



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 3, 2024 – 10:21 PM EST

PDB ID : 1N36
Title : Structure of the *Thermus thermophilus* 30S ribosomal subunit in the presence of crystallographically disordered codon and near-cognate transfer RNA anticodon stem-loop mismatched at the second codon position
Authors : Ogle, J.M.; Murphy IV, F.V.; Tarry, M.J.; Ramakrishnan, V.
Deposited on : 2002-10-25
Resolution : 3.65 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

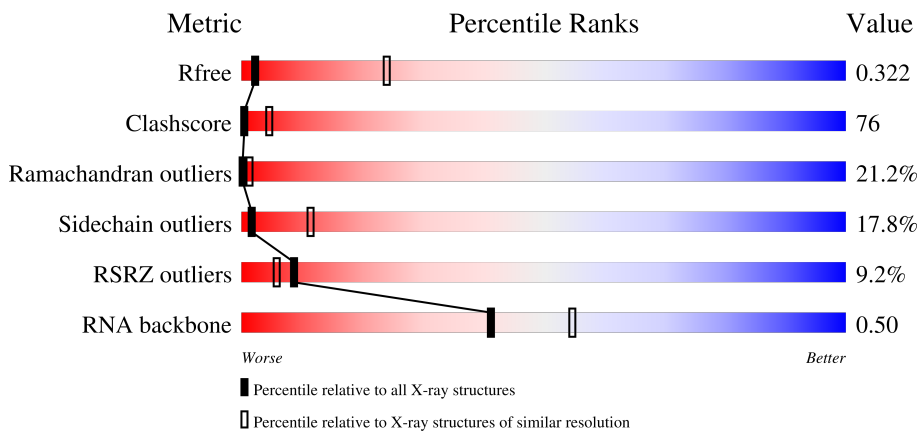
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.65 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1557 (3.82-3.50)
Clashscore	141614	1037 (3.80-3.52)
Ramachandran outliers	138981	1004 (3.80-3.52)
Sidechain outliers	138945	1002 (3.80-3.52)
RSRZ outliers	127900	1441 (3.82-3.50)
RNA backbone	3102	1024 (4.30-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	1522	<div style="display: flex; align-items: center;"> <div style="width: 11%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 9%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 70%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 18%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <div style="display: flex; justify-content: space-between; width: 100%;"> 11% 9% 70% 18% •• </div>
2	B	256	<div style="display: flex; align-items: center;"> <div style="width: 8%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 55%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 23%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 5%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 9%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 8% 55% 23% 5% 9% </div>
3	C	239	<div style="display: flex; align-items: center;"> <div style="width: 8%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 49%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 29%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 14%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <div style="display: flex; justify-content: space-between; width: 100%;"> 8% 6% 49% 29% • 14% </div>
4	D	208	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 12%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 57%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 25%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> </div> <div style="display: flex; justify-content: space-between; width: 100%;"> 3% 12% 57% 25% 6% </div>

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Mol	Chain	Length	Quality of chain
5	E	161	% 14% 56% 22% 7%
6	F	101	5% 17% 63% 19%
7	G	155	4% 14% 59% 26%
8	H	138	% 16% 60% 20%
9	I	128	5% 32% 65% 26%
10	J	104	5% 48% 38% 6%
11	K	129	3% 11% 47% 32% 8%
12	L	135	4% 13% 54% 24% 8%
13	M	126	10% 10% 59% 24% 6%
14	N	60	40% 12% 40% 43% 5%
15	O	88	3% 11% 56% 32%
16	P	88	2% 57% 30% 5% 6%
17	Q	104	3% 12% 66% 19%
18	R	88	3% 11% 49% 18% 5% 17%
19	S	92	24% 10% 57% 20% 13%
20	T	106	% 15% 46% 29% 7%
21	V	26	42% 8% 58% 27% 8%

2 Entry composition i

There are 22 unique types of molecules in this entry. The entry contains 51680 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a RNA chain called 16S RIBOSOMAL RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	A	1513	32508	14472	6016	10509	1511	22	0	0

- Molecule 2 is a protein called 30S RIBOSOMAL PROTEIN S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	234	1900	1213	341	341	5	0	0	0

- Molecule 3 is a protein called 30S RIBOSOMAL PROTEIN S3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	206	1612	1016	314	281	1	0	0	0

- Molecule 4 is a protein called 30S RIBOSOMAL PROTEIN S4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	D	208	1703	1066	339	291	7	0	0	0

- Molecule 5 is a protein called 30S RIBOSOMAL PROTEIN S5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
5	E	150	1146	724	217	201	4	0	0	0

- Molecule 6 is a protein called 30S RIBOSOMAL PROTEIN S6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
6	F	101	843	531	155	154	3	0	0	0

- Molecule 7 is a protein called 30S RIBOSOMAL PROTEIN S7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
7	G	155	1257	781	252	218	6	0	0	0

- Molecule 8 is a protein called 30S RIBOSOMAL PROTEIN S8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
8	H	138	1116	705	215	193	3	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	25	ASP	GLU	conflict	UNP Q5SHQ2
H	37	ARG	LYS	conflict	UNP Q5SHQ2
H	52	ASP	GLU	conflict	UNP Q5SHQ2
H	61	VAL	ILE	conflict	UNP Q5SHQ2
H	62	TYR	HIS	conflict	UNP Q5SHQ2
H	81	HIS	LYS	conflict	UNP Q5SHQ2
H	88	LYS	ARG	conflict	UNP Q5SHQ2
H	115	SER	PRO	conflict	UNP Q5SHQ2

- Molecule 9 is a protein called 30S RIBOSOMAL PROTEIN S9.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
9	I	127	1011	639	198	174	0	0	0

- Molecule 10 is a protein called 30S RIBOSOMAL PROTEIN S10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
10	J	98	792	498	156	137	1	0	0	0

- Molecule 11 is a protein called 30S RIBOSOMAL PROTEIN S11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
11	K	119	885	549	168	165	3	0	0	0

- Molecule 12 is a protein called 30S RIBOSOMAL PROTEIN S12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
12	L	124	970	611	195	163	1	0	0	0

- Molecule 13 is a protein called 30S RIBOSOMAL PROTEIN S13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
13	M	118	937	579	193	163	2	0	0	0

- Molecule 14 is a protein called 30S RIBOSOMAL PROTEIN S14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
14	N	60	492	312	104	72	4	0	0	0

- Molecule 15 is a protein called 30S RIBOSOMAL PROTEIN S15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
15	O	88	734	459	147	126	2	0	0	0

- Molecule 16 is a protein called 30S RIBOSOMAL PROTEIN S16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
16	P	83	700	443	139	117	1	0	0	0

- Molecule 17 is a protein called 30S RIBOSOMAL PROTEIN S17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
17	Q	104	857	547	161	147	2	0	0	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Q	50	LYS	ARG	conflict	UNP Q5SHP7
Q	53	LEU	VAL	conflict	UNP Q5SHP7
Q	62	SER	ALA	conflict	UNP Q5SHP7
Q	79	SER	GLU	conflict	UNP Q5SHP7
Q	82	MET	LEU	conflict	UNP Q5SHP7
Q	90	ILE	VAL	conflict	UNP Q5SHP7

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Chain	Residue	Modelled	Actual	Comment	Reference
Q	96	GLN	ALA	conflict	UNP Q5SHP7

- Molecule 18 is a protein called 30S RIBOSOMAL PROTEIN S18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
18	R	73	597	380	118	99	0	0	0

- Molecule 19 is a protein called 30S RIBOSOMAL PROTEIN S19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
19	S	80	647	414	119	112	2	0	0	0

- Molecule 20 is a protein called 30S RIBOSOMAL PROTEIN S20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
20	T	99	763	470	162	129	2	0	0	0

- Molecule 21 is a protein called 30S RIBOSOMAL PROTEIN THX.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
21	V	24	208	128	50	30	0	0	0

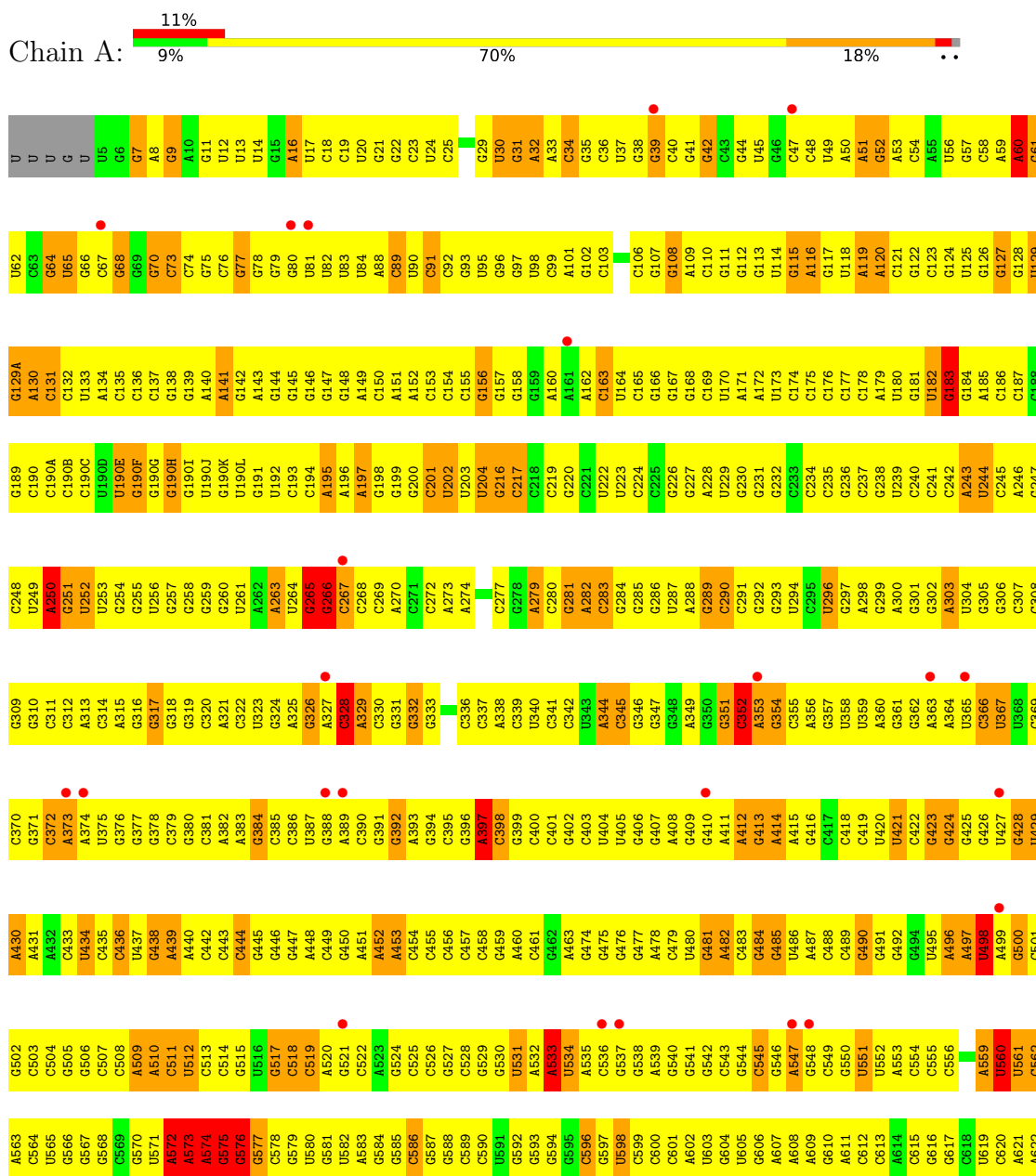
- Molecule 22 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
22	D	1	Total	Zn	0	0
			1	1		
22	N	1	Total	Zn	0	0
			1	1		

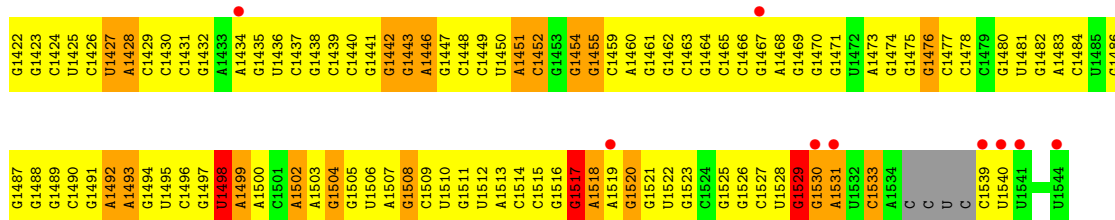
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

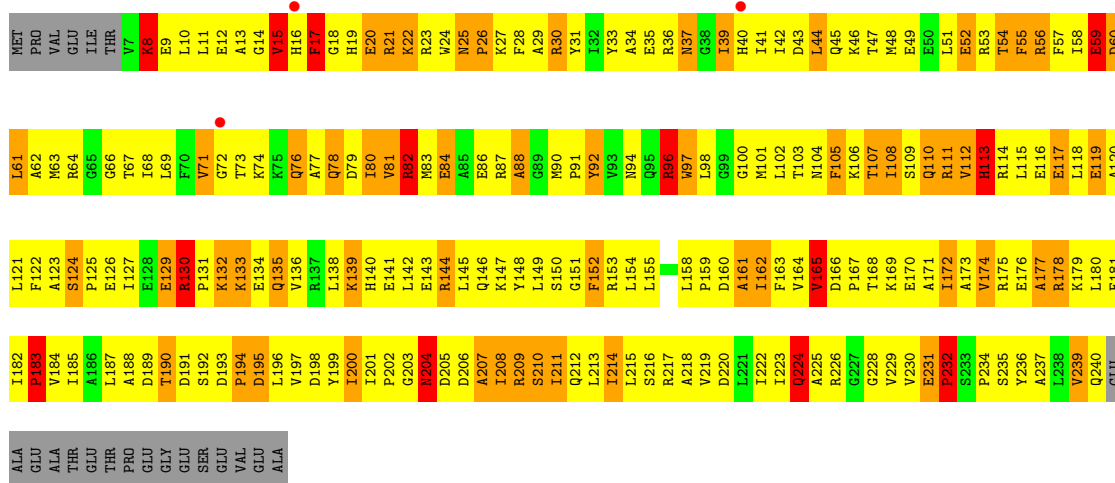
• Molecule 1: 16S RIBOSOMAL RNA



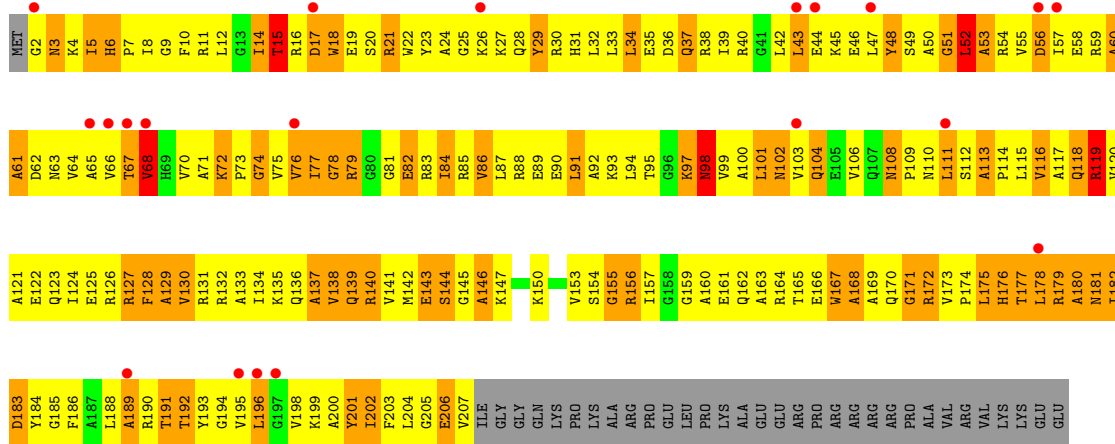
A1360	A1361	C1362	A1363	A1364	A1365	A1366	A1367	A1368	A1369	A1370	A1371	A1372	A1373	A1374	A1375	A1376	A1377	A1378	A1379	A1380	A1381	A1382	A1383	A1384	A1385	A1386	A1387	A1388	A1389	A1390	A1391	A1392	A1393	A1394	A1395	A1396	A1397	A1398	A1399	A1400	A1401	A1402	A1403	A1404	A1405	A1406	A1407	A1408	A1409	A1410	A1411	A1412	A1413	A1414	A1415	A1416	A1417	A1418	A1419	A1420	A1421																	
A1299	G1300	A1301	A1302	A1303	A1304	A1305	A1306	A1307	G1310	A1311	A1312	A1313	A1314	A1315	A1316	A1317	A1318	A1319	A1320	A1321	A1322	A1323	A1324	A1325	A1326	A1327	A1328	A1329	A1330	A1331	A1332	A1333	A1334	A1335	A1336	A1337	A1338	A1339	A1340	A1341	A1342	A1343	A1344	A1345	A1346	A1347	A1348	A1349	A1350	A1351	A1352	A1353	A1354	A1355	A1356	A1357	A1358	A1359																				
A1239	A1240	G1241	A1242	A1243	A1244	A1245	A1246	A1247	A1248	A1249	A1250	A1251	A1252	A1253	A1254	A1255	A1256	A1257	A1258	A1259	A1260	A1261	A1262	A1263	A1264	A1265	A1266	A1267	A1268	A1269	A1270	A1271	A1272	A1273	A1274	A1275	A1276	A1277	A1278	A1279	A1280	A1281	A1282	A1283	A1284	A1285	A1286	A1287	A1288	A1289	A1290	A1291	A1292	A1293	A1294	A1295	A1296	A1297	A1298																			
A1179	A1180	A1181	G1182	A1183	A1184	A1185	A1186	A1187	A1188	A1189	A1190	A1191	A1192	A1193	A1194	A1195	A1196	A1197	A1198	A1199	A1200	A1201	A1202	A1203	A1204	A1205	A1206	A1207	A1208	A1209	A1210	A1211	A1212	A1213	A1214	A1215	A1216	A1217	A1218	A1219	A1220	A1221	A1222	A1223	A1224	A1225	A1226	A1227	A1228	A1229	A1230	A1231	A1232	A1233	A1234	A1235	A1236	A1237	A1238																			
C1116	G1117	C1118	C1119	C1120	C1121	C1122	C1123	C1124	C1125	C1126	C1127	C1128	C1129	C1130	C1131	C1132	C1133	C1134	C1135	C1136	C1137	C1138	C1139	C1140	C1141	C1142	C1143	C1144	C1145	C1146	C1147	C1148	C1149	C1150	C1151	C1152	C1153	C1154	C1155	C1156	C1157	C1158	C1159	C1160	C1161	C1162	C1163	C1164	C1165	C1166	C1167	C1168	C1169	C1170	C1171	C1172	C1173	C1174	C1175	C1176	C1177	C1178																
U1056	G1057	C1058	C1059	C1060	C1061	C1062	C1063	C1064	C1065	C1066	C1067	C1068	C1069	C1070	C1071	C1072	C1073	C1074	C1075	C1076	C1077	C1078	C1079	C1080	C1081	C1082	C1083	C1084	C1085	C1086	C1087	C1088	C1089	C1090	C1091	C1092	C1093	C1094	C1095	C1096	C1097	C1098	C1099	C1100	C1101	C1102	C1103	C1104	C1105	C1106	C1107	C1108	C1109	C1110	C1111	C1112	C1113	C1114	C1115																			
U997	G998	C999	C1000	C1001	C1002	C1003	C1003A	A1004	C1005	C1006	C1007	C1008	C1009	C1010	C1011	C1012	C1013	C1014	C1015	C1016	C1017	C1018	C1019	C1020	C1021	C1022	C1023	C1024	C1025	C1026	C1027	C1028	C1029	C1030	C1030A	C1030B	C1030C	A1030D	C1031	C1032	C1033	C1034	C1035	C1036	C1037	C1038	C1039	C1040	C1041	C1042	C1043	C1044	C1045	C1046	C1047	C1048	C1049	C1050	C1051	C1052	C1053	C1054	C1055															
A937	A938	C939	C940	C941	C942	C943	C944	C945	C946	C947	C948	C949	C950	C951	C952	C953	C954	C955	C956	C957	C958	C959	C960	C961	C962	C963	C964	C965	C966	C967	C968	C969	C970	C971	C972	C973	C974	C975	C976	C977	C978	C979	C980	C981	C982	C983	C984	C985	C986	C987	C988	C989	C990	C991	C992	C993	C994	C995	C996																			
C875	C876	C877	C878	C879	C880	C881	C882	C883	C884	C885	C886	C887	C888	C889	C890	C891	C892	C893	C894	C895	C896	C897	C898	C899	C900	C901	C902	C903	C904	C905	C906	C907	C908	C909	C910	C911	C912	C913	C914	C915	C916	C917	C918	C919	C920	C921	C922	C923	C924	C925	C926	C927	C928	C929	C930	C931	C932	C933	C934	C935	C936																	
C749	C750	C751	C752	C753	C754	C755	C756	C757	C758	C759	C760	C761	C762	C763	C764	C765	C766	C767	C768	C769	C770	C771	C772	C773	C774	C775	C776	C777	C778	C779	C780	C781	C782	C783	C784	C785	C786	C787	C788	C789	C790	C791	C792	C793	C794	C795	C796	C797	C798	C799	C800	C801	C802	C803	C804	C805	C806	C807	C808																			
G609	C610	C611	C612	C613	C614	C615	C616	C617	C618	C619	C620	C621	C622	C623	C624	C625	C626	C627	C628	C629	C630	C631	C632	C633	C634	C635	C636	C637	C638	C639	C640	C641	C642	C643	C644	C645	C646	C647	C648	C649	C650	C651	C652	C653	C654	C655	C656	C657	C658	C659	C660	C661	C662	C663	C664	C665	C666	C667	C668	C669	C670	C671	C672	C673	C674	C675	C676	C677	C678	C679	C680	C681	C682	C683	C684	C685	C686	C687
U686	A687	C688	C689	C690	C691	C692	C693	C694	C695	C696	C697	C698	C699	C700	C701	C702	C703	C704	C705	C706	C707	C708	C709	C710	C711	C712	C713	C714	C715	C716	C717	C718	C719	C720	C721	C722	C723	C724	C725	C726	C727	C728	C729	C730	C731	C732	C733	C734	C735	C736	C737	C738	C739	C740	C741	C742	C743	C744	C745	C746	C747	C748																



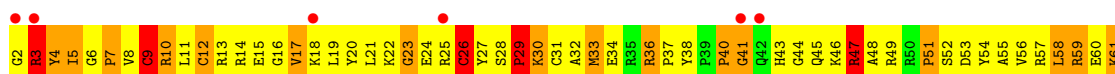
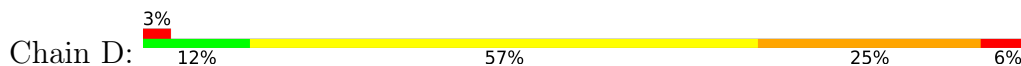
● Molecule 2: 30S RIBOSOMAL PROTEIN S2

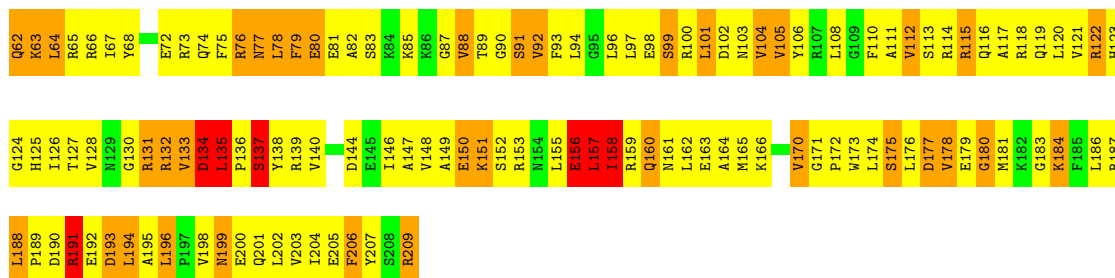


● Molecule 3: 30S RIBOSOMAL PROTEIN S3

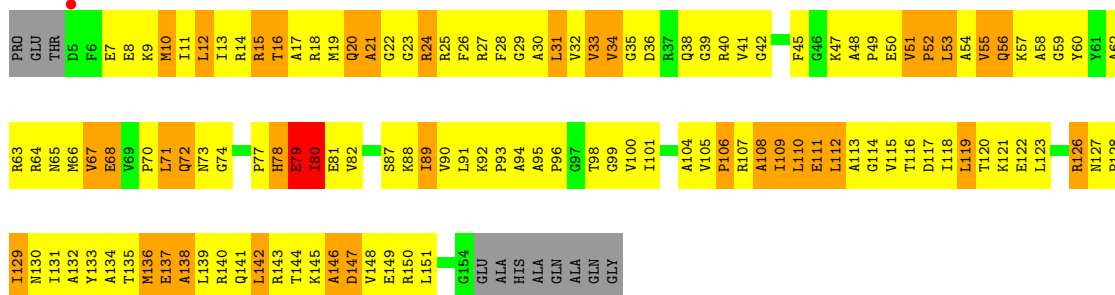
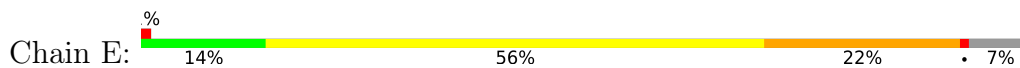


● Molecule 4: 30S RIBOSOMAL PROTEIN S4

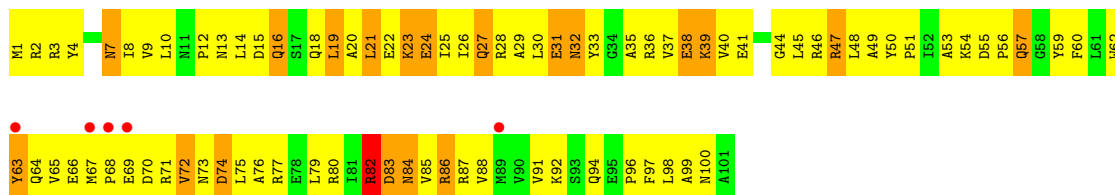




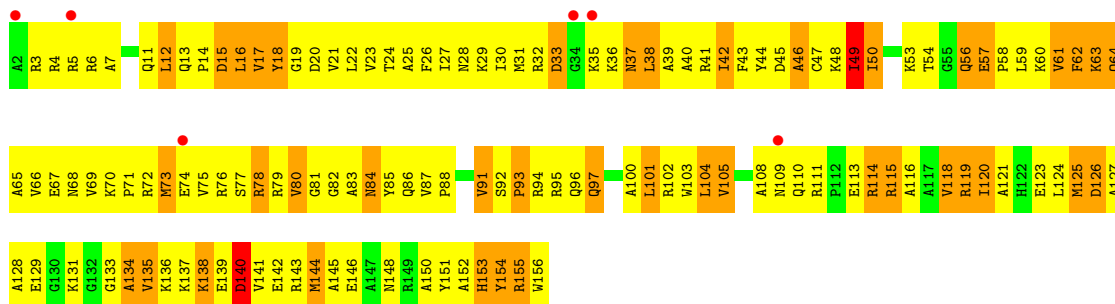
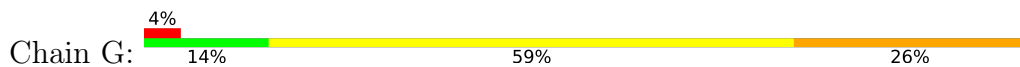
- Molecule 5: 30S RIBOSOMAL PROTEIN S5



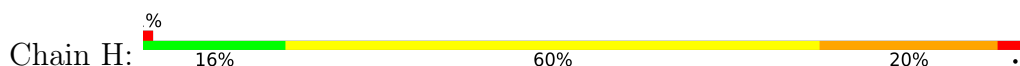
- Molecule 6: 30S RIBOSOMAL PROTEIN S6

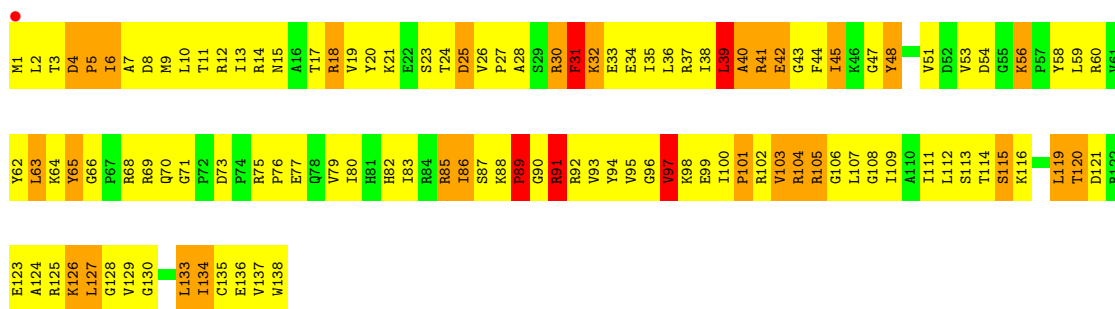


- Molecule 7: 30S RIBOSOMAL PROTEIN S7

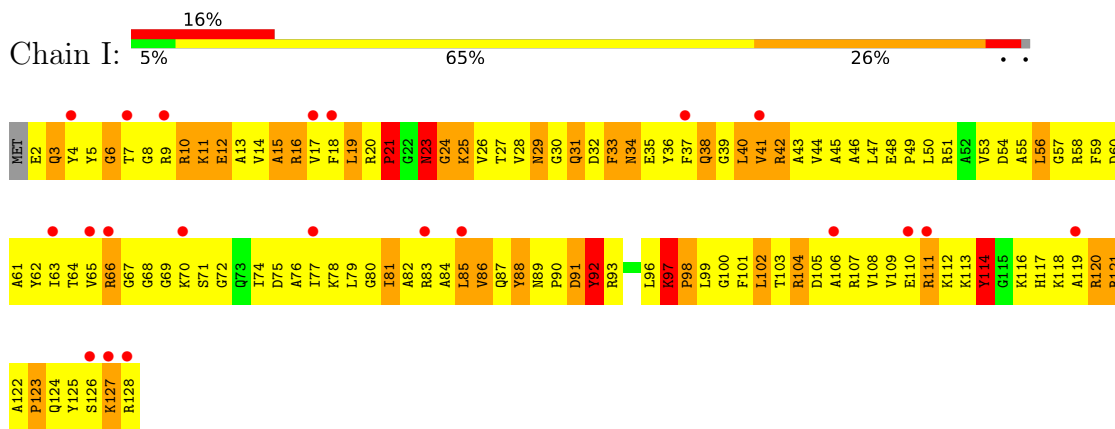


- Molecule 8: 30S RIBOSOMAL PROTEIN S8

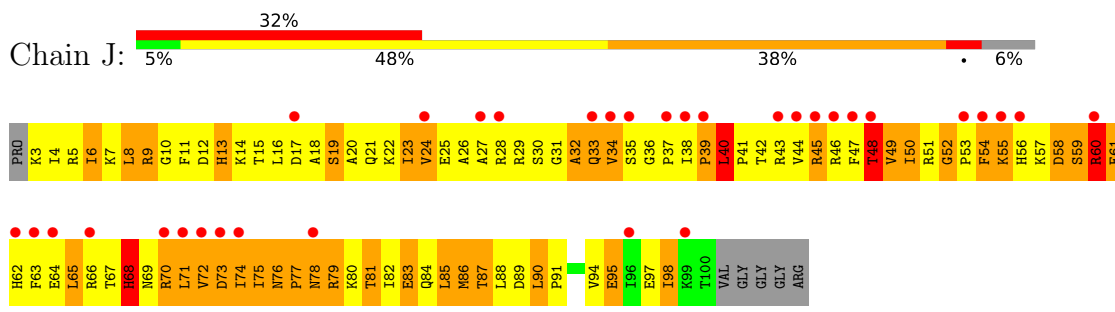




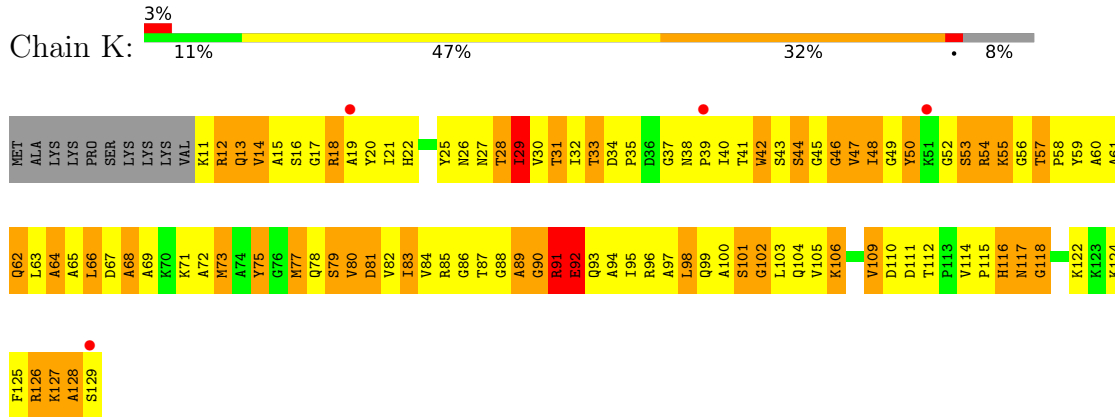
• Molecule 9: 30S RIBOSOMAL PROTEIN S9



• Molecule 10: 30S RIBOSOMAL PROTEIN S10

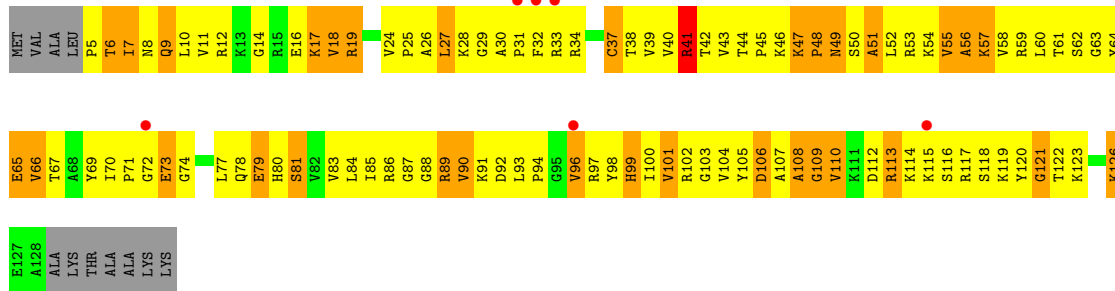


• Molecule 11: 30S RIBOSOMAL PROTEIN S11



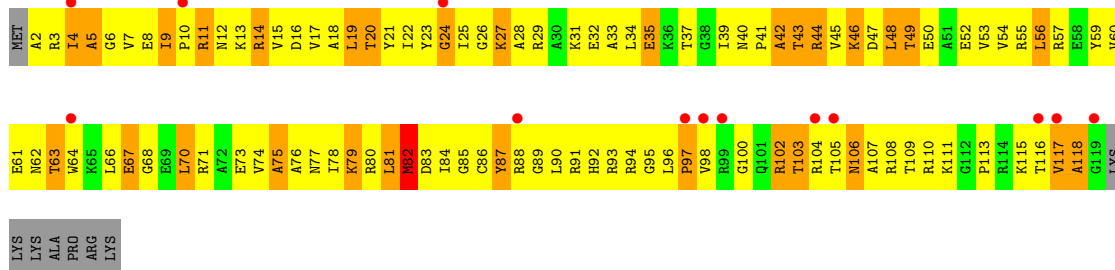
• Molecule 12: 30S RIBOSOMAL PROTEIN S12

Chain L: 4% 13% 54% 24% 8%



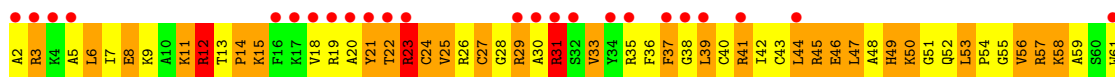
• Molecule 13: 30S RIBOSOMAL PROTEIN S13

Chain M: 10% 10% 59% 24% 6%



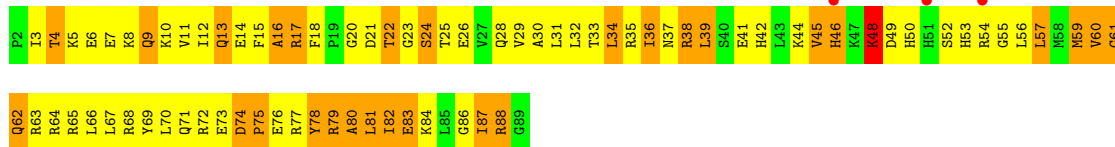
• Molecule 14: 30S RIBOSOMAL PROTEIN S14

Chain N: 40% 12% 40% 43% 5%



• Molecule 15: 30S RIBOSOMAL PROTEIN S15

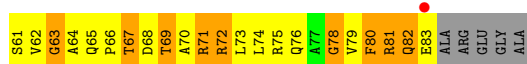
Chain O: 3% 11% 56% 32%



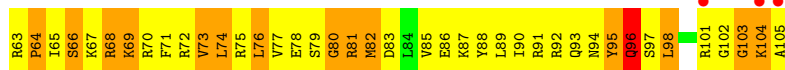
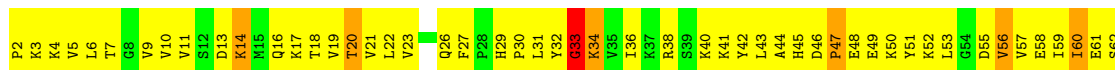
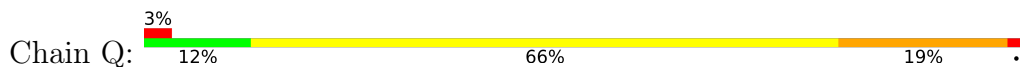
• Molecule 16: 30S RIBOSOMAL PROTEIN S16

Chain P: 2% 57% 30% 5% 6%

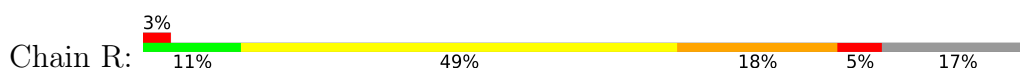




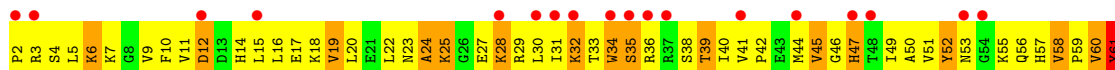
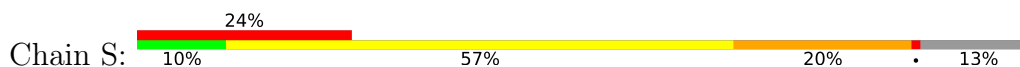
• Molecule 17: 30S RIBOSOMAL PROTEIN S17



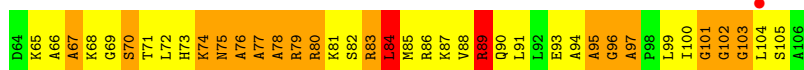
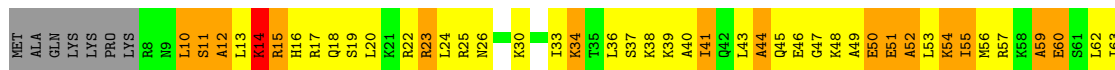
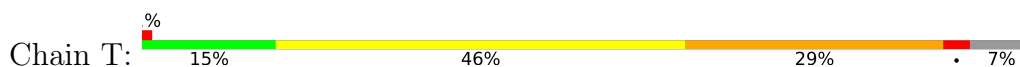
• Molecule 18: 30S RIBOSOMAL PROTEIN S18



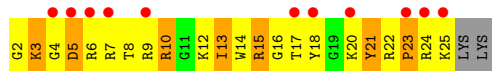
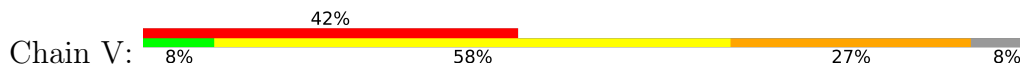
• Molecule 19: 30S RIBOSOMAL PROTEIN S19



• Molecule 20: 30S RIBOSOMAL PROTEIN S20



• Molecule 21: 30S RIBOSOMAL PROTEIN THX



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	402.84Å 402.84Å 174.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	141.42 – 3.65 148.66 – 3.64	Depositor EDS
% Data completeness (in resolution range)	92.6 (141.42-3.65) 89.6 (148.66-3.64)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.44 (at 3.67Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.260 , 0.324 0.259 , 0.322	Depositor DCC
R_{free} test set	7116 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	107.1	Xtrriage
Anisotropy	0.426	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.22 , 121.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	51680	wwPDB-VP
Average B, all atoms (Å ²)	93.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.54% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section:
ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.64	4/36387 (0.0%)	0.78	27/56789 (0.0%)
2	B	0.47	0/1935	0.79	0/2609
3	C	0.46	0/1636	0.77	0/2205
4	D	0.49	0/1733	0.70	0/2318
5	E	0.61	0/1162	0.88	2/1564 (0.1%)
6	F	0.42	0/856	0.72	0/1154
7	G	0.44	0/1276	0.77	1/1709 (0.1%)
8	H	0.65	0/1136	0.87	1/1527 (0.1%)
9	I	0.45	0/1029	0.73	0/1378
10	J	0.47	0/805	0.86	0/1082
11	K	0.49	0/900	0.81	0/1213
12	L	0.51	0/986	0.87	0/1320
13	M	0.40	0/947	0.73	0/1270
14	N	0.46	0/501	0.75	1/664 (0.2%)
15	O	0.51	0/745	0.74	0/992
16	P	0.58	0/716	0.88	1/963 (0.1%)
17	Q	0.58	0/870	0.90	2/1159 (0.2%)
18	R	0.45	0/603	0.75	0/799
19	S	0.47	0/661	0.82	0/890
20	T	0.48	0/765	0.79	0/1007
21	V	0.48	0/212	0.66	0/277
All	All	0.59	4/55861 (0.0%)	0.78	35/82889 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	1	63
8	H	0	1

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Mol	Chain	#Chirality outliers	#Planarity outliers
All	All	1	64

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	858	G	C5-C6	-6.65	1.35	1.42
1	A	1508	G	C5-C6	-5.12	1.37	1.42
1	A	574	A	C5-C6	-5.04	1.36	1.41
1	A	821	G	C5-C6	-5.03	1.37	1.42

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	290	C	N1-C1'-C2'	-8.80	102.33	112.00
1	A	1498	U	C2'-C3'-O3'	8.73	128.71	109.50
1	A	575	G	C2'-C3'-O3'	7.82	126.70	109.50
1	A	60	A	C2'-C3'-O3'	7.67	126.37	109.50
1	A	1454	G	N9-C1'-C2'	-7.34	103.92	112.00
1	A	1517	G	N9-C1'-C2'	-6.94	104.36	112.00
1	A	266	G	C2'-C3'-O3'	6.93	124.79	113.70
1	A	108	G	O4'-C1'-N9	6.84	113.68	108.20
1	A	1529	G	N9-C1'-C2'	6.61	122.59	114.00
1	A	34	C	N1-C1'-C2'	-6.36	105.01	112.00
1	A	141	A	N9-C1'-C2'	-6.31	105.06	112.00
1	A	812	C	N1-C1'-C2'	6.10	121.93	114.00
5	E	110	LEU	CA-CB-CG	-6.10	101.27	115.30
1	A	68	G	N9-C1'-C2'	-5.94	105.47	112.00
1	A	70	G	N9-C1'-C2'	-5.79	105.63	112.00
1	A	1085	U	C2'-C3'-O3'	5.66	122.75	113.70
1	A	190(H)	G	N9-C1'-C2'	-5.58	105.86	112.00
1	A	820	U	N1-C1'-C2'	5.54	121.20	114.00
1	A	265	G	O4'-C1'-N9	5.46	112.57	108.20
1	A	586	C	N1-C1'-C2'	-5.42	106.04	112.00
16	P	51	VAL	N-CA-C	5.32	125.37	111.00
7	G	49	ILE	N-CA-C	-5.30	96.68	111.00
1	A	328	C	C2'-C3'-O3'	5.29	122.17	113.70
1	A	183	G	N9-C1'-C2'	-5.25	106.22	112.00
5	E	119	LEU	CA-CB-CG	-5.25	103.23	115.30
14	N	31	ARG	N-CA-C	5.23	125.13	111.00
1	A	16	A	N9-C1'-C2'	-5.23	106.25	112.00
17	Q	33	GLY	N-CA-C	5.23	126.17	113.10
17	Q	56	VAL	N-CA-C	-5.21	96.94	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	H	89	PRO	N-CA-C	-5.11	98.83	112.10
1	A	1443	G	N9-C1'-C2'	5.10	120.62	114.00
1	A	573	A	N9-C1'-C2'	5.09	120.61	114.00
1	A	818	G	N9-C1'-C2'	5.08	120.61	114.00
1	A	839	U	N1-C1'-C2'	5.08	120.61	114.00
1	A	533	A	C2'-C3'-O3'	5.08	121.82	113.70

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	1498	U	C3'

All (64) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1066	C	Sidechain
1	A	1067	A	Sidechain
1	A	1094	G	Sidechain
1	A	1196	U	Sidechain
1	A	1213	A	Sidechain
1	A	1231	G	Sidechain
1	A	127	G	Sidechain
1	A	1281	U	Sidechain
1	A	129	U	Sidechain
1	A	1301	U	Sidechain
1	A	1370	G	Sidechain
1	A	1381	U	Sidechain
1	A	1401	G	Sidechain
1	A	1454	G	Sidechain
1	A	1455	G	Sidechain
1	A	1502	A	Sidechain
1	A	1531	A	Sidechain
1	A	156	G	Sidechain
1	A	183	G	Sidechain
1	A	239	U	Sidechain
1	A	250	A	Sidechain
1	A	263	A	Sidechain
1	A	265	G	Sidechain
1	A	296	U	Sidechain
1	A	303	A	Sidechain
1	A	305	G	Sidechain
1	A	317	G	Sidechain

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Mol	Chain	Res	Type	Group
1	A	332	G	Sidechain
1	A	352	C	Sidechain
1	A	396	G	Sidechain
1	A	397	A	Sidechain
1	A	434	U	Sidechain
1	A	444	C	Sidechain
1	A	490	G	Sidechain
1	A	498	U	Sidechain
1	A	533	A	Sidechain
1	A	551	U	Sidechain
1	A	560	U	Sidechain
1	A	572	A	Sidechain
1	A	573	A	Sidechain
1	A	574	A	Sidechain
1	A	576	G	Sidechain
1	A	634	C	Sidechain
1	A	666	G	Sidechain
1	A	682	G	Sidechain
1	A	686	U	Sidechain
1	A	691	G	Sidechain
1	A	694	A	Sidechain
1	A	740	U	Sidechain
1	A	756	C	Sidechain
1	A	767	A	Sidechain
1	A	77	G	Sidechain
1	A	777	A	Sidechain
1	A	801	U	Sidechain
1	A	811	C	Sidechain
1	A	819	A	Sidechain
1	A	829	G	Sidechain
1	A	835	U	Sidechain
1	A	870	U	Sidechain
1	A	882	C	Sidechain
1	A	887	G	Sidechain
1	A	898	G	Sidechain
1	A	913	A	Sidechain
8	H	48	TYR	Sidechain

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	32508	0	16414	2582	0
2	B	1900	0	1951	415	0
3	C	1612	0	1677	503	0
4	D	1703	0	1764	410	0
5	E	1146	0	1207	238	0
6	F	843	0	857	150	0
7	G	1257	0	1296	278	0
8	H	1116	0	1177	221	0
9	I	1011	0	1043	277	0
10	J	792	0	835	283	0
11	K	885	0	904	167	0
12	L	970	0	1057	204	0
13	M	937	0	995	225	0
14	N	492	0	532	165	0
15	O	734	0	771	150	0
16	P	700	0	720	199	0
17	Q	857	0	930	177	0
18	R	597	0	668	137	0
19	S	647	0	673	182	0
20	T	763	0	861	206	0
21	V	208	0	221	52	0
22	D	1	0	0	0	0
22	N	1	0	0	0	0
All	All	51680	0	36553	6732	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 76.

All (6732) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:158:ILE:H	4:D:158:ILE:CD1	1.57	1.15
1:A:243:A:H4'	1:A:244:U:H5'	1.22	1.12
4:D:176:LEU:HG	4:D:177:ASP:H	0.96	1.12
1:A:1250:A:H4'	9:I:68:GLY:HA2	1.31	1.12
1:A:1347:G:N2	1:A:1373:G:H2'	1.65	1.12
17:Q:97:SER:HB2	17:Q:103:GLY:HA2	1.21	1.11
14:N:24:CYS:SG	14:N:27:CYS:HB3	1.89	1.11
1:A:1443:G:C5'	1:A:1446:A:H5'	1.80	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:45:ARG:HB3	10:J:45:ARG:HH11	1.01	1.10
4:D:17:VAL:HG12	4:D:18:LYS:H	1.11	1.10
17:Q:70:ARG:HG3	17:Q:70:ARG:HH11	1.17	1.10
10:J:8:LEU:HD13	10:J:16:LEU:HD21	1.34	1.09
20:T:33:ILE:HD11	20:T:63:ILE:HA	1.34	1.09
2:B:101:MET:HA	2:B:108:ILE:HD13	1.31	1.08
1:A:501:C:H2'	1:A:502:G:H8	1.13	1.08
1:A:793:U:H3'	1:A:794:A:C5'	1.83	1.08
3:C:50:ALA:HA	3:C:72:LYS:HD3	1.34	1.08
1:A:1443:G:H5''	1:A:1446:A:C5'	1.82	1.08
11:K:127:LYS:HA	11:K:127:LYS:HE3	1.30	1.08
1:A:1372:U:H5''	9:I:71:SER:HB3	1.28	1.08
15:O:16:ALA:HB1	15:O:21:ASP:HB3	1.34	1.07
15:O:87:ILE:HG22	15:O:88:ARG:H	1.16	1.07
3:C:85:ARG:HA	3:C:88:ARG:HG3	1.33	1.07
9:I:102:LEU:HD23	9:I:102:LEU:H	1.19	1.07
8:H:87:SER:HB2	8:H:93:VAL:HB	1.33	1.06
3:C:71:ALA:HA	3:C:106:VAL:HB	1.10	1.06
1:A:1182:G:H4'	1:A:1183:A:C5'	1.84	1.06
1:A:1123:A:H4'	10:J:37:PRO:HD2	1.37	1.06
3:C:123:GLN:HB3	3:C:128:PHE:HE2	1.19	1.06
1:A:946:A:H2'	1:A:947:G:H8	1.15	1.06
4:D:64:LEU:HB2	4:D:198:VAL:HG21	1.29	1.05
12:L:47:LYS:HG3	12:L:48:PRO:HD3	1.08	1.05
1:A:1182:G:H4'	1:A:1183:A:H5''	1.39	1.05
1:A:1065:U:H4'	1:A:1066:C:H5'	1.35	1.04
8:H:63:LEU:H	8:H:63:LEU:HD12	1.21	1.04
1:A:1150:U:O3'	10:J:41:PRO:HA	1.56	1.04
1:A:1305:G:N2	1:A:1331:G:H2'	1.71	1.04
7:G:135:VAL:HG13	7:G:138:LYS:HZ2	1.20	1.04
9:I:108:VAL:HG12	9:I:109:VAL:H	1.19	1.04
1:A:200:G:H2'	1:A:201:C:H5''	1.38	1.04
4:D:10:ARG:HH11	4:D:10:ARG:HG3	1.21	1.04
7:G:93:PRO:HG2	7:G:94:ARG:H	1.21	1.04
18:R:39:VAL:HG13	18:R:40:LEU:H	1.21	1.04
3:C:76:VAL:HG21	3:C:103:VAL:HG11	1.36	1.03
10:J:86:MET:HG3	10:J:87:THR:H	1.16	1.03
12:L:89:ARG:HH21	12:L:97:ARG:HG2	1.22	1.03
2:B:82:ARG:HA	2:B:92:TYR:CE2	1.93	1.03
7:G:141:VAL:HA	7:G:144:MET:HB2	1.37	1.03
15:O:12:ILE:HD12	15:O:12:ILE:H	1.20	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:S:28:LYS:HD3	19:S:29:ARG:H	1.22	1.03
8:H:44:PHE:O	8:H:45:ILE:HG23	1.58	1.02
4:D:62:GLN:HA	4:D:62:GLN:HE21	1.21	1.02
7:G:111:ARG:HB3	7:G:111:ARG:HH11	1.24	1.02
4:D:22:LYS:HB2	4:D:26:CYS:SG	1.99	1.02
14:N:6:LEU:HD22	14:N:23:ARG:HH21	1.20	1.02
1:A:1127:G:H1'	1:A:1148:U:H3	1.20	1.01
3:C:59:ARG:HG3	3:C:60:ALA:H	1.26	1.01
2:B:28:PHE:CZ	2:B:189:ASP:HA	1.95	1.01
2:B:231:GLU:HB2	2:B:232:PRO:HD2	1.43	1.01
1:A:1305:G:H22	1:A:1331:G:H2'	1.25	1.01
10:J:77:PRO:HA	10:J:81:THR:HG21	1.42	1.01
1:A:64:G:H4'	1:A:65:U:O5'	1.58	1.00
1:A:463:A:H2'	1:A:474:G:H8	1.25	1.00
3:C:182:ILE:HG22	3:C:183:ASP:H	1.25	1.00
10:J:16:LEU:HA	10:J:94:VAL:HG21	1.42	1.00
4:D:158:ILE:HD12	4:D:158:ILE:N	1.71	1.00
11:K:16:SER:O	11:K:35:PRO:HG3	1.59	1.00
4:D:158:ILE:H	4:D:158:ILE:HD12	0.86	1.00
2:B:44:LEU:HD23	2:B:44:LEU:H	1.21	1.00
16:P:82:GLN:NE2	16:P:82:GLN:H	1.60	0.99
3:C:3:ASN:HD22	3:C:3:ASN:H	1.04	0.99
12:L:84:LEU:HB3	12:L:101:VAL:HG21	1.43	0.99
1:A:1129:C:H4'	1:A:1130:A:N7	1.75	0.99
3:C:52:LEU:HD23	3:C:52:LEU:H	1.27	0.99
3:C:190:ARG:HA	3:C:195:VAL:HG22	1.44	0.99
1:A:456:C:H2'	1:A:457:C:H6	1.26	0.99
12:L:84:LEU:HB3	12:L:101:VAL:CG2	1.92	0.99
19:S:22:LEU:HB3	19:S:28:LYS:HB2	1.43	0.99
4:D:176:LEU:CG	4:D:177:ASP:H	1.75	0.99
1:A:518:C:H5''	1:A:530:G:H1'	1.43	0.99
1:A:736:C:H2'	1:A:737:A:C8	1.97	0.99
1:A:1298:C:H4'	1:A:1299:A:H5'	1.45	0.99
4:D:64:LEU:HD23	4:D:198:VAL:HG11	1.45	0.98
4:D:119:GLN:HG3	4:D:123:HIS:HE1	1.24	0.98
1:A:1412:C:H2'	1:A:1413:A:H8	1.26	0.98
4:D:176:LEU:HG	4:D:177:ASP:N	1.76	0.98
3:C:62:ASP:HA	3:C:97:LYS:HB3	1.46	0.98
1:A:269:C:H2'	1:A:270:A:H8	1.27	0.98
16:P:22:THR:HG23	16:P:23:ASP:H	1.28	0.98
4:D:156:GLU:HG2	4:D:157:LEU:HD13	1.42	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:79:ARG:HA	10:J:82:ILE:HB	1.44	0.98
1:A:284:G:H2'	1:A:285:G:H8	1.27	0.98
20:T:40:ALA:HA	20:T:55:ILE:HD11	1.42	0.98
8:H:11:THR:HG23	8:H:14:ARG:NH1	1.79	0.98
13:M:81:LEU:HD13	13:M:88:ARG:HB3	1.46	0.98
19:S:32:LYS:HA	19:S:50:ALA:HB3	1.44	0.98
3:C:66:VAL:HG23	3:C:99:VAL:HG21	1.44	0.97
9:I:19:LEU:HD21	9:I:59:PHE:HB3	1.46	0.97
5:E:55:VAL:HG23	5:E:56:GLN:H	1.29	0.97
3:C:182:ILE:HG22	3:C:183:ASP:N	1.78	0.97
7:G:38:LEU:HG	7:G:42:ILE:HD11	1.43	0.97
9:I:80:GLY:HA2	9:I:83:ARG:HB3	1.46	0.97
12:L:41:ARG:HB3	12:L:41:ARG:HH11	1.26	0.97
7:G:65:ALA:HB2	7:G:128:ALA:HA	1.43	0.97
5:E:8:GLU:HB2	5:E:34:VAL:HG22	1.47	0.97
8:H:65:TYR:HA	8:H:79:VAL:HG23	1.45	0.97
1:A:1280:A:H5''	10:J:40:LEU:HD21	1.43	0.96
2:B:172:ILE:H	2:B:172:ILE:HD12	1.26	0.96
19:S:12:ASP:HB2	19:S:38:SER:HB2	1.47	0.96
4:D:133:VAL:HG12	4:D:134:ASP:H	1.28	0.96
6:F:33:TYR:HB2	6:F:75:LEU:HD23	1.47	0.96
3:C:156:ARG:HH21	3:C:161:GLU:HA	1.29	0.96
1:A:380:G:N2	1:A:382:A:H3'	1.81	0.96
15:O:48:LYS:N	15:O:48:LYS:HZ2	1.63	0.96
3:C:11:ARG:HH12	3:C:179:ARG:H	1.11	0.96
4:D:64:LEU:O	4:D:64:LEU:HD13	1.66	0.96
1:A:201:C:H2'	1:A:202:U:H5''	1.48	0.95
1:A:376:G:H5''	16:P:5:ARG:HG3	1.48	0.95
1:A:501:C:H2'	1:A:502:G:C8	2.01	0.95
3:C:58:GLU:HB2	3:C:65:ALA:HB3	1.47	0.95
3:C:66:VAL:HG12	3:C:68:VAL:HB	1.47	0.95
10:J:30:SER:HB3	10:J:84:GLN:NE2	1.80	0.95
1:A:962:C:H2'	1:A:963:G:H8	1.31	0.95
3:C:88:ARG:HA	3:C:91:LEU:HB3	1.48	0.95
18:R:43:PHE:HA	18:R:51:LEU:HD12	1.48	0.95
13:M:10:PRO:HB2	13:M:18:ALA:HB1	1.48	0.95
1:A:393:A:H2'	1:A:394:G:H8	1.30	0.95
1:A:444:C:H2'	1:A:445:G:H8	1.30	0.95
15:O:82:ILE:HD11	15:O:88:ARG:HG3	1.47	0.95
3:C:83:ARG:HA	3:C:86:VAL:HG23	1.46	0.94
5:E:41:VAL:HG22	5:E:113:ALA:HA	1.49	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:29:LYS:HZ1	7:G:102:ARG:HA	1.31	0.94
5:E:110:LEU:HD13	5:E:118:ILE:HD12	1.48	0.94
10:J:6:ILE:HG13	10:J:73:ASP:HA	1.49	0.94
1:A:1267:C:H1'	21:V:20:LYS:HE3	1.48	0.94
5:E:80:ILE:HG22	8:H:104:ARG:HH21	1.32	0.94
20:T:90:GLN:O	20:T:93:GLU:HG2	1.68	0.94
1:A:243:A:C4'	1:A:244:U:H5'	1.97	0.94
1:A:737:A:H2'	1:A:738:C:H6	1.32	0.94
5:E:35:GLY:H	5:E:112:LEU:HD12	1.30	0.94
15:O:48:LYS:NZ	15:O:48:LYS:H	1.64	0.94
16:P:67:THR:HG22	16:P:69:THR:H	1.32	0.94
1:A:1343:G:H2'	1:A:1344:C:C6	2.01	0.94
7:G:66:VAL:HG22	7:G:104:LEU:HD11	1.49	0.94
3:C:92:ALA:HA	3:C:95:THR:HG22	1.47	0.94
4:D:10:ARG:HG2	4:D:11:LEU:N	1.80	0.94
4:D:62:GLN:HE22	4:D:65:ARG:HH11	1.13	0.94
1:A:737:A:H2'	1:A:738:C:C6	2.03	0.93
19:S:15:LEU:O	19:S:19:VAL:HB	1.69	0.93
12:L:90:VAL:HG11	12:L:93:LEU:HG	1.50	0.93
1:A:131:C:H2'	1:A:132:C:C6	2.04	0.93
3:C:64:VAL:CG2	3:C:97:LYS:HB2	1.98	0.93
14:N:11:LYS:O	14:N:12:ARG:HB3	1.65	0.93
1:A:946:A:H2'	1:A:947:G:C8	2.02	0.93
6:F:47:ARG:HH11	6:F:47:ARG:HA	1.34	0.93
1:A:1203:C:H2'	1:A:1204:A:O4'	1.69	0.93
2:B:36:ARG:HD2	2:B:41:ILE:HD12	1.48	0.93
12:L:66:VAL:HG12	12:L:67:THR:H	1.34	0.92
1:A:256:U:H2'	1:A:257:G:H8	1.30	0.92
1:A:1178:G:H21	1:A:1180:A:H3'	1.32	0.92
2:B:144:ARG:HG3	2:B:145:LEU:N	1.84	0.92
10:J:4:ILE:HB	10:J:74:ILE:HG13	1.51	0.92
1:A:1007:C:H42	1:A:1022:G:H22	1.12	0.92
3:C:71:ALA:CA	3:C:106:VAL:HB	2.00	0.92
5:E:81:GLU:HG2	5:E:90:VAL:HG12	1.46	0.92
12:L:43:VAL:HG12	12:L:44:THR:H	1.34	0.92
1:A:243:A:H4'	1:A:244:U:C5'	1.99	0.92
1:A:420:U:O2'	1:A:421:U:H5''	1.70	0.92
5:E:78:HIS:ND1	8:H:104:ARG:HD2	1.83	0.92
9:I:14:VAL:O	9:I:65:VAL:HG23	1.68	0.92
1:A:397:A:O2'	1:A:398:C:H5''	1.68	0.92
10:J:38:ILE:HB	10:J:71:LEU:HB3	1.50	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:K:40:ILE:HG22	11:K:41:THR:HG23	1.47	0.92
14:N:14:PRO:HG2	14:N:15:LYS:H	1.34	0.92
2:B:219:VAL:HA	2:B:222:ILE:HD12	1.52	0.91
16:P:59:TRP:HA	16:P:59:TRP:CE3	2.05	0.91
1:A:1047:G:H2'	1:A:1048:G:H5''	1.49	0.91
2:B:96:ARG:NE	2:B:96:ARG:H	1.68	0.91
5:E:51:VAL:HB	5:E:52:PRO:HD3	1.52	0.91
7:G:116:ALA:N	7:G:119:ARG:HH21	1.68	0.91
8:H:24:THR:HG22	8:H:63:LEU:HD11	1.51	0.91
10:J:45:ARG:HB3	10:J:45:ARG:NH1	1.85	0.91
1:A:877:C:O2'	1:A:878:G:H5'	1.70	0.91
1:A:1286:A:H2'	1:A:1287:A:H4'	1.52	0.91
1:A:710:G:H5''	6:F:54:LYS:HZ3	1.35	0.91
2:B:20:GLU:HA	2:B:21:ARG:HH21	1.36	0.91
12:L:45:PRO:HB3	12:L:92:ASP:HB3	1.52	0.91
1:A:707:C:H2'	1:A:708:C:H6	1.33	0.90
1:A:1104:G:H4'	2:B:111:ARG:NH2	1.86	0.90
1:A:1329:A:P	13:M:28:ALA:HB3	2.11	0.90
1:A:256:U:H2'	1:A:257:G:C8	2.05	0.90
20:T:67:ALA:HA	20:T:73:HIS:H	1.33	0.90
20:T:75:ASN:HD22	20:T:75:ASN:N	1.64	0.90
1:A:390:C:H2'	1:A:391:G:H8	1.36	0.90
1:A:750:G:N3	15:O:23:GLY:HA3	1.87	0.90
4:D:12:CYS:SG	4:D:19:LEU:HB2	2.12	0.90
7:G:20:ASP:OD2	7:G:22:LEU:HG	1.72	0.90
1:A:982:U:H5	14:N:31:ARG:HH12	1.18	0.90
11:K:33:THR:HA	11:K:39:PRO:HA	1.53	0.90
1:A:405:U:H3'	1:A:406:G:H5'	1.52	0.90
1:A:990:C:H4'	1:A:1018:C:OP1	1.71	0.90
1:A:1236:A:H4'	1:A:1304:G:H4'	1.52	0.90
4:D:62:GLN:HE22	4:D:65:ARG:NH1	1.70	0.90
15:O:48:LYS:HZ2	15:O:48:LYS:H	0.90	0.89
5:E:50:GLU:HG3	5:E:52:PRO:HD2	1.54	0.89
14:N:53:LEU:HD12	14:N:56:VAL:HB	1.51	0.89
18:R:18:ARG:NE	18:R:18:ARG:HA	1.85	0.89
1:A:1090:U:H2'	1:A:1091:U:H6	1.35	0.89
9:I:10:ARG:CZ	9:I:11:LYS:HB2	2.01	0.89
1:A:503:C:H2'	1:A:504:C:H6	1.38	0.89
1:A:560:U:H4'	1:A:561:U:H5''	1.52	0.89
10:J:49:VAL:HG11	14:N:41:ARG:HB2	1.54	0.89
1:A:1107:C:H2'	1:A:1108:G:H5'	1.53	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:65:VAL:HG11	9:I:77:ILE:HD11	1.55	0.89
1:A:167:G:H2'	1:A:168:G:C8	2.08	0.89
13:M:49:THR:HB	13:M:52:GLU:HG3	1.54	0.89
7:G:155:ARG:HA	7:G:155:ARG:HE	1.37	0.89
4:D:104:VAL:HG21	4:D:140:VAL:HG21	1.54	0.89
7:G:65:ALA:HB2	7:G:128:ALA:CA	2.02	0.89
7:G:111:ARG:HB3	7:G:111:ARG:NH1	1.87	0.89
8:H:134:ILE:HG22	8:H:135:CYS:N	1.87	0.89
14:N:56:VAL:HG22	14:N:57:ARG:H	1.38	0.89
1:A:168:G:O2'	1:A:169:C:H5'	1.73	0.89
3:C:150:LYS:HG3	3:C:169:ALA:HB2	1.55	0.89
16:P:21:VAL:HG21	16:P:59:TRP:CD1	2.08	0.88
16:P:74:LEU:HB3	16:P:79:VAL:HG23	1.54	0.88
1:A:894:G:H2'	1:A:895:G:C8	2.07	0.88
1:A:735:C:H2'	1:A:736:C:H6	1.35	0.88
1:A:1117:G:N2	1:A:1180:A:H1'	1.88	0.88
9:I:7:THR:HB	9:I:83:ARG:NH1	1.89	0.88
9:I:48:GLU:HA	9:I:51:ARG:HE	1.39	0.88
13:M:17:VAL:HG22	13:M:27:LYS:HD3	1.55	0.88
1:A:383:A:H2'	1:A:384:G:H5'	1.54	0.88
1:A:793:U:H3'	1:A:794:A:H5'	1.53	0.88
1:A:1519:A:H2'	1:A:1520:G:H5'	1.56	0.88
7:G:38:LEU:HD23	7:G:39:ALA:H	1.34	0.88
11:K:87:THR:HA	11:K:91:ARG:HH12	1.39	0.88
1:A:390:C:H2'	1:A:391:G:C8	2.08	0.88
1:A:436:C:H2'	1:A:437:U:H6	1.37	0.88
1:A:710:G:H5''	6:F:54:LYS:NZ	1.87	0.88
1:A:1412:C:H2'	1:A:1413:A:C8	2.09	0.88
3:C:120:VAL:HA	3:C:123:GLN:HB2	1.56	0.88
1:A:760:G:N2	17:Q:104:LYS:H	1.72	0.88
16:P:4:ILE:HG13	16:P:64:ALA:HB1	1.54	0.88
1:A:1366:C:H2'	1:A:1367:C:H6	1.36	0.88
4:D:187:ARG:HE	4:D:188:LEU:HD12	1.39	0.88
20:T:57:ARG:NH2	20:T:102:GLY:HA3	1.88	0.87
2:B:178:ARG:HH22	8:H:68:ARG:NH2	1.72	0.87
5:E:11:ILE:HD11	5:E:33:VAL:HG22	1.55	0.87
16:P:21:VAL:HG21	16:P:59:TRP:NE1	1.89	0.87
1:A:235:C:H5'	17:Q:70:ARG:HD3	1.53	0.87
7:G:28:ASN:HA	7:G:31:MET:HB2	1.54	0.87
9:I:7:THR:HG22	9:I:8:GLY:H	1.38	0.87
1:A:392:G:H2'	1:A:393:A:C8	2.08	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1008:C:H42	1:A:1021:G:H1	1.18	0.87
10:J:78:ASN:C	10:J:80:LYS:H	1.77	0.87
21:V:17:THR:HG22	21:V:18:TYR:H	1.38	0.87
1:A:373:A:H2'	1:A:374:A:H8	1.38	0.87
13:M:10:PRO:CB	13:M:18:ALA:HB1	2.04	0.87
19:S:11:VAL:HG22	19:S:39:THR:HG22	1.54	0.87
1:A:601:C:O2'	1:A:602:A:H5'	1.74	0.87
1:A:309:G:H2'	1:A:310:G:H8	1.38	0.86
1:A:382:A:H2'	1:A:383:A:H8	1.36	0.86
1:A:1050:G:H22	1:A:1209:C:H1'	1.38	0.86
1:A:1443:G:H5''	1:A:1446:A:H5'	0.91	0.86
2:B:82:ARG:HA	2:B:92:TYR:HE2	1.40	0.86
18:R:58:LEU:HD22	18:R:62:GLU:HB3	1.57	0.86
1:A:1097:C:H2'	1:A:1098:C:H6	1.40	0.86
1:A:1195:C:H3'	1:A:1196:U:H5'	1.55	0.86
4:D:61:LYS:HD3	4:D:62:GLN:N	1.90	0.86
5:E:89:ILE:HD13	5:E:90:VAL:N	1.89	0.86
7:G:136:LYS:O	7:G:140:ASP:N	2.07	0.86
10:J:9:ARG:HH12	10:J:69:ASN:HB3	1.37	0.86
12:L:47:LYS:HG3	12:L:48:PRO:CD	2.01	0.86
14:N:6:LEU:HD22	14:N:23:ARG:NH2	1.91	0.86
1:A:258:G:H2'	1:A:259:G:H8	1.40	0.86
1:A:284:G:H2'	1:A:285:G:C8	2.09	0.86
1:A:986:A:H1'	19:S:55:LYS:HA	1.58	0.86
1:A:1507:A:C2	1:A:1530:G:H1'	2.10	0.86
5:E:26:PHE:O	5:E:27:ARG:HG3	1.75	0.86
16:P:33:ILE:O	16:P:34:GLU:HB3	1.75	0.86
17:Q:98:LEU:HA	17:Q:102:GLY:HA2	1.57	0.86
3:C:123:GLN:HB3	3:C:128:PHE:CE2	2.09	0.86
8:H:134:ILE:HG22	8:H:135:CYS:H	1.38	0.86
10:J:42:THR:HG23	10:J:68:HIS:HA	1.57	0.86
1:A:357:G:O2'	1:A:358:U:H5'	1.74	0.86
1:A:1047:G:C2'	1:A:1048:G:H5''	2.05	0.86
3:C:129:ALA:HB1	3:C:132:ARG:HB2	1.57	0.86
7:G:38:LEU:HD23	7:G:39:ALA:N	1.90	0.86
11:K:16:SER:HA	11:K:79:SER:HB3	1.58	0.86
1:A:1015:A:H2'	1:A:1016:A:C8	2.11	0.86
3:C:18:TRP:HE1	14:N:56:VAL:HG12	1.38	0.86
4:D:157:LEU:HB2	4:D:158:ILE:HD12	1.56	0.86
6:F:94:GLN:HB3	18:R:32:ARG:HH11	1.38	0.86
1:A:894:G:H2'	1:A:895:G:H8	1.39	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:23:GLY:HA2	4:D:112:VAL:O	1.76	0.86
1:A:1215:G:H2'	1:A:1216:G:O4'	1.76	0.86
3:C:64:VAL:HG23	3:C:97:LYS:HB2	1.56	0.86
4:D:52:SER:C	4:D:54:TYR:H	1.77	0.86
10:J:45:ARG:HH11	10:J:45:ARG:CB	1.87	0.86
15:O:56:LEU:HA	15:O:59:MET:HG3	1.58	0.86
15:O:64:ARG:HH22	15:O:68:ARG:HH22	1.20	0.86
4:D:36:ARG:N	4:D:37:PRO:HD3	1.91	0.85
9:I:16:ARG:HD3	9:I:16:ARG:H	1.41	0.85
9:I:102:LEU:HD23	9:I:102:LEU:N	1.89	0.85
11:K:95:ILE:HG22	11:K:99:GLN:HE21	1.41	0.85
2:B:134:GLU:HB3	2:B:138:LEU:HD23	1.58	0.85
4:D:189:PRO:HB2	4:D:194:LEU:HD21	1.58	0.85
10:J:98:ILE:H	10:J:98:ILE:HD12	1.40	0.85
17:Q:97:SER:HB2	17:Q:103:GLY:CA	2.06	0.85
1:A:115:G:H1'	1:A:116:A:N7	1.89	0.85
3:C:3:ASN:HD22	3:C:3:ASN:N	1.69	0.85
6:F:33:TYR:HA	6:F:71:ARG:NH2	1.90	0.85
13:M:79:LYS:HE2	13:M:79:LYS:HA	1.56	0.85
18:R:21:LYS:H	18:R:21:LYS:HD2	1.39	0.85
19:S:38:SER:HB3	19:S:71:LEU:HD11	1.55	0.85
1:A:528:C:H2'	1:A:529:G:H5'	1.58	0.85
5:E:40:ARG:HG2	5:E:68:GLU:OE2	1.76	0.85
1:A:1347:G:H3'	9:I:108:VAL:O	1.76	0.85
14:N:46:GLU:HA	14:N:49:HIS:HD2	1.39	0.85
1:A:1488:G:H2'	1:A:1489:G:C8	2.11	0.85
2:B:130:ARG:HB3	2:B:131:PRO:HD2	1.59	0.85
3:C:76:VAL:O	3:C:83:ARG:HB3	1.77	0.85
4:D:119:GLN:HG3	4:D:123:HIS:CE1	2.11	0.85
5:E:92:LYS:HB3	5:E:119:LEU:HB2	1.59	0.85
15:O:39:LEU:HD13	15:O:56:LEU:HB2	1.57	0.85
17:Q:95:TYR:O	17:Q:97:SER:N	2.10	0.85
1:A:1255:G:O2'	1:A:1258:G:H1'	1.77	0.85
1:A:200:G:C2'	1:A:201:C:H5''	2.06	0.84
20:T:76:ALA:O	20:T:80:ARG:HG2	1.75	0.84
10:J:40:LEU:HD23	10:J:41:PRO:HD2	1.56	0.84
12:L:47:LYS:CG	12:L:48:PRO:HD3	2.02	0.84
21:V:10:ARG:HA	21:V:13:ILE:HG22	1.60	0.84
7:G:17:VAL:HG12	7:G:18:TYR:N	1.92	0.84
7:G:126:ASP:HB3	7:G:131:LYS:O	1.77	0.84
1:A:126:G:H1	1:A:235:C:H42	1.20	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:T:30:LYS:HD3	20:T:72:LEU:HD21	1.57	0.84
1:A:1522:U:O2'	1:A:1523:G:H5'	1.77	0.84
17:Q:86:GLU:O	17:Q:90:ILE:HG13	1.76	0.84
1:A:761:G:H5''	17:Q:101:ARG:O	1.78	0.84
1:A:882:C:O2'	1:A:883:C:H5'	1.78	0.84
14:N:26:ARG:NH2	14:N:47:LEU:HD11	1.92	0.84
1:A:1020:U:H2'	1:A:1021:G:H8	1.42	0.84
3:C:22:TRP:HB3	3:C:59:ARG:HB3	1.60	0.84
3:C:71:ALA:O	3:C:73:PRO:HD3	1.77	0.84
5:E:87:SER:HB3	5:E:131:ILE:HD13	1.60	0.84
18:R:18:ARG:HA	18:R:18:ARG:HE	1.39	0.84
1:A:376:G:H2'	1:A:377:G:H8	1.42	0.84
1:A:1234:C:H1'	1:A:1364:U:O2	1.76	0.84
10:J:14:LYS:HA	10:J:17:ASP:HB3	1.59	0.84
11:K:58:PRO:CB	11:K:93:GLN:HG3	2.06	0.84
3:C:137:ALA:O	3:C:141:VAL:HG22	1.77	0.84
19:S:15:LEU:HB3	19:S:33:THR:HG21	1.60	0.84
1:A:392:G:H2'	1:A:393:A:H8	1.42	0.84
20:T:83:ARG:O	20:T:87:LYS:HG3	1.77	0.84
21:V:6:ARG:HG2	21:V:15:ARG:HH12	1.41	0.83
1:A:1347:G:H21	1:A:1373:G:H2'	1.40	0.83
3:C:23:TYR:CG	3:C:24:ALA:N	2.44	0.83
3:C:112:SER:HB3	3:C:115:LEU:HD12	1.59	0.83
4:D:19:LEU:HB3	4:D:21:LEU:HD13	1.60	0.83
3:C:48:TYR:HB2	3:C:52:LEU:HB3	1.57	0.83
19:S:28:LYS:CD	19:S:29:ARG:H	1.92	0.83
3:C:91:LEU:HD11	3:C:99:VAL:HG22	1.60	0.83
7:G:135:VAL:HG13	7:G:138:LYS:NZ	1.93	0.83
12:L:25:PRO:C	12:L:27:LEU:H	1.81	0.83
20:T:50:GLU:HG2	20:T:100:ILE:HG13	1.60	0.83
1:A:541:G:H2'	1:A:542:G:H8	1.43	0.83
1:A:1154:G:H2'	1:A:1155:G:H8	1.44	0.83
7:G:46:ALA:HB1	7:G:121:ALA:N	1.93	0.83
13:M:98:VAL:HG23	13:M:110:ARG:HH12	1.40	0.83
3:C:121:ALA:O	3:C:124:ILE:HG22	1.78	0.83
12:L:83:VAL:HG22	12:L:84:LEU:H	1.42	0.83
16:P:74:LEU:HB3	16:P:79:VAL:CG2	2.08	0.83
11:K:17:GLY:H	11:K:77:MET:HE1	1.44	0.83
1:A:538:G:H5''	12:L:114:LYS:HB2	1.60	0.83
1:A:858:G:O2'	1:A:859:A:H5''	1.78	0.83
3:C:126:ARG:C	3:C:128:PHE:H	1.80	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:36:ARG:H	4:D:37:PRO:HD3	1.43	0.83
5:E:80:ILE:HD11	5:E:138:ALA:HB1	1.60	0.83
9:I:86:VAL:HG11	9:I:93:ARG:HG3	1.61	0.83
20:T:53:LEU:HD12	20:T:100:ILE:HB	1.60	0.83
1:A:1066:C:H2'	1:A:1067:A:H5'	1.61	0.83
1:A:1134:G:H2'	1:A:1135:U:O4'	1.79	0.83
3:C:110:ASN:O	3:C:111:LEU:HG	1.79	0.83
10:J:62:HIS:HB3	14:N:59:ALA:HB3	1.60	0.83
1:A:875:C:H1'	8:H:15:ASN:OD1	1.77	0.83
13:M:10:PRO:HB3	13:M:18:ALA:O	1.78	0.83
1:A:518:C:H5''	1:A:530:G:C1'	2.08	0.82
4:D:29:PRO:O	4:D:30:LYS:HG3	1.78	0.82
1:A:7:G:H5'	1:A:298:A:H5'	1.61	0.82
1:A:258:G:H2'	1:A:259:G:C8	2.12	0.82
1:A:1048:G:H21	1:A:1214:C:H2'	1.42	0.82
9:I:75:ASP:O	9:I:78:LYS:HB3	1.79	0.82
17:Q:98:LEU:HA	17:Q:102:GLY:CA	2.09	0.82
19:S:53:ASN:ND2	19:S:77:THR:HA	1.94	0.82
3:C:88:ARG:HH12	3:C:101:LEU:HB3	1.42	0.82
4:D:59:ARG:HH21	4:D:62:GLN:HG3	1.43	0.82
8:H:86:ILE:HG22	8:H:87:SER:N	1.93	0.82
10:J:46:ARG:HH12	10:J:64:GLU:HG2	1.42	0.82
1:A:269:C:H2'	1:A:270:A:C8	2.14	0.82
10:J:10:GLY:HA3	10:J:16:LEU:HD12	1.62	0.82
7:G:29:LYS:O	7:G:105:VAL:HG11	1.80	0.82
10:J:40:LEU:HD22	10:J:69:ASN:OD1	1.79	0.82
13:M:49:THR:HG22	13:M:50:GLU:H	1.45	0.82
1:A:726:C:H42	1:A:731:G:H1	1.27	0.82
1:A:382:A:H2'	1:A:383:A:C8	2.13	0.82
2:B:25:ASN:HD22	2:B:25:ASN:C	1.82	0.82
4:D:156:GLU:HG2	4:D:157:LEU:H	1.44	0.82
1:A:148:G:H2'	1:A:149:A:H8	1.44	0.81
1:A:708:C:H2'	1:A:709:G:C8	2.14	0.81
1:A:824:C:H2'	1:A:825:G:H8	1.44	0.81
4:D:9:CYS:HB2	4:D:22:LYS:NZ	1.95	0.81
4:D:15:GLU:HG3	4:D:63:LYS:NZ	1.94	0.81
12:L:6:THR:O	12:L:8:ASN:N	2.12	0.81
15:O:87:ILE:HG22	15:O:88:ARG:N	1.94	0.81
18:R:38:GLU:N	18:R:41:LYS:HG3	1.95	0.81
1:A:200:G:H1	1:A:217:C:H42	1.23	0.81
8:H:11:THR:HG22	8:H:15:ASN:ND2	1.95	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:C:C2'	1:A:202:U:H5''	2.08	0.81
1:A:428:G:H1'	1:A:430:A:C8	2.15	0.81
2:B:21:ARG:HE	2:B:21:ARG:N	1.78	0.81
2:B:71:VAL:HG23	2:B:164:VAL:HA	1.62	0.81
5:E:8:GLU:HA	5:E:34:VAL:HA	1.60	0.81
7:G:17:VAL:HG12	7:G:18:TYR:H	1.45	0.81
13:M:77:ASN:O	13:M:80:ARG:HB3	1.79	0.81
19:S:24:ALA:HB3	19:S:25:LYS:NZ	1.95	0.81
19:S:53:ASN:HD21	19:S:77:THR:HA	1.43	0.81
20:T:51:GLU:HG2	20:T:52:ALA:H	1.43	0.81
21:V:10:ARG:HA	21:V:13:ILE:CG2	2.10	0.81
1:A:1118:C:H1'	1:A:1179:A:C4	2.16	0.81
2:B:231:GLU:HB2	2:B:232:PRO:CD	2.10	0.81
4:D:156:GLU:CG	4:D:157:LEU:H	1.92	0.81
10:J:77:PRO:HA	10:J:81:THR:CG2	2.09	0.81
1:A:1174:G:H2'	1:A:1175:G:H8	1.45	0.81
3:C:19:GLU:H	14:N:51:GLY:HA3	1.46	0.81
12:L:113:ARG:NH1	12:L:115:LYS:HB2	1.95	0.81
14:N:41:ARG:HG2	14:N:41:ARG:HH11	1.46	0.81
1:A:476:G:H2'	1:A:477:G:C8	2.16	0.81
2:B:73:THR:HG23	2:B:96:ARG:NH2	1.96	0.81
3:C:120:VAL:O	3:C:124:ILE:HB	1.80	0.81
5:E:70:PRO:O	5:E:72:GLN:N	2.13	0.81
6:F:1:MET:SD	6:F:66:GLU:HG2	2.20	0.81
10:J:38:ILE:HD12	10:J:71:LEU:HD12	1.61	0.81
9:I:19:LEU:HG	9:I:61:ALA:HB2	1.61	0.81
15:O:64:ARG:HH22	15:O:68:ARG:NH2	1.77	0.81
1:A:107:G:H2'	1:A:108:G:H5'	1.60	0.81
1:A:1065:U:H4'	1:A:1066:C:C5'	2.11	0.81
7:G:71:PRO:HD3	7:G:103:TRP:HZ3	1.46	0.81
8:H:11:THR:HG23	8:H:14:ARG:HH12	1.46	0.81
4:D:9:CYS:SG	4:D:22:LYS:NZ	2.52	0.80
10:J:6:ILE:CG1	10:J:73:ASP:HA	2.10	0.80
12:L:73:GLU:HB2	12:L:110:VAL:HG11	1.62	0.80
1:A:56:U:H2'	1:A:57:G:C8	2.16	0.80
7:G:137:LYS:HA	7:G:140:ASP:HB2	1.63	0.80
12:L:102:ARG:HA	12:L:107:ALA:HB1	1.63	0.80
1:A:129(A):G:HO2'	1:A:190(E):U:H2'	1.46	0.80
3:C:7:PRO:HA	3:C:10:PHE:HB3	1.62	0.80
3:C:196:LEU:N	3:C:196:LEU:HD23	1.97	0.80
1:A:22:G:H2'	1:A:23:C:C6	2.16	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:610:G:H2'	1:A:611:A:C8	2.16	0.80
2:B:43:ASP:O	2:B:46:LYS:HB3	1.81	0.80
3:C:50:ALA:HA	3:C:72:LYS:CD	2.11	0.80
3:C:153:VAL:HA	3:C:198:VAL:HG22	1.62	0.80
5:E:8:GLU:CB	5:E:34:VAL:HG22	2.11	0.80
8:H:89:PRO:O	8:H:91:ARG:N	2.15	0.80
10:J:86:MET:HG3	10:J:87:THR:N	1.95	0.80
19:S:67:VAL:O	19:S:69:HIS:N	2.13	0.80
1:A:1319:A:H61	1:A:1361:G:H21	1.30	0.80
10:J:31:GLY:HA2	10:J:78:ASN:HB2	1.64	0.80
14:N:56:VAL:O	14:N:57:ARG:HB2	1.81	0.80
1:A:436:C:H2'	1:A:437:U:C6	2.16	0.80
8:H:64:LYS:HG3	8:H:79:VAL:HG21	1.62	0.80
9:I:24:GLY:HA2	9:I:60:ASP:HA	1.62	0.80
13:M:78:ILE:H	13:M:78:ILE:HD12	1.47	0.80
1:A:840:C:H5''	1:A:841:U:OP1	1.81	0.80
1:A:1205:U:H4'	3:C:195:VAL:CG2	2.12	0.80
3:C:62:ASP:CA	3:C:97:LYS:HB3	2.12	0.80
3:C:112:SER:O	3:C:115:LEU:HB2	1.81	0.80
6:F:94:GLN:HB3	18:R:32:ARG:NH1	1.96	0.80
18:R:22:VAL:HG12	18:R:23:LYS:N	1.97	0.80
1:A:128:G:H4'	17:Q:3:LYS:HG2	1.63	0.80
1:A:167:G:H2'	1:A:168:G:H8	1.42	0.80
1:A:351:G:H4'	1:A:352:C:OP1	1.82	0.80
1:A:444:C:H2'	1:A:445:G:C8	2.17	0.80
1:A:1130:A:H5''	9:I:20:ARG:HD3	1.62	0.80
1:A:1392:G:O2'	1:A:1393:U:H5'	1.82	0.80
5:E:18:ARG:HH21	5:E:25:ARG:HB3	1.47	0.80
13:M:13:LYS:O	13:M:18:ALA:HB2	1.81	0.80
20:T:30:LYS:NZ	20:T:80:ARG:HH22	1.81	0.80
4:D:25:ARG:HG3	4:D:30:LYS:HB3	1.64	0.79
7:G:15:ASP:OD1	7:G:17:VAL:HB	1.82	0.79
18:R:40:LEU:CD2	18:R:79:LEU:HD21	2.12	0.79
1:A:250:A:H4'	1:A:251:G:O5'	1.81	0.79
1:A:408:A:H5'	4:D:116:GLN:HB2	1.65	0.79
3:C:181:ASN:C	3:C:182:ILE:HD12	2.02	0.79
1:A:182:U:H6	1:A:182:U:H5'	1.46	0.79
1:A:454:C:H5	1:A:478:A:H61	1.30	0.79
1:A:456:C:H2'	1:A:457:C:C6	2.14	0.79
1:A:1261:A:H62	1:A:1274:G:H21	1.28	0.79
2:B:84:GLU:OE1	2:B:216:SER:HA	1.81	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:175:ARG:NH1	2:B:175:ARG:HB2	1.97	0.79
11:K:92:GLU:HA	11:K:95:ILE:HD12	1.65	0.79
13:M:56:LEU:O	13:M:60:VAL:HG23	1.83	0.79
18:R:78:LEU:C	18:R:79:LEU:HD12	2.02	0.79
19:S:49:ILE:HG22	19:S:60:VAL:HB	1.63	0.79
1:A:1090:U:H2'	1:A:1091:U:C6	2.16	0.79
13:M:75:ALA:O	13:M:79:LYS:HB2	1.83	0.79
14:N:12:ARG:O	14:N:14:PRO:HD3	1.81	0.79
19:S:15:LEU:HD13	19:S:44:MET:HE1	1.63	0.79
2:B:111:ARG:HB3	2:B:149:LEU:HD11	1.65	0.79
3:C:32:LEU:O	3:C:36:ASP:HB2	1.82	0.79
12:L:57:LYS:HA	12:L:66:VAL:O	1.83	0.79
1:A:1053:G:HO2'	1:A:1199:U:H5	1.30	0.79
5:E:99:GLY:O	5:E:101:ILE:HG13	1.82	0.79
12:L:70:ILE:HG12	12:L:100:ILE:HD12	1.65	0.79
19:S:11:VAL:HB	19:S:16:LEU:HD22	1.65	0.79
1:A:1244:C:H2'	1:A:1245:A:C8	2.18	0.79
7:G:66:VAL:HG13	7:G:100:ALA:HB1	1.65	0.79
9:I:7:THR:HG22	9:I:8:GLY:N	1.95	0.79
1:A:793:U:H3'	1:A:794:A:H5''	1.65	0.79
1:A:1225:A:H5'	13:M:103:THR:HG23	1.65	0.79
3:C:113:ALA:HB3	3:C:114:PRO:HD3	1.64	0.79
1:A:174:C:H2'	1:A:175:C:H6	1.48	0.78
2:B:97:TRP:CH2	2:B:101:MET:HB2	2.18	0.78
10:J:27:ALA:HA	10:J:30:SER:OG	1.83	0.78
1:A:962:C:H2'	1:A:963:G:C8	2.16	0.78
1:A:1065:U:H1'	1:A:1066:C:OP2	1.83	0.78
4:D:121:VAL:O	4:D:134:ASP:HB2	1.82	0.78
7:G:70:LYS:HE2	7:G:100:ALA:HB2	1.65	0.78
1:A:1333:A:H2'	1:A:1334:G:O4'	1.83	0.78
16:P:22:THR:HG23	16:P:24:ALA:H	1.47	0.78
3:C:27:LYS:HA	3:C:30:ARG:HH12	1.46	0.78
3:C:116:VAL:HA	3:C:119:ARG:HB3	1.66	0.78
8:H:97:VAL:HA	8:H:100:ILE:HD13	1.64	0.78
10:J:26:ALA:HB3	10:J:85:LEU:HG	1.65	0.78
1:A:736:C:H2'	1:A:737:A:H8	1.46	0.78
4:D:52:SER:O	4:D:54:TYR:N	2.16	0.78
7:G:92:SER:O	7:G:96:GLN:HB2	1.83	0.78
10:J:83:GLU:HA	10:J:86:MET:HB2	1.65	0.78
20:T:13:LEU:HD12	20:T:14:LYS:N	1.97	0.78
1:A:237:C:H2'	1:A:238:G:C8	2.19	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1190:G:OP2	3:C:5:ILE:HG23	1.84	0.78
2:B:9:GLU:CD	2:B:217:ARG:HH22	1.87	0.78
2:B:72:GLY:HA3	2:B:81:VAL:HG21	1.66	0.78
6:F:18:GLN:HA	6:F:21:LEU:HD23	1.65	0.78
7:G:39:ALA:HA	7:G:42:ILE:HD12	1.66	0.78
1:A:60:A:H4'	1:A:61:G:O5'	1.83	0.78
1:A:532:A:H2'	1:A:533:A:H5'	1.66	0.78
14:N:46:GLU:HA	14:N:49:HIS:CD2	2.17	0.78
1:A:219:C:H2'	1:A:220:G:H5'	1.66	0.78
2:B:127:ILE:H	2:B:127:ILE:HD12	1.48	0.78
7:G:38:LEU:O	7:G:42:ILE:HG13	1.84	0.78
3:C:26:LYS:NZ	10:J:45:ARG:HE	1.82	0.78
9:I:5:TYR:HE2	9:I:16:ARG:HB2	1.48	0.78
10:J:9:ARG:HB3	10:J:9:ARG:NH1	1.98	0.78
1:A:1228:C:H2'	1:A:1229:A:H8	1.47	0.78
4:D:17:VAL:HG12	4:D:18:LYS:N	1.92	0.78
10:J:34:VAL:HA	10:J:75:ILE:H	1.49	0.78
1:A:517:G:H1'	1:A:519:C:H42	1.48	0.77
1:A:708:C:H2'	1:A:709:G:H8	1.47	0.77
6:F:15:ASP:OD1	6:F:18:GLN:HG3	1.84	0.77
1:A:942:G:N2	1:A:943:U:C2	2.51	0.77
2:B:97:TRP:CZ2	2:B:102:LEU:HD13	2.20	0.77
9:I:108:VAL:HG12	9:I:109:VAL:N	1.98	0.77
1:A:1509:C:O2'	1:A:1510:U:H5'	1.83	0.77
3:C:138:VAL:O	3:C:142:MET:HB2	1.83	0.77
1:A:511:C:O2'	1:A:512:U:H5''	1.84	0.77
1:A:982:U:H5	14:N:31:ARG:NH1	1.82	0.77
1:A:1319:A:H61	1:A:1361:G:N2	1.82	0.77
7:G:93:PRO:HG2	7:G:94:ARG:N	1.98	0.77
1:A:707:C:H2'	1:A:708:C:C6	2.19	0.77
1:A:1014:A:H5'	19:S:14:HIS:HB3	1.67	0.77
4:D:25:ARG:C	4:D:27:TYR:H	1.86	0.77
13:M:81:LEU:O	13:M:89:GLY:HA3	1.85	0.77
1:A:92:C:H2'	1:A:93:G:C8	2.19	0.77
1:A:781:A:H2'	1:A:782:A:H5'	1.67	0.77
12:L:83:VAL:HG21	12:L:100:ILE:HD13	1.67	0.77
12:L:89:ARG:O	12:L:90:VAL:HG23	1.84	0.77
16:P:39:TYR:HE2	16:P:41:PRO:HG3	1.48	0.77
1:A:291:C:O2'	1:A:292:G:H5'	1.84	0.77
2:B:223:ILE:HD13	2:B:230:VAL:HG21	1.65	0.77
15:O:45:VAL:HB	15:O:46:HIS:ND1	1.99	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:R:38:GLU:H	18:R:41:LYS:HE3	1.48	0.77
1:A:608:A:H2'	1:A:609:A:H8	1.50	0.77
1:A:1197:G:OP1	1:A:1197:G:H3'	1.84	0.77
2:B:168:THR:OG1	2:B:192:SER:HA	1.85	0.77
6:F:97:PHE:HD1	18:R:65:ILE:HD12	1.49	0.77
12:L:41:ARG:HH12	12:L:43:VAL:HG22	1.50	0.77
1:A:559:A:H4'	1:A:560:U:O5'	1.85	0.77
3:C:7:PRO:O	3:C:11:ARG:N	2.17	0.77
4:D:17:VAL:CG1	4:D:18:LYS:H	1.94	0.77
12:L:92:ASP:O	12:L:94:PRO:HD3	1.85	0.77
1:A:314:C:O2'	1:A:315:A:H5'	1.85	0.76
1:A:376:G:H5''	16:P:5:ARG:CG	2.15	0.76
1:A:518:C:H4'	1:A:519:C:C6	2.20	0.76
2:B:44:LEU:H	2:B:44:LEU:CD2	1.96	0.76
3:C:62:ASP:HA	3:C:97:LYS:CB	2.14	0.76
14:N:6:LEU:HB3	14:N:23:ARG:HE	1.49	0.76
1:A:99:C:H2'	1:A:101:A:C8	2.20	0.76
1:A:1095:U:OP1	1:A:1108:G:N2	2.19	0.76
2:B:97:TRP:HZ2	2:B:102:LEU:HD13	1.50	0.76
4:D:114:ARG:HA	4:D:117:ALA:HB3	1.66	0.76
1:A:186:C:H2'	1:A:187:C:C6	2.20	0.76
1:A:615:C:H2'	1:A:616:G:H5'	1.68	0.76
6:F:79:LEU:O	6:F:85:VAL:HG11	1.85	0.76
7:G:137:LYS:CA	7:G:140:ASP:HB2	2.16	0.76
8:H:89:PRO:C	8:H:91:ARG:H	1.89	0.76
20:T:80:ARG:HB3	20:T:80:ARG:NH1	2.00	0.76
1:A:222:U:H2'	1:A:223:U:C6	2.21	0.76
1:A:1238:A:N7	1:A:1303:C:H1'	1.99	0.76
2:B:21:ARG:HG2	2:B:22:LYS:N	1.99	0.76
2:B:210:SER:O	2:B:212:GLN:N	2.19	0.76
3:C:130:VAL:O	3:C:134:ILE:HG13	1.86	0.76
6:F:12:PRO:HG3	6:F:57:GLN:HG3	1.65	0.76
12:L:55:VAL:HG12	12:L:56:ALA:H	1.50	0.76
16:P:64:ALA:O	16:P:66:PRO:HD3	1.84	0.76
1:A:186:C:H2'	1:A:187:C:H6	1.49	0.76
3:C:55:VAL:HA	3:C:68:VAL:HG22	1.65	0.76
12:L:43:VAL:HG12	12:L:44:THR:N	2.00	0.76
18:R:22:VAL:HG12	18:R:23:LYS:H	1.50	0.76
1:A:922:G:H5'	5:E:19:MET:O	1.85	0.76
1:A:1020:U:H2'	1:A:1021:G:C8	2.21	0.76
1:A:1046:A:H3'	1:A:1047:G:H8	1.51	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:52:LEU:H	3:C:52:LEU:CD2	1.99	0.76
5:E:70:PRO:C	5:E:72:GLN:H	1.88	0.76
11:K:58:PRO:HB2	11:K:93:GLN:HG3	1.67	0.76
12:L:89:ARG:NH2	12:L:97:ARG:HG2	2.00	0.76
1:A:129(A):G:O2'	1:A:190(E):U:H2'	1.85	0.76
1:A:615:C:C2'	1:A:616:G:H5'	2.16	0.76
1:A:973:G:H3'	1:A:974:A:H5''	1.68	0.76
3:C:59:ARG:HG3	3:C:60:ALA:N	2.01	0.76
10:J:9:ARG:NE	10:J:9:ARG:H	1.83	0.76
12:L:11:VAL:HG13	17:Q:29:HIS:HD2	1.49	0.76
17:Q:10:VAL:HG23	17:Q:55:ASP:O	1.84	0.76
1:A:710:G:OP1	6:F:54:LYS:HD2	1.86	0.76
1:A:1109:C:O2'	1:A:1110:A:H5'	1.86	0.76
1:A:1406:U:H2'	1:A:1407:C:C6	2.21	0.76
19:S:11:VAL:HG11	19:S:16:LEU:HD13	1.67	0.76
1:A:628:G:O2'	1:A:629:G:H5'	1.86	0.76
2:B:69:LEU:HD13	2:B:155:LEU:HD11	1.66	0.76
3:C:47:LEU:HD11	3:C:76:VAL:HG12	1.68	0.76
5:E:55:VAL:HG23	5:E:56:GLN:N	2.01	0.76
10:J:44:VAL:HG12	10:J:45:ARG:H	1.50	0.76
21:V:10:ARG:H	21:V:10:ARG:HD2	1.48	0.76
1:A:265:G:O2'	1:A:266:G:H5'	1.86	0.76
1:A:376:G:O2'	1:A:377:G:H5'	1.86	0.76
3:C:156:ARG:NH2	3:C:161:GLU:HA	1.99	0.76
8:H:63:LEU:H	8:H:63:LEU:CD1	1.98	0.76
19:S:58:VAL:HG12	19:S:59:PRO:HD2	1.66	0.76
1:A:190(H):G:O2'	1:A:190(I):G:H5'	1.85	0.75
1:A:237:C:H2'	1:A:238:G:H8	1.49	0.75
1:A:1343:G:H2'	1:A:1344:C:H6	1.48	0.75
3:C:160:ALA:C	3:C:162:GLN:H	1.88	0.75
9:I:49:PRO:HG2	9:I:50:LEU:HD12	1.68	0.75
14:N:7:ILE:O	14:N:7:ILE:HG22	1.86	0.75
1:A:1150:U:H4'	10:J:41:PRO:HD3	1.66	0.75
1:A:1163:C:H2'	1:A:1164:G:H8	1.50	0.75
3:C:179:ARG:HD2	3:C:180:ALA:N	2.00	0.75
15:O:48:LYS:N	15:O:48:LYS:HD3	2.01	0.75
17:Q:80:GLY:O	17:Q:81:ARG:HB3	1.86	0.75
20:T:60:GLU:HA	20:T:63:ILE:HD12	1.68	0.75
1:A:1326:C:OP1	21:V:12:LYS:HD2	1.86	0.75
4:D:59:ARG:NH2	4:D:62:GLN:HG3	2.01	0.75
11:K:95:ILE:HG22	11:K:99:GLN:NE2	2.00	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:L:7:ILE:O	12:L:10:LEU:N	2.18	0.75
19:S:24:ALA:HB3	19:S:25:LYS:HZ3	1.51	0.75
20:T:15:ARG:O	20:T:19:SER:N	2.18	0.75
1:A:790:A:H2'	1:A:791:G:C8	2.20	0.75
4:D:127:THR:HG22	4:D:147:ALA:O	1.86	0.75
18:R:70:ILE:CG2	18:R:74:ARG:HD2	2.16	0.75
21:V:2:GLY:O	21:V:3:LYS:HG2	1.87	0.75
1:A:597:G:H2'	1:A:598:U:H5'	1.68	0.75
2:B:218:ALA:O	2:B:222:ILE:HG13	1.87	0.75
4:D:201:GLN:HA	4:D:204:ILE:HD12	1.68	0.75
6:F:35:ALA:HB1	6:F:65:VAL:HG21	1.68	0.75
9:I:3:GLN:HE21	9:I:3:GLN:C	1.90	0.75
16:P:21:VAL:O	16:P:21:VAL:HG12	1.87	0.75
17:Q:64:PRO:C	17:Q:65:ILE:HD12	2.07	0.75
1:A:797:C:O2'	1:A:798:G:H5'	1.85	0.75
2:B:74:LYS:HE3	2:B:205:ASP:O	1.87	0.75
8:H:97:VAL:HG22	8:H:98:LYS:N	2.02	0.75
9:I:111:ARG:HH11	9:I:111:ARG:HG3	1.51	0.75
10:J:22:LYS:NZ	10:J:22:LYS:HB2	2.01	0.75
1:A:95:U:H2'	1:A:96:G:C8	2.22	0.75
1:A:144:G:H1	1:A:178:C:H42	1.35	0.75
1:A:514:C:H42	1:A:537:G:H1	1.34	0.75
1:A:1241:G:H2'	1:A:1242:C:C6	2.22	0.75
1:A:1508:G:H2'	1:A:1509:C:H6	1.50	0.75
3:C:88:ARG:NH1	3:C:101:LEU:HB3	2.02	0.75
9:I:85:LEU:O	9:I:89:ASN:HB3	1.87	0.75
14:N:26:ARG:HG3	14:N:27:CYS:H	1.50	0.75
20:T:54:LYS:HA	20:T:57:ARG:HD2	1.69	0.75
20:T:88:VAL:O	20:T:91:LEU:HB2	1.86	0.75
1:A:144:G:H1	1:A:178:C:N4	1.85	0.75
1:A:761:G:C5'	17:Q:102:GLY:HA3	2.17	0.75
8:H:85:ARG:HH11	8:H:85:ARG:HG3	1.52	0.75
1:A:1001:A:C2'	1:A:1002:G:H5'	2.17	0.75
1:A:190(E):U:C2	17:Q:63:ARG:HD3	2.22	0.74
1:A:1369:C:H2'	1:A:1370:G:O4'	1.86	0.74
3:C:90:GLU:HA	3:C:93:LYS:HG3	1.68	0.74
1:A:762:C:H2'	1:A:763:G:H8	1.52	0.74
1:A:1226:C:H5''	13:M:103:THR:OG1	1.87	0.74
3:C:20:SER:C	3:C:21:ARG:HD2	2.08	0.74
3:C:116:VAL:O	3:C:119:ARG:HB3	1.86	0.74
4:D:64:LEU:HB2	4:D:198:VAL:CG2	2.15	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:P:51:VAL:HG12	16:P:52:ASP:N	2.01	0.74
1:A:154:C:O2'	1:A:155:C:H5'	1.88	0.74
1:A:451:A:N6	1:A:481:G:C4	2.55	0.74
1:A:517:G:H1'	1:A:519:C:N4	2.01	0.74
1:A:975:A:O5'	1:A:976:G:H5'	1.87	0.74
1:A:1142:G:H3'	1:A:1143:G:H8	1.52	0.74
3:C:181:ASN:O	3:C:182:ILE:HD12	1.87	0.74
6:F:9:VAL:HB	6:F:87:ARG:HB2	1.68	0.74
7:G:71:PRO:O	7:G:96:GLN:HG3	1.87	0.74
16:P:75:ARG:HA	16:P:80:PHE:CE1	2.23	0.74
1:A:309:G:H2'	1:A:310:G:C8	2.21	0.74
1:A:428:G:H5'	4:D:7:PRO:HB3	1.68	0.74
1:A:709:G:H2'	1:A:710:G:H8	1.51	0.74
1:A:1356:G:H2'	1:A:1357:A:C8	2.22	0.74
4:D:163:GLU:O	4:D:166:LYS:HE2	1.87	0.74
15:O:12:ILE:HD12	15:O:12:ILE:N	2.00	0.74
18:R:31:LEU:HD23	18:R:31:LEU:O	1.86	0.74
1:A:579:G:H2'	1:A:580:U:C6	2.22	0.74
1:A:579:G:H2'	1:A:580:U:H6	1.50	0.74
16:P:21:VAL:HG11	16:P:59:TRP:HE1	1.51	0.74
3:C:180:ALA:HB1	3:C:203:PHE:CE1	2.22	0.74
5:E:139:LEU:HD23	5:E:142:LEU:HD11	1.70	0.74
7:G:12:LEU:HD12	7:G:12:LEU:N	2.02	0.74
13:M:49:THR:HG22	13:M:50:GLU:N	2.03	0.74
13:M:81:LEU:HA	13:M:84:ILE:HG12	1.67	0.74
18:R:39:VAL:HG13	18:R:40:LEU:N	2.00	0.74
20:T:51:GLU:O	20:T:54:LYS:HB3	1.87	0.74
1:A:695:A:H2	1:A:787:A:H1'	1.52	0.74
4:D:64:LEU:CB	4:D:198:VAL:HG21	2.15	0.74
5:E:99:GLY:N	5:E:117:ASP:OD1	2.21	0.74
1:A:707:C:H5''	11:K:20:TYR:HD2	1.52	0.74
1:A:1210:C:H2'	1:A:1211:U:C5'	2.17	0.74
2:B:68:ILE:H	2:B:90:MET:HE3	1.53	0.74
5:E:111:GLU:O	5:E:113:ALA:N	2.21	0.74
9:I:34:ASN:HD22	9:I:34:ASN:N	1.85	0.74
9:I:86:VAL:HA	9:I:89:ASN:O	1.88	0.74
1:A:503:C:H2'	1:A:504:C:C6	2.22	0.74
1:A:538:G:H2'	1:A:539:A:H8	1.50	0.74
1:A:545:C:O2'	1:A:546:G:H5'	1.86	0.74
4:D:161:ASN:O	4:D:165:MET:HG2	1.87	0.74
4:D:200:GLU:O	4:D:204:ILE:HG13	1.88	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:M:20:THR:HA	13:M:25:ILE:HG22	1.68	0.74
13:M:37:THR:HG22	13:M:39:ILE:HG13	1.67	0.74
1:A:89:C:H2'	1:A:90:U:H6	1.53	0.74
1:A:107:G:C2'	1:A:108:G:H5'	2.17	0.74
1:A:112:G:H4'	1:A:389:A:H5''	1.68	0.74
1:A:180:U:C2'	1:A:181:G:H5'	2.18	0.74
1:A:1001:A:H2'	1:A:1002:G:H5'	1.69	0.74
3:C:112:SER:CB	3:C:115:LEU:HD12	2.16	0.74
9:I:48:GLU:N	9:I:49:PRO:HD2	2.02	0.74
1:A:137:C:H42	1:A:226:G:H1	1.35	0.73
1:A:1367:C:H5''	10:J:60:ARG:HH12	1.52	0.73
1:A:1424:C:O2'	1:A:1425:U:H5'	1.88	0.73
1:A:1508:G:H2'	1:A:1509:C:C6	2.23	0.73
2:B:74:LYS:CE	2:B:166:ASP:HB2	2.17	0.73
3:C:183:ASP:OD2	3:C:184:TYR:N	2.21	0.73
10:J:70:ARG:HB3	10:J:70:ARG:HH11	1.52	0.73
17:Q:9:VAL:HG12	17:Q:10:VAL:N	2.02	0.73
1:A:29:G:H5'	1:A:296:U:OP1	1.88	0.73
1:A:1104:G:H4'	2:B:111:ARG:CZ	2.17	0.73
3:C:22:TRP:HH2	3:C:33:LEU:HB2	1.53	0.73
10:J:80:LYS:O	10:J:84:GLN:HB2	1.87	0.73
10:J:86:MET:C	10:J:88:LEU:H	1.90	0.73
17:Q:48:GLU:HG3	17:Q:50:LYS:HB2	1.69	0.73
1:A:491:G:H2'	1:A:492:G:C8	2.23	0.73
2:B:8:LYS:CD	2:B:9:GLU:H	2.01	0.73
20:T:53:LEU:HD21	20:T:101:GLY:O	1.89	0.73
1:A:448:A:H2'	1:A:449:C:C6	2.24	0.73
1:A:1193:G:O2'	1:A:1194:U:H5'	1.88	0.73
2:B:163:PHE:HA	2:B:185:ILE:O	1.89	0.73
3:C:93:LYS:HE2	3:C:93:LYS:HA	1.68	0.73
5:E:9:LYS:O	5:E:33:VAL:HG23	1.88	0.73
5:E:12:LEU:O	5:E:12:LEU:HD13	1.89	0.73
7:G:93:PRO:CG	7:G:94:ARG:H	2.00	0.73
9:I:21:PRO:HA	9:I:59:PHE:HA	1.68	0.73
9:I:26:VAL:HG13	9:I:63:ILE:HB	1.70	0.73
1:A:666:G:H2'	1:A:667:G:H8	1.52	0.73
2:B:208:ILE:HG22	2:B:209:ARG:N	2.00	0.73
3:C:179:ARG:O	3:C:181:ASN:N	2.21	0.73
4:D:13:ARG:HD2	4:D:38:TYR:O	1.88	0.73
8:H:38:ILE:O	8:H:39:LEU:C	2.26	0.73
9:I:42:ARG:NH2	9:I:75:ASP:OD2	2.21	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:16:LEU:HG	10:J:94:VAL:HG11	1.70	0.73
11:K:58:PRO:HB3	11:K:93:GLN:HG3	1.69	0.73
19:S:72:GLY:C	19:S:74:PHE:N	2.40	0.73
1:A:528:C:C2'	1:A:529:G:H5'	2.17	0.73
1:A:673:G:H2'	1:A:674:G:C8	2.23	0.73
1:A:785:G:O2'	1:A:786:G:H5'	1.88	0.73
1:A:1306:A:C2	1:A:1307:U:H1'	2.23	0.73
3:C:59:ARG:HG3	3:C:61:ALA:H	1.53	0.73
3:C:126:ARG:HB3	3:C:128:PHE:HB3	1.71	0.73
7:G:61:VAL:O	7:G:65:ALA:HB3	1.88	0.73
13:M:91:ARG:HB2	13:M:98:VAL:HG22	1.70	0.73
19:S:14:HIS:O	19:S:18:LYS:HG2	1.87	0.73
1:A:375:U:OP1	16:P:69:THR:HG21	1.89	0.73
1:A:476:G:H2'	1:A:477:G:H8	1.49	0.73
1:A:761:G:H4'	17:Q:102:GLY:HA3	1.71	0.73
1:A:1163:C:H2'	1:A:1164:G:C8	2.24	0.73
1:A:1361:G:H8	1:A:1361:G:O5'	1.72	0.73
1:A:1367:C:C5'	10:J:60:ARG:HH12	2.02	0.73
4:D:156:GLU:OE2	4:D:157:LEU:HD22	1.89	0.73
4:D:194:LEU:HD22	4:D:194:LEU:N	2.03	0.73
10:J:30:SER:HB3	10:J:84:GLN:CD	2.08	0.73
16:P:50:LYS:O	16:P:51:VAL:HG23	1.89	0.73
19:S:23:ASN:HA	19:S:27:GLU:HA	1.69	0.73
11:K:87:THR:CA	11:K:91:ARG:HH12	2.00	0.73
13:M:20:THR:O	13:M:20:THR:HG22	1.88	0.73
18:R:43:PHE:C	18:R:44:LEU:HD22	2.09	0.73
1:A:21:G:H2'	1:A:22:G:C8	2.24	0.73
1:A:160:A:H1'	1:A:344:A:N7	2.03	0.73
1:A:538:G:H4'	12:L:114:LYS:HD3	1.69	0.73
1:A:779:C:O2'	1:A:780:A:H5'	1.89	0.73
1:A:1200:C:H2'	1:A:1200:C:O2	1.89	0.73
1:A:1440:C:H2'	1:A:1441:G:H5'	1.71	0.73
3:C:88:ARG:CA	3:C:91:LEU:HB3	2.17	0.73
3:C:129:ALA:HB1	3:C:132:ARG:CB	2.18	0.73
1:A:130:A:H5'	17:Q:63:ARG:HH21	1.54	0.73
2:B:207:ALA:HB3	2:B:210:SER:HB3	1.70	0.73
3:C:83:ARG:C	3:C:85:ARG:H	1.92	0.73
7:G:23:VAL:O	7:G:27:ILE:HG13	1.88	0.73
9:I:112:LYS:HA	9:I:119:ALA:HA	1.71	0.72
13:M:3:ARG:HA	13:M:8:GLU:O	1.89	0.72
20:T:46:GLU:CB	20:T:48:LYS:HE2	2.18	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:694:A:H3'	1:A:695:A:H5''	1.70	0.72
1:A:913:A:H1'	1:A:914:A:O4'	1.89	0.72
1:A:1418:A:H2'	1:A:1419:G:O4'	1.89	0.72
2:B:21:ARG:HE	2:B:21:ARG:H	1.33	0.72
2:B:96:ARG:NE	2:B:96:ARG:N	2.38	0.72
2:B:101:MET:HA	2:B:108:ILE:CD1	2.17	0.72
3:C:67:THR:HA	3:C:102:ASN:HD21	1.52	0.72
4:D:7:PRO:O	4:D:10:ARG:HB3	1.87	0.72
5:E:11:ILE:HB	5:E:12:LEU:HD12	1.69	0.72
5:E:13:ILE:HG22	5:E:30:ALA:HA	1.69	0.72
8:H:82:HIS:HD2	8:H:83:ILE:H	1.37	0.72
11:K:18:ARG:HB2	11:K:33:THR:HG23	1.70	0.72
12:L:41:ARG:HD2	12:L:42:THR:H	1.54	0.72
15:O:12:ILE:H	15:O:12:ILE:CD1	1.97	0.72
21:V:6:ARG:HD2	21:V:15:ARG:HH22	1.53	0.72
1:A:1372:U:H5''	9:I:71:SER:CB	2.15	0.72
1:A:419:C:H5	1:A:425:G:N1	1.87	0.72
1:A:463:A:H2'	1:A:474:G:C8	2.18	0.72
7:G:45:ASP:C	7:G:47:CYS:H	1.92	0.72
9:I:50:LEU:HD12	9:I:50:LEU:H	1.52	0.72
19:S:64:GLU:OE1	19:S:67:VAL:HG21	1.90	0.72
1:A:735:C:H2'	1:A:736:C:C6	2.22	0.72
1:A:1127:G:H1'	1:A:1148:U:N3	2.02	0.72
1:A:1180:A:O2'	1:A:1181:G:H5'	1.89	0.72
1:A:1276:G:C2'	1:A:1277:C:H5'	2.20	0.72
5:E:126:ARG:NH1	5:E:126:ARG:HG2	2.04	0.72
7:G:101:LEU:HD23	7:G:101:LEU:N	2.04	0.72
9:I:19:LEU:HA	9:I:61:ALA:HA	1.71	0.72
13:M:79:LYS:HE2	13:M:79:LYS:CA	2.19	0.72
15:O:16:ALA:CB	15:O:21:ASP:HB3	2.17	0.72
1:A:273:A:C2'	1:A:274:A:H5'	2.19	0.72
1:A:1096:C:O2'	1:A:1097:C:H5'	1.88	0.72
1:A:1250:A:H4'	9:I:68:GLY:CA	2.13	0.72
3:C:85:ARG:CA	3:C:88:ARG:HG3	2.18	0.72
6:F:16:GLN:HA	6:F:19:LEU:HB3	1.70	0.72
6:F:44:GLY:HA2	6:F:59:TYR:CE1	2.24	0.72
9:I:42:ARG:O	9:I:74:ILE:HG21	1.90	0.72
9:I:51:ARG:HG2	9:I:56:LEU:HD12	1.70	0.72
1:A:279:A:OP2	17:Q:95:TYR:OH	2.07	0.72
1:A:397:A:C2'	1:A:398:C:H5''	2.20	0.72
1:A:858:G:H8	1:A:858:G:O5'	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1370:G:H2'	1:A:1371:G:H8	1.55	0.72
4:D:52:SER:C	4:D:54:TYR:N	2.42	0.72
10:J:40:LEU:HD13	10:J:69:ASN:HB2	1.70	0.72
12:L:26:ALA:HB2	12:L:98:TYR:HE2	1.55	0.72
19:S:40:ILE:HD13	19:S:62:ILE:HD11	1.72	0.72
1:A:327:A:O2'	1:A:328:C:O4'	2.07	0.72
1:A:861:G:H2'	1:A:862:C:H6	1.55	0.72
3:C:47:LEU:HD11	3:C:76:VAL:CG1	2.20	0.72
14:N:50:LYS:HG3	14:N:51:GLY:H	1.54	0.72
16:P:43:LYS:HB3	16:P:48:TRP:CG	2.25	0.72
16:P:67:THR:HG22	16:P:69:THR:N	2.03	0.72
1:A:17:U:H2'	1:A:18:C:C6	2.24	0.71
1:A:399:G:O2'	1:A:400:C:H5'	1.89	0.71
1:A:943:U:O2'	1:A:944:G:H5'	1.90	0.71
1:A:1088:G:H2'	1:A:1089:G:H8	1.55	0.71
1:A:1263:C:H2'	1:A:1264:C:O4'	1.90	0.71
2:B:151:GLY:C	2:B:153:ARG:H	1.93	0.71
2:B:165:VAL:O	2:B:167:PRO:HD3	1.89	0.71
3:C:11:ARG:HA	3:C:14:ILE:CG1	2.20	0.71
3:C:34:LEU:O	3:C:38:ARG:HG2	1.90	0.71
4:D:101:LEU:HB2	4:D:138:TYR:HB3	1.71	0.71
6:F:18:GLN:O	6:F:21:LEU:HB3	1.89	0.71
9:I:11:LYS:O	9:I:12:GLU:HB3	1.90	0.71
9:I:21:PRO:HA	9:I:60:ASP:N	2.03	0.71
19:S:15:LEU:HD12	19:S:16:LEU:N	2.05	0.71
1:A:178:C:O2'	1:A:179:A:H5'	1.90	0.71
1:A:1240:U:O4	7:G:30:ILE:HG23	1.89	0.71
17:Q:70:ARG:HH11	17:Q:70:ARG:CG	1.99	0.71
17:Q:88:TYR:HA	17:Q:91:ARG:HE	1.56	0.71
1:A:287:U:O2'	1:A:288:A:H5'	1.90	0.71
1:A:560:U:H4'	1:A:561:U:C5'	2.19	0.71
1:A:1436:U:O2'	1:A:1437:C:H5'	1.91	0.71
3:C:8:ILE:HG21	14:N:50:LYS:HB3	1.70	0.71
3:C:52:LEU:HD23	3:C:52:LEU:N	2.05	0.71
16:P:82:GLN:H	16:P:82:GLN:HE21	1.35	0.71
20:T:15:ARG:HA	20:T:18:GLN:HB2	1.71	0.71
1:A:73:C:H2'	1:A:74:C:C6	2.25	0.71
1:A:518:C:C5'	1:A:530:G:H1'	2.20	0.71
1:A:1266:G:N2	1:A:1268:A:H3'	2.05	0.71
1:A:423:G:H2'	1:A:424:G:H5'	1.71	0.71
1:A:1038:C:H2'	1:A:1039:C:C6	2.26	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:30:ARG:HG3	2:B:31:TYR:CE1	2.26	0.71
3:C:10:PHE:CZ	3:C:178:LEU:HD22	2.26	0.71
5:E:91:LEU:HD23	5:E:120:THR:CG2	2.20	0.71
1:A:977:A:H2'	1:A:978:A:H5''	1.73	0.71
3:C:12:LEU:H	3:C:14:ILE:CD1	2.03	0.71
4:D:9:CYS:HB2	4:D:22:LYS:HZ1	1.53	0.71
4:D:152:SER:O	4:D:155:LEU:HB2	1.90	0.71
4:D:158:ILE:CD1	4:D:158:ILE:N	2.38	0.71
5:E:10:MET:N	5:E:10:MET:SD	2.53	0.71
8:H:104:ARG:HG3	8:H:138:TRP:CE3	2.26	0.71
9:I:5:TYR:CD2	9:I:6:GLY:N	2.57	0.71
10:J:23:ILE:N	10:J:23:ILE:HD12	2.05	0.71
20:T:69:GLY:O	20:T:71:THR:N	2.22	0.71
1:A:434:U:H2'	1:A:435:C:C6	2.26	0.71
1:A:794:A:H2'	1:A:795:C:C6	2.26	0.71
1:A:1260:C:O5'	1:A:1284:C:H4'	1.91	0.71
2:B:188:ALA:HB1	2:B:192:SER:OG	1.90	0.71
10:J:44:VAL:HG12	10:J:45:ARG:N	2.04	0.71
13:M:91:ARG:CB	13:M:98:VAL:HG22	2.20	0.71
2:B:68:ILE:N	2:B:90:MET:HE3	2.05	0.71
4:D:191:ARG:NH2	4:D:198:VAL:O	2.23	0.71
16:P:28:ARG:HG3	16:P:29:ASP:N	2.04	0.71
20:T:75:ASN:N	20:T:75:ASN:ND2	2.36	0.71
1:A:192:U:H5'	20:T:102:GLY:CA	2.21	0.71
1:A:383:A:H2'	1:A:384:G:C5'	2.21	0.71
1:A:853:G:O2'	1:A:854:G:H5'	1.91	0.71
1:A:1215:G:C2	1:A:1216:G:H1'	2.26	0.71
1:A:1305:G:HO2'	1:A:1306:A:H8	1.37	0.71
8:H:123:GLU:O	8:H:126:LYS:HB3	1.91	0.71
13:M:90:LEU:HD23	13:M:93:ARG:HD2	1.72	0.71
17:Q:97:SER:HB3	17:Q:102:GLY:O	1.90	0.71
1:A:624:C:H2'	1:A:625:G:H8	1.55	0.71
1:A:1174:G:H2'	1:A:1175:G:C8	2.24	0.71
5:E:35:GLY:N	5:E:112:LEU:HD12	2.06	0.71
1:A:1007:C:N4	1:A:1022:G:H22	1.88	0.70
2:B:30:ARG:HG3	2:B:31:TYR:CD1	2.26	0.70
18:R:24:ALA:C	18:R:26:LEU:H	1.94	0.70
21:V:21:TYR:O	21:V:23:PRO:HD3	1.91	0.70
1:A:933:G:OP2	7:G:3:ARG:HB3	1.92	0.70
1:A:1102:A:H2'	1:A:1103:C:C6	2.26	0.70
5:E:126:ARG:HG2	5:E:126:ARG:HH11	1.55	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:12:LEU:HD12	7:G:12:LEU:H	1.54	0.70
13:M:62:ASN:O	13:M:63:THR:HB	1.89	0.70
1:A:115:G:H1	1:A:312:C:H42	1.37	0.70
1:A:423:G:C2'	1:A:424:G:H5'	2.21	0.70
1:A:838:G:H2'	1:A:839:U:H5''	1.71	0.70
1:A:994:A:H2'	1:A:994:A:N3	2.06	0.70
1:A:1047:G:C3'	1:A:1048:G:H5''	2.21	0.70
1:A:1339:A:H2'	1:A:1340:A:O4'	1.92	0.70
1:A:1382:C:H2'	1:A:1383:C:C6	2.26	0.70
2:B:91:PRO:HG2	2:B:155:LEU:CD2	2.21	0.70
2:B:91:PRO:HG2	2:B:155:LEU:HD21	1.72	0.70
10:J:3:LYS:HG2	10:J:76:ASN:HD22	1.56	0.70
12:L:90:VAL:HG12	12:L:92:ASP:HB2	1.73	0.70
1:A:352:C:H4'	1:A:354:G:OP1	1.91	0.70
1:A:1218:C:H2'	1:A:1219:U:C6	2.25	0.70
1:A:1355:G:H2'	1:A:1356:G:H8	1.55	0.70
10:J:70:ARG:HH11	10:J:70:ARG:CB	2.04	0.70
11:K:11:LYS:O	11:K:12:ARG:HB2	1.91	0.70
13:M:14:ARG:HB3	13:M:14:ARG:NH1	2.04	0.70
15:O:21:ASP:OD2	15:O:24:SER:HB3	1.90	0.70
16:P:20:VAL:O	16:P:21:VAL:HG23	1.91	0.70
17:Q:51:TYR:CE1	17:Q:73:VAL:HB	2.27	0.70
18:R:40:LEU:HD23	18:R:40:LEU:O	1.92	0.70
1:A:1053:G:H4'	1:A:1054:C:H5'	1.72	0.70
1:A:1142:G:H3'	1:A:1143:G:C8	2.26	0.70
2:B:19:HIS:HB3	2:B:189:ASP:OD2	1.91	0.70
4:D:155:LEU:HB3	4:D:158:ILE:HD13	1.73	0.70
5:E:105:VAL:HG11	5:E:131:ILE:HG22	1.72	0.70
16:P:43:LYS:HB3	16:P:48:TRP:CD1	2.26	0.70
20:T:20:LEU:O	20:T:23:ARG:HB3	1.91	0.70
1:A:185:A:H2'	1:A:186:C:C6	2.25	0.70
1:A:380:G:H21	1:A:382:A:H3'	1.53	0.70
5:E:39:GLY:N	5:E:71:LEU:HD11	2.06	0.70
9:I:18:PHE:O	9:I:61:ALA:HB1	1.91	0.70
12:L:42:THR:HG23	12:L:53:ARG:O	1.90	0.70
15:O:3:ILE:HA	15:O:7:GLU:OE1	1.92	0.70
15:O:86:GLY:C	15:O:87:ILE:HD12	2.12	0.70
19:S:28:LYS:HD3	19:S:29:ARG:N	2.02	0.70
1:A:707:C:H5''	11:K:20:TYR:CD2	2.27	0.70
1:A:1060:C:H2'	1:A:1061:G:H8	1.57	0.70
3:C:180:ALA:CB	3:C:206:GLU:HG3	2.22	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:33:TYR:CB	6:F:75:LEU:HD23	2.21	0.70
10:J:8:LEU:HD13	10:J:16:LEU:CD2	2.18	0.70
10:J:26:ALA:HB3	10:J:85:LEU:CG	2.21	0.70
11:K:99:GLN:HA	11:K:105:VAL:HG21	1.73	0.70
19:S:38:SER:H	19:S:71:LEU:HD12	1.56	0.70
1:A:92:C:H2'	1:A:93:G:H8	1.57	0.70
1:A:1525:G:O2'	1:A:1526:G:H5'	1.91	0.70
2:B:144:ARG:HG3	2:B:145:LEU:H	1.56	0.70
3:C:36:ASP:O	3:C:39:ILE:HG12	1.91	0.70
4:D:120:LEU:O	4:D:125:HIS:HB2	1.92	0.70
8:H:11:THR:HG22	8:H:15:ASN:HD21	1.55	0.70
8:H:12:ARG:NH1	8:H:27:PRO:HD3	2.06	0.70
10:J:64:GLU:HG3	14:N:59:ALA:CB	2.22	0.70
11:K:15:ALA:HA	11:K:77:MET:HA	1.73	0.70
13:M:50:GLU:O	13:M:54:VAL:HG23	1.91	0.70
20:T:50:GLU:O	20:T:53:LEU:N	2.23	0.70
1:A:170:U:O2'	1:A:171:A:H5'	1.91	0.70
1:A:1196:U:H5''	1:A:1197:G:H5'	1.74	0.70
20:T:82:SER:O	20:T:83:ARG:C	2.30	0.70
1:A:292:G:H3'	1:A:293:G:H8	1.55	0.69
1:A:448:A:P	1:A:485:G:H22	2.14	0.69
1:A:792:A:O2'	1:A:793:U:OP2	2.10	0.69
1:A:1093:A:H2	1:A:1109:C:HO2'	1.38	0.69
1:A:1189:C:H2'	1:A:1190:G:H5'	1.74	0.69
9:I:101:PHE:N	9:I:102:LEU:HD23	2.07	0.69
14:N:40:CYS:SG	14:N:42:ILE:HG13	2.32	0.69
18:R:25:THR:C	18:R:26:LEU:HD12	2.12	0.69
2:B:96:ARG:O	2:B:98:LEU:HD23	1.92	0.69
3:C:3:ASN:N	3:C:3:ASN:ND2	2.40	0.69
14:N:11:LYS:HE3	14:N:13:THR:HB	1.74	0.69
20:T:70:SER:HA	20:T:73:HIS:CD2	2.28	0.69
1:A:38:G:H22	1:A:397:A:P	2.15	0.69
1:A:603:U:H2'	1:A:604:G:H8	1.57	0.69
3:C:66:VAL:C	3:C:68:VAL:H	1.95	0.69
3:C:85:ARG:O	3:C:89:GLU:HG2	1.92	0.69
3:C:121:ALA:O	3:C:125:GLU:HG3	1.93	0.69
4:D:79:PHE:HD2	4:D:79:PHE:O	1.74	0.69
6:F:20:ALA:HA	6:F:23:LYS:HB2	1.74	0.69
10:J:40:LEU:HD23	10:J:41:PRO:CD	2.23	0.69
10:J:79:ARG:NH1	10:J:82:ILE:HD12	2.06	0.69
1:A:106:C:O2'	1:A:107:G:H5'	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:499:A:H4'	1:A:500:G:H5'	1.74	0.69
1:A:690:G:OP2	11:K:27:ASN:HB2	1.91	0.69
1:A:1226:C:C5	13:M:104:ARG:HA	2.27	0.69
1:A:1298:C:C4'	1:A:1299:A:H5'	2.21	0.69
7:G:141:VAL:O	7:G:144:MET:N	2.24	0.69
9:I:24:GLY:CA	9:I:60:ASP:HA	2.22	0.69
13:M:19:LEU:C	13:M:21:TYR:H	1.96	0.69
15:O:26:GLU:HA	15:O:81:LEU:HD11	1.75	0.69
17:Q:59:ILE:HG22	17:Q:71:PHE:HB3	1.74	0.69
18:R:39:VAL:CG1	18:R:40:LEU:H	2.00	0.69
20:T:43:LEU:O	20:T:48:LYS:HG2	1.92	0.69
1:A:19:C:H2'	1:A:20:U:H6	1.58	0.69
1:A:163:C:O2	1:A:163:C:H2'	1.92	0.69
1:A:746:A:O2'	1:A:747:C:H5'	1.91	0.69
1:A:1206:G:H2'	1:A:1207:G:C8	2.27	0.69
2:B:20:GLU:O	2:B:39:ILE:HG23	1.91	0.69
4:D:100:ARG:HB3	4:D:102:ASP:OD1	1.91	0.69
4:D:160:GLN:O	4:D:163:GLU:HB3	1.92	0.69
5:E:148:VAL:HG21	8:H:107:LEU:HD13	1.74	0.69
6:F:7:ASN:ND2	18:R:76:LEU:HD11	2.07	0.69
9:I:19:LEU:HG	9:I:60:ASP:O	1.93	0.69
11:K:46:GLY:O	11:K:47:VAL:C	2.31	0.69
12:L:113:ARG:HH12	12:L:115:LYS:HB2	1.52	0.69
13:M:74:VAL:HA	13:M:77:ASN:ND2	2.07	0.69
1:A:135:C:O2	16:P:1:MET:HB2	1.92	0.69
1:A:918:A:H2'	1:A:919:A:O4'	1.92	0.69
1:A:1152:A:H5''	10:J:13:HIS:HB2	1.75	0.69
1:A:1299:A:N3	1:A:1299:A:H2'	2.05	0.69
1:A:1527:C:O2'	1:A:1528:U:H5'	1.93	0.69
2:B:204:ASN:ND2	2:B:206:ASP:H	1.89	0.69
6:F:50:TYR:CE1	18:R:77:GLY:HA2	2.27	0.69
8:H:40:ALA:HA	8:H:45:ILE:CG1	2.23	0.69
14:N:8:GLU:HA	14:N:11:LYS:HD3	1.73	0.69
16:P:2:VAL:O	16:P:64:ALA:HA	1.92	0.69
1:A:404:U:O2'	1:A:405:U:H5'	1.93	0.69
1:A:1036:G:H2'	1:A:1037:C:O4'	1.92	0.69
1:A:1330:U:H4'	13:M:23:TYR:CE1	2.27	0.69
2:B:96:ARG:H	2:B:96:ARG:CZ	2.06	0.69
9:I:19:LEU:CG	9:I:61:ALA:HB2	2.22	0.69
1:A:187:C:O2'	20:T:89:ARG:HD2	1.92	0.69
1:A:600:C:OP1	8:H:97:VAL:HG12	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:627:G:O2'	1:A:628:G:H5'	1.92	0.69
1:A:1104:G:OP1	2:B:111:ARG:HD2	1.93	0.69
1:A:1238:A:H2	1:A:1241:G:N3	1.91	0.69
1:A:1446:A:H5''	1:A:1446:A:N3	2.06	0.69
3:C:90:GLU:HA	3:C:93:LYS:CG	2.23	0.69
3:C:139:GLN:O	3:C:143:GLU:HG3	1.93	0.69
4:D:10:ARG:CG	4:D:11:LEU:N	2.55	0.69
5:E:45:PHE:CE2	5:E:47:LYS:HE3	2.27	0.69
10:J:13:HIS:O	10:J:17:ASP:HB2	1.92	0.69
13:M:8:GLU:HG3	13:M:22:ILE:HG12	1.74	0.69
14:N:12:ARG:O	14:N:12:ARG:HG2	1.92	0.69
16:P:47:ASP:O	16:P:49:LEU:N	2.26	0.69
17:Q:70:ARG:HG3	17:Q:70:ARG:NH1	1.95	0.69
19:S:32:LYS:HA	19:S:50:ALA:CB	2.21	0.69
20:T:14:LYS:O	20:T:16:HIS:N	2.26	0.69
1:A:322:C:O2'	20:T:23:ARG:HD3	1.92	0.69
1:A:541:G:H2'	1:A:542:G:C8	2.28	0.69
1:A:1477:C:H2'	1:A:1478:C:C6	2.28	0.69
3:C:22:TRP:HB3	3:C:59:ARG:CB	2.22	0.69
3:C:125:GLU:O	3:C:127:ARG:HD2	1.93	0.69
6:F:22:GLU:HA	6:F:25:ILE:HG22	1.75	0.69
6:F:97:PHE:HD2	6:F:98:LEU:H	1.41	0.69
8:H:96:GLY:H	8:H:99:GLU:CD	1.95	0.69
10:J:12:ASP:HB3	10:J:15:THR:HB	1.75	0.69
11:K:53:SER:O	11:K:55:LYS:N	2.26	0.69
11:K:82:VAL:O	11:K:83:ILE:HG13	1.93	0.69
15:O:17:ARG:HD3	15:O:77:ARG:HH11	1.58	0.69
1:A:22:G:H2'	1:A:23:C:H6	1.58	0.69
1:A:103:C:OP1	20:T:17:ARG:NH1	2.26	0.69
1:A:1171:G:H2'	1:A:1172:C:C6	2.28	0.69
1:A:1205:U:H4'	3:C:195:VAL:HG21	1.75	0.69
1:A:1315:U:H2'	1:A:1316:G:O4'	1.92	0.69
2:B:36:ARG:CD	2:B:41:ILE:HD12	2.22	0.69
2:B:71:VAL:HG21	2:B:164:VAL:HG22	1.74	0.69
3:C:64:VAL:HG21	3:C:95:THR:OG1	1.93	0.69
4:D:205:GLU:O	4:D:207:TYR:N	2.26	0.69
5:E:13:ILE:HA	5:E:29:GLY:O	1.93	0.69
7:G:116:ALA:CA	7:G:119:ARG:HH21	2.05	0.69
1:A:1425:U:H2'	1:A:1426:C:C6	2.28	0.68
3:C:48:TYR:O	3:C:52:LEU:HD22	1.93	0.68
7:G:155:ARG:HA	7:G:155:ARG:NE	2.08	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:11:LYS:O	9:I:11:LYS:HD3	1.92	0.68
10:J:22:LYS:HB2	10:J:22:LYS:HZ2	1.57	0.68
11:K:127:LYS:HA	11:K:127:LYS:CE	2.12	0.68
17:Q:6:LEU:O	17:Q:58:GLU:HA	1.93	0.68
1:A:1406:U:H1'	1:A:1517:G:N2	2.08	0.68
3:C:11:ARG:HH12	3:C:179:ARG:N	1.88	0.68
3:C:76:VAL:HA	3:C:83:ARG:HD2	1.74	0.68
7:G:44:TYR:O	7:G:47:CYS:HB2	1.93	0.68
18:R:40:LEU:HD23	18:R:40:LEU:C	2.14	0.68
1:A:73:C:H2'	1:A:74:C:H6	1.56	0.68
1:A:478:A:O2'	1:A:479:C:H5'	1.93	0.68
1:A:923:A:O2'	1:A:924:C:H5'	1.93	0.68
1:A:1123:A:H4'	10:J:37:PRO:CD	2.19	0.68
2:B:9:GLU:HB3	2:B:12:GLU:OE2	1.92	0.68
3:C:170:GLN:HG2	3:C:171:GLY:H	1.57	0.68
4:D:24:GLU:HG3	4:D:25:ARG:H	1.58	0.68
6:F:8:ILE:HG12	6:F:88:VAL:HG22	1.74	0.68
10:J:35:SER:HB2	10:J:73:ASP:O	1.93	0.68
12:L:66:VAL:HG11	12:L:98:TYR:CE1	2.28	0.68
13:M:5:ALA:CB	13:M:22:ILE:HD13	2.24	0.68
1:A:114:U:H2'	1:A:115:G:C8	2.28	0.68
1:A:659:U:H3	1:A:746:A:H61	1.38	0.68
1:A:1313:U:H3'	19:S:6:LYS:HD3	1.75	0.68
3:C:37:GLN:HE22	14:N:52:GLN:HE22	1.42	0.68
9:I:6:GLY:HA2	9:I:83:ARG:HD2	1.75	0.68
10:J:49:VAL:HG11	14:N:41:ARG:CB	2.24	0.68
11:K:43:SER:O	11:K:44:SER:HB3	1.92	0.68
12:L:59:ARG:HH11	12:L:59:ARG:HG3	1.58	0.68
13:M:54:VAL:O	13:M:57:ARG:HB3	1.93	0.68
17:Q:76:LEU:C	17:Q:76:LEU:HD12	2.14	0.68
1:A:142:G:H2'	1:A:143:A:H8	1.58	0.68
1:A:376:G:H2'	1:A:377:G:C8	2.27	0.68
1:A:524:G:H2'	1:A:525:C:C6	2.29	0.68
1:A:524:G:H2'	1:A:525:C:H6	1.59	0.68
3:C:12:LEU:H	3:C:14:ILE:HD11	1.58	0.68
3:C:196:LEU:HD23	3:C:196:LEU:H	1.59	0.68
18:R:38:GLU:H	18:R:41:LYS:HG3	1.56	0.68
1:A:824:C:H2'	1:A:825:G:C8	2.28	0.68
1:A:1086:U:O2'	1:A:1087:G:H5'	1.93	0.68
1:A:1361(A):C:C2'	1:A:1362:C:H5''	2.23	0.68
1:A:1507:A:H2	1:A:1530:G:H1'	1.53	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:178:ARG:NE	2:B:196:LEU:O	2.25	0.68
10:J:57:LYS:HE2	10:J:58:ASP:OD2	1.94	0.68
12:L:25:PRO:O	12:L:27:LEU:HD13	1.93	0.68
13:M:53:VAL:O	13:M:57:ARG:HB2	1.94	0.68
1:A:76:C:O2'	1:A:77:G:H5'	1.94	0.68
1:A:254:G:O2'	1:A:255:G:H5'	1.92	0.68
1:A:491:G:H2'	1:A:492:G:H8	1.55	0.68
1:A:769:G:O2'	1:A:770:C:H5'	1.93	0.68
1:A:830:G:O2'	1:A:831:U:H5'	1.94	0.68
1:A:1007:C:H42	1:A:1022:G:N2	1.89	0.68
1:A:1247:U:O2'	1:A:1248:A:H5'	1.94	0.68
1:A:1279:A:H5''	1:A:1280:A:OP1	1.92	0.68
1:A:1316:G:H2'	1:A:1318:A:OP2	1.94	0.68
3:C:18:TRP:CD1	14:N:51:GLY:HA2	2.29	0.68
5:E:139:LEU:CD2	5:E:142:LEU:HD11	2.24	0.68
6:F:28:ARG:HA	6:F:31:GLU:OE1	1.93	0.68
10:J:9:ARG:HH22	10:J:69:ASN:HB2	1.57	0.68
11:K:52:GLY:O	11:K:54:ARG:N	2.27	0.68
12:L:26:ALA:HB2	12:L:98:TYR:CE2	2.29	0.68
14:N:9:LYS:HD3	14:N:9:LYS:C	2.13	0.68
1:A:125:U:H2'	1:A:126:G:C8	2.29	0.68
1:A:782:A:C6	1:A:801:U:C2	2.81	0.68
1:A:930:C:C2'	1:A:931:C:H5'	2.23	0.68
1:A:1158:C:H5''	2:B:133:LYS:HE3	1.76	0.68
2:B:21:ARG:HH12	2:B:23:ARG:NH2	1.92	0.68
2:B:102:LEU:HD21	2:B:162:ILE:HD11	1.76	0.68
3:C:91:LEU:CD1	3:C:99:VAL:HG22	2.23	0.68
8:H:28:ALA:HA	8:H:59:LEU:HD12	1.75	0.68
1:A:1289:A:H3'	1:A:1290:G:H8	1.59	0.68
1:A:1441:G:H4'	1:A:1442:G:C2	2.29	0.68
2:B:13:ALA:C	2:B:15:VAL:H	1.95	0.68
4:D:36:ARG:H	4:D:37:PRO:CD	2.07	0.68
7:G:116:ALA:O	7:G:120:ILE:HG12	1.94	0.68
11:K:73:MET:HA	11:K:77:MET:HB2	1.75	0.68
12:L:70:ILE:HD13	12:L:77:LEU:HD12	1.74	0.68
12:L:89:ARG:O	12:L:90:VAL:CG2	2.41	0.68
17:Q:7:THR:HA	17:Q:57:VAL:O	1.94	0.68
18:R:70:ILE:HG22	18:R:71:LYS:N	2.08	0.68
20:T:85:MET:HB2	20:T:104:LEU:HD21	1.75	0.68
1:A:359:U:H2'	1:A:360:A:H8	1.58	0.68
2:B:105:PHE:HD2	2:B:105:PHE:C	1.98	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:76:VAL:HB	3:C:103:VAL:HG21	1.76	0.68
4:D:2:GLY:O	4:D:3:ARG:HB2	1.92	0.68
5:E:59:GLY:O	5:E:62:ALA:HB3	1.93	0.68
1:A:651:C:O2'	1:A:652:U:H5'	1.94	0.67
1:A:1262:C:H2'	1:A:1263:C:C6	2.29	0.67
1:A:1407:C:O2'	1:A:1408:A:H5'	1.94	0.67
4:D:80:GLU:HA	4:D:80:GLU:OE2	1.93	0.67
6:F:29:ALA:O	6:F:32:ASN:HB3	1.93	0.67
7:G:66:VAL:HG12	7:G:70:LYS:HE3	1.76	0.67
7:G:91:VAL:CG1	7:G:92:SER:N	2.57	0.67
17:Q:67:LYS:HA	17:Q:70:ARG:NH2	2.09	0.67
21:V:6:ARG:O	21:V:12:LYS:HE3	1.94	0.67
1:A:419:C:O2	1:A:419:C:H3'	1.93	0.67
1:A:532:A:H2'	1:A:533:A:C5'	2.24	0.67
1:A:1026:G:H2'	1:A:1027:C:H5'	1.76	0.67
5:E:66:MET:O	5:E:67:VAL:HG13	1.93	0.67
9:I:5:TYR:CE2	9:I:16:ARG:HB2	2.29	0.67
13:M:10:PRO:O	13:M:45:VAL:HG21	1.94	0.67
13:M:33:ALA:O	13:M:37:THR:HB	1.94	0.67
1:A:760:G:H2'	1:A:761:G:O4'	1.94	0.67
1:A:1276:G:O2'	1:A:1277:C:H5'	1.94	0.67
3:C:84:ILE:HG13	3:C:101:LEU:HD22	1.77	0.67
4:D:9:CYS:CB	4:D:22:LYS:NZ	2.57	0.67
4:D:123:HIS:HB2	4:D:125:HIS:CD2	2.30	0.67
6:F:96:PRO:HB2	18:R:30:ASP:OD1	1.93	0.67
7:G:137:LYS:HA	7:G:140:ASP:CB	2.24	0.67
10:J:22:LYS:C	10:J:23:ILE:HD12	2.15	0.67
10:J:77:PRO:O	10:J:78:ASN:HB2	1.95	0.67
15:O:56:LEU:HA	15:O:59:MET:CG	2.24	0.67
1:A:24:U:H2'	1:A:25:C:C6	2.28	0.67
1:A:336:C:H2'	1:A:337:C:C6	2.29	0.67
1:A:538:G:H2'	1:A:539:A:C8	2.29	0.67
1:A:857:C:H2'	1:A:858:G:C8	2.29	0.67
1:A:1048:G:N2	1:A:1214:C:H2'	2.10	0.67
1:A:1428:A:H2'	1:A:1429:C:C6	2.28	0.67
4:D:105:VAL:O	4:D:108:LEU:HB2	1.94	0.67
7:G:103:TRP:NE1	7:G:137:LYS:HD3	2.09	0.67
12:L:33:ARG:HG2	12:L:60:LEU:HD12	1.76	0.67
13:M:90:LEU:O	13:M:93:ARG:HB2	1.95	0.67
14:N:29:ARG:HG2	14:N:30:ALA:H	1.60	0.67
1:A:36:C:H2'	1:A:37:U:H6	1.57	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:C:O2	1:A:352:C:H2'	1.93	0.67
1:A:1030:C:H2'	1:A:1030(A):G:O4'	1.93	0.67
3:C:64:VAL:HG23	3:C:97:LYS:CB	2.25	0.67
7:G:78:ARG:O	7:G:84:ASN:HA	1.94	0.67
8:H:65:TYR:CA	8:H:79:VAL:HG23	2.23	0.67
10:J:7:LYS:HA	10:J:71:LEU:HD22	1.75	0.67
10:J:64:GLU:N	14:N:59:ALA:HB2	2.09	0.67
11:K:87:THR:HA	11:K:91:ARG:NH1	2.09	0.67
13:M:25:ILE:CD1	13:M:66:LEU:HD21	2.25	0.67
15:O:78:TYR:O	15:O:82:ILE:HG22	1.95	0.67
17:Q:17:LYS:O	17:Q:46:ASP:O	2.13	0.67
18:R:38:GLU:HA	18:R:41:LYS:HZ2	1.59	0.67
1:A:243:A:C5'	1:A:244:U:H5'	2.24	0.67
1:A:328:C:O2	1:A:328:C:H2'	1.95	0.67
1:A:849:C:C2'	1:A:850:U:H5'	2.24	0.67
2:B:87:ARG:NH1	2:B:219:VAL:HB	2.09	0.67
2:B:105:PHE:C	2:B:105:PHE:CD2	2.66	0.67
10:J:46:ARG:NH1	10:J:64:GLU:HG2	2.08	0.67
10:J:79:ARG:O	10:J:83:GLU:HB2	1.94	0.67
12:L:66:VAL:HG21	12:L:98:TYR:CE1	2.30	0.67
14:N:46:GLU:O	14:N:49:HIS:HB2	1.93	0.67
18:R:86:VAL:O	18:R:87:ARG:HB2	1.94	0.67
19:S:72:GLY:C	19:S:74:PHE:H	1.96	0.67
1:A:1244:C:H2'	1:A:1245:A:H8	1.57	0.67
1:A:1361:G:O5'	1:A:1361:G:C8	2.47	0.67
3:C:85:ARG:HA	3:C:88:ARG:CG	2.19	0.67
6:F:38:GLU:O	6:F:39:LYS:HB2	1.93	0.67
9:I:77:ILE:HG22	9:I:81:ILE:HD11	1.76	0.67
17:Q:90:ILE:O	17:Q:93:GLN:HB3	1.94	0.67
20:T:80:ARG:HB3	20:T:80:ARG:HH11	1.60	0.67
1:A:625:G:H2'	1:A:626:U:C6	2.29	0.67
1:A:849:C:O2'	1:A:850:U:H5'	1.95	0.67
1:A:1015:A:H2'	1:A:1016:A:H8	1.55	0.67
1:A:1447:G:H2'	1:A:1447:G:N3	2.10	0.67
2:B:27:LYS:O	2:B:194:PRO:HG3	1.95	0.67
3:C:23:TYR:OH	10:J:9:ARG:HB2	1.94	0.67
4:D:15:GLU:HG3	4:D:63:LYS:HZ3	1.57	0.67
13:M:49:THR:O	13:M:53:VAL:HG23	1.95	0.67
14:N:14:PRO:CG	14:N:15:LYS:H	2.08	0.67
16:P:80:PHE:CD1	16:P:80:PHE:N	2.63	0.67
20:T:43:LEU:HD12	20:T:55:ILE:HD13	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1151:A:H5''	10:J:42:THR:OG1	1.95	0.67
1:A:1182:G:H4'	1:A:1183:A:O5'	1.92	0.67
1:A:1222:G:H5''	19:S:78:ARG:NH1	2.10	0.67
9:I:55:ALA:HB1	9:I:59:PHE:CD1	2.30	0.67
12:L:41:ARG:HB3	12:L:41:ARG:NH1	2.04	0.67
13:M:22:ILE:HG21	13:M:66:LEU:HD23	1.77	0.67
15:O:10:LYS:HG3	15:O:11:VAL:N	2.10	0.67
16:P:59:TRP:HA	16:P:59:TRP:HE3	1.55	0.67
1:A:839:U:O2	1:A:839:U:H2'	1.95	0.67
2:B:54:THR:HG23	2:B:199:TYR:HB2	1.77	0.67
4:D:152:SER:OG	4:D:155:LEU:HD12	1.95	0.67
5:E:79:GLU:HA	5:E:91:LEU:O	1.95	0.67
10:J:9:ARG:NE	10:J:9:ARG:N	2.42	0.67
13:M:117:VAL:O	13:M:118:ALA:HB2	1.94	0.67
1:A:452:A:H4'	16:P:72:ARG:NH2	2.10	0.66
1:A:1150:U:O3'	10:J:41:PRO:CA	2.38	0.66
1:A:1372:U:C5'	9:I:71:SER:HB3	2.15	0.66
8:H:45:ILE:HG22	8:H:63:LEU:HA	1.76	0.66
18:R:37:VAL:HG23	18:R:38:GLU:N	2.09	0.66
1:A:322:C:O2'	1:A:323:U:H5'	1.94	0.66
1:A:478:A:H2'	1:A:479:C:O4'	1.94	0.66
1:A:1347:G:H1'	1:A:1348:U:C5	2.30	0.66
1:A:1366:C:H2'	1:A:1367:C:C6	2.24	0.66
4:D:25:ARG:NE	4:D:30:LYS:HB3	2.09	0.66
7:G:92:SER:HB3	7:G:93:PRO:HD2	1.76	0.66
16:P:80:PHE:O	16:P:82:GLN:N	2.28	0.66
19:S:19:VAL:HG22	19:S:47:HIS:HE1	1.59	0.66
20:T:57:ARG:HH21	20:T:100:ILE:HG22	1.60	0.66
1:A:149:A:O2'	1:A:150:C:H5'	1.94	0.66
1:A:174:C:H2'	1:A:175:C:C6	2.30	0.66
1:A:1510:U:H2'	1:A:1511:G:C8	2.30	0.66
5:E:82:VAL:HG21	5:E:138:ALA:H	1.60	0.66
9:I:117:HIS:HB2	9:I:121:ARG:HD2	1.76	0.66
10:J:23:ILE:N	10:J:23:ILE:CD1	2.59	0.66
12:L:46:LYS:HE3	12:L:47:LYS:HE2	1.78	0.66
19:S:63:THR:HB	19:S:65:ASN:OD1	1.95	0.66
1:A:1182:G:C4'	1:A:1183:A:H5''	2.22	0.66
3:C:51:GLY:N	3:C:70:VAL:HG13	2.10	0.66
4:D:25:ARG:HG3	4:D:30:LYS:CB	2.24	0.66
5:E:147:ASP:OD2	5:E:147:ASP:N	2.29	0.66
7:G:46:ALA:HB1	7:G:121:ALA:CA	2.26	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:80:ILE:O	8:H:80:ILE:HG22	1.96	0.66
9:I:7:THR:CG2	9:I:8:GLY:H	2.08	0.66
12:L:66:VAL:HG11	12:L:98:TYR:HE1	1.61	0.66
1:A:590:C:OP1	8:H:30:ARG:N	2.27	0.66
3:C:50:ALA:CA	3:C:72:LYS:HD3	2.19	0.66
3:C:91:LEU:HD12	3:C:101:LEU:HB2	1.76	0.66
8:H:63:LEU:HD12	8:H:63:LEU:N	2.03	0.66
21:V:24:ARG:O	21:V:25:LYS:HB2	1.94	0.66
1:A:129(A):G:HO2'	1:A:130:A:P	2.17	0.66
1:A:525:C:H2'	1:A:526:C:C6	2.30	0.66
1:A:624:C:H2'	1:A:625:G:C8	2.29	0.66
1:A:832:C:O2'	1:A:833:U:H5'	1.96	0.66
1:A:1152:A:C5'	10:J:13:HIS:HB2	2.26	0.66
1:A:1403:C:H2'	1:A:1404:C:C6	2.30	0.66
3:C:180:ALA:HB1	3:C:203:PHE:HE1	1.58	0.66
10:J:16:LEU:CA	10:J:94:VAL:HG21	2.23	0.66
11:K:93:GLN:NE2	11:K:96:ARG:NH2	2.44	0.66
16:P:22:THR:HG23	16:P:23:ASP:N	2.07	0.66
19:S:72:GLY:C	19:S:75:ALA:H	1.98	0.66
5:E:115:VAL:HG12	5:E:116:THR:N	2.10	0.66
8:H:31:PHE:O	8:H:33:GLU:N	2.28	0.66
10:J:78:ASN:C	10:J:80:LYS:N	2.48	0.66
12:L:67:THR:OG1	12:L:96:VAL:HA	1.96	0.66
12:L:79:GLU:O	12:L:79:GLU:HG2	1.95	0.66
18:R:59:SER:H	18:R:62:GLU:HB2	1.60	0.66
1:A:89:C:H2'	1:A:90:U:C6	2.31	0.66
1:A:261:U:O2	1:A:263:A:H8	1.79	0.66
1:A:333:G:H4'	20:T:16:HIS:CE1	2.31	0.66
1:A:489:C:H2'	1:A:490:G:H8	1.59	0.66
1:A:959:A:H3'	1:A:960:U:H5''	1.78	0.66
1:A:1049:U:OP1	14:N:3:ARG:HG2	1.96	0.66
10:J:29:ARG:HG3	10:J:84:GLN:HE22	1.59	0.66
12:L:24:VAL:HG13	12:L:98:TYR:CE2	2.30	0.66
13:M:81:LEU:HA	13:M:84:ILE:CG1	2.25	0.66
15:O:4:THR:OG1	15:O:6:GLU:HG2	1.96	0.66
1:A:254:G:O3'	17:Q:69:LYS:HE2	1.96	0.66
1:A:1277:C:H1'	1:A:1282:C:O2	1.95	0.66
1:A:1360:A:H2'	1:A:1361:G:C8	2.31	0.66
3:C:44:GLU:OE1	3:C:55:VAL:HG22	1.96	0.66
3:C:160:ALA:C	3:C:162:GLN:N	2.49	0.66
4:D:82:ALA:O	4:D:85:LYS:N	2.29	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:2:ARG:O	6:F:66:GLU:HG3	1.96	0.66
1:A:580:U:H2'	1:A:581:G:O4'	1.95	0.66
1:A:741:G:O2'	1:A:742:G:H5'	1.95	0.66
1:A:913:A:O2'	1:A:914:A:P	2.54	0.66
1:A:1201:A:H4'	1:A:1202:G:O5'	1.96	0.66
1:A:1313:U:P	19:S:6:LYS:HG2	2.36	0.66
1:A:1346:A:H5'	9:I:120:ARG:HH12	1.61	0.66
9:I:96:LEU:HG	9:I:102:LEU:HD13	1.77	0.66
12:L:46:LYS:N	12:L:92:ASP:O	2.29	0.66
14:N:41:ARG:HG2	14:N:41:ARG:NH1	2.11	0.66
17:Q:67:LYS:HA	17:Q:70:ARG:HH22	1.61	0.66
18:R:22:VAL:O	18:R:24:ALA:N	2.28	0.66
20:T:75:ASN:O	20:T:76:ALA:C	2.34	0.66
1:A:312:C:H2'	1:A:313:A:C8	2.30	0.65
1:A:376:G:OP1	16:P:5:ARG:HB2	1.95	0.65
1:A:922:G:N2	1:A:1396:A:C4	2.64	0.65
1:A:1293:G:H2'	1:A:1294:G:O4'	1.96	0.65
2:B:178:ARG:NH2	8:H:68:ARG:NH2	2.44	0.65
3:C:132:ARG:O	3:C:136:GLN:HG3	1.95	0.65
4:D:26:CYS:HA	4:D:31:CYS:HB2	1.77	0.65
4:D:153:ARG:HG2	4:D:181:MET:SD	2.36	0.65
13:M:76:ALA:O	13:M:80:ARG:N	2.29	0.65
19:S:70:LYS:H	19:S:73:GLU:HG3	1.61	0.65
20:T:11:SER:O	20:T:13:LEU:N	2.29	0.65
1:A:711:G:H2'	1:A:712:A:C8	2.31	0.65
2:B:96:ARG:O	2:B:98:LEU:CD2	2.45	0.65
2:B:165:VAL:O	2:B:187:LEU:O	2.15	0.65
2:B:204:ASN:C	2:B:204:ASN:HD22	2.00	0.65
4:D:43:HIS:O	4:D:45:GLN:N	2.29	0.65
4:D:144:ASP:O	4:D:184:LYS:HA	1.96	0.65
5:E:87:SER:HB3	5:E:131:ILE:CD1	2.24	0.65
5:E:111:GLU:O	5:E:114:GLY:N	2.25	0.65
7:G:46:ALA:HB2	7:G:120:ILE:HB	1.79	0.65
7:G:47:CYS:C	7:G:49:ILE:H	1.98	0.65
12:L:89:ARG:HA	12:L:96:VAL:O	1.96	0.65
13:M:48:LEU:CD2	13:M:52:GLU:HB2	2.26	0.65
19:S:27:GLU:OE1	19:S:47:HIS:NE2	2.28	0.65
19:S:42:PRO:O	19:S:45:VAL:HG23	1.96	0.65
1:A:534:U:H5''	1:A:535:A:OP2	1.96	0.65
1:A:1064:G:H4'	1:A:1065:U:H5'	1.76	0.65
2:B:77:ALA:HA	2:B:211:ILE:HD11	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:46:ALA:O	9:I:81:ILE:HD12	1.95	0.65
10:J:9:ARG:HA	10:J:69:ASN:HA	1.78	0.65
10:J:44:VAL:HG13	10:J:65:LEU:O	1.95	0.65
10:J:65:LEU:HA	14:N:55:GLY:O	1.95	0.65
16:P:51:VAL:O	16:P:52:ASP:HB3	1.96	0.65
19:S:18:LYS:HE2	19:S:31:ILE:HD12	1.77	0.65
19:S:64:GLU:HA	19:S:67:VAL:HG23	1.77	0.65
20:T:75:ASN:HD22	20:T:75:ASN:H	1.39	0.65
1:A:191:G:C4	20:T:105:SER:HB3	2.32	0.65
1:A:227:G:H2'	1:A:228:A:C8	2.32	0.65
1:A:255:G:H1'	17:Q:16:GLN:OE1	1.97	0.65
1:A:824:C:O2'	1:A:825:G:H5'	1.96	0.65
1:A:1065:U:O4	1:A:1190:G:H1'	1.96	0.65
1:A:1374:A:H2'	1:A:1375:A:H8	1.59	0.65
3:C:101:LEU:HD23	3:C:102:ASN:N	2.11	0.65
9:I:4:TYR:HE1	9:I:21:PRO:HG2	1.60	0.65
16:P:70:ALA:O	16:P:74:LEU:HG	1.96	0.65
1:A:1260:C:H4'	1:A:1284:C:H5'	1.79	0.65
2:B:60:ASP:O	2:B:62:ALA:N	2.28	0.65
4:D:201:GLN:HA	4:D:204:ILE:CD1	2.26	0.65
9:I:24:GLY:N	9:I:60:ASP:HA	2.11	0.65
11:K:79:SER:O	11:K:80:VAL:HB	1.96	0.65
13:M:20:THR:HG23	13:M:26:GLY:HA2	1.76	0.65
1:A:407:G:O2'	4:D:116:GLN:HG3	1.96	0.65
1:A:1066:C:C2'	1:A:1067:A:H5'	2.26	0.65
1:A:1210:C:H2'	1:A:1211:U:H5''	1.78	0.65
4:D:133:VAL:HG12	4:D:134:ASP:N	2.06	0.65
8:H:82:HIS:CD2	8:H:83:ILE:H	2.14	0.65
10:J:7:LYS:HG2	10:J:71:LEU:CD2	2.26	0.65
12:L:90:VAL:HG12	12:L:92:ASP:H	1.61	0.65
19:S:11:VAL:HG21	19:S:41:VAL:HG11	1.78	0.65
1:A:956:U:O2'	1:A:957:U:H5'	1.97	0.65
1:A:1060:C:O2'	1:A:1061:G:H5'	1.96	0.65
1:A:1202:G:O2'	1:A:1203:C:H5'	1.96	0.65
1:A:1343:G:H4'	9:I:122:ALA:O	1.97	0.65
1:A:1425:U:H2'	1:A:1426:C:H6	1.59	0.65
2:B:111:ARG:O	2:B:113:HIS:N	2.27	0.65
3:C:19:GLU:HG2	3:C:54:ARG:HH21	1.60	0.65
3:C:114:PRO:HD3	3:C:184:TYR:O	1.95	0.65
3:C:173:VAL:O	3:C:173:VAL:HG12	1.97	0.65
4:D:25:ARG:CG	4:D:30:LYS:HB3	2.26	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:102:ARG:NH1	8:H:105:ARG:NH1	2.45	0.65
10:J:8:LEU:HA	10:J:95:GLU:OE2	1.97	0.65
16:P:45:THR:OG1	16:P:47:ASP:OD1	2.10	0.65
1:A:76:C:H2'	1:A:77:G:H8	1.62	0.65
1:A:1123:A:O3'	10:J:36:GLY:HA3	1.97	0.65
1:A:1490:C:O2'	1:A:1491:G:H5'	1.97	0.65
4:D:148:VAL:HG12	4:D:149:ALA:N	2.11	0.65
5:E:77:PRO:O	5:E:78:HIS:HB3	1.95	0.65
5:E:135:THR:O	5:E:136:MET:C	2.33	0.65
7:G:93:PRO:O	7:G:97:GLN:HB2	1.97	0.65
18:R:38:GLU:N	18:R:41:LYS:HE3	2.12	0.65
20:T:102:GLY:O	20:T:103:GLY:C	2.35	0.65
1:A:383:A:C2'	1:A:384:G:H5'	2.26	0.65
1:A:794:A:H2'	1:A:795:C:H6	1.61	0.65
1:A:1181:G:O2'	1:A:1182:G:H5'	1.97	0.65
1:A:1251:A:H4'	9:I:12:GLU:OE2	1.96	0.65
1:A:1278:U:H5'	1:A:1279:A:H5'	1.79	0.65
5:E:12:LEU:CD1	5:E:31:LEU:HB2	2.27	0.65
5:E:91:LEU:HA	5:E:120:THR:HG22	1.79	0.65
9:I:43:ALA:HA	9:I:74:ILE:HD13	1.78	0.65
12:L:43:VAL:HG23	12:L:55:VAL:HG21	1.78	0.65
15:O:38:ARG:O	15:O:41:GLU:N	2.30	0.65
15:O:55:GLY:O	15:O:59:MET:HG2	1.96	0.65
16:P:42:ARG:C	16:P:43:LYS:HD2	2.17	0.65
17:Q:101:ARG:HG2	17:Q:101:ARG:HH11	1.62	0.65
1:A:34:C:H42	1:A:550:G:H1	1.44	0.65
1:A:540:G:H2'	1:A:541:G:O4'	1.97	0.65
1:A:608:A:H2'	1:A:609:A:C8	2.32	0.65
15:O:48:LYS:HZ3	15:O:48:LYS:HB2	1.62	0.65
20:T:38:LYS:HA	20:T:41:ILE:HD12	1.77	0.65
1:A:697:U:O2	1:A:798:G:H1'	1.97	0.64
1:A:1413:A:O2'	1:A:1414:U:H5'	1.97	0.64
2:B:184:VAL:N	2:B:198:ASP:OD2	2.30	0.64
2:B:208:ILE:O	2:B:210:SER:N	2.30	0.64
5:E:41:VAL:HG22	5:E:113:ALA:CA	2.26	0.64
5:E:137:GLU:O	5:E:140:ARG:N	2.29	0.64
10:J:37:PRO:HA	10:J:72:VAL:HG13	1.78	0.64
10:J:85:LEU:H	10:J:88:LEU:HD12	1.62	0.64
15:O:71:GLN:HB2	15:O:78:TYR:CD1	2.32	0.64
17:Q:68:ARG:O	17:Q:69:LYS:HB2	1.97	0.64
1:A:371:G:H2'	1:A:372:C:H6	1.63	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:402:G:O2'	1:A:403:C:H5'	1.97	0.64
1:A:634:C:H2'	1:A:635:G:H8	1.62	0.64
1:A:991:U:O4	1:A:1212:U:H2'	1.97	0.64
1:A:1153:C:H2'	1:A:1154:G:C8	2.32	0.64
1:A:1504:G:OP1	1:A:1507:A:H4'	1.96	0.64
1:A:1509:C:C2'	1:A:1510:U:H5'	2.28	0.64
2:B:9:GLU:HG3	2:B:217:ARG:HH12	1.61	0.64
2:B:239:VAL:HG12	2:B:240:GLN:N	2.12	0.64
3:C:137:ALA:HA	3:C:140:ARG:HH11	1.62	0.64
13:M:48:LEU:HD13	13:M:53:VAL:HG22	1.79	0.64
17:Q:66:SER:O	17:Q:69:LYS:HB3	1.97	0.64
1:A:608:A:C4	1:A:609:A:C8	2.85	0.64
1:A:1189:C:H5''	3:C:5:ILE:HD13	1.80	0.64
1:A:1195:C:H3'	1:A:1196:U:C5'	2.25	0.64
1:A:1374:A:H2'	1:A:1375:A:O4'	1.97	0.64
1:A:1467:G:H2'	1:A:1468:A:C8	2.32	0.64
3:C:67:THR:HA	3:C:102:ASN:ND2	2.13	0.64
4:D:104:VAL:HG21	4:D:140:VAL:CG2	2.26	0.64
4:D:104:VAL:HG12	4:D:105:VAL:N	2.11	0.64
10:J:9:ARG:NH1	10:J:69:ASN:HB3	2.11	0.64
16:P:21:VAL:HG21	16:P:59:TRP:HE1	1.59	0.64
16:P:72:ARG:O	16:P:75:ARG:HB2	1.97	0.64
1:A:815:A:O2'	1:A:1527:C:H1'	1.97	0.64
1:A:818:G:O2'	1:A:819:A:H5'	1.97	0.64
1:A:879:C:O2'	1:A:880:C:H5'	1.98	0.64
1:A:979:C:H2'	1:A:980:C:O4'	1.98	0.64
1:A:1024:G:H2'	1:A:1024:G:N3	2.13	0.64
1:A:1240:U:H3'	1:A:1241:G:H5'	1.80	0.64
1:A:1418:A:H8	1:A:1418:A:O5'	1.80	0.64
2:B:69:LEU:HB3	2:B:162:ILE:HG12	1.80	0.64
3:C:91:LEU:HG	3:C:99:VAL:HG13	1.78	0.64
8:H:26:VAL:HG13	8:H:59:LEU:HB2	1.79	0.64
12:L:60:LEU:HD21	12:L:85:ILE:HD12	1.80	0.64
16:P:26:ARG:HD2	16:P:31:LYS:O	1.97	0.64
16:P:28:ARG:HG3	16:P:29:ASP:CG	2.18	0.64
18:R:38:GLU:HA	18:R:41:LYS:NZ	2.11	0.64
1:A:89:C:C6	1:A:90:U:H5	2.16	0.64
1:A:346:G:H2'	1:A:347:G:H5'	1.78	0.64
1:A:575:G:OP1	1:A:575:G:H4'	1.97	0.64
1:A:939:G:H5''	7:G:102:ARG:CZ	2.27	0.64
2:B:208:ILE:HA	2:B:211:ILE:HG12	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:11:ARG:NH2	3:C:175:LEU:O	2.29	0.64
4:D:36:ARG:HA	4:D:38:TYR:CE2	2.32	0.64
8:H:11:THR:O	8:H:12:ARG:C	2.36	0.64
9:I:89:ASN:C	9:I:91:ASP:H	2.01	0.64
12:L:113:ARG:HD3	12:L:114:LYS:H	1.62	0.64
15:O:53:HIS:O	15:O:56:LEU:HB3	1.97	0.64
1:A:248:C:O2'	1:A:249:U:H5'	1.98	0.64
1:A:716:A:H1'	11:K:118:GLY:HA2	1.79	0.64
1:A:1178:G:N2	1:A:1180:A:H3'	2.10	0.64
1:A:1201:A:H5'	1:A:1203:C:OP2	1.96	0.64
1:A:1262:C:H2'	1:A:1263:C:H6	1.62	0.64
3:C:48:TYR:CB	3:C:52:LEU:HB3	2.28	0.64
3:C:138:VAL:CG1	3:C:170:GLN:HB2	2.27	0.64
4:D:120:LEU:O	4:D:120:LEU:HD23	1.98	0.64
6:F:53:ALA:C	6:F:55:ASP:H	2.00	0.64
8:H:97:VAL:HG13	8:H:98:LYS:H	1.62	0.64
12:L:69:TYR:O	12:L:100:ILE:HG13	1.98	0.64
18:R:21:LYS:HG3	18:R:55:ARG:HA	1.79	0.64
21:V:10:ARG:HG2	21:V:10:ARG:HH11	1.60	0.64
1:A:673:G:H5''	6:F:87:ARG:NH1	2.13	0.64
1:A:709:G:H2'	1:A:710:G:C8	2.31	0.64
1:A:1088:G:H8	1:A:1088:G:O5'	1.80	0.64
3:C:37:GLN:NE2	14:N:52:GLN:HE22	1.95	0.64
5:E:19:MET:HE3	5:E:20:GLN:H	1.63	0.64
8:H:18:ARG:HB2	8:H:18:ARG:HH11	1.62	0.64
10:J:6:ILE:HG13	10:J:73:ASP:CA	2.27	0.64
12:L:18:VAL:O	12:L:19:ARG:HB2	1.96	0.64
1:A:162:A:C5	1:A:163:C:H1'	2.33	0.64
1:A:254:G:OP1	17:Q:67:LYS:O	2.16	0.64
1:A:592:G:O2'	1:A:593:G:H5'	1.98	0.64
1:A:1014:A:C2	1:A:1219:U:H1'	2.32	0.64
1:A:1110:A:H8	1:A:1110:A:O5'	1.81	0.64
1:A:1305:G:O2'	1:A:1306:A:H8	1.79	0.64
2:B:168:THR:HG1	2:B:192:SER:HA	1.60	0.64
3:C:76:VAL:HG23	3:C:77:ILE:H	1.62	0.64
4:D:9:CYS:SG	4:D:31:CYS:O	2.56	0.64
4:D:201:GLN:HE22	5:E:99:GLY:CA	2.10	0.64
8:H:11:THR:HA	8:H:14:ARG:NH1	2.12	0.64
16:P:35:LYS:O	16:P:36:ILE:HG23	1.98	0.64
19:S:15:LEU:O	19:S:19:VAL:CB	2.43	0.64
1:A:518:C:H4'	1:A:519:C:C5	2.33	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:694:A:C3'	1:A:695:A:H5''	2.28	0.64
1:A:851:G:H2'	1:A:852:G:H8	1.63	0.64
3:C:42:LEU:HA	3:C:46:GLU:OE2	1.98	0.64
8:H:113:SER:HB2	8:H:134:ILE:HD11	1.79	0.64
1:A:448:A:H2'	1:A:449:C:H6	1.63	0.64
1:A:1365:G:O2'	1:A:1366:C:H5'	1.98	0.64
2:B:59:GLU:O	2:B:62:ALA:HB3	1.98	0.64
5:E:64:ARG:O	5:E:65:ASN:HB3	1.98	0.64
8:H:14:ARG:HB3	8:H:14:ARG:HH11	1.63	0.64
13:M:31:LYS:C	13:M:33:ALA:H	1.98	0.64
13:M:42:ALA:O	13:M:43:THR:O	2.15	0.64
1:A:136:C:O2'	16:P:63:GLY:HA2	1.97	0.63
1:A:317:G:C6	1:A:318:G:N7	2.67	0.63
1:A:318:G:H2'	1:A:319:G:H8	1.62	0.63
1:A:1097:C:H2'	1:A:1098:C:C6	2.28	0.63
1:A:1111:A:H2'	1:A:1112:C:O4'	1.97	0.63
1:A:1196:U:H5''	1:A:1197:G:C5'	2.28	0.63
2:B:118:LEU:HB3	2:B:142:LEU:HD12	1.79	0.63
3:C:198:VAL:HG12	3:C:199:LYS:O	1.98	0.63
5:E:60:TYR:O	5:E:64:ARG:HG2	1.98	0.63
5:E:89:ILE:HD13	5:E:90:VAL:H	1.64	0.63
9:I:13:ALA:O	9:I:76:ALA:HB1	1.98	0.63
14:N:50:LYS:HE3	14:N:52:GLN:HG3	1.80	0.63
1:A:261:U:O2	1:A:263:A:C8	2.50	0.63
1:A:639:G:O2'	1:A:640:A:H5'	1.97	0.63
1:A:1091:U:O2	1:A:1093:A:H8	1.81	0.63
1:A:1114:C:H42	1:A:1186:G:H1	1.44	0.63
1:A:1190:G:H3'	3:C:3:ASN:O	1.99	0.63
1:A:1406:U:H1'	1:A:1517:G:H21	1.61	0.63
4:D:79:PHE:HD2	4:D:79:PHE:C	2.02	0.63
16:P:5:ARG:HG2	16:P:6:LEU:N	2.12	0.63
19:S:5:LEU:O	19:S:6:LYS:HB2	1.97	0.63
1:A:200:G:C3'	1:A:201:C:H5''	2.28	0.63
1:A:285:G:O2'	1:A:286:G:H5'	1.97	0.63
1:A:781:A:C5	1:A:802:A:C2	2.86	0.63
1:A:971:G:H4'	1:A:972:C:C5'	2.28	0.63
2:B:19:HIS:CE1	2:B:206:ASP:HB3	2.34	0.63
2:B:36:ARG:HD2	2:B:41:ILE:CD1	2.26	0.63
3:C:8:ILE:CG2	14:N:50:LYS:HB3	2.28	0.63
4:D:187:ARG:NE	4:D:188:LEU:HD12	2.12	0.63
7:G:62:PHE:CD2	7:G:66:VAL:HG21	2.34	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:63:ILE:HD13	9:I:77:ILE:HG21	1.79	0.63
15:O:48:LYS:N	15:O:48:LYS:CD	2.61	0.63
16:P:40:ASP:HB3	16:P:48:TRP:CB	2.28	0.63
1:A:1369:C:H2'	1:A:1370:G:C8	2.34	0.63
1:A:1488:G:H2'	1:A:1489:G:H8	1.63	0.63
3:C:131:ARG:HA	3:C:134:ILE:HD12	1.81	0.63
4:D:58:LEU:O	4:D:61:LYS:HB3	1.99	0.63
6:F:8:ILE:HG22	6:F:10:LEU:HG	1.81	0.63
9:I:7:THR:HB	9:I:83:ARG:HH12	1.63	0.63
9:I:92:TYR:O	9:I:96:LEU:HB2	1.98	0.63
13:M:106:ASN:O	13:M:107:ALA:HB3	1.97	0.63
20:T:51:GLU:HG2	20:T:52:ALA:N	2.12	0.63
1:A:201:C:C3'	1:A:202:U:H5''	2.29	0.63
1:A:372:C:N3	1:A:387:U:H5	1.97	0.63
1:A:941:G:H1	1:A:1342:C:N4	1.97	0.63
1:A:1481:U:O2'	1:A:1482:G:H5'	1.97	0.63
2:B:21:ARG:H	2:B:21:ARG:NE	1.96	0.63
3:C:27:LYS:HA	3:C:30:ARG:NH1	2.12	0.63
9:I:43:ALA:HA	9:I:74:ILE:HG21	1.78	0.63
9:I:46:ALA:C	9:I:81:ILE:HD12	2.19	0.63
15:O:33:THR:HG23	15:O:63:ARG:NH1	2.13	0.63
1:A:112:G:H5'	1:A:389:A:H4'	1.78	0.63
1:A:452:A:O2'	1:A:453:A:H8	1.81	0.63
1:A:950:U:H2'	1:A:951:G:H8	1.63	0.63
1:A:1016:A:H2'	1:A:1017:G:O4'	1.99	0.63
1:A:1057:G:H5'	3:C:155:GLY:HA3	1.81	0.63
2:B:8:LYS:HD2	2:B:9:GLU:H	1.63	0.63
2:B:60:ASP:C	2:B:62:ALA:N	2.51	0.63
2:B:239:VAL:HG12	2:B:240:GLN:H	1.63	0.63
3:C:48:TYR:HD1	3:C:52:LEU:HD22	1.64	0.63
7:G:70:LYS:HE2	7:G:100:ALA:CB	2.29	0.63
10:J:9:ARG:HH12	10:J:69:ASN:CB	2.09	0.63
11:K:66:LEU:HD23	11:K:69:ALA:CB	2.28	0.63
12:L:55:VAL:O	12:L:56:ALA:HB2	1.98	0.63
12:L:84:LEU:CB	12:L:101:VAL:HG21	2.25	0.63
13:M:14:ARG:HB3	13:M:14:ARG:HH11	1.60	0.63
16:P:64:ALA:O	16:P:66:PRO:CD	2.47	0.63
1:A:361:G:H2'	1:A:362:G:H5'	1.80	0.63
1:A:731:G:OP1	1:A:766:A:H1'	1.99	0.63
2:B:129:GLU:HB3	2:B:130:ARG:HD2	1.81	0.63
3:C:104:GLN:HA	3:C:104:GLN:NE2	2.14	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:126:ARG:C	3:C:128:PHE:N	2.52	0.63
4:D:21:LEU:HD21	4:D:66:ARG:O	1.98	0.63
4:D:96:LEU:N	4:D:96:LEU:HD12	2.14	0.63
5:E:80:ILE:CD1	5:E:138:ALA:HB1	2.28	0.63
6:F:91:VAL:HG11	18:R:72:ARG:NH1	2.14	0.63
20:T:13:LEU:O	20:T:14:LYS:C	2.37	0.63
20:T:30:LYS:NZ	20:T:80:ARG:NH2	2.46	0.63
1:A:371:G:H2'	1:A:372:C:C6	2.34	0.63
1:A:585:G:N3	1:A:879:C:H4'	2.14	0.63
1:A:695:A:H2	1:A:787:A:C1'	2.11	0.63
1:A:811:C:O2'	1:A:901:A:N1	2.30	0.63
1:A:1228:C:H2'	1:A:1229:A:C8	2.32	0.63
3:C:22:TRP:HB3	3:C:59:ARG:HG2	1.81	0.63
4:D:15:GLU:HG3	4:D:63:LYS:HZ2	1.63	0.63
4:D:157:LEU:HB2	4:D:158:ILE:CD1	2.27	0.63
5:E:51:VAL:HB	5:E:52:PRO:CD	2.25	0.63
7:G:115:ARG:HE	7:G:118:VAL:CG2	2.12	0.63
7:G:138:LYS:HD3	7:G:139:GLU:HG3	1.81	0.63
11:K:83:ILE:O	11:K:83:ILE:HG22	1.97	0.63
13:M:87:TYR:CD1	13:M:90:LEU:HD12	2.34	0.63
16:P:47:ASP:C	16:P:49:LEU:H	2.02	0.63
1:A:455:C:O2'	1:A:456:C:H5'	1.99	0.63
1:A:542:G:OP1	4:D:10:ARG:NH2	2.31	0.63
1:A:621:A:C5	1:A:622:A:N7	2.66	0.63
3:C:116:VAL:CA	3:C:119:ARG:HB3	2.29	0.63
8:H:5:PRO:O	8:H:8:ASP:N	2.31	0.63
9:I:85:LEU:HB3	9:I:92:TYR:HD1	1.62	0.63
13:M:48:LEU:HD22	13:M:53:VAL:CG2	2.29	0.63
13:M:74:VAL:O	13:M:76:ALA:N	2.32	0.63
14:N:40:CYS:HB3	14:N:43:CYS:SG	2.38	0.63
18:R:70:ILE:HG22	18:R:74:ARG:HD2	1.79	0.63
1:A:356:A:H2'	1:A:357:G:H8	1.63	0.62
2:B:61:LEU:CD2	2:B:66:GLY:HA3	2.28	0.62
2:B:140:HIS:O	2:B:144:ARG:HB3	1.99	0.62
3:C:21:ARG:HD3	3:C:57:ILE:O	1.99	0.62
3:C:118:GLN:O	3:C:121:ALA:HB3	1.99	0.62
4:D:190:ASP:O	4:D:191:ARG:C	2.36	0.62
5:E:105:VAL:HG21	5:E:128:PRO:HA	1.80	0.62
7:G:22:LEU:O	7:G:25:ALA:HB3	1.98	0.62
11:K:67:ASP:C	11:K:69:ALA:H	2.02	0.62
13:M:88:ARG:HA	13:M:98:VAL:CG1	2.29	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:G:H1	1:A:554:C:H42	1.47	0.62
1:A:386:C:C2'	1:A:387:U:H5'	2.29	0.62
1:A:589:C:O2'	1:A:590:C:H5'	1.99	0.62
1:A:637:G:H2'	1:A:638:G:O4'	1.99	0.62
1:A:877:C:H5''	8:H:88:LYS:HD3	1.80	0.62
1:A:1004:A:H5''	1:A:1025:U:C5	2.34	0.62
3:C:10:PHE:CE2	3:C:178:LEU:HB2	2.34	0.62
4:D:10:ARG:HG2	4:D:11:LEU:H	1.61	0.62
8:H:111:ILE:HD12	8:H:135:CYS:SG	2.38	0.62
9:I:10:ARG:NE	9:I:11:LYS:HB2	2.14	0.62
9:I:108:VAL:CG1	9:I:109:VAL:H	2.04	0.62
10:J:47:PHE:CE2	14:N:37:PHE:HZ	2.17	0.62
11:K:33:THR:CA	11:K:39:PRO:HA	2.27	0.62
1:A:667:G:O2'	15:O:49:ASP:HA	1.99	0.62
1:A:1520:G:H2'	1:A:1521:G:C8	2.34	0.62
2:B:16:HIS:NE2	2:B:214:ILE:HD11	2.13	0.62
2:B:208:ILE:C	2:B:210:SER:H	2.02	0.62
3:C:64:VAL:HG21	3:C:97:LYS:HB2	1.82	0.62
4:D:62:GLN:HA	4:D:62:GLN:NE2	2.01	0.62
5:E:137:GLU:O	5:E:138:ALA:C	2.38	0.62
6:F:71:ARG:O	6:F:73:ASN:N	2.31	0.62
9:I:17:VAL:HG13	9:I:63:ILE:HG12	1.81	0.62
9:I:28:VAL:HG22	9:I:63:ILE:HG22	1.82	0.62
10:J:3:LYS:N	10:J:76:ASN:H	1.98	0.62
12:L:7:ILE:O	12:L:10:LEU:HB2	1.98	0.62
13:M:25:ILE:HD11	13:M:66:LEU:HD21	1.82	0.62
15:O:75:PRO:O	15:O:78:TYR:HB3	2.00	0.62
19:S:27:GLU:HB3	19:S:47:HIS:CD2	2.34	0.62
21:V:10:ARG:HG2	21:V:10:ARG:NH1	2.14	0.62
1:A:370:C:O2'	1:A:371:G:H5'	1.98	0.62
1:A:740:U:O2'	1:A:741:G:H5'	1.99	0.62
1:A:1225:A:H5'	13:M:103:THR:CG2	2.29	0.62
1:A:1237:C:O2'	1:A:1335:C:H5'	2.00	0.62
2:B:131:PRO:O	2:B:133:LYS:N	2.32	0.62
2:B:175:ARG:HB2	2:B:175:ARG:HH11	1.61	0.62
7:G:24:THR:HG22	7:G:28:ASN:HD21	1.64	0.62
9:I:80:GLY:HA2	9:I:83:ARG:CB	2.26	0.62
10:J:46:ARG:HG2	10:J:46:ARG:HH11	1.64	0.62
16:P:41:PRO:O	16:P:43:LYS:HG3	1.97	0.62
1:A:728:A:O2'	1:A:729:A:H5'	1.98	0.62
1:A:1013:G:H1'	1:A:1016:A:H62	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1050:G:N2	1:A:1209:C:H1'	2.12	0.62
1:A:1081:G:H2'	1:A:1082:G:O4'	1.99	0.62
1:A:1168:A:H2'	1:A:1169:A:C8	2.34	0.62
1:A:1337:G:H5''	1:A:1338:G:OP1	1.99	0.62
2:B:178:ARG:HH11	2:B:178:ARG:HG3	1.63	0.62
3:C:203:PHE:CZ	3:C:206:GLU:HB2	2.35	0.62
5:E:51:VAL:O	5:E:54:ALA:HB3	2.00	0.62
5:E:111:GLU:O	5:E:112:LEU:C	2.37	0.62
14:N:46:GLU:CD	14:N:47:LEU:HD23	2.20	0.62
15:O:32:LEU:HD22	15:O:63:ARG:HD3	1.82	0.62
16:P:48:TRP:CE3	16:P:49:LEU:HB2	2.35	0.62
1:A:106:C:O2	1:A:379:C:H4'	2.00	0.62
1:A:148:G:H2'	1:A:149:A:C8	2.32	0.62
1:A:185:A:H2'	1:A:186:C:H6	1.62	0.62
1:A:369:C:O2'	1:A:370:C:H5'	2.00	0.62
1:A:795:C:H5''	1:A:796:C:OP2	2.00	0.62
1:A:1102:A:H2'	1:A:1103:C:H6	1.64	0.62
1:A:1207:G:H2'	1:A:1208:C:C6	2.34	0.62
1:A:1237:C:H3'	1:A:1238:A:H5'	1.81	0.62
4:D:33:MET:O	4:D:37:PRO:HG3	2.00	0.62
4:D:134:ASP:O	4:D:136:PRO:N	2.33	0.62
5:E:79:GLU:O	5:E:80:ILE:HG23	1.99	0.62
5:E:91:LEU:HD12	5:E:138:ALA:HB1	1.80	0.62
6:F:30:LEU:HD23	6:F:35:ALA:HB3	1.81	0.62
7:G:14:PRO:HA	7:G:21:VAL:HG12	1.80	0.62
10:J:19:SER:O	10:J:91:PRO:HG3	1.99	0.62
12:L:84:LEU:HB3	12:L:101:VAL:HG22	1.78	0.62
20:T:96:GLY:O	20:T:97:ALA:HB3	1.99	0.62
1:A:106:C:C2'	1:A:107:G:H5'	2.29	0.62
1:A:372:C:H1'	1:A:373:A:OP2	1.99	0.62
1:A:1189:C:H5''	3:C:5:ILE:HG21	1.80	0.62
1:A:1210:C:H2'	1:A:1211:U:C4'	2.29	0.62
1:A:1494:G:C2'	1:A:1495:U:H5'	2.29	0.62
2:B:112:VAL:HG11	2:B:153:ARG:HA	1.82	0.62
2:B:187:LEU:HD22	2:B:201:ILE:O	2.00	0.62
4:D:10:ARG:HG3	4:D:10:ARG:NH1	1.98	0.62
4:D:176:LEU:H	4:D:176:LEU:HD23	1.64	0.62
4:D:187:ARG:HE	4:D:188:LEU:CD1	2.11	0.62
8:H:4:ASP:OD2	8:H:85:ARG:NH1	2.33	0.62
8:H:136:GLU:O	8:H:136:GLU:HG2	1.99	0.62
9:I:19:LEU:HD23	9:I:21:PRO:HD3	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:90:LEU:N	10:J:91:PRO:HD2	2.15	0.62
12:L:6:THR:O	12:L:7:ILE:C	2.36	0.62
13:M:5:ALA:HB3	13:M:22:ILE:HD13	1.82	0.62
13:M:34:LEU:HD12	13:M:39:ILE:O	1.99	0.62
17:Q:51:TYR:CG	17:Q:73:VAL:HG11	2.34	0.62
18:R:44:LEU:HD12	18:R:48:GLY:O	2.00	0.62
20:T:11:SER:O	20:T:13:LEU:HG	1.99	0.62
1:A:429:U:O4'	1:A:430:A:H5''	2.00	0.62
1:A:549:C:H2'	1:A:550:G:H8	1.65	0.62
1:A:551:U:H2'	1:A:552:U:C6	2.34	0.62
1:A:836:G:H2'	1:A:837:G:H8	1.65	0.62
1:A:945:G:N2	1:A:1334:G:H4'	2.13	0.62
1:A:1314:C:H2'	1:A:1315:U:C6	2.35	0.62
1:A:1323:G:H2'	1:A:1324:A:C8	2.34	0.62
3:C:205:GLY:O	3:C:206:GLU:HB3	2.00	0.62
8:H:17:THR:CG2	8:H:63:LEU:HD23	2.29	0.62
10:J:7:LYS:HG2	10:J:71:LEU:HD21	1.81	0.62
10:J:78:ASN:O	10:J:80:LYS:N	2.32	0.62
12:L:57:LYS:HE2	12:L:67:THR:HG22	1.82	0.62
18:R:64:ARG:O	18:R:67:ALA:HB3	1.99	0.62
1:A:287:U:C2'	1:A:288:A:H5'	2.29	0.62
1:A:399:G:H2'	1:A:400:C:C6	2.35	0.62
1:A:1061:G:O4'	10:J:56:HIS:ND1	2.32	0.62
1:A:1281:U:H5'	1:A:1282:C:C5	2.34	0.62
2:B:112:VAL:HG13	2:B:153:ARG:HG2	1.81	0.62
3:C:156:ARG:HH21	3:C:161:GLU:CA	2.10	0.62
4:D:79:PHE:C	4:D:79:PHE:CD2	2.73	0.62
10:J:15:THR:O	10:J:19:SER:N	2.23	0.62
11:K:18:ARG:HD2	11:K:34:ASP:O	1.99	0.62
12:L:39:VAL:HG12	12:L:41:ARG:H	1.65	0.62
13:M:89:GLY:O	13:M:93:ARG:HG3	1.99	0.62
14:N:24:CYS:O	14:N:28:GLY:HA2	2.00	0.62
19:S:15:LEU:HD13	19:S:44:MET:CE	2.30	0.62
1:A:56:U:H2'	1:A:57:G:H8	1.62	0.62
1:A:325:A:N6	1:A:326:G:N1	2.48	0.62
1:A:573:A:H2'	1:A:574:A:O4'	2.00	0.62
1:A:1088:G:H2'	1:A:1089:G:C8	2.34	0.62
1:A:1269:A:H5'	21:V:18:TYR:O	1.99	0.62
1:A:1424:C:H42	1:A:1476:G:H1	1.47	0.62
7:G:46:ALA:HB1	7:G:121:ALA:H	1.62	0.62
7:G:141:VAL:O	7:G:142:GLU:C	2.38	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:46:ALA:HB1	9:I:81:ILE:CD1	2.30	0.62
10:J:8:LEU:HD12	10:J:70:ARG:O	1.99	0.62
11:K:89:ALA:O	11:K:91:ARG:N	2.33	0.62
13:M:14:ARG:HH11	13:M:14:ARG:CB	2.12	0.62
19:S:22:LEU:HD22	19:S:28:LYS:HB2	1.82	0.62
20:T:51:GLU:O	20:T:55:ILE:HG12	2.00	0.62
1:A:273:A:O2'	1:A:274:A:H5'	2.00	0.61
1:A:745:C:O2'	1:A:746:A:H5'	2.00	0.61
1:A:1150:U:H4'	10:J:41:PRO:CD	2.29	0.61
1:A:1311:G:N7	19:S:2:PRO:HD2	2.15	0.61
1:A:1374:A:C4	1:A:1375:A:C8	2.88	0.61
2:B:92:TYR:CD1	2:B:151:GLY:HA3	2.34	0.61
4:D:130:GLY:C	4:D:132:ARG:H	2.04	0.61
7:G:17:VAL:CG1	7:G:18:TYR:H	2.10	0.61
7:G:135:VAL:O	7:G:139:GLU:HG3	2.00	0.61
11:K:84:VAL:HG22	11:K:110:ASP:HA	1.81	0.61
12:L:11:VAL:HG13	17:Q:29:HIS:CD2	2.31	0.61
12:L:56:ALA:HB2	12:L:70:ILE:HD11	1.81	0.61
13:M:78:ILE:HD12	13:M:78:ILE:N	2.14	0.61
13:M:92:HIS:HA	13:M:110:ARG:NH2	2.14	0.61
19:S:15:LEU:HD11	19:S:38:SER:OG	2.00	0.61
1:A:102:G:H2'	1:A:103:C:H6	1.65	0.61
1:A:411:A:H2'	1:A:413:G:C1'	2.30	0.61
1:A:451:A:O5'	1:A:451:A:H8	1.83	0.61
1:A:1463:C:O2'	1:A:1464:G:H5'	1.99	0.61
1:A:1514:C:H2'	1:A:1515:C:C6	2.35	0.61
2:B:54:THR:HA	2:B:199:TYR:HB3	1.82	0.61
2:B:117:GLU:O	2:B:117:GLU:HG2	2.01	0.61
4:D:4:TYR:O	4:D:5:ILE:HB	2.01	0.61
4:D:120:LEU:HD22	4:D:126:ILE:HD11	1.82	0.61
5:E:71:LEU:HD13	5:E:114:GLY:O	2.00	0.61
7:G:72:ARG:HE	7:G:142:GLU:HB3	1.64	0.61
8:H:44:PHE:O	8:H:45:ILE:CG2	2.40	0.61
20:T:53:LEU:HD11	20:T:101:GLY:O	1.99	0.61
1:A:44:G:OP2	16:P:12:LYS:HB2	2.00	0.61
1:A:344:A:H5''	1:A:345:C:C5	2.36	0.61
1:A:421:U:H3'	1:A:422:C:H5'	1.81	0.61
1:A:529:G:H2'	1:A:530:G:H5'	1.82	0.61
2:B:22:LYS:HE3	2:B:40:HIS:NE2	2.15	0.61
4:D:54:TYR:HE1	4:D:206:PHE:HE1	1.46	0.61
4:D:124:GLY:HA3	4:D:132:ARG:HH21	1.64	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:139:LEU:HA	5:E:142:LEU:CD1	2.31	0.61
9:I:33:PHE:CD2	9:I:34:ASN:ND2	2.67	0.61
12:L:24:VAL:C	12:L:26:ALA:H	2.02	0.61
15:O:20:GLY:O	15:O:22:THR:HG22	2.00	0.61
20:T:40:ALA:CA	20:T:55:ILE:HD11	2.25	0.61
1:A:418:C:O2	1:A:418:C:H2'	2.01	0.61
1:A:445:G:H2'	1:A:446:G:H8	1.63	0.61
1:A:1128:C:O2'	1:A:1129:C:H5''	2.00	0.61
1:A:1465:C:H2'	1:A:1466:C:C6	2.36	0.61
2:B:139:LYS:HD2	2:B:143:GLU:OE2	1.99	0.61
2:B:208:ILE:C	2:B:210:SER:N	2.51	0.61
3:C:87:LEU:O	3:C:91:LEU:N	2.32	0.61
4:D:101:LEU:CB	4:D:138:TYR:HB3	2.30	0.61
4:D:176:LEU:HA	4:D:183:GLY:HA2	1.82	0.61
6:F:99:ALA:O	18:R:28:GLU:HA	2.01	0.61
9:I:47:LEU:C	9:I:49:PRO:HD2	2.20	0.61
9:I:103:THR:HG22	9:I:104:ARG:O	2.00	0.61
11:K:67:ASP:O	11:K:71:LYS:HG3	2.01	0.61
11:K:115:PRO:C	11:K:117:ASN:H	2.02	0.61
13:M:10:PRO:HB2	13:M:18:ALA:CB	2.27	0.61
16:P:74:LEU:HD22	16:P:79:VAL:HG21	1.82	0.61
19:S:22:LEU:O	19:S:25:LYS:HD2	1.99	0.61
1:A:112:G:C2	1:A:113:G:C8	2.88	0.61
1:A:235:C:H1'	17:Q:61:GLU:OE1	1.99	0.61
1:A:292:G:H3'	1:A:293:G:C8	2.36	0.61
1:A:299:G:H2'	1:A:300:A:C8	2.35	0.61
1:A:988:G:H2'	1:A:989:C:O4'	2.00	0.61
1:A:1178:G:N2	1:A:1181:G:OP2	2.33	0.61
2:B:118:LEU:HB3	2:B:142:LEU:CD1	2.31	0.61
2:B:177:ALA:O	2:B:180:LEU:N	2.32	0.61
4:D:47:ARG:HE	4:D:49:ARG:HA	1.64	0.61
7:G:46:ALA:CB	7:G:120:ILE:HB	2.30	0.61
13:M:4:ILE:HG23	13:M:57:ARG:HA	1.83	0.61
15:O:48:LYS:H	15:O:48:LYS:CE	2.13	0.61
19:S:22:LEU:CB	19:S:28:LYS:HB2	2.25	0.61
19:S:49:ILE:HG12	19:S:71:LEU:HD22	1.81	0.61
20:T:30:LYS:HZ3	20:T:80:ARG:HH22	1.48	0.61
20:T:56:MET:CE	20:T:88:VAL:HG11	2.30	0.61
20:T:70:SER:HA	20:T:73:HIS:HD2	1.65	0.61
1:A:36:C:H2'	1:A:37:U:C6	2.35	0.61
1:A:619:U:N3	4:D:135:LEU:HD21	2.16	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1015:A:H2'	1:A:1016:A:O4'	2.01	0.61
1:A:1370:G:C2	1:A:1371:G:N7	2.68	0.61
2:B:54:THR:HG23	2:B:199:TYR:CB	2.31	0.61
3:C:88:ARG:HA	3:C:91:LEU:CB	2.28	0.61
3:C:191:THR:HG23	3:C:194:GLY:N	2.15	0.61
12:L:33:ARG:CG	12:L:60:LEU:HD12	2.31	0.61
18:R:87:ARG:HG3	18:R:87:ARG:NH1	2.14	0.61
1:A:142:G:O2'	1:A:196:A:N1	2.26	0.61
1:A:367:U:O2	1:A:369:C:C5	2.53	0.61
1:A:620:C:C6	4:D:135:LEU:HD13	2.35	0.61
1:A:695:A:C2	1:A:787:A:H1'	2.35	0.61
1:A:927:G:N2	1:A:1391:U:H1'	2.16	0.61
1:A:1201:A:H4'	1:A:1202:G:C5'	2.31	0.61
6:F:33:TYR:HA	6:F:71:ARG:HH21	1.63	0.61
11:K:17:GLY:O	11:K:80:VAL:HA	2.00	0.61
13:M:48:LEU:HD22	13:M:53:VAL:HG23	1.83	0.61
1:A:93:G:O2'	1:A:95:U:H5'	2.00	0.61
1:A:572:A:H5''	1:A:917:G:H4'	1.81	0.61
1:A:1209:C:O2	1:A:1209:C:H2'	2.01	0.61
2:B:71:VAL:CG2	2:B:164:VAL:HG22	2.31	0.61
2:B:134:GLU:C	2:B:136:VAL:H	2.03	0.61
8:H:5:PRO:O	8:H:6:ILE:C	2.38	0.61
8:H:36:LEU:O	8:H:40:ALA:N	2.32	0.61
11:K:30:VAL:HG21	11:K:65:ALA:HB2	1.83	0.61
16:P:50:LYS:HG2	16:P:51:VAL:N	2.16	0.61
16:P:82:GLN:O	16:P:82:GLN:HG2	2.01	0.61
20:T:50:GLU:O	20:T:52:ALA:N	2.34	0.61
1:A:265:G:H2'	1:A:267:C:C5	2.36	0.61
1:A:375:U:H4'	16:P:17:TYR:HE2	1.65	0.61
1:A:385:C:O2'	1:A:386:C:H5'	2.01	0.61
1:A:514:C:N4	1:A:537:G:H1	1.99	0.61
1:A:761:G:H5'	17:Q:102:GLY:HA3	1.82	0.61
1:A:950:U:H5	13:M:102:ARG:NH2	1.99	0.61
1:A:963:G:H1	1:A:972:C:H42	1.47	0.61
2:B:143:GLU:HA	2:B:146:GLN:OE1	2.01	0.61
3:C:74:GLY:C	3:C:76:VAL:H	2.03	0.61
3:C:95:THR:HG23	3:C:98:ASN:HA	1.81	0.61
3:C:113:ALA:HB3	3:C:184:TYR:O	2.01	0.61
3:C:175:LEU:O	3:C:177:THR:N	2.34	0.61
4:D:156:GLU:CG	4:D:157:LEU:N	2.63	0.61
6:F:30:LEU:HD23	6:F:30:LEU:C	2.21	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:111:ARG:HH11	7:G:111:ARG:CB	2.08	0.61
10:J:79:ARG:HH11	10:J:79:ARG:HB2	1.66	0.61
13:M:53:VAL:HA	13:M:56:LEU:HD11	1.83	0.61
16:P:6:LEU:N	16:P:6:LEU:HD12	2.15	0.61
1:A:192:U:H5'	20:T:102:GLY:HA2	1.83	0.61
1:A:665:A:C2	1:A:732:C:C2	2.89	0.61
1:A:1393:U:O2'	1:A:1394:A:H2'	2.00	0.61
2:B:28:PHE:CE1	2:B:189:ASP:HA	2.35	0.61
2:B:132:LYS:C	2:B:134:GLU:H	2.04	0.61
3:C:138:VAL:HG11	3:C:170:GLN:HB2	1.83	0.61
5:E:82:VAL:HG21	5:E:138:ALA:N	2.15	0.61
6:F:3:ARG:HA	6:F:66:GLU:HG3	1.83	0.61
12:L:41:ARG:HH11	12:L:41:ARG:CB	2.08	0.61
18:R:43:PHE:CA	18:R:51:LEU:HD12	2.28	0.61
20:T:11:SER:C	20:T:13:LEU:H	2.04	0.61
1:A:413:G:H2'	1:A:413:G:N3	2.16	0.60
1:A:1128:C:H42	1:A:1143:G:H1	1.49	0.60
1:A:1424:C:H2'	1:A:1425:U:O4'	2.00	0.60
1:A:1476:G:H2'	1:A:1477:C:C6	2.36	0.60
14:N:56:VAL:HG13	14:N:57:ARG:N	2.14	0.60
1:A:255:G:N2	1:A:272:C:H1'	2.16	0.60
1:A:1421:G:H2'	1:A:1422:G:C8	2.37	0.60
1:A:1425:U:O2'	1:A:1426:C:H5'	2.01	0.60
2:B:71:VAL:HG23	2:B:164:VAL:CA	2.31	0.60
3:C:118:GLN:O	3:C:122:GLU:HG3	2.00	0.60
4:D:151:LYS:H	4:D:151:LYS:HD2	1.66	0.60
9:I:31:GLN:HB3	9:I:35:GLU:HB3	1.83	0.60
10:J:31:GLY:HA3	10:J:76:ASN:O	2.01	0.60
14:N:7:ILE:O	14:N:7:ILE:CG2	2.48	0.60
15:O:16:ALA:HB2	15:O:21:ASP:O	2.01	0.60
17:Q:67:LYS:HA	17:Q:70:ARG:NH1	2.16	0.60
1:A:219:C:C2'	1:A:220:G:H5'	2.30	0.60
1:A:378:G:C2	1:A:386:C:O2	2.55	0.60
1:A:1151:A:H5''	10:J:42:THR:N	2.16	0.60
6:F:27:GLN:O	6:F:30:LEU:HB3	2.01	0.60
10:J:42:THR:CG2	10:J:68:HIS:HA	2.30	0.60
11:K:89:ALA:O	11:K:90:GLY:C	2.39	0.60
12:L:74:GLY:O	12:L:110:VAL:HG13	2.02	0.60
16:P:5:ARG:CZ	16:P:22:THR:HG21	2.30	0.60
1:A:129(A):G:O2'	1:A:130:A:P	2.60	0.60
1:A:549:C:H2'	1:A:550:G:C8	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:946:A:C4	1:A:947:G:N7	2.69	0.60
1:A:1137:C:O2'	1:A:1138:G:H5''	2.01	0.60
1:A:1494:G:O2'	1:A:1495:U:H5'	2.01	0.60
2:B:76:GLN:HG3	2:B:206:ASP:OD2	2.01	0.60
3:C:5:ILE:HD12	3:C:10:PHE:HB2	1.82	0.60
4:D:76:ARG:HG3	4:D:77:ASN:N	2.15	0.60
7:G:75:VAL:HG21	7:G:148:ASN:HD22	1.66	0.60
10:J:36:GLY:O	10:J:38:ILE:HG13	1.99	0.60
11:K:82:VAL:HG22	11:K:98:LEU:HD12	1.83	0.60
1:A:229:U:O2'	1:A:230:G:H5'	2.00	0.60
1:A:230:G:C4	1:A:231:G:C8	2.90	0.60
1:A:463:A:H4'	16:P:82:GLN:HE22	1.66	0.60
1:A:486:U:H2'	1:A:487:A:H8	1.65	0.60
1:A:750:G:H1'	15:O:22:THR:HG23	1.81	0.60
1:A:854:G:H3'	1:A:871:U:O4	2.00	0.60
1:A:1063:C:H3'	1:A:1064:G:H2'	1.83	0.60
1:A:1137:C:H4'	1:A:1138:G:C2	2.36	0.60
1:A:1154:G:H2'	1:A:1155:G:C8	2.31	0.60
2:B:55:PHE:HD1	2:B:58:ILE:HD12	1.65	0.60
4:D:25:ARG:CD	4:D:30:LYS:HB3	2.31	0.60
4:D:201:GLN:NE2	5:E:99:GLY:HA2	2.17	0.60
6:F:22:GLU:O	6:F:24:GLU:N	2.34	0.60
9:I:77:ILE:HG22	9:I:81:ILE:CD1	2.31	0.60
9:I:104:ARG:HE	9:I:105:ASP:H	1.49	0.60
11:K:82:VAL:HG12	11:K:83:ILE:N	2.15	0.60
12:L:109:GLY:HA3	12:L:120:TYR:O	2.02	0.60
19:S:69:HIS:HB3	19:S:73:GLU:OE1	2.01	0.60
20:T:36:LEU:O	20:T:39:LYS:HB3	2.02	0.60
20:T:77:ALA:O	20:T:81:LYS:HG3	2.01	0.60
1:A:263:A:OP2	20:T:79:ARG:NH1	2.35	0.60
1:A:770:C:O2'	1:A:771:G:H5'	2.01	0.60
2:B:168:THR:OG1	2:B:192:SER:CA	2.49	0.60
3:C:135:LYS:HZ3	5:E:52:PRO:HG2	1.66	0.60
8:H:126:LYS:C	8:H:128:GLY:H	2.05	0.60
8:H:137:VAL:HG12	8:H:138:TRP:N	2.15	0.60
12:L:78:GLN:HG2	12:L:81:SER:OG	2.00	0.60
13:M:49:THR:CG2	13:M:50:GLU:H	2.14	0.60
13:M:84:ILE:O	13:M:84:ILE:HG13	2.00	0.60
16:P:24:ALA:C	16:P:26:ARG:H	2.03	0.60
19:S:74:PHE:O	19:S:76:PRO:HD3	2.02	0.60
1:A:153:C:N4	1:A:169:C:N4	2.49	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:U:H2'	1:A:181:G:H5'	1.84	0.60
1:A:229:U:O2'	16:P:23:ASP:OD2	2.19	0.60
1:A:944:G:H2'	1:A:945:G:H5''	1.82	0.60
1:A:1461:G:H2'	1:A:1462:G:C8	2.36	0.60
2:B:102:LEU:O	2:B:105:PHE:HB2	2.01	0.60
3:C:182:ILE:CG2	3:C:183:ASP:N	2.51	0.60
5:E:55:VAL:O	5:E:58:ALA:N	2.29	0.60
6:F:62:TRP:C	6:F:63:TYR:HD1	2.04	0.60
8:H:137:VAL:CG1	8:H:138:TRP:N	2.64	0.60
9:I:126:SER:HB2	9:I:127:LYS:NZ	2.17	0.60
16:P:7:ALA:CB	16:P:28:ARG:O	2.50	0.60
20:T:30:LYS:HZ2	20:T:80:ARG:HH22	1.49	0.60
1:A:23:C:H2'	1:A:24:U:H6	1.67	0.60
1:A:260:G:H2'	1:A:261:U:H6	1.67	0.60
1:A:575:G:C6	1:A:821:G:N7	2.70	0.60
1:A:615:C:H2'	1:A:616:G:C5'	2.32	0.60
1:A:1226:C:H5''	13:M:103:THR:HG1	1.67	0.60
1:A:1281:U:H5'	1:A:1282:C:H5	1.67	0.60
2:B:9:GLU:CG	2:B:217:ARG:HH12	2.13	0.60
2:B:44:LEU:HD23	2:B:44:LEU:N	2.06	0.60
3:C:139:GLN:CA	3:C:139:GLN:HE21	2.15	0.60
4:D:80:GLU:OE2	4:D:80:GLU:CA	2.50	0.60
8:H:96:GLY:O	8:H:98:LYS:N	2.35	0.60
9:I:106:ALA:O	9:I:108:VAL:HG23	2.02	0.60
10:J:8:LEU:CD1	10:J:70:ARG:HB2	2.32	0.60
10:J:61:GLU:OE2	14:N:58:LYS:HD3	2.02	0.60
17:Q:104:LYS:HG3	17:Q:105:ALA:H	1.65	0.60
18:R:59:SER:HB3	18:R:62:GLU:OE1	2.01	0.60
1:A:44:G:C2	1:A:45:U:H1'	2.36	0.60
1:A:230:G:H2'	1:A:231:G:O4'	2.02	0.60
1:A:448:A:H62	1:A:486:U:H3	1.48	0.60
1:A:544:G:C5	1:A:545:C:C5	2.90	0.60
1:A:1046:A:H3'	1:A:1047:G:C8	2.34	0.60
1:A:1152:A:H5''	10:J:13:HIS:O	2.01	0.60
1:A:1403:C:O2'	1:A:1404:C:H5'	2.00	0.60
2:B:21:ARG:HH12	2:B:23:ARG:HH22	1.50	0.60
2:B:160:ASP:O	2:B:161:ALA:HB2	2.00	0.60
3:C:12:LEU:HD22	14:N:50:LYS:O	2.02	0.60
9:I:10:ARG:HD3	9:I:105:ASP:HB3	1.84	0.60
9:I:21:PRO:HA	9:I:59:PHE:CA	2.32	0.60
14:N:23:ARG:HA	14:N:29:ARG:O	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:S:22:LEU:HB3	19:S:28:LYS:CB	2.27	0.60
20:T:33:ILE:HG13	20:T:62:LEU:HD22	1.84	0.60
1:A:1515:C:O2'	1:A:1516:G:H5'	2.01	0.60
2:B:52:GLU:O	2:B:56:ARG:HG3	2.02	0.60
3:C:190:ARG:CA	3:C:195:VAL:HG22	2.27	0.60
4:D:36:ARG:HG3	4:D:38:TYR:CZ	2.37	0.60
9:I:19:LEU:CD1	9:I:61:ALA:HB2	2.32	0.60
11:K:15:ALA:HB1	11:K:78:GLN:HB2	1.83	0.60
11:K:86:GLY:HA2	11:K:112:THR:HG23	1.84	0.60
11:K:104:GLN:OE1	11:K:104:GLN:HA	2.01	0.60
14:N:46:GLU:OE1	14:N:47:LEU:HD23	2.02	0.60
1:A:371:G:H21	1:A:374:A:N6	2.00	0.59
1:A:623:C:H2'	1:A:624:C:C6	2.37	0.59
1:A:760:G:N2	17:Q:104:LYS:N	2.47	0.59
1:A:1306:A:H2'	1:A:1307:U:O4'	2.02	0.59
4:D:206:PHE:HD2	4:D:207:TYR:CE2	2.19	0.59
7:G:24:THR:HG22	7:G:28:ASN:ND2	2.17	0.59
1:A:429:U:H4'	1:A:430:A:O5'	2.01	0.59
1:A:1182:G:O2'	1:A:1183:A:OP2	2.19	0.59
3:C:39:ILE:HD11	3:C:57:ILE:HD11	1.84	0.59
3:C:198:VAL:HG12	3:C:199:LYS:N	2.17	0.59
5:E:91:LEU:HD23	5:E:120:THR:HG22	1.83	0.59
6:F:30:LEU:HD23	6:F:31:GLU:N	2.16	0.59
7:G:22:LEU:C	7:G:22:LEU:HD12	2.22	0.59
7:G:91:VAL:HG12	7:G:92:SER:O	2.02	0.59
8:H:34:GLU:OE2	8:H:34:GLU:HA	2.02	0.59
8:H:40:ALA:HA	8:H:45:ILE:HG12	1.84	0.59
9:I:21:PRO:HA	9:I:59:PHE:C	2.23	0.59
13:M:4:ILE:HG22	13:M:5:ALA:N	2.16	0.59
14:N:2:ALA:HA	14:N:27:CYS:O	2.01	0.59
16:P:31:LYS:HG2	16:P:32:TYR:N	2.17	0.59
17:Q:90:ILE:O	17:Q:93:GLN:N	2.34	0.59
1:A:162:A:C4	1:A:163:C:H1'	2.38	0.59
1:A:190:C:H2'	1:A:190(A):C:H6	1.66	0.59
1:A:222:U:H2'	1:A:223:U:C5	2.38	0.59
1:A:502:G:H1'	1:A:550:G:H5'	1.84	0.59
1:A:1130:A:H5''	9:I:20:ARG:CD	2.31	0.59
1:A:1151:A:P	10:J:41:PRO:HA	2.42	0.59
1:A:1184:G:H2'	1:A:1185:G:C8	2.37	0.59
1:A:1355:G:O2'	1:A:1356:G:H5'	2.01	0.59
1:A:1436:U:C4	1:A:1437:C:N3	2.70	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:55:ALA:O	4:D:58:LEU:HB3	2.01	0.59
5:E:13:ILE:HD12	5:E:13:ILE:O	2.02	0.59
7:G:60:LYS:HA	7:G:63:LYS:HB2	1.82	0.59
9:I:11:LYS:O	9:I:12:GLU:CB	2.50	0.59
12:L:46:LYS:HE3	12:L:47:LYS:CE	2.32	0.59
19:S:19:VAL:HG22	19:S:47:HIS:CE1	2.37	0.59
1:A:377:G:OP1	16:P:3:LYS:HD2	2.02	0.59
1:A:1120:G:O2'	1:A:1121:U:H5'	2.02	0.59
1:A:1374:A:H2'	1:A:1375:A:C8	2.38	0.59
1:A:1382:C:H2'	1:A:1383:C:H6	1.67	0.59
1:A:1520:G:H2'	1:A:1521:G:H8	1.67	0.59
2:B:200:ILE:O	2:B:202:PRO:HD3	2.02	0.59
3:C:6:HIS:CD2	3:C:8:ILE:HB	2.36	0.59
3:C:61:ALA:C	3:C:63:ASN:H	2.06	0.59
4:D:25:ARG:C	4:D:27:TYR:N	2.56	0.59
4:D:62:GLN:NE2	4:D:65:ARG:HH11	1.94	0.59
5:E:115:VAL:HG12	5:E:116:THR:H	1.67	0.59
7:G:32:ARG:O	7:G:33:ASP:HB2	2.02	0.59
7:G:39:ALA:HA	7:G:42:ILE:CD1	2.33	0.59
15:O:49:ASP:OD2	15:O:52:SER:HB2	2.02	0.59
16:P:12:LYS:O	16:P:13:HIS:HB2	2.01	0.59
16:P:18:ARG:HD3	16:P:35:LYS:HD3	1.84	0.59
16:P:22:THR:O	16:P:23:ASP:HB2	2.02	0.59
1:A:130:A:OP1	17:Q:63:ARG:HD2	2.02	0.59
1:A:447:G:N1	1:A:485:G:H2'	2.17	0.59
1:A:616:G:O2'	1:A:617:G:H5'	2.03	0.59
1:A:686:U:O2	1:A:687:A:C8	2.56	0.59
1:A:915:A:H2'	1:A:916:G:H5'	1.84	0.59
1:A:1101:A:H4'	1:A:1102:A:O5'	2.02	0.59
1:A:1184:G:H2'	1:A:1185:G:H8	1.67	0.59
2:B:142:LEU:O	2:B:146:GLN:HG3	2.01	0.59
4:D:152:SER:HA	4:D:155:LEU:HD12	1.83	0.59
9:I:28:VAL:HA	9:I:63:ILE:O	2.02	0.59
11:K:93:GLN:HE22	11:K:96:ARG:HH22	1.50	0.59
14:N:26:ARG:HG3	14:N:27:CYS:N	2.15	0.59
1:A:248:C:C2'	1:A:249:U:H5'	2.32	0.59
1:A:1201:A:O2'	1:A:1202:G:OP2	2.17	0.59
3:C:121:ALA:HA	3:C:124:ILE:HB	1.84	0.59
4:D:73:ARG:O	4:D:77:ASN:HB2	2.03	0.59
4:D:117:ALA:O	4:D:121:VAL:HG23	2.02	0.59
5:E:16:THR:HG21	5:E:27:ARG:HB2	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:90:VAL:HG23	5:E:90:VAL:O	2.02	0.59
7:G:100:ALA:O	7:G:104:LEU:HG	2.02	0.59
7:G:139:GLU:O	7:G:141:VAL:N	2.35	0.59
9:I:88:TYR:CD1	9:I:88:TYR:C	2.76	0.59
9:I:97:LYS:HA	9:I:102:LEU:HD11	1.82	0.59
17:Q:21:VAL:O	17:Q:41:LYS:HA	2.02	0.59
18:R:38:GLU:H	18:R:41:LYS:CE	2.14	0.59
19:S:31:ILE:HG22	19:S:33:THR:H	1.66	0.59
1:A:1021:G:H2'	1:A:1022:G:O4'	2.03	0.59
1:A:1053:G:O2'	1:A:1199:U:H5	1.84	0.59
1:A:1118:C:H5'	9:I:104:ARG:CD	2.33	0.59
1:A:1124:G:H2'	1:A:1145:C:N4	2.17	0.59
1:A:1347:G:N7	9:I:107:ARG:HB3	2.17	0.59
3:C:122:GLU:C	3:C:124:ILE:H	2.06	0.59
9:I:16:ARG:H	9:I:16:ARG:CD	2.13	0.59
9:I:50:LEU:HA	9:I:85:LEU:HD11	1.85	0.59
9:I:56:LEU:HD23	9:I:57:GLY:N	2.18	0.59
1:A:1014:A:C5'	19:S:14:HIS:HB3	2.32	0.59
1:A:1226:C:H4'	1:A:1227:A:OP1	2.03	0.59
1:A:1240:U:H3'	1:A:1241:G:C5'	2.31	0.59
2:B:10:LEU:O	2:B:12:GLU:N	2.31	0.59
2:B:60:ASP:C	2:B:62:ALA:H	2.06	0.59
2:B:73:THR:CG2	2:B:169:LYS:HE3	2.33	0.59
3:C:203:PHE:HZ	3:C:206:GLU:HB2	1.67	0.59
4:D:191:ARG:HE	4:D:200:GLU:CD	2.06	0.59
4:D:201:GLN:HE22	5:E:99:GLY:HA2	1.65	0.59
5:E:52:PRO:O	5:E:55:VAL:HG22	2.01	0.59
6:F:82:ARG:HB3	6:F:82:ARG:HH11	1.66	0.59
9:I:21:PRO:CA	9:I:59:PHE:HA	2.33	0.59
12:L:45:PRO:HB2	12:L:49:ASN:O	2.03	0.59
16:P:6:LEU:HD21	16:P:73:LEU:CD1	2.33	0.59
20:T:54:LYS:HB2	20:T:100:ILE:CD1	2.33	0.59
1:A:336:C:H2'	1:A:337:C:H6	1.66	0.59
1:A:615:C:O2'	1:A:616:G:H5'	2.03	0.59
3:C:48:TYR:HB2	3:C:52:LEU:HD13	1.83	0.59
3:C:48:TYR:HA	3:C:52:LEU:HB3	1.83	0.59
3:C:134:ILE:O	3:C:138:VAL:HG23	2.02	0.59
4:D:6:GLY:H	4:D:115:ARG:HH22	1.50	0.59
4:D:9:CYS:O	4:D:12:CYS:HB2	2.02	0.59
4:D:32:ALA:O	4:D:34:GLU:N	2.36	0.59
12:L:90:VAL:CG1	12:L:92:ASP:HB2	2.32	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:P:67:THR:O	16:P:70:ALA:HB3	2.02	0.59
20:T:62:LEU:HD23	20:T:62:LEU:O	2.03	0.59
1:A:587:G:H8	1:A:587:G:O5'	1.86	0.59
1:A:744:C:O2'	1:A:745:C:H5'	2.02	0.59
1:A:838:G:C2'	1:A:839:U:H5''	2.33	0.59
1:A:922:G:H1	1:A:1395:C:N4	2.00	0.59
1:A:956:U:C2'	1:A:957:U:H5'	2.33	0.59
1:A:1003(A):G:H21	1:A:1038:C:H1'	1.67	0.59
1:A:1269:A:C2	1:A:1313:U:H1'	2.38	0.59
2:B:181:PHE:CD2	8:H:70:GLN:HB3	2.38	0.59
3:C:79:ARG:N	3:C:79:ARG:HD3	2.18	0.59
3:C:94:LEU:HD23	3:C:94:LEU:C	2.23	0.59
6:F:62:TRP:CH2	6:F:64:GLN:HB2	2.38	0.59
8:H:104:ARG:HG2	8:H:104:ARG:HH11	1.66	0.59
10:J:49:VAL:CG1	14:N:41:ARG:HB2	2.29	0.59
11:K:68:ALA:O	11:K:72:ALA:HB2	2.03	0.59
13:M:24:GLY:C	13:M:25:ILE:HD12	2.24	0.59
14:N:23:ARG:HD3	14:N:28:GLY:O	2.03	0.59
19:S:22:LEU:HD12	19:S:47:HIS:CE1	2.38	0.59
1:A:147:G:N2	1:A:148:G:C4	2.71	0.58
1:A:180:U:O2'	1:A:181:G:H5'	2.04	0.58
1:A:191:G:H1'	20:T:105:SER:HB3	1.85	0.58
1:A:390:C:H4'	16:P:28:ARG:NH2	2.18	0.58
1:A:687:A:O2'	1:A:688:G:OP2	2.18	0.58
1:A:1019:C:O2'	1:A:1020:U:H5'	2.03	0.58
1:A:1240:U:H1'	7:G:38:LEU:HD11	1.85	0.58
2:B:60:ASP:OD1	2:B:61:LEU:N	2.36	0.58
3:C:65:ALA:N	3:C:99:VAL:HB	2.18	0.58
5:E:12:LEU:CD1	5:E:12:LEU:N	2.66	0.58
5:E:105:VAL:CG1	5:E:131:ILE:HG22	2.33	0.58
7:G:11:GLN:NE2	7:G:11:GLN:HA	2.18	0.58
8:H:82:HIS:ND1	8:H:138:TRP:NE1	2.51	0.58
12:L:70:ILE:CD1	12:L:77:LEU:HD12	2.33	0.58
13:M:102:ARG:NH1	13:M:105:THR:OG1	2.36	0.58
16:P:82:GLN:NE2	16:P:82:GLN:N	2.42	0.58
20:T:44:ALA:O	20:T:47:GLY:N	2.36	0.58
1:A:168:G:C2'	1:A:169:C:H5'	2.32	0.58
1:A:193:C:H2'	1:A:194:C:H6	1.68	0.58
1:A:243:A:C2	1:A:246:A:C8	2.90	0.58
1:A:281:G:O2'	1:A:282:A:P	2.61	0.58
1:A:610:G:H2'	1:A:611:A:H8	1.66	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:763:G:H2'	1:A:764:C:H6	1.68	0.58
1:A:836:G:C6	1:A:851:G:C6	2.91	0.58
1:A:915:A:C2'	1:A:916:G:H5'	2.33	0.58
1:A:1189:C:C2'	1:A:1190:G:H5'	2.33	0.58
1:A:1214:C:H5''	1:A:1215:G:OP2	2.03	0.58
1:A:1319:A:N6	1:A:1361:G:H21	1.98	0.58
1:A:1518:A:OP2	1:A:1518:A:O4'	2.21	0.58
2:B:25:ASN:HD21	2:B:27:LYS:HG3	1.68	0.58
2:B:96:ARG:HA	2:B:96:ARG:HE	1.68	0.58
3:C:137:ALA:N	3:C:140:ARG:NH1	2.51	0.58
4:D:22:LYS:CB	4:D:26:CYS:SG	2.76	0.58
9:I:111:ARG:HG3	9:I:111:ARG:NH1	2.17	0.58
18:R:43:PHE:O	18:R:44:LEU:HD22	2.03	0.58
20:T:33:ILE:HD11	20:T:63:ILE:CA	2.23	0.58
1:A:328:C:O2	1:A:328:C:C2'	2.48	0.58
1:A:754:C:O5'	15:O:72:ARG:NH2	2.36	0.58
1:A:757:U:H2'	1:A:758:G:O4'	2.02	0.58
2:B:19:HIS:N	2:B:39:ILE:HG21	2.18	0.58
3:C:21:ARG:NH2	3:C:56:ASP:OD2	2.36	0.58
3:C:102:ASN:N	3:C:102:ASN:HD22	2.00	0.58
5:E:19:MET:SD	5:E:24:ARG:HD2	2.42	0.58
15:O:11:VAL:HA	15:O:14:GLU:HB3	1.85	0.58
19:S:80:TYR:CG	19:S:81:ARG:N	2.67	0.58
1:A:996:A:H2'	1:A:997:U:C6	2.37	0.58
1:A:1407:C:H2'	1:A:1408:A:H8	1.66	0.58
1:A:1436:U:C5	1:A:1437:C:N3	2.71	0.58
2:B:44:LEU:HA	2:B:47:THR:CB	2.33	0.58
2:B:82:ARG:CA	2:B:92:TYR:HE2	2.13	0.58
2:B:210:SER:C	2:B:212:GLN:H	2.06	0.58
3:C:108:ASN:CG	3:C:111:LEU:HD12	2.23	0.58
4:D:206:PHE:CD2	4:D:207:TYR:CE2	2.91	0.58
6:F:64:GLN:O	6:F:64:GLN:HG2	2.03	0.58
10:J:71:LEU:HD13	10:J:72:VAL:H	1.68	0.58
12:L:58:VAL:HG21	12:L:83:VAL:HG11	1.84	0.58
15:O:48:LYS:NZ	15:O:48:LYS:HB2	2.18	0.58
16:P:10:GLY:HA2	16:P:16:HIS:HB2	1.85	0.58
1:A:438:G:N2	1:A:495:U:H3'	2.18	0.58
1:A:681:C:H2'	1:A:682:G:H8	1.69	0.58
1:A:685:G:C2	1:A:686:U:C5	2.92	0.58
1:A:976:G:N2	1:A:1361(A):C:H2'	2.18	0.58
1:A:1067:A:N3	1:A:1068:G:H1'	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1360:A:H3'	1:A:1361:G:N7	2.18	0.58
2:B:217:ARG:HA	2:B:220:ASP:OD2	2.02	0.58
3:C:167:TRP:O	3:C:168:ALA:CB	2.52	0.58
4:D:52:SER:O	4:D:56:VAL:HG23	2.03	0.58
9:I:48:GLU:HA	9:I:51:ARG:NE	2.15	0.58
9:I:114:TYR:O	9:I:116:LYS:HG2	2.04	0.58
16:P:39:TYR:CE2	16:P:41:PRO:HG3	2.33	0.58
1:A:357:G:C2	1:A:358:U:C6	2.92	0.58
1:A:1226:C:N4	13:M:104:ARG:HG3	2.18	0.58
1:A:1258:G:H1	1:A:1277:C:N4	2.02	0.58
1:A:1367:C:P	9:I:112:LYS:HZ1	2.26	0.58
1:A:1437:C:H6	1:A:1438:G:C8	2.20	0.58
1:A:1489:G:H2'	1:A:1490:C:C6	2.38	0.58
3:C:10:PHE:CD2	3:C:178:LEU:HD13	2.39	0.58
3:C:191:THR:C	3:C:193:TYR:H	2.06	0.58
5:E:78:HIS:ND1	8:H:104:ARG:CD	2.62	0.58
6:F:4:TYR:CZ	6:F:72:VAL:HG21	2.39	0.58
7:G:67:GLU:HA	7:G:70:LYS:HD2	1.84	0.58
10:J:98:ILE:H	10:J:98:ILE:CD1	2.10	0.58
11:K:25:TYR:CE2	11:K:88:GLY:HA2	2.39	0.58
11:K:47:VAL:O	11:K:48:ILE:C	2.40	0.58
12:L:24:VAL:HG13	12:L:98:TYR:HE2	1.67	0.58
13:M:73:GLU:O	13:M:76:ALA:HB3	2.02	0.58
13:M:78:ILE:H	13:M:78:ILE:CD1	2.16	0.58
16:P:50:LYS:HG2	16:P:51:VAL:H	1.66	0.58
20:T:30:LYS:HZ2	20:T:80:ARG:NH2	2.01	0.58
20:T:69:GLY:C	20:T:71:THR:H	2.07	0.58
21:V:10:ARG:H	21:V:10:ARG:CD	2.16	0.58
1:A:58:C:O2'	1:A:59:A:H5'	2.04	0.58
1:A:410:G:OP2	4:D:25:ARG:HD2	2.03	0.58
1:A:580:U:H1'	15:O:57:LEU:HD23	1.86	0.58
1:A:664:G:H22	1:A:741:G:H1	1.52	0.58
1:A:826:C:H2'	1:A:827:U:H6	1.68	0.58
1:A:840:C:OP2	1:A:840:C:H3'	2.03	0.58
1:A:1098:C:O2'	1:A:1099:G:H5'	2.03	0.58
1:A:1404:C:H1'	1:A:1499:A:C2	2.38	0.58
2:B:115:LEU:HD22	2:B:153:ARG:HE	1.67	0.58
3:C:10:PHE:CE1	3:C:178:LEU:HD22	2.38	0.58
3:C:166:GLU:HA	3:C:166:GLU:OE2	2.03	0.58
3:C:203:PHE:O	3:C:204:LEU:HG	2.04	0.58
5:E:11:ILE:HD11	5:E:33:VAL:CG2	2.31	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:85:LEU:HB3	9:I:92:TYR:CD1	2.39	0.58
16:P:42:ARG:O	16:P:43:LYS:HD2	2.03	0.58
1:A:260:G:H2'	1:A:261:U:C6	2.38	0.58
1:A:346:G:C2'	1:A:347:G:H5'	2.33	0.58
1:A:930:C:H2'	1:A:931:C:H5'	1.85	0.58
1:A:1044:A:H2'	1:A:1045:C:O4'	2.04	0.58
1:A:1049:U:O2'	1:A:1050:G:OP2	2.20	0.58
1:A:1241:G:H2'	1:A:1242:C:H6	1.68	0.58
4:D:127:THR:HG1	4:D:130:GLY:C	2.07	0.58
8:H:54:ASP:C	8:H:56:LYS:H	2.07	0.58
10:J:4:ILE:N	10:J:4:ILE:HD12	2.19	0.58
16:P:24:ALA:C	16:P:26:ARG:N	2.57	0.58
17:Q:81:ARG:HG3	17:Q:81:ARG:O	2.03	0.58
1:A:16:A:O2'	1:A:17:U:H5'	2.04	0.58
1:A:21:G:H2'	1:A:22:G:H8	1.69	0.58
1:A:539:A:H2'	1:A:540:G:C8	2.39	0.58
1:A:721:G:C6	1:A:733:A:C2	2.91	0.58
1:A:930:C:O2'	1:A:931:C:H5'	2.02	0.58
1:A:948:C:O2'	1:A:949:A:H5'	2.03	0.58
1:A:1001:A:O2'	1:A:1002:G:H5'	2.02	0.58
1:A:1191:A:C4	1:A:1192:C:H5	2.21	0.58
1:A:1376:U:H2'	1:A:1377:A:C8	2.39	0.58
1:A:1451:A:O2'	1:A:1452:C:OP1	2.21	0.58
4:D:108:LEU:HD22	4:D:176:LEU:HB3	1.86	0.58
5:E:95:ALA:HB1	5:E:96:PRO:CD	2.34	0.58
7:G:15:ASP:C	7:G:17:VAL:H	2.06	0.58
7:G:65:ALA:CB	7:G:128:ALA:HA	2.24	0.58
8:H:17:THR:HG22	8:H:63:LEU:HD23	1.83	0.58
9:I:84:ALA:O	9:I:86:VAL:N	2.37	0.58
10:J:50:ILE:N	10:J:60:ARG:HG3	2.18	0.58
10:J:85:LEU:N	10:J:88:LEU:HD12	2.19	0.58
11:K:30:VAL:HG21	11:K:65:ALA:CA	2.34	0.58
16:P:4:ILE:HG12	16:P:21:VAL:HG22	1.86	0.58
17:Q:9:VAL:CG1	17:Q:10:VAL:N	2.66	0.58
20:T:13:LEU:HD12	20:T:13:LEU:C	2.23	0.58
20:T:50:GLU:O	20:T:51:GLU:C	2.40	0.58
1:A:7:G:H5'	1:A:298:A:C5'	2.34	0.58
1:A:361:G:C2'	1:A:362:G:H5'	2.33	0.58
1:A:502:G:C2	1:A:503:C:C2	2.92	0.58
1:A:509:A:H4'	1:A:510:A:OP1	2.03	0.58
1:A:858:G:C2'	1:A:859:A:H5''	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1028:C:H2'	1:A:1029:C:O4'	2.04	0.58
1:A:1056:U:H5'	3:C:163:ALA:CB	2.34	0.58
1:A:1185:G:O2'	1:A:1186:G:H5'	2.04	0.58
1:A:1187:G:H2'	1:A:1188:A:C8	2.39	0.58
1:A:1347:G:H1'	1:A:1348:U:H5	1.69	0.58
1:A:1465:C:H2'	1:A:1466:C:O4'	2.04	0.58
1:A:1470:G:O2'	1:A:1471:G:H5'	2.04	0.58
4:D:5:ILE:HA	4:D:115:ARG:NH2	2.19	0.58
4:D:152:SER:CA	4:D:155:LEU:HD12	2.33	0.58
7:G:116:ALA:HA	7:G:119:ARG:NH2	2.19	0.58
8:H:104:ARG:C	8:H:106:GLY:H	2.06	0.58
9:I:43:ALA:CA	9:I:74:ILE:HD13	2.33	0.58
9:I:53:VAL:HG21	9:I:85:LEU:CD1	2.32	0.58
10:J:65:LEU:HD23	10:J:66:ARG:N	2.19	0.58
11:K:93:GLN:NE2	11:K:96:ARG:HH22	2.02	0.58
16:P:5:ARG:C	16:P:6:LEU:HD12	2.24	0.58
19:S:22:LEU:HD11	19:S:31:ILE:HD11	1.84	0.58
1:A:102:G:N2	1:A:171:A:H2	2.01	0.57
1:A:444:C:H42	1:A:490:G:H1	1.50	0.57
1:A:989:C:O2'	1:A:990:C:H5'	2.04	0.57
2:B:152:PHE:CD2	2:B:152:PHE:O	2.56	0.57
6:F:2:ARG:C	6:F:66:GLU:HG3	2.24	0.57
7:G:91:VAL:HG12	7:G:92:SER:N	2.16	0.57
12:L:84:LEU:O	12:L:101:VAL:HG13	2.04	0.57
15:O:41:GLU:HA	15:O:44:LYS:HG2	1.86	0.57
1:A:135:C:C2	16:P:1:MET:HB2	2.38	0.57
1:A:496:A:H4'	1:A:497:A:OP1	2.02	0.57
1:A:935:A:O2'	1:A:936:C:H5'	2.05	0.57
1:A:1151:A:H5''	10:J:42:THR:H	1.67	0.57
1:A:1160:G:O2'	1:A:1161:C:H5'	2.04	0.57
1:A:1301:U:O2	1:A:1301:U:H2'	2.05	0.57
1:A:1350:A:OP2	9:I:118:LYS:HD3	2.03	0.57
3:C:67:THR:O	3:C:67:THR:HG22	2.05	0.57
4:D:16:GLY:O	4:D:33:MET:HE1	2.03	0.57
7:G:134:ALA:O	7:G:136:LYS:N	2.37	0.57
19:S:5:LEU:N	19:S:5:LEU:HD12	2.20	0.57
1:A:52:G:O2'	1:A:53:A:H5'	2.04	0.57
1:A:390:C:C3'	16:P:28:ARG:HH22	2.16	0.57
1:A:684:A:O2'	1:A:685:G:H5'	2.04	0.57
1:A:826:C:H2'	1:A:827:U:C6	2.39	0.57
1:A:1291:G:H5''	7:G:41:ARG:HH22	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:18:TRP:HZ2	14:N:56:VAL:O	1.88	0.57
3:C:147:LYS:HD3	3:C:203:PHE:CE2	2.39	0.57
4:D:21:LEU:HD11	4:D:67:ILE:HA	1.85	0.57
4:D:36:ARG:N	4:D:37:PRO:CD	2.63	0.57
8:H:11:THR:O	8:H:14:ARG:N	2.37	0.57
9:I:23:ASN:H	9:I:60:ASP:H	1.52	0.57
9:I:110:GLU:HG2	9:I:111:ARG:H	1.69	0.57
9:I:118:LYS:NZ	9:I:118:LYS:HB3	2.20	0.57
11:K:109:VAL:HG12	18:R:84:LYS:HB3	1.86	0.57
15:O:61:GLY:O	15:O:63:ARG:N	2.36	0.57
19:S:74:PHE:N	19:S:74:PHE:CD1	2.71	0.57
1:A:381:C:H2'	1:A:382:A:O4'	2.04	0.57
1:A:445:G:H2'	1:A:446:G:C8	2.39	0.57
1:A:600:C:OP1	8:H:97:VAL:CG1	2.52	0.57
1:A:650:G:O2'	1:A:651:C:H5'	2.04	0.57
1:A:977:A:C2'	1:A:978:A:H5''	2.33	0.57
1:A:1258:G:H1	1:A:1277:C:H42	1.50	0.57
2:B:12:GLU:C	2:B:14:GLY:H	2.06	0.57
2:B:172:ILE:H	2:B:172:ILE:CD1	2.00	0.57
5:E:71:LEU:HD22	5:E:114:GLY:O	2.04	0.57
6:F:48:LEU:HB3	6:F:50:TYR:O	2.04	0.57
9:I:42:ARG:HH22	9:I:75:ASP:CG	2.07	0.57
11:K:50:TYR:HB3	11:K:54:ARG:HB2	1.86	0.57
12:L:115:LYS:O	12:L:117:ARG:HG3	2.04	0.57
13:M:87:TYR:O	13:M:90:LEU:N	2.38	0.57
14:N:3:ARG:H	14:N:3:ARG:HD3	1.69	0.57
1:A:701:C:H5'	1:A:703:G:O4'	2.04	0.57
1:A:945:G:H21	1:A:1334:G:H4'	1.70	0.57
1:A:967:C:H2'	1:A:968:A:N7	2.20	0.57
1:A:1253:G:N1	1:A:1285:A:N6	2.52	0.57
1:A:1328:C:O3'	13:M:28:ALA:HB3	2.03	0.57
3:C:21:ARG:CZ	3:C:56:ASP:HB3	2.34	0.57
7:G:100:ALA:HB3	7:G:101:LEU:HD23	1.86	0.57
9:I:102:LEU:N	9:I:102:LEU:CD2	2.61	0.57
10:J:3:LYS:N	10:J:75:ILE:HA	2.19	0.57
16:P:57:ARG:NH1	16:P:79:VAL:O	2.37	0.57
16:P:82:GLN:HE21	16:P:82:GLN:N	2.00	0.57
17:Q:85:VAL:HG12	17:Q:89:LEU:HG	1.86	0.57
1:A:411:A:H2'	1:A:413:G:C8	2.39	0.57
1:A:676:A:H2'	1:A:677:U:C6	2.39	0.57
1:A:761:G:C4'	17:Q:102:GLY:HA3	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1435:G:H2'	1:A:1436:U:C6	2.40	0.57
4:D:54:TYR:HE1	4:D:206:PHE:CE1	2.22	0.57
4:D:96:LEU:C	4:D:98:GLU:H	2.08	0.57
7:G:54:THR:HB	7:G:56:GLN:HE22	1.68	0.57
8:H:31:PHE:HZ	8:H:134:ILE:CD1	2.17	0.57
8:H:69:ARG:HD3	8:H:75:ARG:O	2.04	0.57
8:H:134:ILE:CG2	8:H:135:CYS:N	2.58	0.57
14:N:5:ALA:O	14:N:6:LEU:HD12	2.04	0.57
14:N:33:VAL:HA	14:N:39:LEU:O	2.04	0.57
16:P:21:VAL:HG12	16:P:33:ILE:HB	1.85	0.57
17:Q:75:ARG:CG	17:Q:76:LEU:N	2.67	0.57
18:R:26:LEU:HD12	18:R:26:LEU:N	2.19	0.57
18:R:87:ARG:HG3	18:R:87:ARG:HH11	1.69	0.57
1:A:84:U:H2'	1:A:88:A:C8	2.39	0.57
1:A:621:A:H2'	1:A:622:A:H8	1.68	0.57
1:A:1037:C:O2	1:A:1037:C:H2'	2.03	0.57
1:A:1249:C:H3'	1:A:1249:C:H6	1.70	0.57
1:A:1298:C:H4'	1:A:1299:A:C5'	2.28	0.57
1:A:1487:G:C2'	1:A:1488:G:H5'	2.35	0.57
2:B:16:HIS:O	2:B:17:PHE:O	2.22	0.57
11:K:81:ASP:OD1	11:K:106:LYS:HB2	2.04	0.57
12:L:25:PRO:C	12:L:27:LEU:N	2.54	0.57
15:O:50:HIS:O	15:O:53:HIS:N	2.32	0.57
15:O:79:ARG:O	15:O:83:GLU:N	2.34	0.57
17:Q:96:GLN:CD	17:Q:97:SER:N	2.58	0.57
18:R:87:ARG:HH11	18:R:87:ARG:CG	2.17	0.57
1:A:103:C:P	20:T:17:ARG:NH1	2.78	0.57
1:A:596:C:OP2	1:A:597:G:OP2	2.23	0.57
1:A:632:A:H2'	1:A:633:G:H5'	1.85	0.57
1:A:894:G:C2	1:A:895:G:C4	2.92	0.57
2:B:46:LYS:O	2:B:49:GLU:N	2.37	0.57
7:G:45:ASP:C	7:G:47:CYS:N	2.58	0.57
7:G:67:GLU:HA	7:G:70:LYS:HG3	1.86	0.57
9:I:9:ARG:H	9:I:79:LEU:HD12	1.68	0.57
10:J:20:ALA:C	10:J:22:LYS:H	2.08	0.57
11:K:102:GLY:O	11:K:103:LEU:HD23	2.05	0.57
17:Q:40:LYS:HD3	17:Q:42:TYR:OH	2.05	0.57
18:R:86:VAL:O	18:R:87:ARG:CB	2.53	0.57
19:S:30:LEU:HD23	19:S:31:ILE:O	2.05	0.57
1:A:190(F):G:H8	1:A:190(F):G:OP1	1.88	0.57
1:A:219:C:O2'	1:A:381:C:H5'	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:384:G:H2'	1:A:385:C:C6	2.40	0.57
1:A:411:A:C2'	1:A:413:G:H1'	2.34	0.57
1:A:726:C:O2'	1:A:727:G:H5'	2.04	0.57
1:A:1347:G:O2'	1:A:1348:U:P	2.63	0.57
1:A:1374:A:O2'	1:A:1375:A:H5'	2.05	0.57
2:B:12:GLU:C	2:B:14:GLY:N	2.55	0.57
2:B:187:LEU:HD12	2:B:205:ASP:HA	1.87	0.57
3:C:51:GLY:HA3	3:C:70:VAL:HA	1.87	0.57
3:C:76:VAL:HG23	3:C:77:ILE:N	2.20	0.57
4:D:63:LYS:O	4:D:67:ILE:HG13	2.04	0.57
5:E:82:VAL:HB	5:E:89:ILE:HG22	1.87	0.57
7:G:45:ASP:O	7:G:47:CYS:N	2.38	0.57
11:K:79:SER:O	11:K:80:VAL:CB	2.53	0.57
13:M:57:ARG:O	13:M:61:GLU:HG3	2.04	0.57
13:M:94:ARG:HG2	13:M:94:ARG:HH11	1.68	0.57
15:O:21:ASP:CG	15:O:24:SER:HB3	2.25	0.57
17:Q:9:VAL:O	17:Q:21:VAL:HA	2.05	0.57
1:A:1257:U:H4'	1:A:1258:G:OP2	2.03	0.57
2:B:219:VAL:HA	2:B:222:ILE:CD1	2.32	0.57
3:C:11:ARG:HA	3:C:14:ILE:HD11	1.87	0.57
4:D:89:THR:O	4:D:90:GLY:C	2.43	0.57
4:D:96:LEU:H	4:D:96:LEU:CD1	2.18	0.57
4:D:111:ALA:HA	4:D:161:ASN:ND2	2.20	0.57
10:J:40:LEU:HB3	10:J:69:ASN:O	2.05	0.57
16:P:39:TYR:HD1	16:P:49:LEU:HD13	1.69	0.57
19:S:64:GLU:HA	19:S:67:VAL:CG2	2.34	0.57
20:T:11:SER:C	20:T:13:LEU:N	2.56	0.57
1:A:192:U:H5'	20:T:102:GLY:HA3	1.87	0.56
1:A:318:G:H2'	1:A:319:G:C8	2.40	0.56
1:A:1089:G:O2'	1:A:1090:U:H5'	2.04	0.56
1:A:1346:A:O2'	1:A:1347:G:OP2	2.22	0.56
1:A:1509:C:C2	1:A:1510:U:C6	2.93	0.56
4:D:46:LYS:HG2	4:D:47:ARG:H	1.70	0.56
6:F:79:LEU:HD22	6:F:79:LEU:N	2.20	0.56
7:G:5:ARG:HG2	7:G:7:ALA:H	1.71	0.56
7:G:115:ARG:HD3	7:G:115:ARG:H	1.70	0.56
8:H:114:THR:HG22	8:H:130:GLY:O	2.04	0.56
10:J:59:SER:O	10:J:61:GLU:N	2.38	0.56
11:K:66:LEU:O	11:K:69:ALA:HB3	2.05	0.56
12:L:34:ARG:O	12:L:61:THR:HG23	2.05	0.56
19:S:72:GLY:O	19:S:75:ALA:N	2.34	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:C:O2'	1:A:176:C:H5'	2.04	0.56
1:A:355:C:H1'	1:A:388:G:N3	2.20	0.56
1:A:411:A:H2'	1:A:413:G:H1'	1.87	0.56
1:A:428:G:C5'	4:D:7:PRO:HB3	2.35	0.56
1:A:734:G:C6	1:A:735:C:C4	2.92	0.56
1:A:862:C:O2'	1:A:863:U:H5'	2.05	0.56
1:A:1207:G:H2'	1:A:1208:C:H6	1.70	0.56
1:A:1288:A:H2'	1:A:1289:A:C8	2.40	0.56
3:C:184:TYR:CG	3:C:185:GLY:N	2.72	0.56
4:D:148:VAL:CG2	4:D:181:MET:HB3	2.35	0.56
4:D:173:TRP:CD1	4:D:189:PRO:HD3	2.40	0.56
5:E:41:VAL:HG12	5:E:42:GLY:N	2.20	0.56
6:F:19:LEU:HD23	6:F:20:ALA:N	2.20	0.56
7:G:66:VAL:CG1	7:G:100:ALA:HB1	2.35	0.56
7:G:152:ALA:HB1	7:G:155:ARG:CZ	2.35	0.56
11:K:16:SER:HA	11:K:79:SER:CB	2.34	0.56
11:K:30:VAL:HG21	11:K:65:ALA:CB	2.35	0.56
14:N:24:CYS:O	14:N:28:GLY:N	2.39	0.56
15:O:56:LEU:O	15:O:59:MET:N	2.38	0.56
1:A:562:C:H4'	1:A:563:A:H5'	1.86	0.56
1:A:885:G:H1	1:A:912:C:H42	1.53	0.56
1:A:997:U:O2	1:A:1044:A:H2	1.88	0.56
1:A:998:G:O2'	1:A:999:C:H5'	2.06	0.56
1:A:1093:A:H2	1:A:1109:C:O2'	1.88	0.56
1:A:1206:G:H2'	1:A:1207:G:H8	1.70	0.56
1:A:1442:G:C5	1:A:1446:A:N1	2.73	0.56
1:A:1510:U:H1'	1:A:1526:G:N2	2.19	0.56
2:B:22:LYS:HA	2:B:40:HIS:HE1	1.69	0.56
2:B:25:ASN:C	2:B:25:ASN:ND2	2.54	0.56
2:B:61:LEU:HD21	2:B:66:GLY:HA3	1.87	0.56
3:C:182:ILE:CG2	3:C:183:ASP:H	2.00	0.56
5:E:15:ARG:O	5:E:16:THR:HG22	2.05	0.56
6:F:24:GLU:O	6:F:28:ARG:N	2.39	0.56
7:G:71:PRO:HD3	7:G:103:TRP:CZ3	2.35	0.56
7:G:113:GLU:HB3	7:G:118:VAL:HG11	1.87	0.56
8:H:48:TYR:CD1	8:H:48:TYR:C	2.78	0.56
10:J:9:ARG:HB3	10:J:9:ARG:CZ	2.35	0.56
10:J:24:VAL:HG22	10:J:72:VAL:HG11	1.88	0.56
10:J:28:ARG:HA	10:J:33:GLN:HA	1.86	0.56
10:J:40:LEU:HD22	10:J:69:ASN:CG	2.25	0.56
11:K:91:ARG:HD3	11:K:92:GLU:OE1	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:M:63:THR:HG23	13:M:64:TRP:CD2	2.40	0.56
1:A:13:U:C5	1:A:916:G:O6	2.59	0.56
1:A:245:C:O2'	1:A:246:A:H5'	2.05	0.56
1:A:321:A:H2'	1:A:322:C:H6	1.70	0.56
1:A:393:A:H2'	1:A:394:G:C8	2.23	0.56
1:A:433:C:H2'	1:A:434:U:C6	2.40	0.56
1:A:606:G:N2	1:A:631:G:H2'	2.19	0.56
1:A:969:A:O2'	1:A:970:C:H5'	2.06	0.56
1:A:1090:U:O2'	1:A:1091:U:H5'	2.05	0.56
2:B:115:LEU:HD23	2:B:115:LEU:C	2.25	0.56
3:C:48:TYR:C	3:C:50:ALA:H	2.09	0.56
3:C:95:THR:CG2	3:C:98:ASN:HA	2.35	0.56
4:D:14:ARG:HB2	4:D:40:PRO:CG	2.36	0.56
4:D:105:VAL:HG13	4:D:110:PHE:HB2	1.88	0.56
5:E:106:PRO:O	5:E:107:ARG:C	2.43	0.56
15:O:21:ASP:OD1	15:O:24:SER:HB3	2.05	0.56
15:O:64:ARG:NH2	15:O:68:ARG:HH22	1.97	0.56
1:A:649:G:O2'	1:A:650:G:H5'	2.06	0.56
1:A:1091:U:H2'	1:A:1093:A:OP2	2.05	0.56
1:A:1215:G:O2'	1:A:1216:G:H5'	2.05	0.56
1:A:1276:G:H21	1:A:1282:C:H1'	1.70	0.56
3:C:132:ARG:HA	3:C:135:LYS:HD2	1.87	0.56
4:D:125:HIS:HA	4:D:149:ALA:CB	2.35	0.56
4:D:194:LEU:N	4:D:194:LEU:CD2	2.69	0.56
6:F:21:LEU:HD12	6:F:21:LEU:O	2.05	0.56
12:L:97:ARG:C	12:L:98:TYR:CD1	2.79	0.56
13:M:80:ARG:O	13:M:82:MET:N	2.39	0.56
16:P:10:GLY:HA3	16:P:15:PRO:HA	1.86	0.56
1:A:273:A:H2'	1:A:274:A:H5'	1.87	0.56
1:A:359:U:H2'	1:A:360:A:C8	2.39	0.56
1:A:1201:A:H4'	1:A:1202:G:H5''	1.88	0.56
2:B:178:ARG:C	2:B:180:LEU:H	2.08	0.56
4:D:62:GLN:O	4:D:65:ARG:HB3	2.05	0.56
10:J:63:PHE:HE1	14:N:45:ARG:HA	1.71	0.56
10:J:81:THR:HG22	10:J:82:ILE:N	2.19	0.56
11:K:77:MET:SD	11:K:80:VAL:HG23	2.45	0.56
17:Q:74:LEU:O	17:Q:75:ARG:HB2	2.05	0.56
19:S:64:GLU:O	19:S:67:VAL:HB	2.05	0.56
20:T:101:GLY:O	20:T:103:GLY:N	2.39	0.56
1:A:61:G:H2'	1:A:62:U:O4'	2.06	0.56
1:A:519:C:O2'	1:A:520:A:H5'	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:539:A:OP1	12:L:114:LYS:HE2	2.06	0.56
1:A:577:G:H1'	1:A:816:A:N3	2.21	0.56
1:A:781:A:C2'	1:A:782:A:H5'	2.33	0.56
4:D:21:LEU:HD12	4:D:21:LEU:N	2.20	0.56
8:H:114:THR:HG21	8:H:129:VAL:HG23	1.88	0.56
9:I:7:THR:HA	9:I:16:ARG:HA	1.88	0.56
10:J:16:LEU:HD22	10:J:70:ARG:CD	2.36	0.56
15:O:62:GLN:O	15:O:66:LEU:HG	2.05	0.56
16:P:30:GLY:O	16:P:31:LYS:O	2.24	0.56
16:P:39:TYR:CD1	16:P:73:LEU:HD21	2.40	0.56
1:A:96:G:O2'	1:A:97:G:H5'	2.06	0.56
1:A:186:C:C2	1:A:187:C:C5	2.93	0.56
1:A:1207:G:O2'	1:A:1208:C:H5'	2.05	0.56
1:A:1254:C:H41	10:J:43:ARG:CZ	2.18	0.56
3:C:40:ARG:CD	3:C:55:VAL:HG11	2.36	0.56
4:D:8:VAL:O	4:D:10:ARG:N	2.38	0.56
4:D:19:LEU:O	4:D:22:LYS:HG3	2.06	0.56
4:D:116:GLN:HG2	4:D:116:GLN:O	2.05	0.56
5:E:91:LEU:HD23	5:E:120:THR:HG21	1.87	0.56
5:E:92:LYS:O	5:E:118:ILE:HG22	2.06	0.56
5:E:105:VAL:O	5:E:109:ILE:HG12	2.06	0.56
6:F:33:TYR:CA	6:F:71:ARG:HH21	2.19	0.56
7:G:13:GLN:O	7:G:21:VAL:HG12	2.05	0.56
7:G:54:THR:HB	7:G:56:GLN:NE2	2.21	0.56
7:G:141:VAL:C	7:G:143:ARG:N	2.56	0.56
11:K:30:VAL:HG21	11:K:65:ALA:HA	1.86	0.56
12:L:97:ARG:HB2	12:L:98:TYR:CE1	2.41	0.56
14:N:37:PHE:HB2	14:N:39:LEU:HD11	1.88	0.56
14:N:48:ALA:C	14:N:50:LYS:H	2.09	0.56
1:A:168:G:HO2'	1:A:169:C:H5'	1.71	0.56
1:A:428:G:O2'	1:A:429:U:O5'	2.16	0.56
1:A:477:G:O2'	1:A:478:A:H5'	2.05	0.56
1:A:544:G:C4	1:A:545:C:C5	2.94	0.56
1:A:751:U:H1'	15:O:23:GLY:O	2.05	0.56
1:A:1314:C:OP2	19:S:6:LYS:HD2	2.06	0.56
1:A:1370:G:H2'	1:A:1371:G:C8	2.39	0.56
1:A:1461:G:H2'	1:A:1462:G:H8	1.69	0.56
3:C:91:LEU:CD1	3:C:101:LEU:HB2	2.36	0.56
4:D:30:LYS:C	4:D:32:ALA:H	2.09	0.56
4:D:64:LEU:CD2	4:D:198:VAL:HG11	2.30	0.56
4:D:165:MET:C	4:D:166:LYS:HG3	2.27	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:93:PRO:HG2	8:H:105:ARG:HH21	1.71	0.56
7:G:62:PHE:O	7:G:66:VAL:HB	2.05	0.56
7:G:114:ARG:HD3	7:G:114:ARG:H	1.70	0.56
7:G:141:VAL:CA	7:G:144:MET:HB2	2.23	0.56
9:I:79:LEU:HD13	9:I:83:ARG:NH2	2.20	0.56
10:J:60:ARG:O	10:J:61:GLU:HB3	2.06	0.56
11:K:11:LYS:HB2	11:K:11:LYS:NZ	2.21	0.56
12:L:43:VAL:CG1	12:L:44:THR:H	2.14	0.56
13:M:98:VAL:CG2	13:M:110:ARG:HH12	2.16	0.56
15:O:87:ILE:CG2	15:O:88:ARG:H	1.98	0.56
16:P:53:VAL:C	16:P:55:ARG:N	2.55	0.56
20:T:57:ARG:HH22	20:T:102:GLY:HA3	1.68	0.56
1:A:190(L):U:O2	20:T:105:SER:HB2	2.05	0.56
1:A:729:A:H2'	1:A:730:G:H8	1.71	0.56
1:A:742:G:H2'	1:A:743:U:O4'	2.06	0.56
1:A:927:G:H2'	1:A:928:G:H8	1.71	0.56
1:A:1025:U:HO2'	1:A:1026:G:H8	1.54	0.56
1:A:1127:G:H1	1:A:1144:G:N2	2.02	0.56
1:A:1347:G:C6	9:I:107:ARG:CZ	2.89	0.56
3:C:16:ARG:HH22	3:C:183:ASP:HB2	1.71	0.56
3:C:75:VAL:O	3:C:83:ARG:HD2	2.06	0.56
3:C:180:ALA:O	3:C:206:GLU:HA	2.05	0.56
3:C:196:LEU:N	3:C:196:LEU:CD2	2.68	0.56
4:D:79:PHE:CD1	4:D:207:TYR:CD1	2.94	0.56
7:G:69:VAL:HA	7:G:138:LYS:HD2	1.88	0.56
8:H:40:ALA:O	8:H:43:GLY:N	2.39	0.56
8:H:104:ARG:O	8:H:106:GLY:N	2.38	0.56
9:I:96:LEU:HG	9:I:102:LEU:HD22	1.88	0.56
11:K:73:MET:CA	11:K:77:MET:HB2	2.35	0.56
12:L:109:GLY:HA3	12:L:121:GLY:O	2.05	0.56
14:N:37:PHE:HB2	14:N:39:LEU:CD1	2.36	0.56
16:P:6:LEU:HB3	16:P:17:TYR:HD2	1.70	0.56
20:T:84:LEU:O	20:T:87:LYS:HB2	2.06	0.56
1:A:256:U:H5'	17:Q:17:LYS:NZ	2.21	0.55
1:A:376:G:C4	1:A:389:A:C2	2.94	0.55
1:A:407:G:H4'	4:D:116:GLN:HA	1.88	0.55
1:A:601:C:C2'	1:A:602:A:H5'	2.36	0.55
2:B:204:ASN:HD22	2:B:206:ASP:H	1.53	0.55
3:C:91:LEU:O	3:C:95:THR:HG22	2.06	0.55
3:C:113:ALA:O	3:C:114:PRO:C	2.44	0.55
5:E:31:LEU:CD2	5:E:45:PHE:HB2	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:109:ILE:O	5:E:113:ALA:HB2	2.06	0.55
5:E:135:THR:HG22	5:E:136:MET:N	2.20	0.55
7:G:16:LEU:HD22	7:G:16:LEU:N	2.20	0.55
9:I:3:GLN:C	9:I:3:GLN:NE2	2.57	0.55
10:J:16:LEU:HG	10:J:94:VAL:CG1	2.35	0.55
12:L:32:PHE:HA	12:L:85:ILE:O	2.05	0.55
12:L:90:VAL:C	12:L:92:ASP:H	2.09	0.55
14:N:21:TYR:CD1	14:N:21:TYR:N	2.73	0.55
17:Q:98:LEU:CA	17:Q:102:GLY:HA2	2.34	0.55
17:Q:104:LYS:HG3	17:Q:105:ALA:N	2.21	0.55
1:A:189:G:H2'	1:A:190:C:H6	1.71	0.55
1:A:256:U:H5'	17:Q:17:LYS:HZ1	1.71	0.55
1:A:438:G:H22	1:A:495:U:H3'	1.71	0.55
1:A:596:C:O2'	1:A:597:G:H5'	2.05	0.55
1:A:909:A:C8	1:A:910:C:C6	2.94	0.55
1:A:986:A:H1'	19:S:55:LYS:CA	2.34	0.55
1:A:1424:C:C2'	1:A:1425:U:H5'	2.36	0.55
2:B:10:LEU:HG	2:B:48:MET:SD	2.46	0.55
2:B:24:TRP:CZ3	2:B:26:PRO:HA	2.41	0.55
2:B:182:ILE:O	2:B:182:ILE:HG22	2.05	0.55
3:C:83:ARG:HA	3:C:86:VAL:CG2	2.27	0.55
4:D:13:ARG:HD3	4:D:36:ARG:O	2.07	0.55
4:D:117:ALA:O	4:D:120:LEU:N	2.39	0.55
4:D:128:VAL:HG21	4:D:138:TYR:CE2	2.41	0.55
4:D:199:ASN:HD22	4:D:199:ASN:C	2.08	0.55
9:I:55:ALA:HA	9:I:58:ARG:CZ	2.36	0.55
10:J:10:GLY:O	10:J:68:HIS:HD2	1.88	0.55
10:J:47:PHE:O	10:J:48:THR:C	2.44	0.55
11:K:54:ARG:O	11:K:57:THR:HG22	2.05	0.55
14:N:14:PRO:HG2	14:N:15:LYS:N	2.15	0.55
16:P:24:ALA:O	16:P:26:ARG:N	2.40	0.55
16:P:26:ARG:HD3	16:P:31:LYS:HB3	1.88	0.55
20:T:11:SER:O	20:T:14:LYS:NZ	2.39	0.55
20:T:96:GLY:O	20:T:97:ALA:CB	2.54	0.55
1:A:463:A:C4	1:A:474:G:C8	2.94	0.55
1:A:848:C:O2'	1:A:849:C:H5'	2.06	0.55
1:A:895:G:H2'	1:A:896:C:H6	1.71	0.55
1:A:1056:U:C5'	3:C:163:ALA:HB2	2.36	0.55
1:A:1250:A:H2'	1:A:1251:A:C8	2.41	0.55
2:B:61:LEU:HD23	2:B:61:LEU:O	2.07	0.55
2:B:80:ILE:HD12	2:B:208:ILE:HD13	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:5:ILE:HG13	3:C:6:HIS:H	1.70	0.55
3:C:91:LEU:CD2	3:C:99:VAL:HG13	2.37	0.55
4:D:103:ASN:OD1	4:D:114:ARG:NE	2.39	0.55
5:E:51:VAL:CB	5:E:52:PRO:HD3	2.32	0.55
12:L:45:PRO:HA	12:L:93:LEU:HD23	1.88	0.55
16:P:7:ALA:HB2	16:P:28:ARG:O	2.06	0.55
17:Q:67:LYS:HA	17:Q:70:ARG:HH12	1.71	0.55
18:R:40:LEU:HD21	18:R:79:LEU:HD21	1.85	0.55
19:S:51:VAL:HG11	19:S:71:LEU:HB3	1.88	0.55
1:A:13:U:O2	1:A:914:A:C8	2.59	0.55
1:A:39:G:C4	1:A:498:U:O4	2.60	0.55
1:A:820:U:H4'	1:A:821:G:OP2	2.06	0.55
1:A:1119:C:H2'	1:A:1120:G:C8	2.41	0.55
1:A:1255:G:N1	1:A:1283:G:C2	2.75	0.55
3:C:22:TRP:HB3	3:C:59:ARG:CG	2.35	0.55
6:F:83:ASP:C	6:F:85:VAL:H	2.10	0.55
7:G:32:ARG:O	7:G:33:ASP:CB	2.54	0.55
8:H:31:PHE:HZ	8:H:134:ILE:HD13	1.71	0.55
8:H:82:HIS:CD2	8:H:83:ILE:N	2.75	0.55
9:I:88:TYR:CZ	9:I:89:ASN:HB2	2.42	0.55
10:J:34:VAL:C	10:J:75:ILE:HB	2.27	0.55
10:J:50:ILE:HG23	14:N:41:ARG:HD3	1.89	0.55
10:J:86:MET:C	10:J:88:LEU:N	2.60	0.55
11:K:99:GLN:HA	11:K:105:VAL:CG2	2.36	0.55
13:M:8:GLU:OE2	13:M:22:ILE:HA	2.06	0.55
13:M:96:LEU:HB3	13:M:97:PRO:HD2	1.88	0.55
19:S:28:LYS:O	19:S:29:ARG:HG3	2.06	0.55
19:S:49:ILE:HG23	19:S:51:VAL:HG22	1.89	0.55
19:S:70:LYS:HG2	19:S:73:GLU:OE2	2.06	0.55
1:A:101:A:C2'	1:A:102:G:H5'	2.37	0.55
1:A:193:C:H2'	1:A:194:C:C6	2.41	0.55
1:A:309:G:C4	1:A:310:G:N7	2.75	0.55
1:A:503:C:C2	1:A:504:C:C5	2.94	0.55
1:A:596:C:O2	1:A:596:C:H2'	2.06	0.55
1:A:1278:U:H5'	1:A:1279:A:C5'	2.35	0.55
1:A:1451:A:H4'	1:A:1452:C:OP2	2.06	0.55
2:B:58:ILE:O	2:B:59:GLU:C	2.45	0.55
3:C:6:HIS:HD2	3:C:8:ILE:H	1.53	0.55
3:C:11:ARG:HH22	3:C:177:THR:C	2.10	0.55
6:F:60:PHE:CZ	18:R:78:LEU:HD21	2.41	0.55
8:H:24:THR:HG23	8:H:24:THR:O	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:30:ARG:O	8:H:31:PHE:O	2.25	0.55
9:I:72:GLY:O	9:I:75:ASP:N	2.34	0.55
16:P:23:ASP:O	16:P:25:ARG:N	2.30	0.55
17:Q:59:ILE:CG2	17:Q:71:PHE:HB3	2.35	0.55
18:R:36:ASN:HD22	18:R:36:ASN:C	2.08	0.55
18:R:59:SER:OG	18:R:62:GLU:HG3	2.06	0.55
1:A:254:G:H5'	17:Q:69:LYS:HD3	1.89	0.55
1:A:411:A:C1'	1:A:413:G:H1'	2.37	0.55
1:A:854:G:H3'	1:A:871:U:C4	2.41	0.55
1:A:1251:A:H8	1:A:1251:A:O5'	1.90	0.55
1:A:1327:C:OP1	21:V:21:TYR:CE1	2.59	0.55
1:A:1442:G:C6	1:A:1446:A:N6	2.74	0.55
2:B:97:TRP:CZ2	2:B:101:MET:HB2	2.41	0.55
3:C:66:VAL:C	3:C:68:VAL:N	2.59	0.55
3:C:85:ARG:C	3:C:87:LEU:N	2.60	0.55
4:D:8:VAL:O	4:D:11:LEU:N	2.40	0.55
4:D:74:GLN:HE22	4:D:137:SER:HB3	1.71	0.55
4:D:206:PHE:CD2	4:D:207:TYR:HE2	2.24	0.55
6:F:97:PHE:HD1	18:R:65:ILE:CD1	2.20	0.55
6:F:97:PHE:CD2	6:F:98:LEU:N	2.74	0.55
7:G:46:ALA:HA	7:G:121:ALA:HB2	1.88	0.55
11:K:125:PHE:N	11:K:125:PHE:CD1	2.73	0.55
15:O:41:GLU:HA	15:O:44:LYS:CG	2.37	0.55
16:P:11:SER:O	16:P:13:HIS:N	2.39	0.55
20:T:75:ASN:ND2	20:T:75:ASN:H	2.01	0.55
21:V:17:THR:HG22	21:V:18:TYR:N	2.14	0.55
1:A:29:G:O2'	1:A:30:U:H5'	2.06	0.55
1:A:101:A:O2'	1:A:102:G:H5'	2.07	0.55
1:A:400:C:H2'	1:A:401:C:C6	2.42	0.55
1:A:706:A:O2'	11:K:29:ILE:HD11	2.07	0.55
1:A:1225:A:H5'	13:M:103:THR:OG1	2.07	0.55
1:A:1477:C:H2'	1:A:1478:C:H6	1.72	0.55
1:A:1497:G:O2'	1:A:1498:U:H5'	2.06	0.55
1:A:1526:G:H2'	1:A:1527:C:H6	1.70	0.55
2:B:12:GLU:HB3	2:B:213:LEU:HD11	1.89	0.55
2:B:13:ALA:C	2:B:15:VAL:N	2.60	0.55
2:B:29:ALA:C	2:B:31:TYR:H	2.10	0.55
3:C:182:ILE:HG22	3:C:183:ASP:O	2.07	0.55
4:D:32:ALA:C	4:D:34:GLU:N	2.60	0.55
4:D:122:ARG:NE	4:D:134:ASP:OD2	2.37	0.55
5:E:94:ALA:HB3	5:E:117:ASP:O	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:16:LEU:HD22	7:G:16:LEU:H	1.71	0.55
7:G:29:LYS:HZ1	7:G:102:ARG:CA	2.11	0.55
7:G:75:VAL:HG11	7:G:86:GLN:HB3	1.88	0.55
9:I:2:GLU:OE1	9:I:20:ARG:HG2	2.06	0.55
11:K:62:GLN:HE22	11:K:97:ALA:HA	1.71	0.55
13:M:55:ARG:HG3	13:M:55:ARG:HH11	1.70	0.55
14:N:8:GLU:O	14:N:11:LYS:HE2	2.06	0.55
15:O:56:LEU:HD12	15:O:59:MET:HB2	1.87	0.55
17:Q:29:HIS:ND1	17:Q:30:PRO:HD2	2.22	0.55
17:Q:97:SER:CB	17:Q:103:GLY:HA2	2.14	0.55
19:S:25:LYS:HD2	19:S:25:LYS:N	2.22	0.55
1:A:402:G:C2'	1:A:403:C:H5'	2.37	0.55
1:A:1102:A:C4	1:A:1103:C:C5	2.94	0.55
1:A:1127:G:H1	1:A:1144:G:H22	1.54	0.55
1:A:1176:A:H2'	1:A:1177:G:C1'	2.36	0.55
1:A:1520:G:O2'	1:A:1521:G:H5'	2.07	0.55
2:B:71:VAL:CG2	2:B:164:VAL:HG13	2.36	0.55
3:C:93:LYS:HA	3:C:93:LYS:CE	2.36	0.55
3:C:122:GLU:O	3:C:126:ARG:HG3	2.07	0.55
4:D:102:ASP:O	4:D:103:ASN:C	2.43	0.55
6:F:1:MET:CE	6:F:36:ARG:HH21	2.20	0.55
7:G:54:THR:HB	7:G:56:GLN:OE1	2.07	0.55
12:L:90:VAL:CG1	12:L:93:LEU:HG	2.32	0.55
12:L:113:ARG:HB2	12:L:122:THR:HG21	1.89	0.55
17:Q:29:HIS:CG	17:Q:30:PRO:HD2	2.42	0.55
18:R:53:ARG:HD3	18:R:63:GLN:HG2	1.89	0.55
20:T:53:LEU:CD1	20:T:100:ILE:HB	2.33	0.55
20:T:62:LEU:O	20:T:65:LYS:HB3	2.07	0.55
1:A:327:A:C3'	1:A:328:C:H5''	2.37	0.55
1:A:880:C:OP2	12:L:6:THR:OG1	2.23	0.55
1:A:1089:G:H1	1:A:1096:C:H42	1.54	0.55
1:A:1150:U:H4'	10:J:41:PRO:HB3	1.88	0.55
1:A:1347:G:H22	1:A:1373:G:H2'	1.62	0.55
1:A:1351:U:O2'	1:A:1352:C:H5'	2.07	0.55
3:C:25:GLY:HA2	3:C:29:TYR:H	1.72	0.55
3:C:133:ALA:O	3:C:136:GLN:HB2	2.07	0.55
3:C:141:VAL:HG23	3:C:142:MET:N	2.21	0.55
4:D:20:TYR:C	4:D:22:LYS:H	2.09	0.55
4:D:49:ARG:HG2	4:D:49:ARG:HH11	1.72	0.55
6:F:82:ARG:O	6:F:84:ASN:N	2.40	0.55
9:I:34:ASN:N	9:I:34:ASN:ND2	2.51	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:N:22:THR:HB	14:N:33:VAL:HG11	1.88	0.55
19:S:63:THR:HG22	19:S:64:GLU:N	2.22	0.55
1:A:191:G:N9	20:T:105:SER:HB3	2.21	0.55
1:A:293:G:O2'	1:A:294:U:H5'	2.07	0.55
1:A:505:G:H1	1:A:526:C:H42	1.54	0.55
1:A:1157:A:H4'	1:A:1158:C:O5'	2.07	0.55
1:A:1349:A:O2'	1:A:1350:A:H5'	2.06	0.55
1:A:1390:U:H2'	1:A:1391:U:H6	1.70	0.55
2:B:8:LYS:HD3	2:B:9:GLU:H	1.72	0.55
2:B:67:THR:HA	2:B:90:MET:HE1	1.88	0.55
2:B:116:GLU:C	2:B:118:LEU:H	2.10	0.55
3:C:143:GLU:OE1	3:C:144:SER:N	2.39	0.55
5:E:26:PHE:C	5:E:27:ARG:HG3	2.26	0.55
9:I:25:LYS:HD3	9:I:25:LYS:N	2.22	0.55
10:J:39:PRO:HA	10:J:70:ARG:HH22	1.72	0.55
13:M:22:ILE:CG2	13:M:66:LEU:HD23	2.37	0.55
16:P:53:VAL:O	16:P:54:GLU:C	2.44	0.55
19:S:22:LEU:CD1	19:S:31:ILE:HD11	2.37	0.55
19:S:70:LYS:H	19:S:73:GLU:CG	2.20	0.55
1:A:153:C:N3	1:A:169:C:N3	2.55	0.54
1:A:293:G:C2	1:A:294:U:C6	2.95	0.54
1:A:321:A:HO2'	1:A:322:C:H6	1.55	0.54
1:A:418:C:O2	1:A:418:C:C2'	2.55	0.54
1:A:782:A:H2'	1:A:783:C:O4'	2.07	0.54
1:A:861:G:C5	1:A:862:C:C5	2.95	0.54
1:A:1197:G:OP1	1:A:1198:G:OP2	2.25	0.54
1:A:1281:U:H4'	1:A:1282:C:OP2	2.07	0.54
1:A:1503:A:H5'	1:A:1531:A:C1'	2.37	0.54
2:B:60:ASP:CG	2:B:61:LEU:N	2.61	0.54
2:B:74:LYS:HE3	2:B:166:ASP:HB2	1.89	0.54
2:B:167:PRO:O	2:B:171:ALA:HB2	2.07	0.54
2:B:200:ILE:HD12	2:B:200:ILE:N	2.22	0.54
4:D:68:TYR:N	4:D:68:TYR:CD1	2.74	0.54
5:E:40:ARG:NH1	5:E:68:GLU:OE2	2.39	0.54
7:G:59:LEU:O	7:G:63:LYS:HG3	2.07	0.54
16:P:9:PHE:O	16:P:10:GLY:O	2.24	0.54
17:Q:34:LYS:HB2	17:Q:34:LYS:NZ	2.22	0.54
17:Q:67:LYS:HA	17:Q:70:ARG:CZ	2.37	0.54
1:A:44:G:H2'	1:A:45:U:O4'	2.07	0.54
1:A:397:A:H3'	1:A:397:A:N3	2.22	0.54
1:A:1001:A:H2'	1:A:1002:G:C5'	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1135:U:H4'	1:A:1136:U:H5	1.71	0.54
1:A:1215:G:C2'	1:A:1216:G:H5'	2.37	0.54
1:A:1346:A:O2'	1:A:1347:G:O4'	2.25	0.54
3:C:50:ALA:O	3:C:72:LYS:N	2.31	0.54
3:C:150:LYS:HB2	3:C:173:VAL:HG21	1.90	0.54
5:E:7:GLU:O	5:E:34:VAL:HA	2.07	0.54
7:G:28:ASN:OD1	7:G:36:LYS:HE3	2.07	0.54
9:I:55:ALA:HA	9:I:58:ARG:NH2	2.22	0.54
13:M:20:THR:HG23	13:M:25:ILE:O	2.07	0.54
15:O:80:ALA:O	15:O:84:LYS:HG3	2.07	0.54
17:Q:62:SER:OG	17:Q:63:ARG:N	2.40	0.54
17:Q:85:VAL:CG1	17:Q:89:LEU:HG	2.38	0.54
1:A:184:G:C4'	1:A:224:C:H4'	2.37	0.54
1:A:191:G:N2	20:T:103:GLY:O	2.36	0.54
1:A:322:C:C2'	1:A:323:U:H5'	2.37	0.54
1:A:498:U:H2'	1:A:498:U:O2	2.07	0.54
1:A:886:G:H2'	1:A:887:G:O4'	2.07	0.54
1:A:949:A:C2	1:A:1233:G:N3	2.75	0.54
2:B:200:ILE:HD12	2:B:200:ILE:H	1.71	0.54
5:E:8:GLU:HA	5:E:34:VAL:CA	2.36	0.54
6:F:38:GLU:O	6:F:39:LYS:CB	2.55	0.54
8:H:19:VAL:CG2	8:H:21:LYS:HG2	2.38	0.54
8:H:27:PRO:HG3	8:H:58:TYR:CE2	2.43	0.54
10:J:64:GLU:O	10:J:65:LEU:HB2	2.08	0.54
12:L:56:ALA:O	12:L:58:VAL:HG23	2.06	0.54
13:M:90:LEU:HA	13:M:93:ARG:HD2	1.90	0.54
20:T:53:LEU:O	20:T:56:MET:HB3	2.07	0.54
1:A:9:G:H5''	5:E:126:ARG:CZ	2.37	0.54
1:A:410:G:OP2	4:D:30:LYS:HD3	2.07	0.54
1:A:442:C:O2'	1:A:443:C:H5'	2.08	0.54
1:A:828:A:H2'	1:A:829:G:O4'	2.08	0.54
1:A:1240:U:C4	7:G:30:ILE:HG23	2.42	0.54
2:B:55:PHE:CD1	2:B:58:ILE:HD12	2.42	0.54
3:C:153:VAL:HG22	3:C:198:VAL:HG21	1.90	0.54
5:E:129:ILE:O	5:E:129:ILE:HG22	2.06	0.54
9:I:72:GLY:O	9:I:75:ASP:HB2	2.08	0.54
12:L:28:LYS:O	12:L:30:ALA:N	2.41	0.54
13:M:94:ARG:O	13:M:96:LEU:HD12	2.08	0.54
19:S:15:LEU:HB3	19:S:33:THR:CG2	2.34	0.54
1:A:451:A:H4'	16:P:72:ARG:HH12	1.72	0.54
1:A:924:C:H2'	1:A:925:G:H8	1.71	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:988:G:HO2'	1:A:1016:A:H2	1.52	0.54
1:A:1187:G:H2'	1:A:1188:A:H8	1.72	0.54
2:B:44:LEU:HA	2:B:47:THR:HB	1.89	0.54
3:C:66:VAL:O	3:C:68:VAL:N	2.40	0.54
5:E:88:LYS:HD3	5:E:123:LEU:HB2	1.89	0.54
6:F:32:ASN:O	6:F:32:ASN:ND2	2.41	0.54
13:M:31:LYS:C	13:M:33:ALA:N	2.61	0.54
19:S:36:ARG:HH21	19:S:75:ALA:CB	2.20	0.54
19:S:38:SER:H	19:S:71:LEU:CD1	2.21	0.54
19:S:44:MET:O	19:S:46:GLY:N	2.40	0.54
20:T:90:GLN:HG3	20:T:91:LEU:N	2.21	0.54
1:A:123:C:H42	1:A:238:G:H1	1.56	0.54
1:A:323:U:H2'	1:A:324:G:O4'	2.08	0.54
1:A:444:C:N4	1:A:491:G:C6	2.75	0.54
1:A:481:G:H5'	1:A:482:A:OP1	2.08	0.54
1:A:508:C:H4'	1:A:509:A:O5'	2.08	0.54
1:A:707:C:O2'	1:A:708:C:H5'	2.08	0.54
1:A:833:U:H3	1:A:853:G:H1	1.55	0.54
2:B:151:GLY:O	2:B:153:ARG:N	2.40	0.54
3:C:64:VAL:HG12	3:C:65:ALA:N	2.22	0.54
3:C:139:GLN:O	3:C:142:MET:HB3	2.07	0.54
4:D:54:TYR:CE1	4:D:206:PHE:HE1	2.26	0.54
4:D:158:ILE:O	4:D:162:LEU:HD13	2.08	0.54
8:H:116:LYS:NZ	8:H:127:LEU:HB3	2.23	0.54
9:I:48:GLU:CD	9:I:51:ARG:HH21	2.11	0.54
14:N:11:LYS:NZ	14:N:13:THR:O	2.25	0.54
17:Q:101:ARG:HG2	17:Q:101:ARG:NH1	2.23	0.54
20:T:10:LEU:O	20:T:12:ALA:N	2.41	0.54
1:A:13:U:O2	1:A:914:A:H3'	2.07	0.54
1:A:412:A:H4'	1:A:413:G:OP1	2.06	0.54
1:A:425:G:O2'	1:A:426:G:H5'	2.08	0.54
1:A:815:A:O2'	1:A:1527:C:C1'	2.54	0.54
1:A:823:G:O2'	1:A:824:C:H5'	2.08	0.54
1:A:1074:G:C2	1:A:1102:A:C5	2.96	0.54
1:A:1122:U:H2'	1:A:1123:A:O4'	2.08	0.54
1:A:1192:C:O2	1:A:1192:C:H2'	2.06	0.54
2:B:116:GLU:O	2:B:118:LEU:N	2.41	0.54
3:C:23:TYR:HD1	10:J:11:PHE:CE1	2.25	0.54
3:C:88:ARG:HA	3:C:91:LEU:HD13	1.90	0.54
5:E:109:ILE:O	5:E:113:ALA:CB	2.55	0.54
5:E:144:THR:O	5:E:148:VAL:HG23	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:33:TYR:C	6:F:71:ARG:HH21	2.11	0.54
10:J:79:ARG:NH1	10:J:79:ARG:HB2	2.23	0.54
11:K:63:LEU:O	11:K:67:ASP:N	2.41	0.54
13:M:109:THR:HG23	13:M:110:ARG:N	2.22	0.54
14:N:11:LYS:CE	14:N:13:THR:HB	2.37	0.54
14:N:26:ARG:O	14:N:27:CYS:C	2.46	0.54
14:N:37:PHE:N	14:N:37:PHE:CD1	2.75	0.54
15:O:56:LEU:O	15:O:59:MET:HB2	2.08	0.54
17:Q:18:THR:HG21	17:Q:69:LYS:HD2	1.89	0.54
21:V:6:ARG:HD2	21:V:15:ARG:NH2	2.21	0.54
1:A:66:G:N3	1:A:66:G:H2'	2.23	0.54
1:A:265:G:H2'	1:A:267:C:H5	1.72	0.54
1:A:399:G:C5	1:A:400:C:C4	2.96	0.54
1:A:603:U:H3	1:A:635:G:H1	1.56	0.54
1:A:1094:G:O2'	1:A:1108:G:N2	2.41	0.54
2:B:131:PRO:C	2:B:133:LYS:H	2.11	0.54
4:D:126:ILE:HG22	4:D:127:THR:H	1.73	0.54
5:E:19:MET:HE3	5:E:23:GLY:O	2.08	0.54
7:G:74:GLU:HG2	7:G:91:VAL:HG22	1.90	0.54
8:H:38:ILE:HG22	8:H:39:LEU:N	2.22	0.54
12:L:93:LEU:HB2	12:L:96:VAL:HG21	1.89	0.54
19:S:18:LYS:HG3	19:S:18:LYS:O	2.07	0.54
1:A:129(A):G:N3	1:A:190(E):U:H5''	2.23	0.54
1:A:316:G:H2'	1:A:317:G:C8	2.43	0.54
1:A:374:A:H2'	1:A:375:U:C6	2.43	0.54
1:A:950:U:H5	13:M:102:ARG:CZ	2.21	0.54
1:A:1177:G:H2'	1:A:1178:G:H8	1.72	0.54
1:A:1249:C:O2	9:I:70:LYS:HE3	2.08	0.54
1:A:1275:A:H2'	1:A:1276:G:C8	2.43	0.54
1:A:1352:C:H42	1:A:1370:G:H1	1.56	0.54
1:A:1419:G:C6	1:A:1482:G:H1'	2.43	0.54
2:B:79:ASP:O	2:B:80:ILE:C	2.47	0.54
3:C:23:TYR:HA	10:J:11:PHE:CE1	2.42	0.54
3:C:84:ILE:HG23	3:C:84:ILE:O	2.08	0.54
4:D:28:SER:O	4:D:29:PRO:C	2.46	0.54
4:D:100:ARG:O	4:D:101:LEU:C	2.45	0.54
5:E:47:LYS:N	5:E:47:LYS:HD2	2.23	0.54
6:F:97:PHE:C	6:F:98:LEU:HG	2.26	0.54
8:H:100:ILE:HG13	8:H:112:LEU:HD21	1.88	0.54
12:L:41:ARG:HD2	12:L:42:THR:N	2.22	0.54
12:L:78:GLN:O	12:L:80:HIS:N	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:L:101:VAL:O	12:L:102:ARG:C	2.45	0.54
12:L:103:GLY:HA2	12:L:108:ALA:HA	1.90	0.54
13:M:45:VAL:C	13:M:47:ASP:H	2.11	0.54
13:M:81:LEU:HA	13:M:84:ILE:CD1	2.38	0.54
13:M:91:ARG:HB3	13:M:97:PRO:O	2.08	0.54
14:N:14:PRO:O	14:N:15:LYS:HB2	2.06	0.54
14:N:44:LEU:HD12	14:N:44:LEU:O	2.08	0.54
16:P:76:GLN:C	16:P:78:GLY:H	2.10	0.54
18:R:75:ILE:C	18:R:77:GLY:H	2.11	0.54
20:T:15:ARG:O	20:T:16:HIS:C	2.46	0.54
20:T:84:LEU:O	20:T:87:LYS:N	2.41	0.54
1:A:99:C:H2'	1:A:101:A:O4'	2.08	0.54
1:A:124:G:C6	1:A:125:U:C4	2.96	0.54
1:A:227:G:H2'	1:A:228:A:H8	1.71	0.54
1:A:267:C:N4	1:A:268:C:N4	2.56	0.54
1:A:341:C:O2	1:A:349:A:C2	2.61	0.54
1:A:428:G:OP2	4:D:7:PRO:HG3	2.07	0.54
1:A:715:A:H2'	1:A:716:A:C8	2.43	0.54
1:A:1004:A:H5''	1:A:1025:U:H5	1.73	0.54
1:A:1442:G:C4	1:A:1446:A:N1	2.76	0.54
2:B:9:GLU:HG3	2:B:217:ARG:NH1	2.23	0.54
2:B:87:ARG:HH11	2:B:219:VAL:HB	1.70	0.54
4:D:2:GLY:N	4:D:3:ARG:NH2	2.56	0.54
5:E:70:PRO:C	5:E:72:GLN:N	2.57	0.54
9:I:10:ARG:HD3	9:I:105:ASP:CB	2.38	0.54
10:J:47:PHE:CE2	14:N:37:PHE:CZ	2.96	0.54
11:K:90:GLY:O	11:K:91:ARG:C	2.45	0.54
15:O:71:GLN:HB2	15:O:78:TYR:CG	2.42	0.54
20:T:67:ALA:HA	20:T:73:HIS:N	2.15	0.54
21:V:7:ARG:HB2	21:V:21:TYR:CZ	2.43	0.54
1:A:115:G:C2	1:A:313:A:C2	2.96	0.53
1:A:152:A:N6	1:A:170:U:C2	2.76	0.53
1:A:290:C:H2'	1:A:291:C:H5'	1.88	0.53
1:A:303:A:O2'	1:A:304:U:H5'	2.08	0.53
1:A:340:U:H2'	1:A:341:C:C6	2.44	0.53
1:A:429:U:C1'	1:A:430:A:H5''	2.38	0.53
1:A:529:G:C2'	1:A:530:G:H5'	2.38	0.53
1:A:864:A:H2	1:A:917:G:N3	2.05	0.53
1:A:874:G:O2'	1:A:875:C:H5'	2.07	0.53
1:A:877:C:HO2'	1:A:878:G:H5'	1.70	0.53
1:A:1078:U:O2'	1:A:1079:G:H5'	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1200:C:N3	1:A:1206:G:O6	2.40	0.53
1:A:1276:G:H21	1:A:1282:C:C1'	2.20	0.53
1:A:1347:G:HO2'	1:A:1348:U:H5	1.52	0.53
7:G:67:GLU:HA	7:G:70:LYS:CD	2.38	0.53
9:I:97:LYS:NZ	9:I:102:LEU:CD1	2.71	0.53
13:M:49:THR:CG2	13:M:50:GLU:N	2.72	0.53
20:T:13:LEU:CD1	20:T:14:LYS:N	2.70	0.53
20:T:51:GLU:O	20:T:55:ILE:HG23	2.07	0.53
20:T:76:ALA:O	20:T:77:ALA:C	2.46	0.53
1:A:291:C:C2'	1:A:292:G:H5'	2.38	0.53
1:A:298:A:H2'	1:A:299:G:O4'	2.08	0.53
1:A:868:C:H2'	1:A:869:G:H5'	1.89	0.53
1:A:913:A:O2'	1:A:914:A:OP2	2.26	0.53
1:A:931:C:H42	1:A:1386:G:H1	1.57	0.53
2:B:96:ARG:NE	2:B:96:ARG:CA	2.72	0.53
2:B:130:ARG:HB3	2:B:131:PRO:CD	2.34	0.53
7:G:20:ASP:OD2	7:G:23:VAL:HG13	2.08	0.53
8:H:51:VAL:HG21	8:H:60:ARG:HG3	1.89	0.53
11:K:101:SER:OG	11:K:102:GLY:N	2.40	0.53
12:L:53:ARG:HG3	12:L:93:LEU:HD21	1.90	0.53
14:N:22:THR:HB	14:N:33:VAL:HB	1.89	0.53
15:O:7:GLU:O	15:O:10:LYS:HB3	2.08	0.53
17:Q:6:LEU:O	17:Q:59:ILE:N	2.40	0.53
19:S:40:ILE:HD12	19:S:69:HIS:HB2	1.90	0.53
1:A:187:C:O2'	20:T:89:ARG:CD	2.56	0.53
1:A:358:U:H2'	1:A:359:U:C6	2.43	0.53
1:A:623:C:H2'	1:A:624:C:H6	1.71	0.53
1:A:723:U:O2	1:A:723:U:H2'	2.08	0.53
1:A:1089:G:C2'	1:A:1090:U:H5'	2.39	0.53
1:A:1109:C:OP2	3:C:176:HIS:CD2	2.61	0.53
1:A:1259:C:O2	1:A:1259:C:H2'	2.07	0.53
2:B:19:HIS:CD2	2:B:20:GLU:HG2	2.44	0.53
3:C:16:ARG:NH2	3:C:183:ASP:HB2	2.22	0.53
3:C:115:LEU:HA	3:C:118:GLN:OE1	2.09	0.53
5:E:108:ALA:O	5:E:109:ILE:C	2.47	0.53
7:G:21:VAL:O	7:G:24:THR:N	2.41	0.53
7:G:70:LYS:HG2	7:G:100:ALA:HB2	1.88	0.53
8:H:104:ARG:C	8:H:106:GLY:N	2.61	0.53
9:I:4:TYR:CE1	9:I:21:PRO:HG2	2.43	0.53
9:I:16:ARG:NH2	9:I:64:THR:HG22	2.24	0.53
12:L:60:LEU:HD11	12:L:85:ILE:HD12	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:T:94:ALA:O	20:T:95:ALA:HB2	2.08	0.53
1:A:20:U:O2'	1:A:21:G:H5'	2.08	0.53
1:A:167:G:C2'	1:A:168:G:H8	2.17	0.53
1:A:1519:A:H2'	1:A:1520:G:C5'	2.34	0.53
2:B:51:LEU:HD22	2:B:55:PHE:CE2	2.44	0.53
2:B:178:ARG:NH2	2:B:196:LEU:O	2.41	0.53
16:P:39:TYR:OH	16:P:41:PRO:HA	2.09	0.53
17:Q:5:VAL:HG12	17:Q:6:LEU:N	2.23	0.53
19:S:7:LYS:HG3	19:S:7:LYS:O	2.08	0.53
19:S:11:VAL:HG22	19:S:39:THR:CG2	2.31	0.53
19:S:17:GLU:C	19:S:19:VAL:H	2.12	0.53
1:A:182:U:H3'	1:A:182:U:OP2	2.08	0.53
1:A:327:A:H3'	1:A:328:C:H5''	1.91	0.53
1:A:520:A:H61	1:A:529:G:H1'	1.74	0.53
1:A:538:G:C4'	12:L:114:LYS:HD3	2.37	0.53
1:A:987:G:N2	1:A:1219:U:N3	2.57	0.53
1:A:1109:C:OP1	3:C:176:HIS:NE2	2.41	0.53
1:A:1206:G:C6	1:A:1207:G:C6	2.97	0.53
1:A:1225:A:H2'	1:A:1226:C:C6	2.44	0.53
1:A:1325:C:O3'	21:V:17:THR:HG21	2.09	0.53
1:A:1376:U:H2'	1:A:1377:A:H8	1.72	0.53
1:A:1407:C:H2'	1:A:1408:A:C8	2.43	0.53
1:A:1514:C:O2'	1:A:1515:C:H5'	2.09	0.53
2:B:47:THR:O	2:B:48:MET:C	2.47	0.53
2:B:151:GLY:C	2:B:153:ARG:N	2.61	0.53
2:B:170:GLU:OE2	2:B:172:ILE:HD13	2.07	0.53
3:C:7:PRO:HB2	3:C:11:ARG:HD2	1.90	0.53
5:E:18:ARG:HE	5:E:25:ARG:HB2	1.73	0.53
5:E:41:VAL:HG12	5:E:42:GLY:H	1.73	0.53
7:G:139:GLU:O	7:G:140:ASP:C	2.47	0.53
12:L:89:ARG:HH21	12:L:97:ARG:CG	2.07	0.53
13:M:74:VAL:C	13:M:76:ALA:N	2.62	0.53
20:T:65:LYS:O	20:T:68:LYS:CB	2.56	0.53
1:A:44:G:H1'	1:A:399:G:N2	2.24	0.53
1:A:52:G:H2'	1:A:53:A:H8	1.74	0.53
1:A:300:A:H8	1:A:300:A:O5'	1.91	0.53
1:A:439:A:C4	1:A:497:A:C2	2.96	0.53
1:A:612:C:O2'	1:A:613:C:H5'	2.08	0.53
1:A:855:G:H2'	1:A:856:C:C6	2.44	0.53
1:A:960:U:H2'	1:A:1225:A:N6	2.24	0.53
1:A:1105:A:H2'	1:A:1106:G:H8	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1204:A:C6	1:A:1205:U:C2	2.96	0.53
3:C:108:ASN:OD1	3:C:111:LEU:HD12	2.09	0.53
3:C:118:GLN:HA	3:C:121:ALA:HB3	1.91	0.53
3:C:150:LYS:HA	3:C:169:ALA:HA	1.89	0.53
4:D:21:LEU:HD12	4:D:21:LEU:H	1.74	0.53
4:D:103:ASN:O	4:D:104:VAL:C	2.46	0.53
4:D:156:GLU:O	4:D:157:LEU:C	2.47	0.53
5:E:18:ARG:HH21	5:E:25:ARG:CB	2.19	0.53
6:F:74:ASP:HA	6:F:77:ARG:NE	2.24	0.53
9:I:18:PHE:HB2	9:I:62:TYR:HB3	1.90	0.53
9:I:40:LEU:O	9:I:42:ARG:N	2.41	0.53
9:I:89:ASN:O	9:I:91:ASP:N	2.38	0.53
11:K:91:ARG:O	11:K:92:GLU:C	2.47	0.53
13:M:113:PRO:O	13:M:115:LYS:HG3	2.08	0.53
15:O:62:GLN:O	15:O:65:ARG:HB3	2.07	0.53
16:P:52:ASP:OD2	16:P:54:GLU:HB3	2.08	0.53
16:P:53:VAL:HB	16:P:79:VAL:HG13	1.89	0.53
16:P:71:ARG:HA	16:P:74:LEU:HD12	1.91	0.53
19:S:10:PHE:O	19:S:39:THR:HB	2.07	0.53
20:T:46:GLU:HB2	20:T:48:LYS:HE2	1.91	0.53
1:A:428:G:C1'	1:A:430:A:C8	2.90	0.53
1:A:490:G:C2	1:A:491:G:C8	2.96	0.53
1:A:849:C:H2'	1:A:850:U:H5'	1.89	0.53
1:A:1179:A:O2'	9:I:104:ARG:HB2	2.07	0.53
1:A:1243:C:O2'	1:A:1244:C:H5'	2.08	0.53
2:B:22:LYS:HG3	2:B:40:HIS:CE1	2.44	0.53
4:D:15:GLU:O	4:D:63:LYS:NZ	2.41	0.53
4:D:162:LEU:HD23	4:D:181:MET:HG2	1.90	0.53
8:H:54:ASP:O	8:H:54:ASP:CG	2.47	0.53
8:H:116:LYS:HZ3	8:H:127:LEU:HB3	1.72	0.53
9:I:18:PHE:HD1	9:I:62:TYR:HD2	1.57	0.53
10:J:90:LEU:H	10:J:91:PRO:HD2	1.72	0.53
12:L:34:ARG:HB3	12:L:61:THR:HG21	1.89	0.53
12:L:101:VAL:O	12:L:101:VAL:HG23	2.08	0.53
13:M:87:TYR:O	13:M:88:ARG:C	2.46	0.53
17:Q:92:ARG:O	17:Q:95:TYR:HB2	2.09	0.53
18:R:36:ASN:ND2	18:R:37:VAL:O	2.42	0.53
20:T:88:VAL:HA	20:T:91:LEU:HD12	1.90	0.53
1:A:379:C:O2'	1:A:380:G:H5'	2.08	0.53
1:A:419:C:O2	1:A:419:C:C3'	2.57	0.53
1:A:715:A:O2'	1:A:716:A:H5'	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1190:G:OP1	3:C:4:LYS:HA	2.09	0.53
1:A:1359:C:OP2	14:N:22:THR:HG21	2.09	0.53
1:A:1407:C:H2'	1:A:1408:A:O4'	2.09	0.53
1:A:1526:G:O2'	1:A:1527:C:H5'	2.09	0.53
2:B:225:ALA:O	2:B:226:ARG:C	2.46	0.53
3:C:12:LEU:C	3:C:14:ILE:H	2.11	0.53
3:C:77:ILE:HD13	3:C:84:ILE:HD12	1.90	0.53
3:C:126:ARG:C	3:C:127:ARG:HD2	2.28	0.53
3:C:167:TRP:O	3:C:168:ALA:HB2	2.08	0.53
3:C:175:LEU:HD22	3:C:175:LEU:N	2.24	0.53
3:C:176:HIS:O	3:C:178:LEU:N	2.37	0.53
4:D:21:LEU:H	4:D:21:LEU:CD1	2.22	0.53
6:F:68:PRO:O	6:F:72:VAL:HG22	2.08	0.53
7:G:67:GLU:HA	7:G:70:LYS:CG	2.39	0.53
9:I:118:LYS:NZ	9:I:118:LYS:CB	2.72	0.53
10:J:31:GLY:O	10:J:32:ALA:CB	2.56	0.53
10:J:80:LYS:O	10:J:84:GLN:CB	2.55	0.53
12:L:117:ARG:HG3	12:L:117:ARG:HH11	1.74	0.53
13:M:78:ILE:C	13:M:80:ARG:N	2.62	0.53
14:N:39:LEU:HD22	14:N:43:CYS:HB3	1.90	0.53
15:O:3:ILE:HD12	15:O:3:ILE:N	2.23	0.53
15:O:6:GLU:H	15:O:6:GLU:CD	2.12	0.53
17:Q:40:LYS:HD3	17:Q:42:TYR:CZ	2.44	0.53
19:S:55:LYS:HE2	19:S:56:GLN:HE21	1.72	0.53
20:T:33:ILE:O	20:T:37:SER:N	2.36	0.53
1:A:59:A:N6	1:A:331:G:H1'	2.23	0.53
1:A:103:C:O2'	1:A:172:A:N1	2.35	0.53
1:A:198:G:H1	1:A:219:C:H42	1.56	0.53
1:A:377:G:C2	1:A:387:U:O2	2.62	0.53
1:A:443:C:O2'	1:A:444:C:H5'	2.09	0.53
1:A:597:G:H2'	1:A:598:U:C5'	2.38	0.53
1:A:877:C:H1'	8:H:3:THR:CG2	2.39	0.53
1:A:1230:C:H2'	1:A:1231:G:H8	1.74	0.53
1:A:1418:A:H61	1:A:1482:G:C2'	2.21	0.53
2:B:100:GLY:O	2:B:101:MET:C	2.46	0.53
3:C:27:LYS:HA	3:C:30:ARG:HH22	1.73	0.53
3:C:64:VAL:HG12	3:C:65:ALA:H	1.73	0.53
3:C:104:GLN:HA	3:C:104:GLN:HE21	1.74	0.53
3:C:179:ARG:CD	3:C:180:ALA:N	2.71	0.53
6:F:100:ASN:H	18:R:23:LYS:HE3	1.72	0.53
8:H:120:THR:HG23	8:H:123:GLU:HB2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:30:GLY:O	9:I:31:GLN:C	2.47	0.53
9:I:96:LEU:O	9:I:102:LEU:HD21	2.09	0.53
10:J:49:VAL:C	10:J:60:ARG:HG3	2.29	0.53
14:N:22:THR:HB	14:N:33:VAL:CG1	2.39	0.53
15:O:60:VAL:O	15:O:64:ARG:HG2	2.09	0.53
16:P:75:ARG:C	16:P:78:GLY:H	2.12	0.53
18:R:36:ASN:C	18:R:36:ASN:ND2	2.61	0.53
20:T:33:ILE:HG22	20:T:34:LYS:N	2.22	0.53
20:T:46:GLU:HB3	20:T:48:LYS:HE2	1.91	0.53
1:A:67:C:H2'	1:A:68:G:C8	2.44	0.53
1:A:597:G:N2	8:H:94:TYR:CE2	2.77	0.53
1:A:1026:G:H2'	1:A:1026:G:N3	2.23	0.53
1:A:1119:C:H2'	1:A:1120:G:H8	1.74	0.53
1:A:1240:U:OP1	7:G:119:ARG:NH2	2.33	0.53
1:A:1426:C:H2'	1:A:1427:U:C6	2.43	0.53
2:B:181:PHE:CE2	8:H:70:GLN:HB3	2.44	0.53
3:C:131:ARG:O	3:C:134:ILE:HB	2.09	0.53
4:D:100:ARG:O	4:D:102:ASP:N	2.42	0.53
4:D:101:LEU:O	4:D:104:VAL:HB	2.09	0.53
6:F:75:LEU:C	6:F:75:LEU:HD13	2.30	0.53
7:G:16:LEU:H	7:G:16:LEU:CD2	2.22	0.53
7:G:156:TRP:CD1	7:G:156:TRP:O	2.61	0.53
8:H:96:GLY:N	8:H:99:GLU:HB2	2.23	0.53
12:L:69:TYR:CZ	12:L:70:ILE:O	2.62	0.53
13:M:8:GLU:CG	13:M:22:ILE:HG12	2.37	0.53
13:M:88:ARG:HA	13:M:98:VAL:HG13	1.91	0.53
15:O:16:ALA:C	15:O:18:PHE:H	2.12	0.53
15:O:29:VAL:HG11	15:O:67:LEU:HD21	1.91	0.53
18:R:30:ASP:C	18:R:32:ARG:H	2.12	0.53
1:A:619:U:O2	4:D:133:VAL:HG13	2.09	0.52
1:A:646:U:H2'	1:A:647:C:C6	2.44	0.52
1:A:1355:G:C4	1:A:1356:G:C8	2.97	0.52
1:A:1473:A:H2'	1:A:1474:G:O4'	2.09	0.52
3:C:191:THR:HG23	3:C:194:GLY:H	1.73	0.52
4:D:9:CYS:SG	4:D:32:ALA:HB3	2.49	0.52
8:H:28:ALA:HA	8:H:59:LEU:CD1	2.38	0.52
9:I:112:LYS:HE2	9:I:116:LYS:O	2.09	0.52
13:M:33:ALA:C	13:M:35:GLU:H	2.11	0.52
16:P:40:ASP:HB3	16:P:48:TRP:HB2	1.89	0.52
1:A:390:C:H6	1:A:390:C:O5'	1.91	0.52
1:A:423:G:C3'	1:A:424:G:H5'	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:836:G:H2'	1:A:837:G:C8	2.45	0.52
1:A:941:G:H1	1:A:1342:C:H42	1.57	0.52
1:A:986:A:H2'	1:A:987:G:O4'	2.09	0.52
1:A:1065:U:O4	1:A:1190:G:C1'	2.56	0.52
5:E:126:ARG:HH11	5:E:126:ARG:CG	2.22	0.52
7:G:151:TYR:HA	7:G:153:HIS:CE1	2.43	0.52
8:H:10:LEU:CD2	8:H:83:ILE:HD11	2.39	0.52
8:H:86:ILE:HG22	8:H:87:SER:H	1.72	0.52
9:I:19:LEU:HD23	9:I:21:PRO:CD	2.39	0.52
10:J:62:HIS:C	14:N:59:ALA:HB3	2.30	0.52
10:J:68:HIS:H	10:J:68:HIS:CD2	2.26	0.52
15:O:8:LYS:O	15:O:9:GLN:C	2.47	0.52
16:P:17:TYR:N	16:P:17:TYR:CD1	2.78	0.52
18:R:76:LEU:O	18:R:78:LEU:HG	2.09	0.52
20:T:14:LYS:O	20:T:17:ARG:N	2.39	0.52
1:A:39:G:H1'	1:A:498:U:C5	2.44	0.52
1:A:191:G:C1'	20:T:105:SER:HB3	2.40	0.52
1:A:328:C:H4'	1:A:329:A:H5'	1.90	0.52
1:A:377:G:H5''	16:P:24:ALA:HB1	1.90	0.52
1:A:568:G:N2	1:A:883:C:C6	2.77	0.52
1:A:756:C:O2'	1:A:757:U:H5'	2.10	0.52
1:A:774:G:N2	1:A:775:G:H1'	2.24	0.52
1:A:1078:U:C2'	1:A:1079:G:H5'	2.39	0.52
1:A:1203:C:H5'	14:N:2:ALA:HB1	1.90	0.52
1:A:1312:G:N2	1:A:1326:C:N3	2.57	0.52
1:A:1437:C:H2'	1:A:1438:G:C8	2.45	0.52
3:C:91:LEU:HD21	3:C:99:VAL:H	1.74	0.52
3:C:139:GLN:HA	3:C:142:MET:HB3	1.91	0.52
4:D:125:HIS:O	4:D:126:ILE:HD13	2.09	0.52
7:G:46:ALA:CA	7:G:121:ALA:HB2	2.39	0.52
7:G:101:LEU:C	7:G:103:TRP:N	2.63	0.52
7:G:115:ARG:C	7:G:119:ARG:HH21	2.13	0.52
11:K:21:ILE:HD11	11:K:95:ILE:HA	1.91	0.52
11:K:98:LEU:N	11:K:98:LEU:HD23	2.24	0.52
13:M:64:TRP:HB2	13:M:66:LEU:HD11	1.91	0.52
14:N:23:ARG:N	14:N:33:VAL:HG21	2.25	0.52
16:P:39:TYR:CD1	16:P:73:LEU:CD2	2.93	0.52
1:A:38:G:N2	1:A:397:A:H5''	2.24	0.52
1:A:91:C:H2'	1:A:92:C:H6	1.73	0.52
1:A:102:G:O2'	1:A:103:C:H5'	2.10	0.52
1:A:123:C:OP1	1:A:312:C:H5'	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:718:G:H4'	11:K:117:ASN:HD21	1.73	0.52
1:A:922:G:N2	1:A:1396:A:C5	2.78	0.52
1:A:1142:G:H2'	1:A:1143:G:O4'	2.09	0.52
1:A:1219:U:H2'	1:A:1220:G:H8	1.74	0.52
1:A:1316:G:H4'	14:N:18:VAL:CG1	2.40	0.52
2:B:210:SER:C	2:B:212:GLN:N	2.62	0.52
4:D:105:VAL:HG12	4:D:106:TYR:N	2.24	0.52
4:D:121:VAL:HG12	4:D:134:ASP:O	2.09	0.52
4:D:190:ASP:O	4:D:192:GLU:N	2.42	0.52
6:F:97:PHE:CD1	18:R:65:ILE:HD12	2.37	0.52
8:H:114:THR:HG22	8:H:130:GLY:C	2.29	0.52
9:I:50:LEU:HD11	9:I:81:ILE:HG22	1.91	0.52
9:I:97:LYS:O	9:I:99:LEU:N	2.42	0.52
12:L:81:SER:HB3	12:L:106:ASP:HB2	1.91	0.52
14:N:43:CYS:O	14:N:47:LEU:HB2	2.10	0.52
15:O:48:LYS:H	15:O:48:LYS:CD	2.21	0.52
17:Q:51:TYR:CD2	17:Q:73:VAL:HG11	2.43	0.52
19:S:22:LEU:HD22	19:S:28:LYS:HG3	1.91	0.52
20:T:39:LYS:O	20:T:43:LEU:HG	2.09	0.52
20:T:79:ARG:O	20:T:83:ARG:N	2.43	0.52
1:A:437:U:H2'	1:A:438:G:O4'	2.10	0.52
1:A:450:G:N2	1:A:482:A:H61	2.08	0.52
1:A:858:G:O6	1:A:869:G:C8	2.62	0.52
1:A:1276:G:H2'	1:A:1277:C:H5'	1.92	0.52
2:B:73:THR:HG21	2:B:169:LYS:HE3	1.92	0.52
2:B:204:ASN:ND2	2:B:204:ASN:C	2.62	0.52
3:C:191:THR:OG1	3:C:193:TYR:CD2	2.63	0.52
5:E:101:ILE:HD12	5:E:119:LEU:HD21	1.92	0.52
5:E:139:LEU:HA	5:E:142:LEU:HG	1.90	0.52
5:E:144:THR:C	5:E:146:ALA:N	2.63	0.52
7:G:111:ARG:HD2	7:G:123:GLU:OE2	2.10	0.52
8:H:100:ILE:HG23	8:H:101:PRO:HD2	1.90	0.52
12:L:88:GLY:C	12:L:89:ARG:HG2	2.30	0.52
16:P:2:VAL:O	16:P:64:ALA:CA	2.57	0.52
18:R:56:THR:O	18:R:58:LEU:N	2.43	0.52
1:A:145:G:H2'	1:A:146:G:H8	1.75	0.52
1:A:665:A:H2'	1:A:732:C:O2	2.10	0.52
1:A:799:G:C2'	1:A:800:G:H5'	2.39	0.52
1:A:1046:A:H2'	1:A:1046:A:N3	2.25	0.52
1:A:1487:G:H2'	1:A:1488:G:H5'	1.91	0.52
1:A:1511:G:C6	1:A:1512:U:N3	2.78	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1526:G:H2'	1:A:1527:C:C6	2.44	0.52
2:B:44:LEU:HA	2:B:47:THR:OG1	2.09	0.52
7:G:14:PRO:O	7:G:15:ASP:C	2.47	0.52
7:G:65:ALA:CB	7:G:128:ALA:CA	2.84	0.52
7:G:135:VAL:O	7:G:138:LYS:HD3	2.09	0.52
8:H:87:SER:HB2	8:H:93:VAL:CB	2.24	0.52
9:I:53:VAL:HG21	9:I:85:LEU:HD13	1.90	0.52
16:P:21:VAL:HG11	16:P:59:TRP:NE1	2.23	0.52
16:P:28:ARG:HD2	16:P:29:ASP:OD2	2.10	0.52
18:R:35:ARG:O	18:R:37:VAL:N	2.42	0.52
19:S:23:ASN:O	19:S:25:LYS:N	2.43	0.52
20:T:67:ALA:CA	20:T:73:HIS:H	2.15	0.52
1:A:277:C:H5'	17:Q:68:ARG:NH1	2.24	0.52
1:A:316:G:C2	1:A:317:G:C5	2.98	0.52
1:A:415:A:H2'	1:A:416:G:C8	2.44	0.52
1:A:529:G:H4'	1:A:533:A:C2	2.44	0.52
1:A:986:A:O2'	1:A:987:G:H5'	2.09	0.52
1:A:1118:C:H5'	9:I:104:ARG:HD2	1.92	0.52
1:A:1121:U:H2'	1:A:1122:U:C6	2.45	0.52
1:A:1129:C:H1'	1:A:1131:G:C8	2.44	0.52
1:A:1240:U:C1'	7:G:38:LEU:HD11	2.40	0.52
1:A:1372:U:O2'	1:A:1373:G:H5'	2.09	0.52
1:A:1427:U:O4	1:A:1428:A:N6	2.43	0.52
2:B:161:ALA:O	2:B:162:ILE:HG13	2.09	0.52
3:C:92:ALA:CA	3:C:95:THR:HG22	2.32	0.52
3:C:113:ALA:HB3	3:C:183:ASP:OD2	2.09	0.52
4:D:8:VAL:C	4:D:10:ARG:N	2.61	0.52
6:F:18:GLN:C	6:F:21:LEU:HB3	2.30	0.52
7:G:61:VAL:O	7:G:65:ALA:CB	2.57	0.52
9:I:64:THR:CG2	9:I:66:ARG:HH21	2.23	0.52
10:J:14:LYS:HA	10:J:17:ASP:CB	2.35	0.52
10:J:31:GLY:HA3	10:J:81:THR:OG1	2.10	0.52
12:L:40:VAL:O	12:L:40:VAL:HG12	2.10	0.52
15:O:73:GLU:O	15:O:74:ASP:HB2	2.10	0.52
16:P:74:LEU:C	16:P:80:PHE:HE1	2.12	0.52
17:Q:90:ILE:O	17:Q:91:ARG:C	2.47	0.52
18:R:37:VAL:HB	18:R:41:LYS:CG	2.39	0.52
1:A:906:G:H8	1:A:906:G:O5'	1.92	0.52
1:A:909:A:H2'	1:A:910:C:O4'	2.10	0.52
1:A:915:A:H2'	1:A:916:G:C5'	2.40	0.52
1:A:1114:C:H2'	1:A:1115:C:H5'	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1267:C:O2'	21:V:20:LYS:HG3	2.10	0.52
2:B:78:GLN:HG3	2:B:94:ASN:OD1	2.10	0.52
4:D:176:LEU:O	4:D:177:ASP:HB3	2.10	0.52
4:D:201:GLN:OE1	4:D:204:ILE:HD12	2.10	0.52
4:D:203:VAL:O	4:D:204:ILE:C	2.47	0.52
5:E:137:GLU:OE1	5:E:141:GLN:NE2	2.37	0.52
6:F:69:GLU:O	6:F:72:VAL:HG23	2.10	0.52
7:G:39:ALA:HA	7:G:42:ILE:CG1	2.40	0.52
7:G:54:THR:CB	7:G:56:GLN:HE22	2.23	0.52
8:H:85:ARG:HG3	8:H:85:ARG:NH1	2.25	0.52
9:I:97:LYS:HZ1	9:I:102:LEU:HG	1.75	0.52
14:N:9:LYS:C	14:N:11:LYS:N	2.63	0.52
16:P:5:ARG:HH21	16:P:28:ARG:HA	1.74	0.52
18:R:18:ARG:C	18:R:19:LYS:HG2	2.29	0.52
19:S:12:ASP:H	19:S:38:SER:CB	2.22	0.52
19:S:49:ILE:CG1	19:S:51:VAL:HG13	2.40	0.52
1:A:14:U:O2	1:A:17:U:H5	1.93	0.52
1:A:552:U:H4'	12:L:86:ARG:O	2.09	0.52
1:A:553:A:O2'	1:A:554:C:H5'	2.09	0.52
1:A:567:G:H2'	1:A:568:G:O4'	2.10	0.52
1:A:575:G:O2'	1:A:576:G:OP2	2.26	0.52
1:A:628:G:C2'	1:A:629:G:H5'	2.39	0.52
1:A:895:G:H2'	1:A:896:C:C6	2.44	0.52
1:A:916:G:C2	1:A:917:G:N7	2.78	0.52
1:A:941:G:H2'	1:A:942:G:H8	1.75	0.52
1:A:1514:C:H2'	1:A:1515:C:H6	1.75	0.52
3:C:76:VAL:HA	3:C:83:ARG:CD	2.39	0.52
3:C:83:ARG:C	3:C:85:ARG:N	2.62	0.52
4:D:92:VAL:O	4:D:96:LEU:HD13	2.10	0.52
7:G:116:ALA:CA	7:G:119:ARG:NH2	2.73	0.52
8:H:54:ASP:C	8:H:56:LYS:N	2.63	0.52
14:N:46:GLU:OE1	14:N:47:LEU:N	2.41	0.52
16:P:67:THR:HB	16:P:70:ALA:HB2	1.91	0.52
20:T:51:GLU:O	20:T:55:ILE:N	2.43	0.52
1:A:742:G:O2'	1:A:743:U:H5'	2.10	0.52
1:A:961:U:OP1	1:A:1223:C:H4'	2.10	0.52
1:A:1130:A:O3'	9:I:3:GLN:HG3	2.10	0.52
1:A:1291:G:O2'	9:I:38:GLN:HG3	2.10	0.52
1:A:1328:C:H5''	13:M:28:ALA:HB1	1.92	0.52
1:A:1329:A:H2'	1:A:1330:U:O4'	2.10	0.52
1:A:1391:U:H2'	1:A:1392:G:C8	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1467:G:H8	1:A:1467:G:O5'	1.93	0.52
2:B:14:GLY:C	2:B:15:VAL:HG22	2.30	0.52
2:B:21:ARG:NH1	2:B:23:ARG:NH2	2.57	0.52
2:B:97:TRP:HZ3	2:B:176:GLU:OE2	1.92	0.52
3:C:6:HIS:NE2	3:C:8:ILE:HB	2.24	0.52
3:C:24:ALA:HB1	3:C:28:GLN:HE22	1.75	0.52
3:C:127:ARG:HD2	3:C:127:ARG:N	2.24	0.52
5:E:50:GLU:OE1	5:E:51:VAL:HG23	2.10	0.52
7:G:41:ARG:O	7:G:42:ILE:C	2.48	0.52
10:J:31:GLY:O	10:J:32:ALA:HB2	2.10	0.52
12:L:6:THR:O	12:L:9:GLN:N	2.42	0.52
12:L:50:SER:O	12:L:51:ALA:HB2	2.09	0.52
19:S:23:ASN:OD1	19:S:27:GLU:OE1	2.29	0.52
20:T:66:ALA:C	20:T:68:LYS:H	2.12	0.52
1:A:180:U:H3	1:A:196:A:N6	2.08	0.51
1:A:539:A:H2'	1:A:540:G:H8	1.74	0.51
1:A:942:G:O2'	1:A:943:U:H5'	2.10	0.51
1:A:1180:A:OP1	9:I:103:THR:HG23	2.10	0.51
1:A:1414:U:O2'	1:A:1415:G:H5'	2.09	0.51
2:B:116:GLU:HG2	2:B:153:ARG:NH2	2.25	0.51
2:B:118:LEU:O	2:B:120:ALA:N	2.43	0.51
2:B:134:GLU:HB3	2:B:138:LEU:CD2	2.36	0.51
2:B:178:ARG:O	2:B:180:LEU:N	2.43	0.51
2:B:187:LEU:CD1	2:B:205:ASP:HA	2.40	0.51
3:C:135:LYS:NZ	5:E:52:PRO:HG2	2.26	0.51
5:E:48:ALA:O	5:E:50:GLU:N	2.42	0.51
8:H:103:VAL:HG21	8:H:109:ILE:O	2.10	0.51
9:I:85:LEU:O	9:I:92:TYR:HB2	2.10	0.51
10:J:22:LYS:HZ2	10:J:23:ILE:CD1	2.23	0.51
11:K:66:LEU:HD21	11:K:101:SER:HA	1.92	0.51
13:M:35:GLU:C	13:M:37:THR:H	2.13	0.51
19:S:28:LYS:CG	19:S:29:ARG:H	2.19	0.51
1:A:204:U:O2'	1:A:216:G:O5'	2.24	0.51
1:A:489:C:H2'	1:A:490:G:C8	2.41	0.51
1:A:521:G:C6	1:A:529:G:N2	2.79	0.51
1:A:949:A:C5	1:A:950:U:N3	2.78	0.51
1:A:1014:A:H4'	19:S:14:HIS:CG	2.45	0.51
1:A:1112:C:O2	3:C:179:ARG:CB	2.58	0.51
1:A:1124:G:H5''	10:J:35:SER:O	2.10	0.51
1:A:1243:C:OP2	21:V:10:ARG:NE	2.43	0.51
2:B:55:PHE:O	2:B:58:ILE:N	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:131:PRO:C	2:B:133:LYS:N	2.64	0.51
7:G:95:ARG:O	7:G:96:GLN:C	2.48	0.51
8:H:113:SER:HB2	8:H:134:ILE:CD1	2.40	0.51
9:I:7:THR:CG2	9:I:8:GLY:N	2.65	0.51
12:L:78:GLN:C	12:L:80:HIS:H	2.13	0.51
13:M:96:LEU:O	13:M:97:PRO:O	2.29	0.51
15:O:46:HIS:ND1	15:O:46:HIS:N	2.57	0.51
16:P:50:LYS:O	16:P:51:VAL:CG2	2.58	0.51
19:S:11:VAL:CG1	19:S:16:LEU:HD13	2.39	0.51
19:S:34:TRP:O	19:S:51:VAL:HA	2.10	0.51
20:T:43:LEU:O	20:T:46:GLU:HB2	2.10	0.51
21:V:14:TRP:C	21:V:16:GLY:N	2.62	0.51
1:A:286:G:H2'	1:A:287:U:C6	2.45	0.51
1:A:308:C:H2'	1:A:309:G:C8	2.46	0.51
1:A:318:G:C2	1:A:319:G:C5	2.99	0.51
1:A:380:G:H22	1:A:382:A:H3'	1.67	0.51
1:A:690:G:H2'	1:A:691:G:O4'	2.10	0.51
1:A:718:G:O4'	11:K:117:ASN:ND2	2.43	0.51
1:A:1215:G:H2'	1:A:1216:G:C4'	2.39	0.51
1:A:1276:G:H21	1:A:1282:C:C2'	2.23	0.51
2:B:92:TYR:CD1	2:B:92:TYR:C	2.82	0.51
2:B:235:SER:C	2:B:237:ALA:N	2.63	0.51
3:C:11:ARG:HG2	3:C:11:ARG:HH11	1.76	0.51
3:C:75:VAL:O	3:C:83:ARG:NH1	2.43	0.51
3:C:91:LEU:HD23	3:C:91:LEU:C	2.31	0.51
4:D:96:LEU:HD12	4:D:96:LEU:H	1.75	0.51
7:G:53:LYS:O	7:G:54:THR:HG23	2.10	0.51
8:H:6:ILE:HB	8:H:85:ARG:NH1	2.26	0.51
8:H:91:ARG:NH1	17:Q:33:GLY:HA3	2.26	0.51
8:H:125:ARG:H	8:H:125:ARG:HD2	1.76	0.51
12:L:62:SER:O	12:L:64:TYR:N	2.43	0.51
15:O:63:ARG:C	15:O:65:ARG:N	2.62	0.51
17:Q:80:GLY:O	17:Q:81:ARG:CB	2.55	0.51
19:S:34:TRP:CD2	19:S:57:HIS:HE1	2.28	0.51
1:A:39:G:O6	1:A:547:A:H5''	2.10	0.51
1:A:400:C:O5'	1:A:400:C:H6	1.94	0.51
1:A:476:G:C2'	1:A:477:G:H8	2.21	0.51
1:A:603:U:O2'	1:A:604:G:H5'	2.10	0.51
1:A:807:A:H2'	1:A:808:C:C6	2.46	0.51
1:A:1074:G:O2'	2:B:103:THR:HG21	2.10	0.51
1:A:1152:A:H5''	10:J:13:HIS:CB	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1255:G:N2	1:A:1259:C:O2	2.35	0.51
1:A:1518:A:OP2	1:A:1518:A:C4	2.63	0.51
2:B:98:LEU:HD23	2:B:98:LEU:N	2.24	0.51
3:C:21:ARG:NH2	3:C:56:ASP:HB3	2.26	0.51
3:C:31:HIS:HA	3:C:34:LEU:HB2	1.92	0.51
3:C:79:ARG:NH2	3:C:82:GLU:HG3	2.25	0.51
3:C:102:ASN:N	3:C:102:ASN:ND2	2.58	0.51
4:D:20:TYR:C	4:D:22:LYS:N	2.64	0.51
5:E:144:THR:HB	5:E:147:ASP:OD2	2.10	0.51
7:G:5:ARG:O	7:G:7:ALA:N	2.42	0.51
8:H:4:ASP:OD2	8:H:7:ALA:HB2	2.10	0.51
11:K:122:LYS:O	11:K:125:PHE:N	2.40	0.51
13:M:20:THR:O	13:M:20:THR:CG2	2.59	0.51
14:N:48:ALA:O	14:N:50:LYS:N	2.41	0.51
17:Q:98:LEU:HD23	17:Q:98:LEU:H	1.76	0.51
1:A:138:G:O2'	1:A:139:G:H5'	2.11	0.51
1:A:146:G:C2	1:A:147:G:C8	2.98	0.51
1:A:290:C:C4	1:A:291:C:C5	2.98	0.51
1:A:701:C:OP1	1:A:703:G:H5'	2.10	0.51
1:A:949:A:N1	1:A:1233:G:N3	2.58	0.51
1:A:1107:C:OP1	3:C:174:PRO:HD3	2.11	0.51
1:A:1347:G:C6	9:I:107:ARG:NH2	2.78	0.51
1:A:1440:C:C2'	1:A:1441:G:H5'	2.39	0.51
3:C:116:VAL:HA	3:C:119:ARG:CB	2.40	0.51
4:D:157:LEU:CB	4:D:158:ILE:HD12	2.33	0.51
6:F:62:TRP:C	6:F:63:TYR:CD1	2.83	0.51
10:J:19:SER:OG	10:J:91:PRO:HB2	2.10	0.51
11:K:56:GLY:O	11:K:57:THR:C	2.49	0.51
12:L:37:CYS:HA	12:L:58:VAL:HA	1.93	0.51
13:M:17:VAL:HG12	13:M:21:TYR:HE1	1.76	0.51
13:M:42:ALA:O	13:M:43:THR:C	2.48	0.51
16:P:18:ARG:HD3	16:P:35:LYS:CD	2.39	0.51
16:P:78:GLY:C	16:P:80:PHE:N	2.62	0.51
19:S:18:LYS:O	19:S:22:LEU:HD11	2.09	0.51
1:A:74:C:H2'	1:A:75:G:O4'	2.10	0.51
1:A:251:G:H4'	1:A:252:U:O5'	2.10	0.51
1:A:255:G:O2'	17:Q:16:GLN:NE2	2.43	0.51
1:A:769:G:H4'	1:A:1513:A:H4'	1.93	0.51
1:A:1228:C:OP2	13:M:111:LYS:HE2	2.11	0.51
1:A:1298:C:C4	7:G:114:ARG:HD2	2.46	0.51
1:A:1409:C:H2'	1:A:1410:G:H8	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1430:C:O2'	1:A:1431:C:H5'	2.10	0.51
1:A:1529:G:H4'	1:A:1530:G:OP2	2.09	0.51
2:B:102:LEU:HD21	2:B:162:ILE:CD1	2.40	0.51
3:C:132:ARG:HA	3:C:135:LYS:CD	2.40	0.51
3:C:139:GLN:HE21	3:C:139:GLN:C	2.14	0.51
3:C:177:THR:O	3:C:179:ARG:N	2.44	0.51
4:D:82:ALA:O	4:D:83:SER:C	2.49	0.51
4:D:175:SER:OG	4:D:186:LEU:HD21	2.11	0.51
5:E:79:GLU:O	5:E:79:GLU:OE1	2.28	0.51
7:G:41:ARG:O	7:G:44:TYR:N	2.43	0.51
7:G:91:VAL:CG1	7:G:92:SER:H	2.24	0.51
8:H:28:ALA:H	8:H:58:TYR:HA	1.74	0.51
13:M:116:THR:HG22	13:M:117:VAL:N	2.25	0.51
16:P:22:THR:OG1	16:P:26:ARG:HB3	2.10	0.51
17:Q:94:ASN:HD22	17:Q:94:ASN:N	2.08	0.51
1:A:53:A:C2	1:A:54:C:C2	2.98	0.51
1:A:393:A:O2'	1:A:394:G:H5'	2.10	0.51
1:A:642:A:N3	8:H:113:SER:O	2.43	0.51
1:A:735:C:H1'	18:R:75:ILE:HD11	1.93	0.51
1:A:835:U:H3	1:A:851:G:H1	1.59	0.51
1:A:875:C:O2'	8:H:14:ARG:HD2	2.11	0.51
1:A:1245:A:C2	1:A:1293:G:N3	2.79	0.51
1:A:1434:A:O2'	1:A:1435:G:H5'	2.10	0.51
2:B:21:ARG:HH22	2:B:23:ARG:NH2	2.08	0.51
2:B:164:VAL:HG12	2:B:165:VAL:N	2.25	0.51
2:B:208:ILE:CG2	2:B:209:ARG:N	2.72	0.51
3:C:84:ILE:O	3:C:88:ARG:HD2	2.10	0.51
3:C:85:ARG:C	3:C:87:LEU:H	2.12	0.51
4:D:8:VAL:O	4:D:9:CYS:C	2.49	0.51
6:F:9:VAL:CB	6:F:87:ARG:HB2	2.40	0.51
6:F:71:ARG:O	6:F:72:VAL:C	2.49	0.51
6:F:74:ASP:HB3	6:F:77:ARG:HH21	1.76	0.51
10:J:86:MET:H	10:J:88:LEU:HG	1.76	0.51
10:J:89:ASP:O	10:J:90:LEU:HB2	2.11	0.51
12:L:62:SER:C	12:L:64:TYR:H	2.14	0.51
12:L:66:VAL:HG12	12:L:67:THR:N	2.15	0.51
13:M:15:VAL:O	13:M:19:LEU:HD12	2.11	0.51
20:T:79:ARG:O	20:T:80:ARG:C	2.48	0.51
21:V:6:ARG:CD	21:V:15:ARG:HH22	2.24	0.51
1:A:409:G:H2'	1:A:410:G:O4'	2.11	0.51
1:A:597:G:C5	1:A:598:U:C6	2.98	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:891:U:O2'	1:A:892:A:H5'	2.09	0.51
1:A:925:G:N1	1:A:927:G:C5	2.79	0.51
1:A:991:U:O2'	1:A:993:G:C8	2.63	0.51
1:A:1144:G:H21	1:A:1146:A:H61	1.56	0.51
1:A:1153:C:OP1	10:J:13:HIS:CE1	2.64	0.51
1:A:1318:A:H4'	19:S:10:PHE:CD2	2.45	0.51
1:A:1474:G:H2'	1:A:1475:G:H8	1.75	0.51
3:C:39:ILE:CD1	3:C:57:ILE:HD11	2.41	0.51
3:C:191:THR:CG2	3:C:194:GLY:O	2.58	0.51
4:D:113:SER:O	4:D:117:ALA:N	2.42	0.51
4:D:149:ALA:O	4:D:150:GLU:C	2.49	0.51
6:F:19:LEU:HD23	6:F:23:LYS:HG3	1.91	0.51
6:F:71:ARG:C	6:F:73:ASN:N	2.62	0.51
8:H:100:ILE:HG21	8:H:125:ARG:HE	1.76	0.51
20:T:99:LEU:C	20:T:101:GLY:N	2.61	0.51
1:A:185:A:H1'	20:T:81:LYS:HZ3	1.76	0.51
1:A:264:U:C2'	1:A:265:G:H5'	2.41	0.51
1:A:407:G:H2'	1:A:408:A:C8	2.46	0.51
1:A:763:G:H2'	1:A:764:C:C6	2.46	0.51
1:A:1021:G:H2'	1:A:1022:G:C8	2.46	0.51
1:A:1361(A):C:H2'	1:A:1362:C:H5''	1.91	0.51
2:B:91:PRO:HG3	2:B:154:LEU:HB2	1.93	0.51
3:C:44:GLU:OE1	3:C:68:VAL:HG22	2.11	0.51
3:C:50:ALA:HA	3:C:72:LYS:HB2	1.92	0.51
4:D:43:HIS:HB3	4:D:46:LYS:HE3	1.92	0.51
4:D:96:LEU:N	4:D:96:LEU:CD1	2.73	0.51
5:E:25:ARG:C	5:E:26:PHE:CD1	2.84	0.51
8:H:97:VAL:C	8:H:99:GLU:H	2.14	0.51
9:I:122:ALA:HB1	9:I:123:PRO:HD2	1.93	0.51
11:K:67:ASP:C	11:K:69:ALA:N	2.64	0.51
13:M:5:ALA:HB2	13:M:22:ILE:HD13	1.93	0.51
14:N:24:CYS:O	14:N:28:GLY:CA	2.58	0.51
14:N:36:PHE:HB3	14:N:37:PHE:CE1	2.46	0.51
15:O:16:ALA:O	15:O:18:PHE:N	2.44	0.51
16:P:43:LYS:CB	16:P:48:TRP:CD1	2.92	0.51
18:R:69:THR:O	18:R:70:ILE:C	2.49	0.51
1:A:186:C:H4'	20:T:82:SER:HB2	1.93	0.51
1:A:321:A:H2'	1:A:322:C:C6	2.46	0.51
1:A:877:C:C2'	1:A:878:G:H5'	2.40	0.51
1:A:1418:A:O2'	1:A:1419:G:H5'	2.11	0.51
1:A:1474:G:H2'	1:A:1475:G:C8	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:76:GLN:O	2:B:208:ILE:HD11	2.10	0.51
2:B:173:ALA:O	2:B:174:VAL:C	2.49	0.51
3:C:201:TYR:O	3:C:202:ILE:HG13	2.10	0.51
4:D:200:GLU:O	4:D:203:VAL:HB	2.11	0.51
5:E:9:LYS:O	5:E:33:VAL:N	2.43	0.51
7:G:36:LYS:O	7:G:40:ALA:N	2.43	0.51
8:H:4:ASP:OD2	8:H:7:ALA:CB	2.59	0.51
8:H:53:VAL:O	8:H:54:ASP:HB3	2.10	0.51
11:K:33:THR:OG1	11:K:34:ASP:N	2.42	0.51
13:M:56:LEU:HD12	13:M:57:ARG:H	1.74	0.51
15:O:28:GLN:O	15:O:32:LEU:HB2	2.11	0.51
15:O:29:VAL:O	15:O:32:LEU:HB3	2.11	0.51
16:P:4:ILE:CG1	16:P:64:ALA:HB1	2.35	0.51
16:P:12:LYS:O	16:P:13:HIS:CB	2.59	0.51
19:S:19:VAL:HG12	19:S:20:LEU:N	2.26	0.51
1:A:24:U:H2'	1:A:25:C:H6	1.74	0.50
1:A:277:C:P	17:Q:68:ARG:HH12	2.33	0.50
1:A:363:A:O2'	1:A:364:A:H5'	2.10	0.50
1:A:399:G:N7	1:A:400:C:N4	2.59	0.50
1:A:436:C:O2'	1:A:437:U:H5'	2.11	0.50
1:A:665:A:H3'	1:A:725:G:N2	2.27	0.50
1:A:841:U:H5'	1:A:848:C:C5	2.46	0.50
1:A:1080:A:H5'	5:E:14:ARG:NH2	2.25	0.50
1:A:1144:G:H21	1:A:1146:A:N6	2.09	0.50
1:A:1176:A:H2'	1:A:1177:G:O4'	2.11	0.50
1:A:1224:G:N2	1:A:1362:C:N3	2.59	0.50
1:A:1492:A:O2'	1:A:1493:A:C8	2.65	0.50
2:B:10:LEU:HG	2:B:48:MET:CE	2.40	0.50
2:B:142:LEU:O	2:B:143:GLU:C	2.50	0.50
3:C:11:ARG:HA	3:C:14:ILE:CD1	2.41	0.50
5:E:150:ARG:O	5:E:151:LEU:C	2.48	0.50
7:G:74:GLU:N	7:G:91:VAL:HG23	2.26	0.50
7:G:110:GLN:OE1	7:G:110:GLN:HA	2.10	0.50
10:J:60:ARG:H	10:J:60:ARG:HD3	1.75	0.50
11:K:58:PRO:O	11:K:59:TYR:C	2.49	0.50
14:N:8:GLU:CA	14:N:11:LYS:HD3	2.41	0.50
16:P:80:PHE:H	16:P:80:PHE:HD1	1.55	0.50
17:Q:11:VAL:HG11	17:Q:22:LEU:HB2	1.93	0.50
17:Q:65:ILE:HD12	17:Q:65:ILE:N	2.25	0.50
19:S:36:ARG:HH21	19:S:75:ALA:HB1	1.75	0.50
20:T:59:ALA:O	20:T:60:GLU:C	2.49	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:C:H2'	1:A:352:C:H41	1.76	0.50
1:A:59:A:H2'	1:A:331:G:N2	2.26	0.50
1:A:75:G:O2'	1:A:76:C:H5'	2.11	0.50
1:A:101:A:C2	1:A:102:G:C8	2.99	0.50
1:A:290:C:H42	1:A:310:G:H1	1.58	0.50
1:A:424:G:H2'	1:A:425:G:H8	1.76	0.50
1:A:620:C:N1	4:D:135:LEU:HD13	2.26	0.50
1:A:687:A:O2'	1:A:688:G:P	2.69	0.50
1:A:722:A:H3'	1:A:722:A:N3	2.25	0.50
1:A:1036:G:H2'	1:A:1037:C:C4'	2.41	0.50
1:A:1118:C:H1'	1:A:1179:A:N9	2.27	0.50
1:A:1371:G:C4	1:A:1372:U:C5	3.00	0.50
2:B:25:ASN:O	2:B:27:LYS:N	2.45	0.50
3:C:126:ARG:CB	3:C:128:PHE:HB3	2.41	0.50
3:C:179:ARG:C	3:C:181:ASN:H	2.14	0.50
4:D:150:GLU:O	4:D:152:SER:N	2.45	0.50
4:D:163:GLU:O	4:D:165:MET:N	2.44	0.50
7:G:65:ALA:HB2	7:G:128:ALA:CB	2.41	0.50
7:G:108:ALA:O	7:G:119:ARG:HB3	2.12	0.50
8:H:4:ASP:CG	8:H:85:ARG:HH12	2.15	0.50
8:H:104:ARG:HG2	8:H:104:ARG:NH1	2.26	0.50
10:J:55:LYS:HB2	10:J:56:HIS:CD2	2.47	0.50
12:L:56:ALA:O	12:L:57:LYS:C	2.49	0.50
14:N:9:LYS:HD3	14:N:9:LYS:O	2.09	0.50
15:O:15:PHE:O	15:O:16:ALA:C	2.49	0.50
18:R:37:VAL:HB	18:R:41:LYS:HD3	1.93	0.50
19:S:35:SER:O	19:S:71:LEU:HD13	2.11	0.50
19:S:65:ASN:OD1	19:S:66:MET:HG3	2.12	0.50
1:A:21:G:C2	1:A:22:G:C6	2.98	0.50
1:A:57:G:C2	1:A:58:C:O2	2.65	0.50
1:A:102:G:H22	1:A:171:A:H2	1.57	0.50
1:A:235:C:H5'	17:Q:70:ARG:CD	2.33	0.50
1:A:290:C:C2'	1:A:291:C:H5'	2.41	0.50
1:A:342:C:O5'	1:A:342:C:H6	1.95	0.50
1:A:564:C:H5'	17:Q:32:TYR:CE2	2.47	0.50
1:A:834:C:O2'	1:A:835:U:H5'	2.12	0.50
1:A:922:G:H1	1:A:1395:C:H42	1.58	0.50
1:A:939:G:H5''	7:G:102:ARG:NH2	2.26	0.50
1:A:1064:G:OP1	1:A:1386:G:H4'	2.11	0.50
1:A:1084:G:OP1	1:A:1086:U:C4	2.65	0.50
1:A:1291:G:H5''	7:G:41:ARG:NH2	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1298:C:O4'	1:A:1299:A:C6	2.64	0.50
1:A:1313:U:OP2	19:S:6:LYS:HA	2.11	0.50
2:B:213:LEU:HD23	2:B:213:LEU:C	2.32	0.50
3:C:8:ILE:HG22	14:N:50:LYS:HA	1.93	0.50
5:E:95:ALA:HB1	5:E:96:PRO:HD2	1.92	0.50
6:F:12:PRO:CG	6:F:57:GLN:HG3	2.37	0.50
7:G:73:MET:HA	7:G:91:VAL:HG23	1.92	0.50
8:H:10:LEU:HD22	8:H:83:ILE:HD11	1.93	0.50
8:H:21:LYS:O	8:H:65:TYR:OH	2.22	0.50
9:I:49:PRO:HG2	9:I:50:LEU:CD1	2.38	0.50
11:K:48:ILE:HG13	11:K:64:ALA:HA	1.93	0.50
13:M:67:GLU:O	13:M:68:GLY:C	2.50	0.50
14:N:14:PRO:CG	14:N:15:LYS:N	2.73	0.50
14:N:37:PHE:CD2	14:N:53:LEU:HD21	2.46	0.50
16:P:74:LEU:HB3	16:P:79:VAL:HG21	1.92	0.50
18:R:38:GLU:H	18:R:41:LYS:CG	2.25	0.50
18:R:79:LEU:HD12	18:R:79:LEU:N	2.26	0.50
20:T:54:LYS:HB2	20:T:100:ILE:HD11	1.93	0.50
20:T:100:ILE:O	20:T:102:GLY:N	2.44	0.50
1:A:137:C:N3	1:A:226:G:N2	2.44	0.50
1:A:140:A:H2'	1:A:141:A:O4'	2.11	0.50
1:A:570:G:C6	1:A:571:U:O4	2.65	0.50
1:A:772:U:O2'	1:A:773:G:H5'	2.12	0.50
1:A:960:U:H4'	1:A:961:U:C5'	2.42	0.50
1:A:1346:A:O2'	1:A:1347:G:P	2.69	0.50
2:B:74:LYS:HE2	2:B:166:ASP:HB2	1.90	0.50
3:C:50:ALA:HA	3:C:72:LYS:CG	2.41	0.50
3:C:116:VAL:C	3:C:119:ARG:HB3	2.31	0.50
3:C:143:GLU:C	3:C:145:GLY:H	2.14	0.50
3:C:191:THR:OG1	3:C:192:THR:N	2.45	0.50
4:D:23:GLY:HA2	4:D:113:SER:HB2	1.92	0.50
5:E:12:LEU:HD13	5:E:12:LEU:C	2.32	0.50
7:G:13:GLN:C	7:G:21:VAL:HG12	2.32	0.50
8:H:97:VAL:HA	8:H:100:ILE:CD1	2.37	0.50
12:L:104:VAL:O	12:L:105:TYR:HB2	2.11	0.50
13:M:22:ILE:O	13:M:24:GLY:N	2.41	0.50
14:N:59:ALA:HB1	14:N:61:TRP:HZ3	1.75	0.50
16:P:21:VAL:CG1	16:P:33:ILE:HB	2.42	0.50
16:P:75:ARG:CA	16:P:80:PHE:CE1	2.94	0.50
17:Q:18:THR:HG23	17:Q:69:LYS:NZ	2.26	0.50
18:R:88:LYS:OXT	18:R:88:LYS:HG3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:S:49:ILE:HG22	19:S:60:VAL:CB	2.39	0.50
1:A:189:G:H2'	1:A:190:C:C6	2.46	0.50
1:A:272:C:O2'	1:A:273:A:H5'	2.11	0.50
1:A:386:C:H2'	1:A:387:U:H5'	1.92	0.50
1:A:407:G:C6	1:A:408:A:N6	2.80	0.50
1:A:438:G:H4'	1:A:439:A:OP1	2.10	0.50
1:A:691:G:OP2	11:K:26:ASN:ND2	2.42	0.50
1:A:777:A:C6	1:A:778:G:C5	3.00	0.50
1:A:961:U:H1'	1:A:984:C:O4'	2.12	0.50
1:A:1361(A):C:O2'	1:A:1362:C:H5''	2.12	0.50
1:A:1403:C:H2'	1:A:1404:C:H6	1.72	0.50
2:B:42:ILE:CD1	2:B:203:GLY:HA2	2.41	0.50
2:B:81:VAL:O	2:B:83:MET:N	2.44	0.50
2:B:92:TYR:HE1	2:B:151:GLY:N	2.10	0.50
2:B:112:VAL:CG1	2:B:153:ARG:HA	2.41	0.50
3:C:141:VAL:O	3:C:146:ALA:HB3	2.11	0.50
3:C:188:LEU:O	3:C:189:ALA:HB2	2.11	0.50
5:E:9:LYS:O	5:E:32:VAL:HA	2.11	0.50
5:E:20:GLN:O	5:E:21:ALA:C	2.50	0.50
5:E:80:ILE:C	5:E:81:GLU:HG3	2.31	0.50
5:E:130:ASN:O	5:E:133:TYR:N	2.45	0.50
5:E:143:ARG:HH11	5:E:143:ARG:HG3	1.75	0.50
8:H:86:ILE:CG2	8:H:87:SER:N	2.65	0.50
10:J:44:VAL:CG1	10:J:45:ARG:H	2.23	0.50
10:J:65:LEU:HD23	10:J:65:LEU:C	2.31	0.50
13:M:14:ARG:HH12	13:M:16:ASP:HB2	1.77	0.50
13:M:17:VAL:C	13:M:19:LEU:N	2.64	0.50
13:M:23:TYR:CE2	13:M:70:LEU:HD22	2.46	0.50
13:M:88:ARG:HD2	19:S:3:ARG:NH2	2.26	0.50
17:Q:63:ARG:HH11	17:Q:63:ARG:HG3	1.76	0.50
18:R:22:VAL:HG13	18:R:26:LEU:HD11	1.93	0.50
20:T:79:ARG:HG2	20:T:83:ARG:HH11	1.76	0.50
1:A:39:G:C2	1:A:40:C:C6	3.00	0.50
1:A:53:A:C6	1:A:54:C:C4	2.99	0.50
1:A:112:G:O2'	1:A:113:G:H5'	2.11	0.50
1:A:298:A:C6	1:A:299:G:C2	3.00	0.50
1:A:392:G:C2	1:A:393:A:C5	3.00	0.50
1:A:602:A:H2'	1:A:603:U:O4'	2.12	0.50
1:A:619:U:H3	4:D:135:LEU:HD21	1.76	0.50
1:A:665:A:N3	1:A:732:C:H2'	2.26	0.50
1:A:886:G:H1	1:A:911:U:H3	1.60	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:959:A:C2	1:A:1222:G:O4'	2.65	0.50
1:A:1162:C:H2'	1:A:1163:C:C6	2.47	0.50
1:A:1509:C:H2'	1:A:1510:U:O4'	2.10	0.50
1:A:1533:C:O2	1:A:1533:C:H3'	2.12	0.50
2:B:88:ALA:HB2	2:B:219:VAL:CG1	2.42	0.50
2:B:223:ILE:HG21	2:B:230:VAL:HG23	1.93	0.50
3:C:137:ALA:CA	3:C:140:ARG:HH11	2.25	0.50
3:C:183:ASP:OD2	3:C:184:TYR:O	2.30	0.50
4:D:134:ASP:O	4:D:135:LEU:C	2.49	0.50
4:D:192:GLU:O	4:D:193:ASP:C	2.50	0.50
7:G:63:LYS:O	7:G:67:GLU:HG2	2.12	0.50
7:G:152:ALA:CB	7:G:155:ARG:CZ	2.89	0.50
8:H:38:ILE:O	8:H:40:ALA:N	2.44	0.50
8:H:40:ALA:O	8:H:41:ARG:C	2.50	0.50
9:I:97:LYS:NZ	9:I:102:LEU:HD11	2.27	0.50
10:J:22:LYS:HZ1	10:J:91:PRO:HD3	1.76	0.50
12:L:24:VAL:O	12:L:26:ALA:N	2.41	0.50
1:A:142:G:H1'	1:A:196:A:C2	2.47	0.50
1:A:366:C:O2'	1:A:367:U:H5''	2.11	0.50
1:A:626:U:O2'	1:A:627:G:H5'	2.11	0.50
1:A:635:G:H2'	1:A:636:U:C6	2.47	0.50
1:A:687:A:H4'	1:A:688:G:O5'	2.12	0.50
1:A:762:C:H2'	1:A:763:G:C8	2.40	0.50
1:A:1317:C:O2'	1:A:1318:A:H5'	2.11	0.50
1:A:1326:C:OP1	21:V:17:THR:HB	2.12	0.50
1:A:1413:A:C2	1:A:1414:U:N1	2.80	0.50
2:B:35:GLU:HA	2:B:40:HIS:HA	1.93	0.50
2:B:220:ASP:O	2:B:224:GLN:HB2	2.11	0.50
4:D:32:ALA:C	4:D:34:GLU:H	2.14	0.50
4:D:110:PHE:N	4:D:110:PHE:CD1	2.80	0.50
8:H:31:PHE:CZ	8:H:134:ILE:HD13	2.47	0.50
10:J:33:GLN:C	10:J:75:ILE:HG22	2.31	0.50
14:N:53:LEU:HD12	14:N:56:VAL:CB	2.34	0.50
15:O:48:LYS:C	15:O:50:HIS:N	2.63	0.50
20:T:39:LYS:HG2	20:T:55:ILE:HD12	1.94	0.50
20:T:60:GLU:O	20:T:63:ILE:HB	2.11	0.50
1:A:16:A:C2	1:A:920:U:O2	2.65	0.50
1:A:157:G:H2'	1:A:158:G:H8	1.75	0.50
1:A:597:G:C2'	1:A:598:U:H5'	2.40	0.50
1:A:624:C:H6	1:A:624:C:O5'	1.94	0.50
1:A:679:C:H2'	1:A:680:C:C6	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:711:G:H2'	1:A:712:A:H8	1.73	0.50
1:A:923:A:H2	1:A:1395:C:N3	2.10	0.50
1:A:927:G:N2	1:A:1391:U:C1'	2.75	0.50
1:A:1245:A:C2	1:A:1293:G:C2	3.00	0.50
1:A:1392:G:C2'	1:A:1393:U:H5'	2.42	0.50
1:A:1418:A:C2'	1:A:1419:G:H5'	2.42	0.50
2:B:71:VAL:HG23	2:B:164:VAL:HG13	1.93	0.50
3:C:21:ARG:HD2	3:C:21:ARG:N	2.26	0.50
3:C:71:ALA:O	3:C:73:PRO:CD	2.56	0.50
4:D:12:CYS:HG	4:D:19:LEU:HB2	1.75	0.50
9:I:6:GLY:CA	9:I:83:ARG:HD2	2.40	0.50
13:M:11:ARG:HD3	13:M:12:ASN:N	2.27	0.50
13:M:13:LYS:O	13:M:14:ARG:C	2.49	0.50
16:P:75:ARG:HA	16:P:80:PHE:HE1	1.75	0.50
19:S:49:ILE:HG12	19:S:71:LEU:CD2	2.42	0.50
20:T:75:ASN:O	20:T:78:ALA:N	2.45	0.50
20:T:85:MET:O	20:T:86:ARG:C	2.48	0.50
21:V:9:ARG:O	21:V:13:ILE:HB	2.12	0.50
1:A:16:A:C2'	1:A:17:U:H5'	2.42	0.50
1:A:190(B):C:H2'	1:A:190(C):C:O4'	2.11	0.50
1:A:376:G:HO2'	1:A:377:G:H5'	1.75	0.50
1:A:452:A:C4'	16:P:72:ARG:NH2	2.74	0.50
1:A:1240:U:O2	7:G:38:LEU:HD21	2.11	0.50
1:A:1399:C:C2	1:A:1502:A:N6	2.80	0.50
1:A:1448:C:N3	1:A:1449:C:C5	2.80	0.50
1:A:1533:C:O2	1:A:1533:C:C2'	2.60	0.50
2:B:30:ARG:O	2:B:31:TYR:CD1	2.65	0.50
2:B:91:PRO:HG3	2:B:154:LEU:CB	2.42	0.50
3:C:47:LEU:HD21	3:C:76:VAL:CG1	2.41	0.50
3:C:83:ARG:O	3:C:85:ARG:N	2.43	0.50
3:C:119:ARG:O	3:C:123:GLN:HG3	2.11	0.50
4:D:24:GLU:O	4:D:25:ARG:HB3	2.12	0.50
5:E:132:ALA:O	5:E:135:THR:N	2.39	0.50
6:F:22:GLU:C	6:F:24:GLU:N	2.65	0.50
8:H:25:ASP:OD1	8:H:25:ASP:N	2.36	0.50
8:H:97:VAL:C	8:H:99:GLU:N	2.64	0.50
8:H:108:GLY:HA3	8:H:138:TRP:CB	2.42	0.50
9:I:118:LYS:H	9:I:121:ARG:HB3	1.77	0.50
10:J:10:GLY:CA	10:J:16:LEU:HD12	2.37	0.50
10:J:26:ALA:O	10:J:84:GLN:OE1	2.28	0.50
10:J:48:THR:HG21	10:J:62:HIS:CE1	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:L:60:LEU:HD23	12:L:65:GLU:HA	1.92	0.50
14:N:40:CYS:SG	14:N:41:ARG:N	2.85	0.50
16:P:75:ARG:N	16:P:80:PHE:HE1	2.09	0.50
17:Q:90:ILE:HA	17:Q:93:GLN:CB	2.42	0.50
19:S:24:ALA:HB3	19:S:25:LYS:HZ2	1.73	0.50
20:T:99:LEU:O	20:T:100:ILE:C	2.50	0.50
21:V:8:THR:HG22	21:V:9:ARG:H	1.77	0.50
1:A:39:G:C2	1:A:40:C:C5	2.99	0.49
1:A:115:G:H1	1:A:312:C:N4	2.09	0.49
1:A:115:G:C1'	1:A:116:A:N7	2.71	0.49
1:A:924:C:H2'	1:A:925:G:C8	2.47	0.49
1:A:943:U:C2'	1:A:944:G:H5'	2.42	0.49
1:A:1077:G:N2	1:A:1080:A:OP2	2.39	0.49
1:A:1483:A:H2'	1:A:1484:C:O4'	2.12	0.49
1:A:1483:A:C2'	1:A:1484:C:H5'	2.41	0.49
2:B:210:SER:OG	2:B:211:ILE:N	2.43	0.49
3:C:27:LYS:CA	3:C:30:ARG:HH12	2.21	0.49
4:D:159:ARG:O	4:D:163:GLU:HB2	2.11	0.49
7:G:17:VAL:O	7:G:19:GLY:N	2.45	0.49
7:G:80:VAL:HG12	7:G:81:GLY:N	2.27	0.49
8:H:34:GLU:O	8:H:37:ARG:N	2.45	0.49
12:L:59:ARG:HH11	12:L:59:ARG:CG	2.25	0.49
12:L:115:LYS:O	12:L:117:ARG:N	2.34	0.49
13:M:71:ARG:HG2	13:M:71:ARG:HH11	1.77	0.49
14:N:46:GLU:OE2	14:N:47:LEU:HD23	2.11	0.49
17:Q:11:VAL:CG1	17:Q:22:LEU:HB2	2.42	0.49
1:A:44:G:N2	1:A:45:U:H1'	2.27	0.49
1:A:67:C:O2'	1:A:171:A:H1'	2.11	0.49
1:A:89:C:OP2	1:A:89:C:H6	1.95	0.49
1:A:182:U:H2'	1:A:183:G:H5'	1.93	0.49
1:A:788:U:O2'	1:A:789:U:H5'	2.12	0.49
1:A:886:G:C2	1:A:887:G:C4	3.00	0.49
1:A:1128:C:OP1	9:I:66:ARG:NH2	2.45	0.49
1:A:1349:A:OP2	9:I:118:LYS:NZ	2.45	0.49
1:A:1455:G:O5'	1:A:1455:G:H8	1.95	0.49
2:B:46:LYS:O	2:B:47:THR:C	2.50	0.49
2:B:54:THR:O	2:B:58:ILE:HG13	2.12	0.49
3:C:22:TRP:HD1	3:C:59:ARG:HB2	1.76	0.49
3:C:55:VAL:HG13	3:C:68:VAL:CG2	2.42	0.49
3:C:150:LYS:CG	3:C:169:ALA:HB2	2.37	0.49
5:E:30:ALA:HB1	5:E:55:VAL:HG12	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:80:ILE:HD11	5:E:91:LEU:HD12	1.95	0.49
6:F:30:LEU:CD2	6:F:35:ALA:HB3	2.42	0.49
7:G:118:VAL:O	7:G:120:ILE:N	2.45	0.49
10:J:4:ILE:O	10:J:6:ILE:HG12	2.12	0.49
10:J:47:PHE:CZ	14:N:37:PHE:HZ	2.29	0.49
10:J:64:GLU:HG3	14:N:59:ALA:HB2	1.94	0.49
11:K:32:ILE:HG22	11:K:33:THR:N	2.27	0.49
13:M:11:ARG:O	13:M:12:ASN:HB3	2.12	0.49
13:M:70:LEU:O	13:M:70:LEU:HD23	2.13	0.49
14:N:6:LEU:CD2	14:N:23:ARG:HE	2.25	0.49
14:N:21:TYR:O	14:N:23:ARG:N	2.45	0.49
16:P:60:LEU:HD21	16:P:66:PRO:HD3	1.94	0.49
16:P:74:LEU:O	16:P:78:GLY:N	2.45	0.49
17:Q:48:GLU:OE1	17:Q:50:LYS:HD2	2.11	0.49
19:S:34:TRP:HE3	19:S:34:TRP:H	1.59	0.49
1:A:414:A:OP2	1:A:428:G:N2	2.45	0.49
1:A:597:G:C6	1:A:644:G:C6	3.00	0.49
1:A:606:G:H2'	1:A:631:G:N2	2.27	0.49
1:A:971:G:OP1	1:A:971:G:H3'	2.12	0.49
1:A:1497:G:C5	1:A:1498:U:C4	3.00	0.49
3:C:120:VAL:CA	3:C:123:GLN:HB2	2.34	0.49
3:C:131:ARG:O	3:C:135:LYS:HG3	2.12	0.49
3:C:139:GLN:O	3:C:140:ARG:C	2.50	0.49
4:D:156:GLU:OE1	4:D:156:GLU:N	2.45	0.49
5:E:12:LEU:HD11	5:E:31:LEU:HB2	1.93	0.49
7:G:22:LEU:HD12	7:G:23:VAL:N	2.27	0.49
7:G:24:THR:HA	7:G:27:ILE:HD12	1.93	0.49
7:G:155:ARG:HE	7:G:155:ARG:CA	2.16	0.49
10:J:6:ILE:HD11	10:J:74:ILE:N	2.26	0.49
10:J:86:MET:N	10:J:88:LEU:HG	2.27	0.49
16:P:5:ARG:HB2	16:P:67:THR:OG1	2.11	0.49
18:R:26:LEU:HD21	18:R:42:ARG:HD2	1.93	0.49
1:A:171:A:O2'	1:A:172:A:H5'	2.12	0.49
1:A:192:U:O3'	20:T:57:ARG:NH1	2.45	0.49
1:A:356:A:O2'	1:A:357:G:H5'	2.12	0.49
1:A:562:C:H2'	1:A:562:C:OP2	2.12	0.49
1:A:580:U:C1'	15:O:57:LEU:HD23	2.41	0.49
1:A:855:G:H2'	1:A:856:C:H6	1.76	0.49
1:A:1175:G:O2'	1:A:1176:A:H5'	2.12	0.49
1:A:1306:A:C5	1:A:1332:A:C8	3.01	0.49
1:A:1371:G:H2'	1:A:1372:U:H6	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:212:GLN:O	2:B:213:LEU:C	2.50	0.49
3:C:91:LEU:CG	3:C:99:VAL:HG13	2.43	0.49
3:C:203:PHE:CG	3:C:204:LEU:N	2.79	0.49
5:E:92:LYS:O	5:E:118:ILE:CG2	2.61	0.49
6:F:22:GLU:C	6:F:24:GLU:H	2.15	0.49
6:F:28:ARG:O	6:F:31:GLU:OE2	2.30	0.49
8:H:19:VAL:HG23	8:H:19:VAL:O	2.12	0.49
10:J:16:LEU:HA	10:J:19:SER:HB3	1.93	0.49
10:J:22:LYS:NZ	10:J:23:ILE:CD1	2.75	0.49
10:J:84:GLN:O	10:J:85:LEU:CB	2.59	0.49
10:J:95:GLU:CD	10:J:95:GLU:C	2.70	0.49
16:P:19:ILE:O	16:P:36:ILE:HG13	2.12	0.49
18:R:53:ARG:HD3	18:R:63:GLN:CG	2.42	0.49
19:S:19:VAL:HG12	19:S:20:LEU:HD23	1.94	0.49
1:A:375:U:O2'	1:A:376:G:H5'	2.12	0.49
1:A:399:G:C5	1:A:400:C:N4	2.81	0.49
1:A:499:A:H4'	1:A:500:G:OP1	2.12	0.49
1:A:582:U:H2'	1:A:583:A:O4'	2.13	0.49
1:A:684:A:H2'	1:A:685:G:H8	1.77	0.49
1:A:893:C:C4	1:A:894:G:N7	2.80	0.49
1:A:971:G:H4'	1:A:972:C:H5'	1.94	0.49
1:A:1152:A:H5''	10:J:13:HIS:CA	2.42	0.49
1:A:1343:G:C4	1:A:1344:C:C5	3.00	0.49
2:B:24:TRP:HA	2:B:190:THR:CG2	2.43	0.49
2:B:204:ASN:OD1	2:B:207:ALA:HB2	2.12	0.49
3:C:26:LYS:C	3:C:28:GLN:N	2.66	0.49
3:C:91:LEU:HD12	3:C:101:LEU:HD12	1.94	0.49
4:D:148:VAL:HG12	4:D:149:ALA:H	1.76	0.49
7:G:111:ARG:O	7:G:119:ARG:HG2	2.13	0.49
8:H:39:LEU:O	8:H:44:PHE:HB2	2.12	0.49
8:H:120:THR:OG1	8:H:121:ASP:N	2.46	0.49
9:I:3:GLN:NE2	9:I:4:TYR:N	2.60	0.49
10:J:54:PHE:O	10:J:55:LYS:O	2.30	0.49
11:K:34:ASP:OD1	11:K:38:ASN:HB2	2.12	0.49
12:L:28:LYS:C	12:L:30:ALA:H	2.15	0.49
12:L:33:ARG:C	12:L:84:LEU:HD12	2.32	0.49
12:L:109:GLY:N	12:L:121:GLY:O	2.46	0.49
13:M:70:LEU:HD23	13:M:70:LEU:C	2.32	0.49
14:N:6:LEU:O	14:N:23:ARG:HD2	2.12	0.49
15:O:17:ARG:HG3	15:O:17:ARG:HH11	1.77	0.49
15:O:26:GLU:HG3	15:O:81:LEU:HD12	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
18:R:72:ARG:O	18:R:75:ILE:N	2.45	0.49
20:T:99:LEU:O	20:T:101:GLY:N	2.46	0.49
1:A:68:G:H8	1:A:68:G:O5'	1.95	0.49
1:A:84:U:C4	1:A:88:A:C6	3.01	0.49
1:A:411:A:N9	1:A:413:G:H1'	2.27	0.49
1:A:707:C:O3'	11:K:20:TYR:HE2	1.95	0.49
1:A:1438:G:H2'	1:A:1439:C:C6	2.48	0.49
2:B:73:THR:HG23	2:B:96:ARG:HH21	1.72	0.49
2:B:135:GLN:O	2:B:139:LYS:HB2	2.13	0.49
3:C:37:GLN:HE22	14:N:52:GLN:NE2	2.08	0.49
3:C:198:VAL:CG1	3:C:199:LYS:N	2.75	0.49
4:D:56:VAL:HG12	4:D:202:LEU:HD13	1.93	0.49
4:D:177:ASP:OD2	4:D:178:VAL:N	2.45	0.49
4:D:201:GLN:O	4:D:205:GLU:HG3	2.13	0.49
5:E:25:ARG:C	5:E:26:PHE:HD1	2.15	0.49
5:E:93:PRO:CG	8:H:105:ARG:HH21	2.26	0.49
5:E:105:VAL:HB	5:E:106:PRO:HD3	1.95	0.49
6:F:80:ARG:HG2	6:F:80:ARG:HH11	1.77	0.49
7:G:16:LEU:HG	9:I:41:VAL:O	2.12	0.49
7:G:16:LEU:HD12	9:I:44:VAL:HG13	1.95	0.49
7:G:124:LEU:O	7:G:127:ALA:HB3	2.11	0.49
9:I:7:THR:N	9:I:83:ARG:HD2	2.28	0.49
9:I:8:GLY:HA2	9:I:79:LEU:HD13	1.94	0.49
10:J:51:ARG:O	10:J:52:GLY:O	2.31	0.49
13:M:48:LEU:HD21	13:M:52:GLU:HB2	1.94	0.49
13:M:76:ALA:O	13:M:77:ASN:C	2.50	0.49
14:N:22:THR:O	14:N:23:ARG:O	2.29	0.49
16:P:4:ILE:HG12	16:P:21:VAL:CG2	2.43	0.49
16:P:53:VAL:O	16:P:55:ARG:N	2.45	0.49
18:R:73:ALA:CB	18:R:79:LEU:HD13	2.42	0.49
20:T:43:LEU:HD12	20:T:55:ILE:CD1	2.42	0.49
20:T:50:GLU:HB3	20:T:100:ILE:HD11	1.95	0.49
21:V:18:TYR:CE2	21:V:23:PRO:O	2.65	0.49
1:A:54:C:H2'	1:A:352:C:N4	2.28	0.49
1:A:376:G:OP2	16:P:67:THR:HG21	2.12	0.49
1:A:681:C:H2'	1:A:682:G:C8	2.48	0.49
1:A:1347:G:O2'	1:A:1348:U:OP2	2.31	0.49
2:B:74:LYS:O	2:B:77:ALA:N	2.44	0.49
4:D:9:CYS:HB2	4:D:22:LYS:CE	2.43	0.49
4:D:101:LEU:O	4:D:101:LEU:HD12	2.13	0.49
4:D:171:GLY:O	4:D:173:TRP:N	2.46	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:61:VAL:HG13	7:G:128:ALA:CB	2.41	0.49
7:G:68:ASN:C	7:G:70:LYS:H	2.16	0.49
7:G:145:ALA:O	7:G:146:GLU:HB3	2.11	0.49
8:H:35:ILE:O	8:H:39:LEU:HB2	2.13	0.49
8:H:91:ARG:HH12	17:Q:33:GLY:HA3	1.77	0.49
10:J:47:PHE:CB	14:N:44:LEU:HD21	2.43	0.49
13:M:70:LEU:C	13:M:70:LEU:CD2	2.81	0.49
17:Q:98:LEU:HB3	17:Q:103:GLY:H	1.77	0.49
19:S:36:ARG:NE	19:S:52:TYR:O	2.46	0.49
20:T:62:LEU:C	20:T:65:LYS:HB3	2.33	0.49
1:A:248:C:H2'	1:A:249:U:H5'	1.95	0.49
1:A:403:C:H2'	1:A:404:U:H6	1.76	0.49
1:A:606:G:H21	1:A:631:G:H2'	1.77	0.49
1:A:766:A:C8	1:A:814:A:N6	2.81	0.49
1:A:766:A:H2'	1:A:767:A:O4'	2.12	0.49
1:A:865:A:N6	1:A:866:C:N4	2.61	0.49
1:A:935:A:C2	7:G:3:ARG:NH2	2.78	0.49
1:A:1060:C:HO2'	10:J:56:HIS:HD1	1.61	0.49
1:A:1110:A:H2'	1:A:1111:A:H5'	1.95	0.49
1:A:1313:U:C5	19:S:4:SER:OG	2.64	0.49
1:A:1379:G:H2'	1:A:1380:U:H5'	1.94	0.49
2:B:213:LEU:O	2:B:216:SER:OG	2.29	0.49
3:C:95:THR:HG23	3:C:95:THR:O	2.13	0.49
3:C:127:ARG:HG2	3:C:127:ARG:HH11	1.78	0.49
4:D:88:VAL:O	4:D:91:SER:HB3	2.12	0.49
6:F:83:ASP:O	6:F:85:VAL:N	2.45	0.49
8:H:21:LYS:N	8:H:65:TYR:OH	2.46	0.49
10:J:94:VAL:HG12	10:J:95:GLU:N	2.28	0.49
11:K:82:VAL:CG1	11:K:83:ILE:N	2.76	0.49
13:M:84:ILE:CG1	13:M:84:ILE:O	2.61	0.49
14:N:25:VAL:HG12	14:N:26:ARG:N	2.27	0.49
18:R:87:ARG:O	18:R:88:LYS:HB2	2.13	0.49
20:T:23:ARG:O	20:T:26:ASN:N	2.45	0.49
20:T:52:ALA:O	20:T:53:LEU:C	2.50	0.49
21:V:4:GLY:O	21:V:6:ARG:N	2.45	0.49
1:A:36:C:O2'	1:A:501:C:OP1	2.31	0.49
1:A:119:A:H4'	1:A:120:A:C8	2.48	0.49
1:A:255:G:C2	1:A:272:C:C2	3.00	0.49
1:A:421:U:H3'	1:A:422:C:C5'	2.41	0.49
1:A:781:A:C3'	1:A:782:A:H5'	2.42	0.49
1:A:881:G:OP2	12:L:12:ARG:NH2	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:908:A:C2	1:A:909:A:C4	3.00	0.49
1:A:1210:C:C2'	1:A:1211:U:H5''	2.42	0.49
1:A:1317:C:H2'	1:A:1318:A:O4'	2.13	0.49
1:A:1468:A:H3'	1:A:1469:G:C8	2.48	0.49
1:A:1509:C:C2'	1:A:1510:U:C5'	2.90	0.49
2:B:34:ALA:O	2:B:41:ILE:N	2.45	0.49
3:C:37:GLN:NE2	14:N:52:GLN:NE2	2.59	0.49
3:C:48:TYR:CA	3:C:52:LEU:HB3	2.43	0.49
4:D:9:CYS:HA	4:D:22:LYS:HD2	1.94	0.49
4:D:61:LYS:NZ	4:D:72:GLU:OE2	2.44	0.49
4:D:104:VAL:HG11	4:D:146:ILE:HD13	1.95	0.49
5:E:36:ASP:C	5:E:38:GLN:N	2.66	0.49
7:G:32:ARG:O	7:G:32:ARG:HG3	2.12	0.49
7:G:66:VAL:O	7:G:70:LYS:HG3	2.13	0.49
7:G:82:GLY:HA3	7:G:85:TYR:OH	2.13	0.49
12:L:18:VAL:O	12:L:19:ARG:CB	2.60	0.49
12:L:41:ARG:HH11	12:L:42:THR:H	1.60	0.49
12:L:46:LYS:CE	12:L:47:LYS:HE2	2.43	0.49
12:L:93:LEU:O	12:L:96:VAL:HG23	2.13	0.49
13:M:81:LEU:HA	13:M:84:ILE:HD11	1.95	0.49
13:M:91:ARG:HB3	13:M:98:VAL:HG22	1.91	0.49
14:N:8:GLU:O	14:N:8:GLU:HG2	2.12	0.49
15:O:5:LYS:O	15:O:8:LYS:HB3	2.12	0.49
16:P:19:ILE:O	16:P:36:ILE:CG1	2.61	0.49
18:R:21:LYS:O	18:R:22:VAL:O	2.31	0.49
19:S:11:VAL:HG21	19:S:41:VAL:CG1	2.40	0.49
19:S:55:LYS:HE2	19:S:56:GLN:NE2	2.28	0.49
20:T:80:ARG:HH11	20:T:80:ARG:CB	2.24	0.49
1:A:190(K):G:O2'	1:A:190(L):U:H5'	2.12	0.49
1:A:241:C:O2'	1:A:242:C:H5'	2.12	0.49
1:A:245:C:O2	1:A:283:C:N3	2.46	0.49
1:A:325:A:H2'	1:A:326:G:O4'	2.12	0.49
1:A:357:G:C2'	1:A:358:U:H5'	2.42	0.49
1:A:433:C:H2'	1:A:434:U:H6	1.77	0.49
1:A:695:A:H2'	1:A:696:A:C8	2.48	0.49
1:A:860:A:H2'	1:A:861:G:O4'	2.13	0.49
1:A:979:C:H3'	1:A:980:C:C6	2.48	0.49
1:A:1423:G:O2'	1:A:1424:C:H5'	2.13	0.49
1:A:1483:A:C2	1:A:1484:C:H1'	2.47	0.49
1:A:1509:C:H2'	1:A:1510:U:C5'	2.43	0.49
2:B:114:ARG:O	2:B:115:LEU:C	2.51	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:123:ALA:O	2:B:124:SER:C	2.50	0.49
2:B:141:GLU:O	2:B:145:LEU:HG	2.12	0.49
2:B:178:ARG:C	2:B:180:LEU:N	2.65	0.49
3:C:44:GLU:HG3	3:C:52:LEU:HA	1.94	0.49
4:D:148:VAL:HG23	4:D:181:MET:HB3	1.93	0.49
4:D:163:GLU:C	4:D:165:MET:N	2.65	0.49
5:E:36:ASP:OD2	5:E:40:ARG:HB2	2.13	0.49
7:G:57:GLU:O	7:G:59:LEU:N	2.46	0.49
8:H:37:ARG:O	8:H:41:ARG:HB2	2.13	0.49
8:H:100:ILE:HG21	8:H:121:ASP:HB2	1.95	0.49
9:I:80:GLY:O	9:I:84:ALA:N	2.39	0.49
10:J:8:LEU:HD13	10:J:70:ARG:HB2	1.94	0.49
10:J:42:THR:HG23	10:J:68:HIS:CA	2.36	0.49
10:J:85:LEU:H	10:J:88:LEU:CD1	2.25	0.49
13:M:81:LEU:HD13	13:M:88:ARG:CB	2.31	0.49
15:O:8:LYS:HG2	15:O:12:ILE:HD11	1.94	0.49
15:O:9:GLN:HA	15:O:12:ILE:HD13	1.95	0.49
15:O:78:TYR:O	15:O:79:ARG:C	2.50	0.49
16:P:78:GLY:HA2	16:P:80:PHE:HD1	1.78	0.49
18:R:70:ILE:O	18:R:71:LYS:C	2.51	0.49
19:S:41:VAL:HG23	19:S:44:MET:HG3	1.93	0.49
20:T:65:LYS:O	20:T:68:LYS:N	2.40	0.49
20:T:78:ALA:O	20:T:82:SER:N	2.43	0.49
21:V:14:TRP:O	21:V:16:GLY:N	2.46	0.49
1:A:35:G:H2'	1:A:36:C:C6	2.48	0.48
1:A:176:C:H2'	1:A:177:C:H6	1.78	0.48
1:A:267:C:C4	1:A:268:C:N4	2.81	0.48
1:A:296:U:H2'	1:A:297:G:C8	2.48	0.48
1:A:337:C:H2'	1:A:338:A:C8	2.48	0.48
1:A:502:G:OP1	12:L:118:SER:N	2.46	0.48
1:A:658:G:O2'	1:A:659:U:H5'	2.12	0.48
1:A:791:G:N1	1:A:792:A:N6	2.61	0.48
1:A:988:G:O2'	1:A:1016:A:H2	1.96	0.48
1:A:1215:G:H2'	1:A:1216:G:C5'	2.43	0.48
1:A:1464:G:O2'	1:A:1465:C:H5'	2.13	0.48
2:B:24:TRP:HA	2:B:190:THR:HG23	1.94	0.48
2:B:79:ASP:O	2:B:82:ARG:N	2.46	0.48
2:B:172:ILE:HD12	2:B:172:ILE:N	2.10	0.48
3:C:51:GLY:C	3:C:53:ALA:H	2.15	0.48
3:C:58:GLU:HG2	3:C:67:THR:OG1	2.14	0.48
3:C:113:ALA:HB3	3:C:114:PRO:CD	2.40	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:38:GLU:O	6:F:38:GLU:HG2	2.13	0.48
9:I:10:ARG:HD2	9:I:11:LYS:N	2.28	0.48
9:I:45:ALA:O	9:I:48:GLU:HB3	2.13	0.48
11:K:44:SER:O	11:K:46:GLY:N	2.46	0.48
15:O:10:LYS:CG	15:O:11:VAL:N	2.75	0.48
15:O:61:GLY:O	15:O:65:ARG:N	2.41	0.48
15:O:82:ILE:O	15:O:82:ILE:HG12	2.13	0.48
19:S:15:LEU:HD21	19:S:38:SER:HB3	1.95	0.48
19:S:16:LEU:O	19:S:20:LEU:HG	2.13	0.48
1:A:11:G:C5	1:A:12:U:C4	3.01	0.48
1:A:635:G:H2'	1:A:636:U:H6	1.78	0.48
1:A:657:G:O2'	1:A:658:G:H5'	2.13	0.48
1:A:814:A:O2'	1:A:815:A:H3'	2.14	0.48
1:A:986:A:C2	19:S:52:TYR:HE2	2.30	0.48
1:A:1124:G:H2'	1:A:1145:C:H41	1.78	0.48
2:B:33:TYR:O	2:B:34:ALA:HB2	2.12	0.48
2:B:107:THR:O	2:B:110:GLN:HG3	2.13	0.48
2:B:126:GLU:OE2	2:B:129:GLU:HB3	2.13	0.48
4:D:61:LYS:O	4:D:62:GLN:C	2.52	0.48
5:E:79:GLU:OE1	8:H:104:ARG:HD3	2.13	0.48
10:J:28:ARG:HG2	10:J:28:ARG:O	2.13	0.48
11:K:60:ALA:O	11:K:61:ALA:C	2.50	0.48
11:K:81:ASP:OD1	11:K:106:LYS:CB	2.60	0.48
16:P:57:ARG:O	16:P:61:SER:N	2.47	0.48
17:Q:59:ILE:HG22	17:Q:71:PHE:CD1	2.47	0.48
19:S:31:ILE:HG22	19:S:32:LYS:N	2.28	0.48
19:S:33:THR:HG23	19:S:33:THR:O	2.13	0.48
1:A:142:G:N3	1:A:196:A:H2	2.09	0.48
1:A:231:G:C2	1:A:232:G:C8	3.01	0.48
1:A:255:G:O3'	17:Q:17:LYS:HD2	2.12	0.48
1:A:455:C:H2'	1:A:456:C:C6	2.48	0.48
1:A:502:G:H2'	1:A:503:C:C6	2.49	0.48
2:B:44:LEU:O	2:B:45:GLN:C	2.49	0.48
2:B:118:LEU:C	2:B:120:ALA:N	2.66	0.48
2:B:223:ILE:O	2:B:224:GLN:C	2.51	0.48
4:D:176:LEU:CG	4:D:177:ASP:N	2.49	0.48
6:F:91:VAL:HG11	18:R:72:ARG:CZ	2.43	0.48
7:G:79:ARG:HA	7:G:83:ALA:O	2.12	0.48
11:K:52:GLY:O	11:K:53:SER:C	2.52	0.48
13:M:60:VAL:HG13	13:M:64:TRP:CZ3	2.48	0.48
16:P:17:TYR:N	16:P:17:TYR:HD1	2.11	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
19:S:42:PRO:HG3	19:S:64:GLU:OE1	2.13	0.48
20:T:48:LYS:HB3	20:T:51:GLU:CG	2.43	0.48
1:A:393:A:C4	1:A:394:G:C8	3.01	0.48
1:A:518:C:N4	12:L:49:ASN:HD21	2.12	0.48
1:A:644:G:C5	1:A:645:C:C5	3.00	0.48
1:A:696:A:O5'	1:A:696:A:H8	1.97	0.48
1:A:913:A:H1'	1:A:914:A:C1'	2.43	0.48
1:A:1025:U:H2'	1:A:1026:G:H8	1.79	0.48
1:A:1075:C:O2'	1:A:1076:C:H5'	2.14	0.48
1:A:1316:G:C4	1:A:1318:A:OP2	2.65	0.48
1:A:1440:C:H2'	1:A:1441:G:C5'	2.41	0.48
3:C:66:VAL:CG1	3:C:68:VAL:HB	2.33	0.48
5:E:128:PRO:O	5:E:130:ASN:N	2.47	0.48
6:F:55:ASP:C	6:F:57:GLN:H	2.15	0.48
6:F:73:ASN:O	6:F:75:LEU:N	2.47	0.48
7:G:26:PHE:O	7:G:29:LYS:N	2.39	0.48
7:G:46:ALA:CB	7:G:121:ALA:H	2.26	0.48
8:H:20:TYR:HA	8:H:65:TYR:CE2	2.49	0.48
8:H:20:TYR:HA	8:H:65:TYR:CZ	2.48	0.48
9:I:15:ALA:HB2	9:I:65:VAL:CG2	2.43	0.48
9:I:91:ASP:N	9:I:91:ASP:OD2	2.47	0.48
9:I:97:LYS:HZ1	9:I:102:LEU:CD1	2.26	0.48
9:I:97:LYS:HZ1	9:I:102:LEU:CG	2.27	0.48
10:J:40:LEU:HD13	10:J:69:ASN:CB	2.39	0.48
10:J:46:ARG:NH1	10:J:46:ARG:HG2	2.27	0.48
15:O:17:ARG:CD	15:O:77:ARG:HH11	2.24	0.48
16:P:20:VAL:O	16:P:21:VAL:CB	2.61	0.48
18:R:29:PHE:HE1	18:R:31:LEU:HA	1.79	0.48
18:R:87:ARG:NH1	18:R:87:ARG:CG	2.75	0.48
20:T:94:ALA:O	20:T:95:ALA:CB	2.61	0.48
1:A:490:G:O2'	1:A:491:G:H5'	2.13	0.48
1:A:508:C:P	4:D:209:ARG:HH22	2.36	0.48
1:A:544:G:H2'	1:A:545:C:H6	1.79	0.48
1:A:605:U:O2'	1:A:606:G:H5'	2.12	0.48
1:A:617:G:H1	1:A:623:C:H42	1.62	0.48
1:A:892:A:C6	1:A:893:C:C4	3.02	0.48
1:A:914:A:O2'	1:A:915:A:H5'	2.14	0.48
1:A:1110:A:O5'	1:A:1110:A:C8	2.65	0.48
1:A:1152:A:H5''	10:J:13:HIS:C	2.33	0.48
1:A:1305:G:O2'	1:A:1306:A:P	2.72	0.48
2:B:139:LYS:O	2:B:143:GLU:HG3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:14:ILE:HD13	3:C:14:ILE:N	2.28	0.48
3:C:91:LEU:HG	3:C:99:VAL:CG1	2.44	0.48
4:D:7:PRO:HB2	4:D:10:ARG:HB3	1.96	0.48
4:D:57:ARG:HB3	4:D:206:PHE:HB2	1.94	0.48
4:D:63:LYS:C	4:D:65:ARG:H	2.17	0.48
4:D:205:GLU:C	4:D:207:TYR:N	2.64	0.48
5:E:151:LEU:HD11	8:H:77:GLU:OE2	2.13	0.48
7:G:29:LYS:NZ	7:G:102:ARG:HA	2.16	0.48
11:K:92:GLU:CA	11:K:95:ILE:HD12	2.38	0.48
12:L:60:LEU:HD21	12:L:85:ILE:CD1	2.43	0.48
18:R:21:LYS:HD2	18:R:21:LYS:N	2.19	0.48
18:R:24:ALA:C	18:R:26:LEU:N	2.60	0.48
18:R:65:ILE:O	18:R:66:LEU:C	2.51	0.48
1:A:337:C:H2'	1:A:338:A:H8	1.78	0.48
1:A:345:C:H4'	1:A:346:G:H5''	1.96	0.48
1:A:509:A:H2'	1:A:510:A:C8	2.48	0.48
1:A:642:A:C8	8:H:115:SER:HA	2.48	0.48
1:A:768:A:C2'	1:A:769:G:H5'	2.43	0.48
1:A:939:G:H4'	7:G:102:ARG:HH21	1.78	0.48
1:A:948:C:C5	13:M:106:ASN:OD1	2.67	0.48
1:A:1114:C:N4	1:A:1115:C:C4	2.81	0.48
1:A:1231:G:C2	1:A:1232:U:C2	3.02	0.48
1:A:1360:A:C3'	1:A:1361:G:C8	2.97	0.48
1:A:1379:G:C2'	1:A:1380:U:H5'	2.43	0.48
2:B:61:LEU:HD23	2:B:66:GLY:HA3	1.96	0.48
2:B:111:ARG:C	2:B:113:HIS:N	2.66	0.48
3:C:5:ILE:CG1	3:C:6:HIS:H	2.26	0.48
3:C:48:TYR:HB2	3:C:52:LEU:CB	2.36	0.48
3:C:126:ARG:O	3:C:128:PHE:N	2.46	0.48
3:C:162:GLN:HG3	3:C:164:ARG:CD	2.43	0.48
3:C:205:GLY:O	3:C:206:GLU:CB	2.60	0.48
4:D:136:PRO:C	4:D:138:TYR:H	2.16	0.48
4:D:151:LYS:H	4:D:151:LYS:CD	2.27	0.48
4:D:173:TRP:CD1	4:D:187:ARG:O	2.66	0.48
7:G:156:TRP:O	7:G:156:TRP:HD1	1.96	0.48
10:J:60:ARG:H	10:J:60:ARG:CD	2.27	0.48
11:K:124:LYS:HB3	11:K:125:PHE:CD1	2.49	0.48
14:N:24:CYS:SG	14:N:27:CYS:CB	2.83	0.48
15:O:70:LEU:HD12	15:O:78:TYR:HA	1.95	0.48
1:A:102:G:N2	1:A:171:A:C2	2.81	0.48
1:A:841:U:H5'	1:A:848:C:C6	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1296:C:H5'	13:M:44:ARG:HH22	1.79	0.48
1:A:1385:G:O2'	1:A:1386:G:H5'	2.14	0.48
1:A:1397:C:HO2'	1:A:1398:A:P	2.36	0.48
1:A:1437:C:O2'	1:A:1438:G:H5'	2.13	0.48
1:A:1518:A:C8	1:A:1518:A:H3'	2.48	0.48
2:B:73:THR:HG23	2:B:96:ARG:CZ	2.42	0.48
2:B:109:SER:C	2:B:111:ARG:H	2.16	0.48
2:B:184:VAL:HG11	2:B:197:VAL:HG22	1.95	0.48
3:C:139:GLN:CA	3:C:139:GLN:NE2	2.75	0.48
3:C:153:VAL:HG22	3:C:198:VAL:CG2	2.44	0.48
3:C:184:TYR:HA	3:C:200:ALA:O	2.12	0.48
5:E:32:VAL:HG12	5:E:33:VAL:N	2.29	0.48
6:F:40:VAL:HG23	6:F:41:GLU:N	2.27	0.48
7:G:65:ALA:HB1	7:G:127:ALA:C	2.34	0.48
7:G:131:LYS:O	7:G:131:LYS:HG3	2.12	0.48
8:H:95:VAL:HB	8:H:99:GLU:HB3	1.95	0.48
11:K:17:GLY:H	11:K:77:MET:CE	2.22	0.48
16:P:68:ASP:O	16:P:71:ARG:HG3	2.13	0.48
18:R:37:VAL:HG23	18:R:38:GLU:H	1.77	0.48
19:S:22:LEU:CD2	19:S:31:ILE:HD11	2.43	0.48
20:T:76:ALA:C	20:T:80:ARG:HG2	2.34	0.48
1:A:184:G:O2'	1:A:185:A:H5'	2.14	0.48
1:A:219:C:H2'	1:A:220:G:C5'	2.41	0.48
1:A:371:G:H21	1:A:374:A:H62	1.62	0.48
1:A:458:C:N3	1:A:459:G:C2	2.82	0.48
1:A:922:G:C6	1:A:923:A:N1	2.82	0.48
1:A:925:G:C2	1:A:927:G:C8	3.02	0.48
1:A:1176:A:C6	1:A:1177:G:C2	3.02	0.48
1:A:1288:A:N6	1:A:1289:A:N6	2.62	0.48
1:A:1367:C:H5'	10:J:60:ARG:HH12	1.76	0.48
1:A:1397:C:H4'	1:A:1398:A:OP2	2.14	0.48
4:D:4:TYR:HE2	4:D:6:GLY:C	2.17	0.48
4:D:31:CYS:O	4:D:31:CYS:SG	2.72	0.48
6:F:74:ASP:HA	6:F:77:ARG:HE	1.78	0.48
7:G:137:LYS:O	7:G:138:LYS:C	2.50	0.48
8:H:23:SER:HA	8:H:63:LEU:HD13	1.96	0.48
10:J:3:LYS:HB2	10:J:76:ASN:HA	1.96	0.48
10:J:13:HIS:ND1	10:J:14:LYS:N	2.61	0.48
12:L:89:ARG:C	12:L:90:VAL:HG23	2.34	0.48
13:M:80:ARG:C	13:M:82:MET:H	2.17	0.48
15:O:68:ARG:CB	15:O:68:ARG:HH11	2.27	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:Q:76:LEU:HD12	17:Q:77:VAL:N	2.28	0.48
18:R:22:VAL:CG1	18:R:23:LYS:N	2.68	0.48
21:V:5:ASP:HB3	21:V:8:THR:OG1	2.13	0.48
1:A:73:C:C2	1:A:74:C:C5	3.02	0.48
1:A:152:A:N6	1:A:170:U:N3	2.62	0.48
1:A:321:A:H2	1:A:332:G:H22	1.61	0.48
1:A:325:A:N6	1:A:326:G:C6	2.82	0.48
1:A:551:U:H2'	1:A:552:U:H6	1.77	0.48
1:A:975:A:H4'	1:A:976:G:O5'	2.14	0.48
1:A:1075:C:O5'	1:A:1075:C:H6	1.97	0.48
1:A:1162:C:C2	1:A:1175:G:C2	3.02	0.48
1:A:1247:U:C2'	1:A:1248:A:H5'	2.44	0.48
1:A:1397:C:O2'	1:A:1398:A:P	2.71	0.48
3:C:99:VAL:HG23	3:C:101:LEU:N	2.29	0.48
4:D:79:PHE:O	4:D:79:PHE:CD2	2.62	0.48
4:D:156:GLU:N	4:D:156:GLU:CD	2.67	0.48
5:E:129:ILE:O	5:E:129:ILE:CG2	2.62	0.48
7:G:68:ASN:O	7:G:138:LYS:HE3	2.14	0.48
7:G:70:LYS:HB3	7:G:96:GLN:HG2	1.96	0.48
11:K:97:ALA:O	11:K:100:ALA:N	2.44	0.48
15:O:16:ALA:C	15:O:18:PHE:N	2.67	0.48
17:Q:70:ARG:CG	17:Q:70:ARG:NH1	2.62	0.48
18:R:21:LYS:NZ	18:R:54:ARG:O	2.39	0.48
18:R:37:VAL:HG23	18:R:41:LYS:HE3	1.95	0.48
18:R:43:PHE:CD1	18:R:43:PHE:N	2.82	0.48
1:A:67:C:HO2'	1:A:171:A:H1'	1.79	0.48
1:A:236:G:H2'	1:A:237:C:O4'	2.14	0.48
1:A:287:U:H2'	1:A:288:A:H5'	1.94	0.48
1:A:376:G:C2	1:A:389:A:C2	3.02	0.48
1:A:405:U:C3'	1:A:406:G:H5'	2.36	0.48
1:A:602:A:O2'	1:A:603:U:H5'	2.14	0.48
1:A:607:A:C4	1:A:608:A:C8	3.02	0.48
1:A:947:G:C2	1:A:948:C:C2	3.02	0.48
1:A:1151:A:C2	10:J:39:PRO:HG3	2.49	0.48
1:A:1226:C:OP2	13:M:103:THR:HG21	2.13	0.48
1:A:1254:C:N4	10:J:43:ARG:CZ	2.77	0.48
1:A:1437:C:C6	1:A:1438:G:N7	2.81	0.48
3:C:44:GLU:CG	3:C:55:VAL:HG22	2.43	0.48
3:C:51:GLY:HA3	3:C:71:ALA:H	1.79	0.48
3:C:162:GLN:O	3:C:164:ARG:HG3	2.14	0.48
3:C:180:ALA:HA	3:C:207:VAL:H	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:118:ARG:O	4:D:121:VAL:HB	2.14	0.48
6:F:13:ASN:ND2	6:F:57:GLN:OE1	2.46	0.48
7:G:21:VAL:HA	7:G:24:THR:OG1	2.14	0.48
8:H:11:THR:CG2	8:H:14:ARG:NH1	2.67	0.48
10:J:20:ALA:C	10:J:22:LYS:N	2.67	0.48
10:J:29:ARG:HG3	10:J:29:ARG:O	2.13	0.48
13:M:77:ASN:O	13:M:81:LEU:HG	2.14	0.48
13:M:81:LEU:CD1	13:M:88:ARG:HB3	2.33	0.48
13:M:117:VAL:O	13:M:118:ALA:CB	2.60	0.48
16:P:20:VAL:O	16:P:21:VAL:CG2	2.59	0.48
16:P:66:PRO:O	16:P:67:THR:C	2.52	0.48
17:Q:94:ASN:O	17:Q:95:TYR:C	2.52	0.48
21:V:8:THR:HG22	21:V:9:ARG:N	2.29	0.48
1:A:341:C:O5'	1:A:341:C:H6	1.96	0.47
1:A:960:U:H4'	1:A:961:U:H5''	1.96	0.47
1:A:1327:C:H5''	21:V:20:LYS:HB2	1.96	0.47
2:B:97:TRP:CE3	2:B:98:LEU:O	2.67	0.47
3:C:48:TYR:CD1	3:C:52:LEU:HD22	2.46	0.47
6:F:86:ARG:H	6:F:86:ARG:HG2	1.46	0.47
7:G:105:VAL:O	7:G:105:VAL:HG12	2.14	0.47
8:H:40:ALA:HA	8:H:45:ILE:HG13	1.95	0.47
11:K:13:GLN:HA	11:K:75:TYR:O	2.14	0.47
12:L:113:ARG:HD3	12:L:114:LYS:N	2.29	0.47
13:M:55:ARG:O	13:M:59:TYR:N	2.41	0.47
15:O:71:GLN:C	15:O:71:GLN:OE1	2.52	0.47
15:O:79:ARG:O	15:O:83:GLU:HB3	2.13	0.47
21:V:6:ARG:HG2	21:V:15:ARG:NH1	2.21	0.47
1:A:151:A:H2'	1:A:152:A:C8	2.49	0.47
1:A:164:U:H2'	1:A:165:C:C6	2.50	0.47
1:A:253:U:H2'	1:A:254:G:H8	1.78	0.47
1:A:255:G:N2	1:A:272:C:C2	2.83	0.47
1:A:438:G:H22	1:A:496:A:P	2.37	0.47
1:A:688:G:H2'	1:A:689:C:C6	2.49	0.47
1:A:707:C:H4'	11:K:20:TYR:CE2	2.49	0.47
1:A:778:G:O2'	1:A:779:C:H5'	2.14	0.47
1:A:825:G:C6	1:A:876:G:C6	3.02	0.47
1:A:1127:G:H8	1:A:1127:G:OP2	1.97	0.47
1:A:1369:C:C2'	1:A:1370:G:O4'	2.58	0.47
1:A:1432:G:H1'	1:A:1468:A:H61	1.79	0.47
1:A:1435:G:H2'	1:A:1436:U:C5	2.49	0.47
2:B:138:LEU:HD22	2:B:138:LEU:N	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:47:ARG:HE	4:D:49:ARG:CA	2.26	0.47
4:D:91:SER:O	4:D:92:VAL:C	2.52	0.47
7:G:69:VAL:O	7:G:138:LYS:HB2	2.14	0.47
8:H:9:MET:HG3	8:H:26:VAL:HG21	1.95	0.47
8:H:40:ALA:HB1	8:H:47:GLY:HA2	1.96	0.47
8:H:88:LYS:C	8:H:89:PRO:O	2.50	0.47
14:N:22:THR:HB	14:N:33:VAL:CB	2.44	0.47
14:N:47:LEU:C	14:N:49:HIS:N	2.67	0.47
15:O:30:ALA:O	15:O:31:LEU:C	2.51	0.47
15:O:76:GLU:O	15:O:77:ARG:C	2.52	0.47
16:P:31:LYS:CG	16:P:32:TYR:N	2.77	0.47
16:P:71:ARG:HG2	16:P:71:ARG:HH11	1.79	0.47
16:P:78:GLY:C	16:P:80:PHE:H	2.17	0.47
17:Q:18:THR:HG23	17:Q:69:LYS:HZ2	1.79	0.47
20:T:17:ARG:O	20:T:20:LEU:HB2	2.14	0.47
1:A:134:A:H8	1:A:134:A:O5'	1.96	0.47
1:A:264:U:O2'	17:Q:64:PRO:HB2	2.14	0.47
1:A:750:G:C2	15:O:23:GLY:HA3	2.49	0.47
1:A:787:A:O2'	1:A:788:U:H5'	2.14	0.47
1:A:877:C:O2	8:H:3:THR:HG21	2.13	0.47
1:A:956:U:H2'	1:A:957:U:H5'	1.96	0.47
1:A:1064:G:H5'	1:A:1066:C:O4'	2.14	0.47
1:A:1102:A:C6	1:A:1103:C:N4	2.82	0.47
1:A:1210:C:H2'	1:A:1211:U:H4'	1.97	0.47
1:A:1288:A:H2'	1:A:1289:A:H8	1.79	0.47
1:A:1392:G:H2'	1:A:1393:U:H6	1.79	0.47
3:C:113:ALA:O	3:C:116:VAL:N	2.47	0.47
3:C:153:VAL:HG13	3:C:198:VAL:HG22	1.95	0.47
3:C:175:LEU:N	3:C:175:LEU:CD2	2.77	0.47
6:F:53:ALA:C	6:F:55:ASP:N	2.67	0.47
7:G:5:ARG:C	7:G:7:ALA:H	2.18	0.47
7:G:25:ALA:O	7:G:29:LYS:N	2.47	0.47
7:G:91:VAL:HG13	7:G:92:SER:H	1.78	0.47
8:H:126:LYS:C	8:H:128:GLY:N	2.68	0.47
12:L:38:THR:HG22	12:L:39:VAL:HG23	1.95	0.47
13:M:19:LEU:C	13:M:21:TYR:N	2.64	0.47
15:O:35:ARG:O	15:O:36:ILE:C	2.51	0.47
15:O:49:ASP:CG	15:O:49:ASP:O	2.53	0.47
1:A:128:G:O2'	17:Q:3:LYS:HE3	2.14	0.47
1:A:132:C:H2'	1:A:133:U:O5'	2.14	0.47
1:A:231:G:H2'	1:A:231:G:N3	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:419:C:C5	1:A:425:G:C2	3.02	0.47
1:A:425:G:C2	1:A:426:G:C8	3.02	0.47
1:A:487:A:H2'	1:A:488:C:O4'	2.15	0.47
1:A:506:G:N7	1:A:507:C:C5	2.82	0.47
1:A:720:C:N3	1:A:721:G:C6	2.82	0.47
1:A:942:G:N2	1:A:943:U:N1	2.63	0.47
1:A:1056:U:C5'	3:C:163:ALA:CB	2.92	0.47
1:A:1074:G:O4'	2:B:104:ASN:HB2	2.15	0.47
1:A:1103:C:H2'	1:A:1104:G:O4'	2.14	0.47
1:A:1365:G:C2'	1:A:1366:C:H5'	2.45	0.47
2:B:96:ARG:HE	2:B:96:ARG:CA	2.27	0.47
2:B:112:VAL:HG21	2:B:152:PHE:O	2.14	0.47
4:D:130:GLY:O	4:D:132:ARG:N	2.47	0.47
5:E:16:THR:CG2	5:E:27:ARG:HB2	2.44	0.47
5:E:133:TYR:O	5:E:137:GLU:HB2	2.14	0.47
6:F:48:LEU:O	6:F:49:ALA:C	2.53	0.47
7:G:68:ASN:C	7:G:70:LYS:N	2.68	0.47
11:K:69:ALA:O	11:K:72:ALA:HB3	2.14	0.47
12:L:42:THR:HG22	12:L:52:LEU:HD22	1.97	0.47
16:P:39:TYR:HA	16:P:49:LEU:HD12	1.97	0.47
19:S:19:VAL:HA	19:S:22:LEU:CD1	2.43	0.47
19:S:34:TRP:O	19:S:35:SER:C	2.52	0.47
19:S:70:LYS:H	19:S:73:GLU:CD	2.17	0.47
1:A:79:G:C2	1:A:80:G:C8	3.02	0.47
1:A:519:C:O2'	1:A:520:A:C5'	2.63	0.47
1:A:1056:U:O5'	3:C:163:ALA:HB2	2.15	0.47
1:A:1210:C:C3'	1:A:1211:U:H5''	2.44	0.47
1:A:1229:A:O2'	1:A:1230:C:H5'	2.13	0.47
1:A:1305:G:O2'	1:A:1306:A:O5'	2.32	0.47
1:A:1382:C:H2'	1:A:1383:C:C5	2.49	0.47
1:A:1419:G:N7	1:A:1420:C:C4	2.82	0.47
1:A:1513:A:H2'	1:A:1514:C:C6	2.49	0.47
2:B:210:SER:O	2:B:213:LEU:N	2.48	0.47
3:C:179:ARG:C	3:C:181:ASN:N	2.66	0.47
4:D:128:VAL:HG11	4:D:138:TYR:HE2	1.80	0.47
4:D:202:LEU:O	4:D:205:GLU:HB2	2.13	0.47
5:E:33:VAL:C	5:E:34:VAL:HG23	2.33	0.47
5:E:146:ALA:O	5:E:149:GLU:HB2	2.14	0.47
6:F:10:LEU:HD23	6:F:85:VAL:HA	1.96	0.47
7:G:63:LYS:HB3	7:G:63:LYS:NZ	2.30	0.47
11:K:67:ASP:O	11:K:69:ALA:N	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:K:97:ALA:O	11:K:100:ALA:HB3	2.14	0.47
12:L:83:VAL:HG22	12:L:84:LEU:N	2.20	0.47
17:Q:11:VAL:HA	17:Q:85:VAL:HG22	1.96	0.47
18:R:33:ASP:OD2	18:R:36:ASN:HB3	2.15	0.47
19:S:15:LEU:HD12	19:S:16:LEU:H	1.77	0.47
19:S:22:LEU:HD21	19:S:31:ILE:HD11	1.96	0.47
20:T:62:LEU:HA	20:T:65:LYS:HB3	1.96	0.47
1:A:142:G:O2'	1:A:143:A:H5'	2.13	0.47
1:A:306:G:C4	1:A:307:C:C5	3.03	0.47
1:A:447:G:H2'	1:A:485:G:N2	2.30	0.47
1:A:560:U:O2'	1:A:561:U:OP2	2.30	0.47
1:A:588:G:C6	1:A:589:C:N4	2.83	0.47
1:A:632:A:C2'	1:A:633:G:H5'	2.45	0.47
1:A:913:A:H4'	1:A:914:A:O5'	2.15	0.47
1:A:1413:A:H2'	1:A:1414:U:O4'	2.15	0.47
2:B:21:ARG:HG2	2:B:22:LYS:H	1.76	0.47
2:B:25:ASN:O	2:B:26:PRO:C	2.53	0.47
2:B:73:THR:O	2:B:73:THR:HG22	2.13	0.47
2:B:108:ILE:HG13	2:B:108:ILE:O	2.13	0.47
3:C:191:THR:HG21	3:C:193:TYR:CE1	2.49	0.47
4:D:38:TYR:CG	4:D:45:GLN:HG2	2.49	0.47
5:E:28:PHE:CE1	5:E:50:GLU:HA	2.49	0.47
10:J:22:LYS:HD2	10:J:90:LEU:HD13	1.97	0.47
10:J:64:GLU:HG3	14:N:59:ALA:HA	1.97	0.47
11:K:53:SER:O	11:K:54:ARG:C	2.52	0.47
13:M:23:TYR:HE2	13:M:70:LEU:HD22	1.78	0.47
16:P:14:ASN:N	16:P:15:PRO:CD	2.78	0.47
16:P:48:TRP:HE3	16:P:49:LEU:HB2	1.76	0.47
16:P:80:PHE:O	16:P:82:GLN:NE2	2.45	0.47
17:Q:96:GLN:NE2	17:Q:97:SER:OG	2.44	0.47
1:A:11:G:C6	1:A:12:U:C4	3.03	0.47
1:A:243:A:O2'	1:A:244:U:OP2	2.29	0.47
1:A:391:G:C6	1:A:392:G:C5	3.02	0.47
1:A:393:A:OP2	16:P:12:LYS:HE2	2.14	0.47
1:A:430:A:P	4:D:7:PRO:HA	2.54	0.47
1:A:533:A:C6	1:A:536:C:C4	3.03	0.47
1:A:579:G:H5'	1:A:728:A:H1'	1.96	0.47
1:A:608:A:C2'	1:A:609:A:H8	2.23	0.47
1:A:616:G:C2'	1:A:617:G:H5'	2.44	0.47
1:A:665:A:H1'	1:A:733:A:O4'	2.15	0.47
1:A:666:G:H5'	1:A:726:C:H1'	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:894:G:C2	1:A:895:G:C5	3.03	0.47
1:A:951:G:O2'	1:A:952:U:H5'	2.14	0.47
1:A:986:A:H2	19:S:52:TYR:HE2	1.62	0.47
1:A:987:G:N2	1:A:1219:U:H3	2.12	0.47
1:A:1188:A:O3'	14:N:58:LYS:HE2	2.15	0.47
1:A:1267:C:O2	21:V:20:LYS:HD2	2.15	0.47
1:A:1287:A:C5	1:A:1288:A:C6	3.02	0.47
1:A:1313:U:OP2	19:S:6:LYS:HG2	2.15	0.47
1:A:1322:C:H6	1:A:1322:C:OP1	1.98	0.47
1:A:1326:C:H2'	1:A:1327:C:H6	1.78	0.47
1:A:1356:G:C2	1:A:1357:A:C5	3.02	0.47
1:A:1360:A:C2'	1:A:1361:G:C8	2.98	0.47
1:A:1402:C:H2'	1:A:1403:C:C6	2.50	0.47
1:A:1443:G:C4'	1:A:1446:A:H5'	2.42	0.47
2:B:120:ALA:C	2:B:122:PHE:H	2.17	0.47
2:B:134:GLU:O	2:B:136:VAL:N	2.46	0.47
3:C:14:ILE:H	3:C:14:ILE:HD13	1.79	0.47
3:C:19:GLU:N	14:N:51:GLY:HA3	2.24	0.47
3:C:23:TYR:CD1	3:C:24:ALA:N	2.79	0.47
3:C:128:PHE:O	3:C:129:ALA:C	2.52	0.47
3:C:178:LEU:O	3:C:179:ARG:HB2	2.13	0.47
4:D:10:ARG:HH12	4:D:40:PRO:CB	2.27	0.47
4:D:63:LYS:O	4:D:65:ARG:N	2.48	0.47
4:D:131:ARG:O	4:D:132:ARG:O	2.32	0.47
4:D:178:VAL:C	4:D:180:GLY:H	2.17	0.47
5:E:135:THR:O	5:E:137:GLU:N	2.48	0.47
6:F:14:LEU:HD11	6:F:84:ASN:OD1	2.14	0.47
6:F:18:GLN:HA	6:F:21:LEU:CD2	2.41	0.47
6:F:19:LEU:CD2	6:F:20:ALA:N	2.77	0.47
8:H:11:THR:CG2	8:H:14:ARG:HH12	2.21	0.47
8:H:14:ARG:NH1	8:H:14:ARG:HB3	2.28	0.47
8:H:37:ARG:O	8:H:41:ARG:CB	2.63	0.47
9:I:99:LEU:N	9:I:99:LEU:HD22	2.29	0.47
10:J:16:LEU:HD13	10:J:70:ARG:HG3	1.97	0.47
10:J:80:LYS:O	10:J:84:GLN:N	2.46	0.47
12:L:16:GLU:O	12:L:17:LYS:O	2.33	0.47
12:L:84:LEU:HD23	12:L:101:VAL:HG11	1.96	0.47
16:P:15:PRO:O	16:P:41:PRO:HD2	2.15	0.47
16:P:21:VAL:CG1	16:P:59:TRP:HE1	2.25	0.47
16:P:71:ARG:O	16:P:75:ARG:N	2.48	0.47
16:P:76:GLN:C	16:P:78:GLY:N	2.68	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:Q:20:THR:HA	17:Q:43:LEU:HD23	1.96	0.47
20:T:102:GLY:O	20:T:103:GLY:O	2.32	0.47
1:A:77:G:O2'	1:A:78:G:H5'	2.15	0.47
1:A:99:C:H2'	1:A:101:A:H8	1.72	0.47
1:A:131:C:C2'	1:A:132:C:C6	2.89	0.47
1:A:346:G:H2'	1:A:347:G:C5'	2.45	0.47
1:A:372:C:N3	1:A:387:U:C5	2.79	0.47
1:A:390:C:O3'	16:P:28:ARG:NH2	2.43	0.47
1:A:787:A:H2'	1:A:788:U:H6	1.80	0.47
1:A:936:C:H2'	1:A:937:A:O4'	2.14	0.47
1:A:1109:C:C2'	1:A:1110:A:H5'	2.45	0.47
1:A:1148:U:OP1	9:I:7:THR:HG21	2.14	0.47
1:A:1220:G:HO2'	19:S:52:TYR:HD2	1.63	0.47
1:A:1261:A:H2'	1:A:1262:C:H5'	1.95	0.47
1:A:1284:C:H3'	1:A:1285:A:H8	1.79	0.47
1:A:1290:G:H2'	1:A:1291:G:H8	1.79	0.47
1:A:1326:C:OP2	21:V:6:ARG:NH2	2.47	0.47
1:A:1355:G:H2'	1:A:1356:G:C8	2.44	0.47
1:A:1539:C:H2'	1:A:1540:U:C6	2.50	0.47
2:B:191:ASP:OD1	2:B:191:ASP:N	2.48	0.47
3:C:162:GLN:HG3	3:C:164:ARG:HD3	1.96	0.47
4:D:31:CYS:C	4:D:33:MET:H	2.16	0.47
6:F:7:ASN:O	6:F:8:ILE:HG13	2.14	0.47
6:F:53:ALA:O	6:F:55:ASP:N	2.47	0.47
7:G:24:THR:O	7:G:28:ASN:ND2	2.48	0.47
9:I:59:PHE:HE2	9:I:88:TYR:HH	1.61	0.47
9:I:79:LEU:HD22	9:I:83:ARG:HH21	1.80	0.47
11:K:127:LYS:HE3	11:K:127:LYS:CA	2.22	0.47
12:L:126:LYS:O	12:L:126:LYS:CD	2.63	0.47
14:N:56:VAL:HG22	14:N:57:ARG:N	2.17	0.47
15:O:11:VAL:O	15:O:14:GLU:HB3	2.14	0.47
16:P:67:THR:CG2	16:P:68:ASP:N	2.78	0.47
17:Q:66:SER:HB3	17:Q:69:LYS:HD3	1.96	0.47
17:Q:76:LEU:HD12	17:Q:78:GLU:H	1.79	0.47
1:A:38:G:N2	1:A:397:A:C5'	2.78	0.47
1:A:344:A:H5''	1:A:345:C:H5	1.76	0.47
1:A:345:C:H4'	1:A:346:G:C5'	2.45	0.47
1:A:443:C:H42	1:A:491:G:H1	1.61	0.47
1:A:673:G:H5''	6:F:87:ARG:CZ	2.45	0.47
1:A:992:U:H3	1:A:1044:A:H62	1.62	0.47
1:A:1061:G:O2'	1:A:1062:U:H5'	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1126:U:H2'	1:A:1127:G:O4'	2.15	0.47
1:A:1313:U:O5'	19:S:6:LYS:HG2	2.15	0.47
1:A:1374:A:N3	1:A:1375:A:C8	2.82	0.47
3:C:43:LEU:HD11	3:C:68:VAL:HG21	1.97	0.47
4:D:47:ARG:HH11	4:D:47:ARG:HG3	1.80	0.47
5:E:130:ASN:O	5:E:133:TYR:HB2	2.15	0.47
6:F:3:ARG:CA	6:F:66:GLU:HG3	2.44	0.47
6:F:51:PRO:HB3	6:F:56:PRO:HA	1.96	0.47
8:H:31:PHE:O	8:H:32:LYS:C	2.52	0.47
10:J:62:HIS:CB	14:N:59:ALA:HB3	2.40	0.47
11:K:44:SER:C	11:K:46:GLY:N	2.68	0.47
12:L:41:ARG:HH11	12:L:42:THR:N	2.13	0.47
12:L:88:GLY:O	12:L:89:ARG:HG2	2.14	0.47
17:Q:76:LEU:HD12	17:Q:78:GLU:N	2.29	0.47
1:A:59:A:H1'	1:A:354:G:C2	2.50	0.47
1:A:559:A:O2'	1:A:560:U:OP2	2.32	0.47
1:A:740:U:H1'	15:O:42:HIS:CE1	2.50	0.47
1:A:1013:G:H1'	1:A:1016:A:N6	2.29	0.47
1:A:1450:U:H2'	1:A:1452:C:C5	2.50	0.47
2:B:239:VAL:CG1	2:B:240:GLN:N	2.76	0.47
4:D:74:GLN:O	4:D:78:LEU:N	2.43	0.47
4:D:157:LEU:HD13	4:D:157:LEU:H	1.80	0.47
7:G:67:GLU:O	7:G:70:LYS:HB2	2.15	0.47
7:G:148:ASN:C	7:G:150:ALA:H	2.18	0.47
8:H:13:ILE:O	8:H:14:ARG:C	2.51	0.47
9:I:10:ARG:HD3	9:I:105:ASP:CG	2.36	0.47
9:I:93:ARG:HA	9:I:96:LEU:CD2	2.44	0.47
9:I:96:LEU:O	9:I:102:LEU:CD2	2.63	0.47
10:J:79:ARG:HH11	10:J:82:ILE:HD12	1.76	0.47
13:M:17:VAL:C	13:M:19:LEU:H	2.17	0.47
13:M:110:ARG:HG2	13:M:110:ARG:HH11	1.80	0.47
16:P:28:ARG:HG3	16:P:29:ASP:OD1	2.14	0.47
1:A:97:G:H2'	1:A:98:U:H5'	1.97	0.46
1:A:190(B):C:N3	1:A:190(H):G:C2	2.82	0.46
1:A:636:U:H5'	17:Q:2:PRO:HG2	1.97	0.46
1:A:687:A:HO2'	1:A:688:G:P	2.38	0.46
1:A:721:G:OP2	18:R:53:ARG:HG2	2.14	0.46
1:A:992:U:H1'	1:A:993:G:OP2	2.15	0.46
1:A:1057:G:H2'	1:A:1058:G:C8	2.50	0.46
1:A:1115:C:O4'	14:N:61:TRP:HA	2.15	0.46
1:A:1141:C:H2'	1:A:1142:G:O4'	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1150:U:H4'	10:J:41:PRO:CB	2.44	0.46
1:A:1451:A:O3'	1:A:1452:C:H6	1.99	0.46
1:A:1500:A:OP2	1:A:1505:G:OP2	2.32	0.46
2:B:14:GLY:O	2:B:15:VAL:HG22	2.14	0.46
2:B:20:GLU:HA	2:B:21:ARG:NH2	2.18	0.46
2:B:97:TRP:CZ3	2:B:176:GLU:OE2	2.69	0.46
3:C:117:ALA:O	3:C:121:ALA:HB2	2.14	0.46
3:C:138:VAL:O	3:C:142:MET:CB	2.60	0.46
4:D:124:GLY:CA	4:D:132:ARG:HE	2.27	0.46
5:E:26:PHE:O	5:E:27:ARG:CG	2.57	0.46
6:F:68:PRO:O	6:F:72:VAL:CG2	2.63	0.46
7:G:15:ASP:C	7:G:17:VAL:N	2.69	0.46
8:H:85:ARG:NE	8:H:87:SER:O	2.48	0.46
12:L:5:PRO:O	12:L:7:ILE:N	2.48	0.46
13:M:34:LEU:HD13	13:M:39:ILE:HB	1.96	0.46
14:N:36:PHE:O	14:N:36:PHE:CD1	2.68	0.46
16:P:45:THR:OG1	16:P:46:PRO:N	2.47	0.46
16:P:57:ARG:CZ	16:P:79:VAL:O	2.63	0.46
16:P:66:PRO:O	16:P:67:THR:O	2.32	0.46
18:R:47:THR:HG22	18:R:48:GLY:H	1.80	0.46
19:S:38:SER:HB3	19:S:71:LEU:CD1	2.38	0.46
19:S:40:ILE:CG2	19:S:62:ILE:HG12	2.45	0.46
20:T:78:ALA:O	20:T:81:LYS:HB2	2.15	0.46
1:A:9:G:OP1	5:E:122:GLU:HB2	2.15	0.46
1:A:70:G:H2'	1:A:73:C:O4'	2.16	0.46
1:A:190(J):U:H2'	1:A:190(K):G:C8	2.50	0.46
1:A:204:U:H4'	1:A:216:G:OP1	2.14	0.46
1:A:341:C:O2'	1:A:342:C:H5'	2.15	0.46
1:A:379:C:C2'	1:A:380:G:H5'	2.46	0.46
1:A:391:G:P	16:P:28:ARG:HH12	2.38	0.46
1:A:452:A:O2'	1:A:453:A:C8	2.66	0.46
1:A:807:A:C2	1:A:808:C:C2	3.03	0.46
1:A:852:G:O2'	1:A:853:G:H5'	2.16	0.46
1:A:1070:U:O2'	1:A:1071:C:H5'	2.15	0.46
1:A:1120:G:C2	1:A:1121:U:C4	3.03	0.46
1:A:1326:C:OP2	21:V:6:ARG:CZ	2.63	0.46
1:A:1326:C:N4	1:A:1327:C:H41	2.13	0.46
1:A:1343:G:C6	1:A:1344:C:N4	2.84	0.46
1:A:1352:C:H2'	1:A:1353:G:C8	2.51	0.46
1:A:1522:U:C2'	1:A:1523:G:H5'	2.44	0.46
2:B:15:VAL:HB	2:B:16:HIS:H	1.55	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:239:VAL:O	2:B:240:GLN:C	2.52	0.46
3:C:51:GLY:CA	3:C:70:VAL:HG13	2.45	0.46
3:C:124:ILE:HD11	3:C:153:VAL:HG11	1.98	0.46
4:D:110:PHE:N	4:D:110:PHE:HD1	2.14	0.46
4:D:135:LEU:O	4:D:138:TYR:HB2	2.15	0.46
6:F:62:TRP:CE2	18:R:35:ARG:NH2	2.83	0.46
8:H:5:PRO:O	8:H:8:ASP:HB3	2.14	0.46
8:H:31:PHE:HE1	8:H:35:ILE:HG13	1.80	0.46
10:J:84:GLN:O	10:J:85:LEU:HB2	2.14	0.46
12:L:45:PRO:HD3	12:L:51:ALA:O	2.14	0.46
12:L:53:ARG:HH12	12:L:92:ASP:CB	2.28	0.46
13:M:3:ARG:CD	13:M:7:VAL:HA	2.45	0.46
13:M:14:ARG:HG3	13:M:44:ARG:CZ	2.45	0.46
14:N:23:ARG:HH12	14:N:30:ALA:HB2	1.80	0.46
15:O:8:LYS:O	15:O:10:LYS:N	2.49	0.46
18:R:18:ARG:O	18:R:19:LYS:O	2.33	0.46
20:T:13:LEU:HD12	20:T:14:LYS:CA	2.44	0.46
20:T:13:LEU:CG	20:T:14:LYS:N	2.79	0.46
21:V:10:ARG:CA	21:V:13:ILE:HG22	2.38	0.46
21:V:17:THR:O	21:V:22:ARG:HG2	2.15	0.46
1:A:132:C:O2'	1:A:133:U:H5'	2.15	0.46
1:A:142:G:H1'	1:A:196:A:H2	1.80	0.46
1:A:357:G:HO2'	1:A:358:U:H5'	1.79	0.46
1:A:389:A:C6	1:A:390:C:H1'	2.51	0.46
1:A:768:A:H2'	1:A:769:G:H5'	1.97	0.46
1:A:927:G:C5	1:A:928:G:N7	2.83	0.46
1:A:976:G:C8	1:A:1358:U:C2	3.03	0.46
1:A:1004:A:H8	1:A:1037:C:HO2'	1.63	0.46
1:A:1057:G:O2'	1:A:1058:G:H5'	2.15	0.46
1:A:1061:G:C2'	1:A:1062:U:H5'	2.45	0.46
1:A:1125:U:O4	10:J:5:ARG:CD	2.63	0.46
1:A:1201:A:HO2'	1:A:1202:G:P	2.36	0.46
1:A:1345:U:C2	1:A:1377:A:C2	3.02	0.46
1:A:1390:U:H2'	1:A:1391:U:C6	2.49	0.46
2:B:21:ARG:O	2:B:39:ILE:HA	2.16	0.46
2:B:44:LEU:C	2:B:46:LYS:N	2.66	0.46
2:B:51:LEU:O	2:B:52:GLU:C	2.54	0.46
2:B:111:ARG:CB	2:B:149:LEU:HD11	2.41	0.46
2:B:116:GLU:C	2:B:118:LEU:N	2.68	0.46
4:D:205:GLU:C	4:D:207:TYR:H	2.19	0.46
5:E:78:HIS:CG	5:E:78:HIS:O	2.68	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:115:VAL:CG1	5:E:116:THR:N	2.79	0.46
8:H:9:MET:O	8:H:10:LEU:C	2.53	0.46
9:I:16:ARG:HH21	9:I:64:THR:HG22	1.80	0.46
16:P:1:MET:HE1	16:P:65:GLN:HB2	1.97	0.46
20:T:56:MET:HE2	20:T:88:VAL:HG11	1.97	0.46
1:A:44:G:OP2	16:P:12:LYS:HD2	2.15	0.46
1:A:144:G:N2	1:A:178:C:N3	2.57	0.46
1:A:167:G:HO2'	1:A:168:G:H8	1.62	0.46
1:A:245:C:C6	1:A:284:G:N2	2.84	0.46
1:A:318:G:N2	1:A:319:G:C4	2.83	0.46
1:A:401:C:P	4:D:73:ARG:HH21	2.38	0.46
1:A:676:A:H2'	1:A:677:U:H6	1.79	0.46
1:A:866:C:H2'	1:A:867:G:O4'	2.14	0.46
1:A:1073:U:H3	1:A:1102:A:H61	1.63	0.46
1:A:1089:G:H2'	1:A:1090:U:H5'	1.98	0.46
1:A:1189:C:C5'	3:C:5:ILE:HD13	2.45	0.46
1:A:1286:A:H8	1:A:1287:A:H4'	1.81	0.46
2:B:92:TYR:C	2:B:92:TYR:HD1	2.19	0.46
2:B:109:SER:C	2:B:111:ARG:N	2.68	0.46
3:C:126:ARG:HD2	3:C:128:PHE:CD2	2.51	0.46
4:D:61:LYS:HD3	4:D:62:GLN:CA	2.45	0.46
4:D:148:VAL:CG1	4:D:149:ALA:N	2.77	0.46
5:E:48:ALA:HB3	5:E:54:ALA:HA	1.97	0.46
6:F:85:VAL:O	6:F:85:VAL:HG12	2.14	0.46
10:J:65:LEU:C	10:J:65:LEU:CD2	2.83	0.46
10:J:82:ILE:O	10:J:86:MET:SD	2.73	0.46
10:J:86:MET:CG	10:J:87:THR:H	1.97	0.46
12:L:98:TYR:CD1	12:L:98:TYR:N	2.84	0.46
17:Q:47:PRO:C	17:Q:49:GLU:H	2.18	0.46
18:R:24:ALA:O	18:R:26:LEU:N	2.49	0.46
19:S:44:MET:O	19:S:45:VAL:C	2.54	0.46
20:T:33:ILE:CD1	20:T:63:ILE:HA	2.24	0.46
20:T:51:GLU:HA	20:T:54:LYS:HB3	1.97	0.46
1:A:112:G:C5'	1:A:389:A:H4'	2.45	0.46
1:A:370:C:H2'	1:A:371:G:H8	1.81	0.46
1:A:374:A:C4	1:A:375:U:C5	3.04	0.46
1:A:402:G:H1'	1:A:620:C:H42	1.80	0.46
1:A:463:A:H4'	16:P:82:GLN:NE2	2.31	0.46
1:A:501:C:O2'	1:A:502:G:H5'	2.16	0.46
1:A:754:C:H1'	15:O:69:TYR:CG	2.51	0.46
1:A:927:G:C4	1:A:928:G:C8	3.04	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:939:G:H4'	7:G:102:ARG:NH2	2.30	0.46
1:A:955:U:H2'	1:A:956:U:H6	1.81	0.46
1:A:1158:C:H5''	2:B:133:LYS:CE	2.44	0.46
1:A:1250:A:C4'	9:I:68:GLY:HA2	2.21	0.46
1:A:1279:A:O2'	1:A:1282:C:N4	2.48	0.46
1:A:1320:C:O4'	19:S:73:GLU:HG2	2.15	0.46
1:A:1334:G:H8	1:A:1334:G:O5'	1.99	0.46
3:C:26:LYS:C	3:C:28:GLN:H	2.18	0.46
3:C:77:ILE:CD1	3:C:84:ILE:HD12	2.45	0.46
3:C:111:LEU:HD21	3:C:146:ALA:N	2.31	0.46
3:C:179:ARG:HD2	3:C:180:ALA:CA	2.45	0.46
4:D:99:SER:O	4:D:140:VAL:HG23	2.15	0.46
4:D:149:ALA:O	4:D:152:SER:N	2.42	0.46
4:D:152:SER:HA	4:D:155:LEU:CD1	2.45	0.46
4:D:173:TRP:HB2	4:D:187:ARG:O	2.16	0.46
5:E:10:MET:HA	5:E:31:LEU:O	2.15	0.46
5:E:93:PRO:HG2	8:H:105:ARG:NH2	2.29	0.46
7:G:43:PHE:HD2	7:G:44:TYR:HD2	1.62	0.46
9:I:97:LYS:CB	9:I:98:PRO:HD3	2.45	0.46
10:J:9:ARG:O	10:J:95:GLU:OE1	2.34	0.46
13:M:19:LEU:O	13:M:21:TYR:N	2.48	0.46
14:N:21:TYR:H	14:N:21:TYR:HD1	1.54	0.46
15:O:25:THR:HG21	15:O:70:LEU:HD23	1.98	0.46
15:O:35:ARG:HB3	15:O:59:MET:HE1	1.98	0.46
15:O:74:ASP:OD2	15:O:77:ARG:HG3	2.15	0.46
16:P:49:LEU:CD1	16:P:73:LEU:HD22	2.45	0.46
18:R:30:ASP:O	18:R:32:ARG:N	2.49	0.46
18:R:76:LEU:O	18:R:77:GLY:C	2.51	0.46
19:S:23:ASN:CA	19:S:27:GLU:HA	2.43	0.46
20:T:48:LYS:O	20:T:52:ALA:HB3	2.16	0.46
20:T:65:LYS:O	20:T:68:LYS:HB2	2.14	0.46
1:A:248:C:H2'	1:A:249:U:C5'	2.46	0.46
1:A:319:G:O2'	1:A:320:C:H5'	2.16	0.46
1:A:428:G:O3'	4:D:36:ARG:NH2	2.47	0.46
1:A:541:G:O2'	1:A:542:G:H5'	2.16	0.46
1:A:655:A:H2'	1:A:656:C:O4'	2.16	0.46
1:A:793:U:C3'	1:A:794:A:C5'	2.74	0.46
1:A:860:A:N6	1:A:861:G:C2	2.83	0.46
1:A:1076:C:C2	1:A:1077:G:C8	3.04	0.46
1:A:1262:C:H42	1:A:1273:G:H1	1.62	0.46
1:A:1340:A:O2'	1:A:1341:U:H5'	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1512:U:O2'	1:A:1513:A:H5'	2.16	0.46
2:B:126:GLU:OE2	2:B:129:GLU:CB	2.64	0.46
4:D:104:VAL:O	4:D:105:VAL:C	2.52	0.46
5:E:19:MET:SD	5:E:24:ARG:HA	2.56	0.46
5:E:50:GLU:O	5:E:51:VAL:C	2.54	0.46
6:F:99:ALA:O	6:F:100:ASN:C	2.54	0.46
9:I:50:LEU:HD12	9:I:50:LEU:N	2.28	0.46
10:J:64:GLU:HG3	14:N:59:ALA:CA	2.46	0.46
12:L:67:THR:HG1	12:L:96:VAL:HA	1.81	0.46
15:O:87:ILE:HD12	15:O:87:ILE:N	2.31	0.46
17:Q:75:ARG:HG2	17:Q:76:LEU:N	2.31	0.46
1:A:7:G:C6	1:A:298:A:C2	3.04	0.46
1:A:131:C:N3	1:A:232:G:C6	2.84	0.46
1:A:308:C:H2'	1:A:309:G:H8	1.81	0.46
1:A:382:A:C2	1:A:383:A:C4	3.03	0.46
1:A:439:A:H3'	1:A:440:A:H8	1.80	0.46
1:A:451:A:C2	1:A:480:U:C5	3.03	0.46
1:A:575:G:C4	1:A:881:G:N2	2.83	0.46
1:A:594:G:N1	1:A:596:C:N4	2.64	0.46
1:A:760:G:H21	17:Q:103:GLY:HA3	1.81	0.46
1:A:1432:G:O2'	1:A:1468:A:N6	2.49	0.46
1:A:1518:A:H3'	1:A:1518:A:H8	1.81	0.46
3:C:16:ARG:O	3:C:17:ASP:C	2.54	0.46
4:D:74:GLN:O	4:D:78:LEU:HB2	2.16	0.46
7:G:153:HIS:HD2	7:G:154:TYR:CZ	2.33	0.46
9:I:19:LEU:HD21	9:I:59:PHE:O	2.15	0.46
9:I:19:LEU:HD12	9:I:61:ALA:HB2	1.97	0.46
10:J:68:HIS:CD2	10:J:68:HIS:N	2.82	0.46
11:K:98:LEU:HD23	11:K:98:LEU:H	1.80	0.46
12:L:28:LYS:C	12:L:30:ALA:N	2.68	0.46
12:L:42:THR:CG2	12:L:52:LEU:HD22	2.45	0.46
13:M:26:GLY:O	13:M:28:ALA:N	2.49	0.46
14:N:41:ARG:HG3	14:N:42:ILE:N	2.31	0.46
16:P:6:LEU:N	16:P:6:LEU:CD1	2.78	0.46
16:P:14:ASN:OD1	16:P:14:ASN:O	2.34	0.46
16:P:71:ARG:O	16:P:75:ARG:HG3	2.15	0.46
19:S:52:TYR:CD1	19:S:56:GLN:O	2.68	0.46
19:S:63:THR:O	19:S:67:VAL:HG23	2.16	0.46
21:V:18:TYR:HE2	21:V:23:PRO:O	1.97	0.46
1:A:390:C:O2'	16:P:28:ARG:NH2	2.49	0.46
1:A:696:A:H2'	1:A:697:U:H6	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:718:G:C4'	11:K:117:ASN:ND2	2.79	0.46
1:A:760:G:H22	17:Q:104:LYS:H	1.55	0.46
1:A:799:G:H2'	1:A:800:G:H5'	1.98	0.46
1:A:1077:G:N2	1:A:1079:G:H3'	2.30	0.46
1:A:1120:G:H1	1:A:1153:C:H42	1.64	0.46
1:A:1437:C:H6	1:A:1438:G:N7	2.14	0.46
3:C:139:GLN:HA	3:C:139:GLN:NE2	2.31	0.46
3:C:170:GLN:O	3:C:172:ARG:N	2.49	0.46
3:C:180:ALA:HB2	3:C:206:GLU:HG3	1.97	0.46
3:C:191:THR:HG23	3:C:194:GLY:C	2.37	0.46
4:D:19:LEU:HB3	4:D:21:LEU:CD1	2.39	0.46
4:D:46:LYS:HG2	4:D:47:ARG:N	2.30	0.46
5:E:12:LEU:HD13	5:E:31:LEU:HB2	1.98	0.46
5:E:45:PHE:CD2	5:E:47:LYS:HE3	2.50	0.46
5:E:137:GLU:CD	5:E:141:GLN:HE21	2.18	0.46
7:G:29:LYS:HG3	7:G:101:LEU:HD12	1.98	0.46
7:G:71:PRO:CD	7:G:103:TRP:HZ3	2.21	0.46
10:J:48:THR:HG1	10:J:62:HIS:CE1	2.29	0.46
11:K:19:ALA:HB2	11:K:80:VAL:HG13	1.97	0.46
13:M:20:THR:HA	13:M:25:ILE:O	2.16	0.46
14:N:24:CYS:HA	14:N:33:VAL:HG13	1.98	0.46
18:R:33:ASP:OD2	18:R:36:ASN:CB	2.64	0.46
18:R:85:LEU:HD12	18:R:85:LEU:C	2.36	0.46
19:S:67:VAL:C	19:S:69:HIS:N	2.69	0.46
1:A:507:C:H3'	1:A:508:C:H2'	1.97	0.46
1:A:562:C:C4'	1:A:563:A:H5'	2.45	0.46
1:A:676:A:C2	1:A:677:U:C4	3.04	0.46
1:A:705:U:C5	1:A:706:A:C5	3.04	0.46
1:A:1186:G:C2	1:A:1187:G:C4	3.04	0.46
2:B:10:LEU:C	2:B:12:GLU:H	2.16	0.46
3:C:11:ARG:HA	3:C:14:ILE:HG12	1.95	0.46
4:D:59:ARG:O	4:D:62:GLN:HB2	2.16	0.46
4:D:170:VAL:HG13	4:D:174:LEU:HB2	1.98	0.46
7:G:75:VAL:CG1	7:G:86:GLN:HB3	2.45	0.46
9:I:93:ARG:HA	9:I:96:LEU:HD23	1.96	0.46
10:J:49:VAL:O	10:J:60:ARG:HA	2.16	0.46
12:L:78:GLN:C	12:L:80:HIS:N	2.69	0.46
13:M:43:THR:O	13:M:44:ARG:O	2.33	0.46
14:N:14:PRO:HB2	14:N:19:ARG:HB2	1.97	0.46
15:O:9:GLN:HA	15:O:9:GLN:OE1	2.15	0.46
15:O:33:THR:HG23	15:O:63:ARG:HH11	1.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:O:48:LYS:NZ	15:O:48:LYS:CB	2.78	0.46
16:P:10:GLY:HA3	16:P:16:HIS:N	2.31	0.46
19:S:53:ASN:N	19:S:56:GLN:O	2.49	0.46
21:V:17:THR:CG2	21:V:18:TYR:H	2.19	0.46
1:A:373:A:H2'	1:A:374:A:C8	2.31	0.46
1:A:670:G:O2'	1:A:671:G:H5'	2.16	0.46
1:A:799:G:O2'	1:A:800:G:H5'	2.16	0.46
1:A:977:A:C8	1:A:1223:C:C2	3.03	0.46
1:A:1006:C:H2'	1:A:1007:C:O4'	2.15	0.46
1:A:1225:A:H5'	13:M:103:THR:CB	2.46	0.46
1:A:1242:C:C4	1:A:1243:C:C5	3.04	0.46
1:A:1298:C:C4'	1:A:1299:A:C5'	2.92	0.46
1:A:1326:C:OP1	21:V:12:LYS:CD	2.62	0.46
1:A:1349:A:H2'	1:A:1350:A:C5'	2.45	0.46
1:A:1470:G:H2'	1:A:1471:G:C8	2.51	0.46
6:F:54:LYS:O	6:F:56:PRO:HD3	2.16	0.46
6:F:74:ASP:O	6:F:77:ARG:HG2	2.16	0.46
7:G:47:CYS:C	7:G:49:ILE:N	2.69	0.46
8:H:31:PHE:CE1	8:H:35:ILE:HG13	2.51	0.46
8:H:119:LEU:CD1	8:H:124:ALA:HA	2.46	0.46
10:J:26:ALA:HB3	10:J:85:LEU:CD1	2.45	0.46
10:J:31:GLY:O	10:J:76:ASN:HB2	2.16	0.46
20:T:78:ALA:O	20:T:79:ARG:C	2.53	0.46
1:A:266:G:O2'	1:A:267:C:OP2	2.26	0.45
1:A:309:G:H1'	1:A:608:A:C2	2.51	0.45
1:A:321:A:N3	1:A:322:C:C5	2.84	0.45
1:A:355:C:C2	1:A:356:A:C8	3.04	0.45
1:A:451:A:H1'	1:A:452:A:C8	2.51	0.45
1:A:568:G:O6	12:L:5:PRO:HD3	2.16	0.45
1:A:585:G:C6	1:A:586:C:C4	3.04	0.45
1:A:940:C:O2'	1:A:941:G:H5'	2.17	0.45
1:A:979:C:OP2	1:A:980:C:H5	1.98	0.45
1:A:1049:U:H4'	1:A:1050:G:O5'	2.17	0.45
1:A:1095:U:P	1:A:1108:G:H1	2.38	0.45
1:A:1228:C:OP1	13:M:115:LYS:HD3	2.15	0.45
1:A:1275:A:H2'	1:A:1276:G:H8	1.81	0.45
1:A:1516:G:N2	1:A:1520:G:C4	2.84	0.45
2:B:111:ARG:HD3	2:B:111:ARG:HA	1.60	0.45
2:B:135:GLN:C	2:B:139:LYS:HB2	2.37	0.45
3:C:24:ALA:HB1	3:C:28:GLN:NE2	2.31	0.45
3:C:38:ARG:HD3	3:C:38:ARG:HA	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:61:ALA:O	3:C:62:ASP:HB2	2.16	0.45
3:C:76:VAL:O	3:C:78:GLY:N	2.49	0.45
7:G:114:ARG:HD3	7:G:114:ARG:N	2.22	0.45
8:H:66:GLY:O	8:H:76:PRO:HB3	2.16	0.45
9:I:15:ALA:HB1	9:I:77:ILE:HG12	1.98	0.45
9:I:18:PHE:CD1	9:I:62:TYR:HD2	2.35	0.45
9:I:63:ILE:HD13	9:I:77:ILE:CG2	2.46	0.45
9:I:80:GLY:C	9:I:83:ARG:H	2.19	0.45
9:I:114:TYR:HE2	10:J:61:GLU:H	1.64	0.45
10:J:57:LYS:HG2	10:J:60:ARG:NE	2.30	0.45
10:J:82:ILE:O	10:J:86:MET:HA	2.17	0.45
10:J:90:LEU:C	10:J:90:LEU:HD23	2.36	0.45
11:K:41:THR:O	11:K:42:TRP:HB3	2.15	0.45
11:K:46:GLY:O	11:K:48:ILE:N	2.49	0.45
12:L:49:ASN:HD22	12:L:49:ASN:HA	1.54	0.45
12:L:58:VAL:HG12	12:L:59:ARG:N	2.31	0.45
18:R:33:ASP:OD2	18:R:36:ASN:N	2.47	0.45
1:A:75:G:N2	1:A:96:G:N2	2.64	0.45
1:A:153:C:N4	1:A:169:C:H42	2.12	0.45
1:A:245:C:C2'	1:A:246:A:H5'	2.46	0.45
1:A:427:U:C4'	1:A:541:G:H5''	2.47	0.45
1:A:507:C:OP2	1:A:508:C:O2'	2.32	0.45
1:A:667:G:N2	1:A:668:G:C4	2.84	0.45
1:A:677:U:H2'	1:A:678:U:O4'	2.16	0.45
1:A:690:G:C6	1:A:691:G:C2	3.04	0.45
1:A:861:G:C4	1:A:862:C:C5	3.04	0.45
1:A:867:G:N2	1:A:868:C:C2	2.85	0.45
1:A:950:U:H2'	1:A:951:G:C8	2.47	0.45
1:A:1112:C:O2	3:C:179:ARG:HB3	2.15	0.45
1:A:1350:A:C6	1:A:1351:U:C4	3.04	0.45
1:A:1380:U:C5	7:G:3:ARG:HA	2.50	0.45
1:A:1486:G:H2'	1:A:1487:G:C8	2.51	0.45
2:B:118:LEU:O	2:B:119:GLU:C	2.55	0.45
3:C:21:ARG:N	3:C:21:ARG:CD	2.79	0.45
3:C:27:LYS:HA	3:C:30:ARG:NH2	2.32	0.45
4:D:127:THR:HA	4:D:132:ARG:HA	1.99	0.45
6:F:33:TYR:CG	6:F:75:LEU:HD23	2.51	0.45
7:G:14:PRO:CA	7:G:21:VAL:HG12	2.45	0.45
9:I:23:ASN:O	9:I:56:LEU:O	2.35	0.45
10:J:4:ILE:HD12	10:J:74:ILE:O	2.16	0.45
10:J:6:ILE:CD1	10:J:73:ASP:HA	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:20:ALA:O	10:J:23:ILE:N	2.32	0.45
11:K:18:ARG:HD3	11:K:33:THR:HG23	1.98	0.45
11:K:87:THR:C	11:K:91:ARG:HH12	2.18	0.45
13:M:41:PRO:O	13:M:42:ALA:O	2.33	0.45
13:M:92:HIS:CE1	13:M:98:VAL:HG21	2.50	0.45
14:N:26:ARG:CZ	14:N:47:LEU:HD21	2.46	0.45
15:O:36:ILE:CD1	15:O:63:ARG:HD3	2.46	0.45
17:Q:4:LYS:O	17:Q:60:ILE:HA	2.16	0.45
17:Q:87:LYS:O	17:Q:88:TYR:C	2.55	0.45
20:T:45:GLN:C	20:T:45:GLN:CD	2.74	0.45
1:A:142:G:N3	1:A:196:A:C2	2.85	0.45
1:A:356:A:H2'	1:A:357:G:C8	2.48	0.45
1:A:507:C:C4	1:A:508:C:H5	2.35	0.45
1:A:541:G:C4	1:A:542:G:C8	3.04	0.45
1:A:1213:A:C6	1:A:1215:G:C8	3.04	0.45
1:A:1250:A:H61	1:A:1354:C:C1'	2.29	0.45
1:A:1340:A:H2'	1:A:1341:U:O4'	2.16	0.45
1:A:1361:G:H8	1:A:1361:G:P	2.39	0.45
2:B:74:LYS:HD3	2:B:206:ASP:HB2	1.98	0.45
2:B:142:LEU:C	2:B:144:ARG:N	2.68	0.45
2:B:231:GLU:CB	2:B:232:PRO:CD	2.87	0.45
3:C:76:VAL:CG2	3:C:77:ILE:N	2.79	0.45
3:C:153:VAL:HG13	3:C:198:VAL:CG2	2.46	0.45
3:C:162:GLN:HG3	3:C:164:ARG:HG3	1.97	0.45
3:C:206:GLU:O	3:C:207:VAL:HB	2.16	0.45
4:D:100:ARG:HG2	4:D:100:ARG:HH11	1.82	0.45
7:G:11:GLN:HA	7:G:11:GLN:HE21	1.82	0.45
8:H:11:THR:O	8:H:13:ILE:N	2.49	0.45
8:H:11:THR:HA	8:H:14:ARG:HH12	1.78	0.45
9:I:85:LEU:HD22	9:I:92:TYR:CE1	2.51	0.45
10:J:16:LEU:HD21	10:J:70:ARG:HB2	1.97	0.45
11:K:46:GLY:O	11:K:49:GLY:N	2.49	0.45
13:M:81:LEU:CA	13:M:84:ILE:HG12	2.41	0.45
15:O:17:ARG:CZ	15:O:77:ARG:NH1	2.80	0.45
15:O:74:ASP:O	15:O:76:GLU:N	2.50	0.45
17:Q:85:VAL:O	17:Q:89:LEU:N	2.48	0.45
18:R:38:GLU:O	18:R:39:VAL:C	2.54	0.45
19:S:41:VAL:HA	19:S:42:PRO:HD3	1.83	0.45
20:T:79:ARG:CG	20:T:83:ARG:HH11	2.29	0.45
1:A:38:G:H22	1:A:397:A:C5'	2.29	0.45
1:A:98:U:H2'	1:A:99:C:C6	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:G:C2	1:A:168:G:C5	3.04	0.45
1:A:192:U:H4'	20:T:57:ARG:NH1	2.31	0.45
1:A:223:U:C4	1:A:224:C:C5	3.05	0.45
1:A:319:G:N2	1:A:320:C:C2	2.85	0.45
1:A:456:C:C6	1:A:457:C:H5	2.34	0.45
1:A:580:U:H1'	15:O:57:LEU:CD2	2.46	0.45
1:A:665:A:N3	1:A:732:C:C2	2.84	0.45
1:A:811:C:H4'	1:A:900:A:N6	2.32	0.45
1:A:944:G:C2'	1:A:945:G:H5''	2.46	0.45
1:A:955:U:O2'	1:A:956:U:H5'	2.16	0.45
1:A:1313:U:H3'	19:S:6:LYS:CD	2.45	0.45
2:B:109:SER:O	2:B:111:ARG:N	2.49	0.45
2:B:158:LEU:CD2	2:B:159:PRO:HD2	2.46	0.45
4:D:10:ARG:CG	4:D:10:ARG:NH1	2.70	0.45
4:D:52:SER:OG	4:D:55:ALA:N	2.36	0.45
5:E:141:GLN:O	5:E:142:LEU:C	2.54	0.45
7:G:125:MET:O	7:G:128:ALA:N	2.50	0.45
9:I:75:ASP:C	9:I:78:LYS:HB3	2.36	0.45
9:I:97:LYS:O	9:I:100:GLY:N	2.36	0.45
10:J:3:LYS:HB2	10:J:77:PRO:HD3	1.98	0.45
10:J:22:LYS:NZ	10:J:23:ILE:HD11	2.31	0.45
11:K:59:TYR:O	11:K:63:LEU:HG	2.17	0.45
12:L:55:VAL:HG12	12:L:56:ALA:N	2.24	0.45
12:L:84:LEU:HG	12:L:101:VAL:HG11	1.98	0.45
13:M:2:ALA:O	13:M:10:PRO:HD2	2.17	0.45
14:N:28:GLY:O	14:N:29:ARG:O	2.34	0.45
14:N:43:CYS:O	14:N:46:GLU:OE1	2.35	0.45
18:R:29:PHE:C	18:R:29:PHE:CD1	2.90	0.45
1:A:44:G:C2	1:A:399:G:N1	2.85	0.45
1:A:191:G:O2'	20:T:102:GLY:HA2	2.16	0.45
1:A:411:A:C8	1:A:413:G:C8	3.04	0.45
1:A:773:G:C5	1:A:774:G:N7	2.84	0.45
1:A:778:G:C5	1:A:779:C:C5	3.04	0.45
1:A:913:A:C1'	1:A:914:A:O4'	2.63	0.45
1:A:1114:C:C4	1:A:1115:C:C5	3.05	0.45
1:A:1342:C:O2'	1:A:1343:G:H5'	2.17	0.45
1:A:1411:C:H2'	1:A:1412:C:O4'	2.17	0.45
2:B:142:LEU:HD23	2:B:146:GLN:CG	2.47	0.45
2:B:178:ARG:HG3	2:B:178:ARG:NH1	2.29	0.45
3:C:14:ILE:O	3:C:15:THR:HG23	2.17	0.45
3:C:79:ARG:HH21	3:C:82:GLU:HG3	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:201:TYR:N	3:C:201:TYR:CD1	2.85	0.45
4:D:4:TYR:HE2	4:D:6:GLY:O	1.99	0.45
4:D:10:ARG:O	4:D:11:LEU:C	2.53	0.45
4:D:150:GLU:OE1	4:D:150:GLU:HA	2.16	0.45
5:E:15:ARG:CZ	5:E:26:PHE:HE2	2.29	0.45
5:E:73:ASN:O	5:E:73:ASN:ND2	2.48	0.45
6:F:1:MET:HE1	6:F:36:ARG:HH21	1.81	0.45
6:F:14:LEU:HD22	6:F:18:GLN:NE2	2.31	0.45
13:M:44:ARG:N	13:M:44:ARG:HD2	2.32	0.45
14:N:8:GLU:O	14:N:11:LYS:CE	2.64	0.45
15:O:78:TYR:CZ	15:O:82:ILE:HD12	2.52	0.45
17:Q:6:LEU:N	17:Q:59:ILE:O	2.50	0.45
17:Q:67:LYS:CA	17:Q:70:ARG:HH12	2.29	0.45
1:A:68:G:O4'	1:A:171:A:H1'	2.17	0.45
1:A:281:G:HO2'	1:A:282:A:P	2.39	0.45
1:A:327:A:C6	1:A:329:A:C5	3.04	0.45
1:A:621:A:C4	1:A:622:A:N7	2.85	0.45
1:A:636:U:H5'	17:Q:2:PRO:CG	2.46	0.45
1:A:639:G:N2	1:A:640:A:C4	2.85	0.45
1:A:657:G:H2'	1:A:658:G:H8	1.82	0.45
1:A:659:U:H3	1:A:746:A:N6	2.10	0.45
1:A:686:U:O2	1:A:687:A:C4	2.70	0.45
1:A:849:C:H2'	1:A:850:U:C5'	2.46	0.45
1:A:864:A:C6	1:A:865:A:N1	2.85	0.45
1:A:880:C:O2'	1:A:881:G:H5'	2.17	0.45
1:A:977:A:C8	1:A:1223:C:N3	2.85	0.45
1:A:1014:A:N6	1:A:1015:A:N6	2.64	0.45
1:A:1128:C:H2'	1:A:1139:G:N7	2.31	0.45
1:A:1220:G:H1'	19:S:52:TYR:HD2	1.81	0.45
1:A:1249:C:H3'	1:A:1249:C:C6	2.51	0.45
3:C:76:VAL:CG2	3:C:77:ILE:H	2.17	0.45
4:D:191:ARG:O	4:D:192:GLU:C	2.54	0.45
6:F:25:ILE:HG23	6:F:26:ILE:N	2.32	0.45
7:G:123:GLU:O	7:G:124:LEU:C	2.53	0.45
7:G:136:LYS:O	7:G:140:ASP:HB2	2.17	0.45
8:H:60:ARG:HB3	8:H:62:TYR:HE2	1.81	0.45
9:I:49:PRO:HB2	9:I:82:ALA:HA	1.98	0.45
9:I:50:LEU:HD11	9:I:81:ILE:CG2	2.46	0.45
10:J:9:ARG:CZ	10:J:9:ARG:CB	2.94	0.45
11:K:91:ARG:HG2	11:K:95:ILE:HD11	1.99	0.45
15:O:13:GLN:HE21	15:O:13:GLN:HB3	1.48	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:Q:10:VAL:HG21	17:Q:55:ASP:HB2	1.98	0.45
17:Q:81:ARG:C	17:Q:83:ASP:H	2.19	0.45
19:S:22:LEU:HD22	19:S:28:LYS:CB	2.46	0.45
19:S:36:ARG:NH2	19:S:75:ALA:HB3	2.32	0.45
19:S:49:ILE:HG12	19:S:51:VAL:HG13	1.98	0.45
19:S:68:GLY:O	19:S:69:HIS:ND1	2.49	0.45
1:A:30:U:O2'	1:A:31:G:OP1	2.27	0.45
1:A:96:G:C2	1:A:97:G:C8	3.05	0.45
1:A:444:C:N3	1:A:491:G:C2	2.85	0.45
1:A:559:A:H4'	1:A:560:U:C5'	2.46	0.45
1:A:806:C:H2'	1:A:807:A:H8	1.81	0.45
1:A:865:A:N6	1:A:866:C:H42	2.14	0.45
1:A:901:A:C5	1:A:902:G:H1'	2.52	0.45
1:A:967:C:O2'	9:I:128:ARG:NH2	2.50	0.45
1:A:1371:G:O3'	9:I:69:GLY:HA3	2.17	0.45
2:B:36:ARG:O	2:B:37:ASN:O	2.35	0.45
2:B:71:VAL:O	2:B:165:VAL:HG22	2.15	0.45
3:C:12:LEU:C	3:C:14:ILE:N	2.70	0.45
4:D:195:ALA:O	4:D:196:LEU:C	2.54	0.45
5:E:39:GLY:CA	5:E:71:LEU:HD11	2.45	0.45
8:H:18:ARG:HH11	8:H:18:ARG:CB	2.28	0.45
8:H:87:SER:CB	8:H:93:VAL:HB	2.24	0.45
8:H:100:ILE:CG2	8:H:101:PRO:N	2.79	0.45
9:I:25:LYS:HD3	9:I:25:LYS:H	1.80	0.45
9:I:88:TYR:O	9:I:90:PRO:HD2	2.16	0.45
10:J:16:LEU:HD22	10:J:70:ARG:HD3	1.99	0.45
10:J:24:VAL:HG12	10:J:25:GLU:N	2.31	0.45
14:N:48:ALA:HA	14:N:53:LEU:HB2	1.98	0.45
18:R:61:LYS:O	18:R:62:GLU:C	2.55	0.45
1:A:178:C:O2'	1:A:179:A:C5'	2.62	0.45
1:A:281:G:C2'	1:A:282:A:OP2	2.64	0.45
1:A:478:A:C2'	1:A:479:C:H5'	2.47	0.45
1:A:529:G:C4'	1:A:533:A:C2	3.00	0.45
1:A:538:G:H5''	12:L:114:LYS:CB	2.41	0.45
1:A:599:C:O2'	1:A:600:C:H5'	2.16	0.45
1:A:639:G:H2'	1:A:640:A:H8	1.82	0.45
1:A:644:G:H2'	1:A:645:C:O4'	2.17	0.45
1:A:839:U:O2	1:A:839:U:C2'	2.64	0.45
1:A:1081:G:OP1	5:E:17:ALA:O	2.35	0.45
1:A:1219:U:H2'	1:A:1220:G:C8	2.52	0.45
1:A:1329:A:OP1	13:M:28:ALA:HB3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1356:G:H2'	1:A:1357:A:H8	1.76	0.45
1:A:1492:A:HO2'	1:A:1493:A:H8	1.60	0.45
2:B:43:ASP:C	2:B:43:ASP:OD1	2.55	0.45
2:B:177:ALA:O	2:B:178:ARG:C	2.55	0.45
2:B:182:ILE:O	2:B:183:PRO:C	2.55	0.45
5:E:132:ALA:O	5:E:135:THR:HB	2.17	0.45
7:G:38:LEU:CG	7:G:42:ILE:HD11	2.32	0.45
9:I:6:GLY:O	9:I:80:GLY:HA3	2.16	0.45
9:I:32:ASP:O	9:I:33:PHE:C	2.55	0.45
9:I:48:GLU:N	9:I:49:PRO:CD	2.73	0.45
12:L:108:ALA:O	12:L:109:GLY:O	2.34	0.45
12:L:126:LYS:HD2	12:L:126:LYS:C	2.37	0.45
13:M:17:VAL:HG22	13:M:27:LYS:CD	2.37	0.45
14:N:42:ILE:O	14:N:45:ARG:N	2.49	0.45
16:P:10:GLY:HA3	16:P:16:HIS:H	1.82	0.45
20:T:67:ALA:CB	20:T:77:ALA:HB2	2.47	0.45
1:A:56:U:O2'	1:A:57:G:H5'	2.17	0.45
1:A:131:C:O2	1:A:232:G:C2	2.70	0.45
1:A:307:C:H2'	1:A:308:C:H6	1.82	0.45
1:A:307:C:C6	1:A:308:C:H5	2.35	0.45
1:A:407:G:H1	1:A:435:C:H42	1.64	0.45
1:A:419:C:O2	1:A:419:C:C2'	2.65	0.45
1:A:568:G:O2'	1:A:574:A:N1	2.38	0.45
1:A:734:G:C5	1:A:735:C:C4	3.05	0.45
1:A:1095:U:C5'	1:A:1109:C:O2	2.65	0.45
1:A:1112:C:O2	3:C:179:ARG:HB2	2.16	0.45
1:A:1260:C:P	1:A:1284:C:H4'	2.57	0.45
1:A:1353:G:C2	1:A:1354:C:C5	3.05	0.45
3:C:180:ALA:HB1	3:C:206:GLU:HG3	1.99	0.45
3:C:201:TYR:O	3:C:202:ILE:CG1	2.65	0.45
4:D:201:GLN:NE2	5:E:99:GLY:CA	2.76	0.45
5:E:16:THR:HG22	5:E:27:ARG:O	2.16	0.45
6:F:47:ARG:HA	6:F:47:ARG:NH1	2.17	0.45
6:F:67:MET:SD	6:F:72:VAL:HA	2.57	0.45
7:G:70:LYS:HA	7:G:71:PRO:HD2	1.78	0.45
7:G:125:MET:HB3	7:G:126:ASP:H	1.59	0.45
9:I:110:GLU:OE2	9:I:113:LYS:NZ	2.49	0.45
9:I:113:LYS:N	9:I:113:LYS:HD2	2.32	0.45
10:J:9:ARG:HB3	10:J:9:ARG:HH11	1.80	0.45
10:J:94:VAL:HG12	10:J:95:GLU:H	1.81	0.45
11:K:91:ARG:O	11:K:93:GLN:N	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:N:25:VAL:HG23	14:N:38:GLY:O	2.17	0.45
17:Q:5:VAL:O	17:Q:6:LEU:HG	2.17	0.45
17:Q:67:LYS:O	17:Q:68:ARG:C	2.54	0.45
19:S:11:VAL:HG12	19:S:16:LEU:HB2	1.99	0.45
1:A:12:U:H4'	1:A:526:C:O2'	2.17	0.45
1:A:127:G:O3'	17:Q:2:PRO:HD2	2.16	0.45
1:A:197:A:H1'	1:A:198:G:H1'	1.97	0.45
1:A:377:G:C2	1:A:378:G:N7	2.85	0.45
1:A:555:C:H2'	1:A:556:C:C6	2.52	0.45
1:A:765:G:O6	1:A:812:C:C6	2.70	0.45
1:A:913:A:O2'	1:A:914:A:C5'	2.65	0.45
1:A:1076:C:N3	1:A:1077:G:N7	2.65	0.45
1:A:1260:C:C5'	1:A:1284:C:H4'	2.47	0.45
1:A:1312:G:C2	1:A:1326:C:N3	2.85	0.45
1:A:1343:G:C5	1:A:1344:C:C4	3.05	0.45
1:A:1483:A:H2'	1:A:1484:C:H5'	1.99	0.45
1:A:1521:G:C2	1:A:1522:U:N3	2.85	0.45
2:B:21:ARG:HA	2:B:39:ILE:CG1	2.47	0.45
2:B:55:PHE:HE1	2:B:218:ALA:HA	1.82	0.45
2:B:148:TYR:CD2	2:B:148:TYR:N	2.84	0.45
2:B:194:PRO:HG2	2:B:195:ASP:H	1.82	0.45
3:C:141:VAL:HG23	3:C:142:MET:H	1.81	0.45
3:C:182:ILE:O	3:C:183:ASP:HB2	2.17	0.45
4:D:194:LEU:HB3	4:D:196:LEU:HG	1.97	0.45
5:E:39:GLY:HA2	5:E:71:LEU:HD21	1.98	0.45
5:E:122:GLU:OE1	5:E:131:ILE:HG21	2.17	0.45
5:E:148:VAL:O	5:E:149:GLU:C	2.55	0.45
7:G:61:VAL:HG13	7:G:128:ALA:HB2	1.98	0.45
8:H:65:TYR:CD1	8:H:65:TYR:N	2.84	0.45
9:I:118:LYS:HB2	9:I:121:ARG:HB2	1.99	0.45
10:J:22:LYS:HZ2	10:J:23:ILE:HD13	1.81	0.45
13:M:48:LEU:HD22	13:M:53:VAL:HG22	1.98	0.45
13:M:57:ARG:HH11	13:M:57:ARG:HG2	1.82	0.45
17:Q:44:ALA:HB2	17:Q:59:ILE:HD12	1.99	0.45
17:Q:88:TYR:O	17:Q:89:LEU:C	2.56	0.45
17:Q:95:TYR:C	17:Q:97:SER:N	2.70	0.45
20:T:18:GLN:O	20:T:22:ARG:HG3	2.17	0.45
20:T:101:GLY:O	20:T:102:GLY:C	2.54	0.45
1:A:293:G:N2	1:A:294:U:C2	2.85	0.44
1:A:499:A:C4'	1:A:500:G:H5'	2.45	0.44
1:A:547:A:OP2	4:D:2:GLY:N	2.49	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:658:G:OP1	15:O:31:LEU:HD11	2.16	0.44
1:A:722:A:O2'	1:A:723:U:C2	2.69	0.44
1:A:922:G:O2'	1:A:1398:A:N1	2.47	0.44
1:A:1036:G:C2'	1:A:1037:C:H5'	2.46	0.44
1:A:1238:A:N7	1:A:1301:U:O4	2.50	0.44
1:A:1241:G:C6	1:A:1242:C:N4	2.85	0.44
1:A:1269:A:C2	1:A:1313:U:C1'	3.00	0.44
1:A:1270:C:O2'	1:A:1271:G:H5'	2.16	0.44
1:A:1347:G:N1	9:I:107:ARG:NH2	2.66	0.44
1:A:1405:G:C2	1:A:1497:G:C2	3.04	0.44
1:A:1409:C:H2'	1:A:1410:G:C8	2.52	0.44
1:A:1507:A:N6	1:A:1508:G:O6	2.50	0.44
1:A:1518:A:O2'	1:A:1519:A:H5'	2.17	0.44
3:C:62:ASP:C	3:C:97:LYS:HB3	2.38	0.44
3:C:173:VAL:O	3:C:175:LEU:HD22	2.17	0.44
4:D:58:LEU:HD13	4:D:58:LEU:C	2.37	0.44
4:D:111:ALA:HB1	4:D:116:GLN:OE1	2.17	0.44
9:I:79:LEU:O	9:I:83:ARG:N	2.50	0.44
10:J:44:VAL:CG1	10:J:45:ARG:N	2.75	0.44
11:K:54:ARG:C	11:K:56:GLY:H	2.20	0.44
12:L:41:ARG:NH1	12:L:42:THR:H	2.15	0.44
14:N:50:LYS:CG	14:N:51:GLY:H	2.21	0.44
15:O:26:GLU:HA	15:O:81:LEU:CD1	2.45	0.44
15:O:48:LYS:O	15:O:49:ASP:C	2.55	0.44
16:P:9:PHE:CE2	16:P:18:ARG:HD2	2.52	0.44
17:Q:23:VAL:O	17:Q:23:VAL:HG12	2.16	0.44
17:Q:66:SER:O	17:Q:70:ARG:NH1	2.50	0.44
20:T:73:HIS:HB3	20:T:74:LYS:H	1.39	0.44
1:A:21:G:C2	1:A:22:G:C5	3.05	0.44
1:A:42:G:C2	1:A:401:C:O2	2.70	0.44
1:A:282:A:C4	1:A:283:C:C6	3.06	0.44
1:A:303:A:C2	1:A:304:U:O2	2.70	0.44
1:A:419:C:H5	1:A:425:G:C2	2.34	0.44
1:A:586:C:H1'	1:A:878:G:O2'	2.16	0.44
1:A:882:C:C2'	1:A:883:C:H5'	2.47	0.44
1:A:1214:C:H4'	1:A:1215:G:OP1	2.18	0.44
1:A:1310:G:OP2	13:M:88:ARG:NH1	2.50	0.44
1:A:1494:G:H2'	1:A:1495:U:H5'	1.98	0.44
2:B:149:LEU:O	2:B:153:ARG:HB2	2.17	0.44
2:B:175:ARG:HH11	2:B:175:ARG:CB	2.28	0.44
2:B:214:ILE:O	2:B:215:LEU:C	2.55	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:97:LYS:N	3:C:97:LYS:HD2	2.31	0.44
4:D:125:HIS:HA	4:D:149:ALA:HB2	2.00	0.44
4:D:149:ALA:O	4:D:152:SER:HB3	2.17	0.44
5:E:31:LEU:HD21	5:E:45:PHE:HB2	1.99	0.44
5:E:52:PRO:O	5:E:53:LEU:C	2.52	0.44
6:F:65:VAL:HG22	6:F:66:GLU:N	2.32	0.44
9:I:46:ALA:HB1	9:I:81:ILE:HD11	1.98	0.44
10:J:9:ARG:NH1	10:J:69:ASN:CB	2.77	0.44
11:K:22:HIS:CG	11:K:22:HIS:O	2.69	0.44
11:K:124:LYS:C	11:K:125:PHE:HD1	2.21	0.44
13:M:55:ARG:HG3	13:M:55:ARG:NH1	2.31	0.44
14:N:27:CYS:SG	14:N:29:ARG:CB	3.05	0.44
14:N:47:LEU:O	14:N:49:HIS:N	2.50	0.44
15:O:9:GLN:CA	15:O:12:ILE:HD13	2.47	0.44
16:P:53:VAL:HA	16:P:56:ALA:HB3	1.98	0.44
18:R:37:VAL:HB	18:R:41:LYS:CD	2.47	0.44
18:R:41:LYS:HZ2	18:R:41:LYS:HB2	1.81	0.44
19:S:6:LYS:HB3	19:S:7:LYS:H	1.45	0.44
19:S:16:LEU:O	19:S:19:VAL:HG12	2.17	0.44
1:A:131:C:H2'	1:A:132:C:C5	2.50	0.44
1:A:190:C:H2'	1:A:190(A):C:C6	2.49	0.44
1:A:253:U:H2'	1:A:254:G:C8	2.52	0.44
1:A:353:A:H5'	1:A:353:A:H8	1.83	0.44
1:A:448:A:N7	1:A:486:U:O4	2.50	0.44
1:A:541:G:N3	1:A:542:G:C8	2.86	0.44
1:A:564:C:C4	1:A:565:U:C4	3.05	0.44
1:A:740:U:H2'	1:A:741:G:H8	1.81	0.44
1:A:775:G:O2'	1:A:776:G:H5'	2.17	0.44
1:A:813:U:O2	1:A:813:U:H2'	2.17	0.44
1:A:902:G:O2'	1:A:903:G:H5'	2.17	0.44
1:A:913:A:O2'	1:A:914:A:O5'	2.35	0.44
1:A:1000:U:H2'	1:A:1001:A:C8	2.52	0.44
1:A:1064:G:N3	1:A:1066:C:N4	2.66	0.44
2:B:116:GLU:HG2	2:B:153:ARG:CZ	2.47	0.44
3:C:177:THR:O	3:C:177:THR:HG23	2.18	0.44
5:E:107:ARG:O	5:E:108:ALA:C	2.56	0.44
9:I:10:ARG:CD	9:I:105:ASP:HB3	2.46	0.44
9:I:23:ASN:HD22	9:I:23:ASN:HA	1.59	0.44
9:I:55:ALA:O	9:I:58:ARG:N	2.51	0.44
11:K:85:ARG:HD3	11:K:111:ASP:O	2.18	0.44
12:L:8:ASN:O	12:L:9:GLN:C	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:M:9:ILE:HA	13:M:10:PRO:HD2	1.81	0.44
18:R:22:VAL:HG12	18:R:26:LEU:HD13	1.99	0.44
1:A:182:U:H6	1:A:182:U:C5'	2.26	0.44
1:A:250:A:H5'	1:A:252:U:H5'	1.99	0.44
1:A:332:G:H2'	1:A:333:G:H8	1.82	0.44
1:A:436:C:C2	1:A:437:U:C5	3.05	0.44
1:A:455:C:H2'	1:A:456:C:H6	1.82	0.44
1:A:707:C:C5'	11:K:20:TYR:CD2	2.98	0.44
1:A:1029:C:C2'	1:A:1030:C:H5'	2.48	0.44
1:A:1128:C:N4	1:A:1143:G:H1	2.15	0.44
1:A:1149:C:OP1	9:I:9:ARG:HD3	2.17	0.44
1:A:1258:G:H2'	1:A:1259:C:H6	1.82	0.44
1:A:1438:G:C6	1:A:1439:C:N4	2.85	0.44
1:A:1468:A:H2'	1:A:1469:G:O4'	2.17	0.44
1:A:1521:G:C2	1:A:1522:U:C2	3.05	0.44
3:C:126:ARG:C	3:C:127:ARG:CD	2.86	0.44
3:C:162:GLN:HG3	3:C:164:ARG:CG	2.47	0.44
4:D:96:LEU:C	4:D:98:GLU:N	2.69	0.44
5:E:15:ARG:CZ	5:E:26:PHE:CE2	3.01	0.44
5:E:139:LEU:HA	5:E:142:LEU:CG	2.48	0.44
7:G:58:PRO:HA	7:G:61:VAL:CG2	2.47	0.44
10:J:3:LYS:CB	10:J:77:PRO:HD3	2.48	0.44
10:J:10:GLY:O	10:J:67:THR:HA	2.16	0.44
12:L:90:VAL:HG12	12:L:92:ASP:N	2.31	0.44
13:M:39:ILE:HD12	13:M:56:LEU:CD2	2.48	0.44
15:O:79:ARG:O	15:O:80:ALA:C	2.55	0.44
17:Q:56:VAL:HB	17:Q:77:VAL:HB	1.99	0.44
17:Q:94:ASN:O	17:Q:95:TYR:O	2.36	0.44
19:S:10:PHE:HZ	19:S:12:ASP:OD2	1.99	0.44
20:T:19:SER:OG	20:T:20:LEU:N	2.50	0.44
1:A:97:G:C2	1:A:98:U:H1'	2.53	0.44
1:A:167:G:C2'	1:A:168:G:C8	2.91	0.44
1:A:306:G:H2'	1:A:307:C:H6	1.83	0.44
1:A:355:C:C4	1:A:356:A:N7	2.86	0.44
1:A:369:C:C2'	1:A:370:C:H5'	2.48	0.44
1:A:482:A:C2	1:A:483:C:H1'	2.53	0.44
1:A:593:G:H2'	1:A:594:G:O4'	2.18	0.44
1:A:957:U:H3	1:A:960:U:H5''	1.82	0.44
1:A:1110:A:C2'	1:A:1111:A:H5'	2.48	0.44
1:A:1349:A:C2'	1:A:1350:A:H5'	2.47	0.44
1:A:1487:G:N2	1:A:1488:G:H1'	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:48:TYR:HD1	3:C:52:LEU:CD2	2.30	0.44
4:D:7:PRO:HG2	4:D:10:ARG:HD2	1.99	0.44
4:D:24:GLU:C	4:D:26:CYS:H	2.21	0.44
5:E:138:ALA:O	5:E:139:LEU:C	2.55	0.44
5:E:144:THR:HG22	5:E:145:LYS:N	2.31	0.44
6:F:1:MET:SD	6:F:66:GLU:O	2.76	0.44
7:G:36:LYS:HA	7:G:39:ALA:HB3	1.99	0.44
7:G:41:ARG:O	7:G:43:PHE:N	2.50	0.44
7:G:60:LYS:HG2	7:G:64:GLN:HB3	1.99	0.44
9:I:5:TYR:CG	9:I:6:GLY:N	2.83	0.44
10:J:84:GLN:O	10:J:85:LEU:HG	2.17	0.44
15:O:79:ARG:O	15:O:82:ILE:CG2	2.65	0.44
15:O:87:ILE:CG2	15:O:88:ARG:N	2.66	0.44
18:R:41:LYS:NZ	18:R:41:LYS:HB2	2.32	0.44
18:R:59:SER:N	18:R:62:GLU:HB2	2.29	0.44
19:S:22:LEU:O	19:S:25:LYS:CD	2.64	0.44
19:S:30:LEU:HD23	19:S:30:LEU:C	2.38	0.44
1:A:616:G:C6	1:A:617:G:N7	2.85	0.44
1:A:740:U:O2'	1:A:741:G:C5'	2.64	0.44
1:A:823:G:C6	1:A:824:C:N4	2.85	0.44
1:A:922:G:C5'	5:E:19:MET:O	2.63	0.44
1:A:971:G:H5''	1:A:972:C:H5''	2.00	0.44
1:A:1036:G:C6	1:A:1037:C:H1'	2.53	0.44
1:A:1149:C:C2'	1:A:1150:U:O5'	2.66	0.44
1:A:1285:A:O2'	1:A:1286:A:P	2.76	0.44
1:A:1291:G:OP1	7:G:41:ARG:NH2	2.51	0.44
1:A:1302:U:H3'	1:A:1303:C:C5'	2.48	0.44
2:B:72:GLY:HA2	2:B:165:VAL:CG2	2.47	0.44
3:C:127:ARG:HG2	3:C:127:ARG:NH1	2.32	0.44
4:D:21:LEU:N	4:D:21:LEU:CD1	2.81	0.44
4:D:79:PHE:CD1	4:D:207:TYR:HD1	2.33	0.44
4:D:126:ILE:HG22	4:D:127:THR:N	2.32	0.44
7:G:29:LYS:HB3	7:G:105:VAL:HG21	2.00	0.44
9:I:20:ARG:O	9:I:21:PRO:O	2.36	0.44
11:K:25:TYR:OH	11:K:87:THR:HB	2.18	0.44
11:K:66:LEU:HD23	11:K:69:ALA:HB3	1.99	0.44
11:K:98:LEU:H	11:K:98:LEU:CD2	2.27	0.44
12:L:46:LYS:HG2	12:L:47:LYS:N	2.32	0.44
12:L:83:VAL:HG13	12:L:84:LEU:N	2.33	0.44
12:L:119:LYS:O	12:L:120:TYR:HB2	2.17	0.44
13:M:35:GLU:C	13:M:37:THR:N	2.70	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:Q:22:LEU:HD12	17:Q:22:LEU:HA	1.73	0.44
18:R:21:LYS:H	18:R:21:LYS:CD	2.20	0.44
18:R:40:LEU:CD2	18:R:40:LEU:C	2.86	0.44
19:S:23:ASN:HA	19:S:27:GLU:CA	2.44	0.44
20:T:23:ARG:O	20:T:24:LEU:C	2.56	0.44
20:T:33:ILE:O	20:T:34:LYS:C	2.56	0.44
20:T:69:GLY:C	20:T:71:THR:N	2.70	0.44
1:A:142:G:H2'	1:A:143:A:C8	2.46	0.44
1:A:146:G:C4	1:A:147:G:C8	3.06	0.44
1:A:172:A:O2'	1:A:173:U:H5'	2.18	0.44
1:A:364:A:C2	1:A:365:U:O4	2.71	0.44
1:A:373:A:C2	1:A:482:A:C6	3.05	0.44
1:A:411:A:C8	1:A:413:G:N9	2.86	0.44
1:A:517:G:C6	1:A:531:U:O4'	2.71	0.44
1:A:735:C:C2	1:A:736:C:C5	3.06	0.44
1:A:936:C:H2'	1:A:937:A:C8	2.53	0.44
1:A:1070:U:H5'	5:E:18:ARG:HH12	1.83	0.44
1:A:1135:U:H4'	1:A:1136:U:C5	2.50	0.44
1:A:1210:C:H3'	1:A:1211:U:H5''	2.00	0.44
1:A:1272:G:H2'	1:A:1273:G:H8	1.83	0.44
3:C:160:ALA:O	3:C:162:GLN:N	2.51	0.44
4:D:152:SER:CB	4:D:155:LEU:HD12	2.47	0.44
5:E:20:GLN:O	5:E:21:ALA:O	2.35	0.44
5:E:80:ILE:HG22	8:H:104:ARG:NH2	2.15	0.44
5:E:119:LEU:HA	5:E:119:LEU:HD23	1.53	0.44
6:F:9:VAL:HA	6:F:59:TYR:O	2.18	0.44
7:G:148:ASN:C	7:G:150:ALA:N	2.69	0.44
8:H:44:PHE:C	8:H:45:ILE:HG12	2.37	0.44
8:H:102:ARG:HH12	8:H:105:ARG:NH1	2.15	0.44
10:J:39:PRO:O	10:J:70:ARG:NH1	2.51	0.44
10:J:62:HIS:HB3	14:N:59:ALA:CB	2.41	0.44
10:J:67:THR:O	10:J:68:HIS:C	2.56	0.44
12:L:24:VAL:O	12:L:24:VAL:HG12	2.16	0.44
12:L:110:VAL:O	12:L:122:THR:HG21	2.18	0.44
13:M:71:ARG:HG2	13:M:71:ARG:NH1	2.33	0.44
14:N:37:PHE:O	14:N:39:LEU:HG	2.18	0.44
15:O:36:ILE:HG22	15:O:37:ASN:N	2.32	0.44
18:R:35:ARG:O	18:R:36:ASN:C	2.56	0.44
1:A:53:A:C2'	1:A:54:C:O5'	2.65	0.44
1:A:106:C:H2'	1:A:107:G:O4'	2.17	0.44
1:A:134:A:H2'	1:A:135:C:O4'	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:G:C6	1:A:192:U:C4	3.05	0.44
1:A:229:U:H2'	1:A:230:G:H8	1.83	0.44
1:A:521:G:O2'	1:A:522:C:H5'	2.18	0.44
1:A:543:C:H2'	1:A:544:G:O4'	2.18	0.44
1:A:1064:G:H4'	1:A:1065:U:C5'	2.47	0.44
1:A:1250:A:C2	1:A:1287:A:C2	3.06	0.44
1:A:1513:A:O2'	1:A:1514:C:H5'	2.18	0.44
2:B:63:MET:HB3	2:B:225:ALA:HB1	1.99	0.44
3:C:91:LEU:O	3:C:95:THR:CG2	2.66	0.44
3:C:108:ASN:OD1	3:C:108:ASN:N	2.49	0.44
3:C:113:ALA:O	3:C:116:VAL:HG23	2.18	0.44
3:C:116:VAL:O	3:C:117:ALA:C	2.57	0.44
3:C:154:SER:HA	3:C:165:THR:HA	1.98	0.44
5:E:118:ILE:HG22	5:E:119:LEU:H	1.83	0.44
6:F:19:LEU:HD23	6:F:19:LEU:C	2.38	0.44
7:G:76:ARG:O	7:G:86:GLN:HA	2.18	0.44
7:G:88:PRO:CG	7:G:151:TYR:O	2.66	0.44
7:G:114:ARG:HD3	7:G:114:ARG:HA	1.81	0.44
9:I:100:GLY:C	9:I:102:LEU:H	2.21	0.44
9:I:121:ARG:O	9:I:121:ARG:HD3	2.18	0.44
13:M:33:ALA:C	13:M:35:GLU:N	2.71	0.44
13:M:34:LEU:CD1	13:M:41:PRO:HA	2.48	0.44
13:M:39:ILE:CD1	13:M:52:GLU:HB3	2.47	0.44
13:M:80:ARG:C	13:M:82:MET:N	2.71	0.44
13:M:80:ARG:HG2	13:M:81:LEU:N	2.31	0.44
13:M:85:GLY:O	13:M:86:CYS:C	2.56	0.44
16:P:35:LYS:HG2	16:P:36:ILE:N	2.33	0.44
17:Q:102:GLY:O	17:Q:103:GLY:O	2.36	0.44
18:R:28:GLU:OE1	18:R:28:GLU:N	2.51	0.44
18:R:37:VAL:O	18:R:38:GLU:HB2	2.18	0.44
19:S:18:LYS:O	19:S:22:LEU:HD21	2.18	0.44
1:A:32:A:H2'	1:A:33:A:O4'	2.16	0.44
1:A:75:G:N2	1:A:96:G:C2	2.85	0.44
1:A:197:A:H1'	1:A:198:G:C1'	2.47	0.44
1:A:290:C:C3'	1:A:290:C:C6	3.01	0.44
1:A:290:C:N4	1:A:291:C:C5	2.86	0.44
1:A:597:G:C5	1:A:644:G:C6	3.06	0.44
1:A:833:U:H2'	1:A:834:C:C6	2.53	0.44
1:A:1039:C:H2'	1:A:1040:U:H6	1.82	0.44
1:A:1048:G:H21	1:A:1214:C:C2'	2.22	0.44
1:A:1129:C:OP1	9:I:18:PHE:HE1	2.01	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1273:G:H2'	1:A:1274:G:C8	2.53	0.44
1:A:1406:U:H2'	1:A:1407:C:C5	2.53	0.44
1:A:1502:A:H5'	1:A:1504:G:N7	2.32	0.44
2:B:119:GLU:O	2:B:123:ALA:HB2	2.18	0.44
3:C:38:ARG:HH11	3:C:38:ARG:HG3	1.82	0.44
3:C:99:VAL:HG23	3:C:100:ALA:N	2.33	0.44
3:C:157:ILE:HB	3:C:164:ARG:NH2	2.33	0.44
4:D:61:LYS:HA	4:D:203:VAL:HG22	2.00	0.44
5:E:55:VAL:CG2	5:E:56:GLN:H	2.13	0.44
6:F:83:ASP:C	6:F:85:VAL:N	2.70	0.44
7:G:101:LEU:N	7:G:101:LEU:CD2	2.71	0.44
8:H:32:LYS:O	8:H:33:GLU:C	2.55	0.44
8:H:41:ARG:C	8:H:43:GLY:H	2.20	0.44
9:I:43:ALA:CB	9:I:74:ILE:HD13	2.48	0.44
9:I:50:LEU:HD11	9:I:81:ILE:HB	1.99	0.44
11:K:58:PRO:HB2	11:K:93:GLN:CG	2.43	0.44
12:L:84:LEU:CG	12:L:101:VAL:HG11	2.47	0.44
14:N:9:LYS:C	14:N:11:LYS:H	2.22	0.44
14:N:24:CYS:SG	14:N:27:CYS:N	2.91	0.44
15:O:31:LEU:C	15:O:31:LEU:HD23	2.39	0.44
16:P:7:ALA:HB1	16:P:28:ARG:O	2.18	0.44
16:P:49:LEU:HD13	16:P:73:LEU:HD22	2.00	0.44
17:Q:56:VAL:C	17:Q:57:VAL:HG13	2.38	0.44
18:R:34:TYR:CE1	18:R:35:ARG:HG3	2.53	0.44
1:A:110:C:H2'	1:A:111:G:C5'	2.48	0.43
1:A:160:A:H61	1:A:347:G:H1'	1.82	0.43
1:A:182:U:H2'	1:A:183:G:C5'	2.47	0.43
1:A:190(E):U:N3	17:Q:72:ARG:CZ	2.81	0.43
1:A:229:U:H2'	1:A:230:G:C8	2.53	0.43
1:A:542:G:H5'	4:D:41:GLY:HA2	2.00	0.43
1:A:639:G:C2	1:A:640:A:C5	3.06	0.43
1:A:750:G:N2	1:A:751:U:C2	2.86	0.43
1:A:835:U:OP1	18:R:64:ARG:NH2	2.51	0.43
1:A:1003(A):G:N2	1:A:1038:C:C2	2.85	0.43
1:A:1055:A:C8	1:A:1206:G:C2	3.06	0.43
1:A:1095:U:H2'	1:A:1096:C:C6	2.52	0.43
1:A:1117:G:N2	1:A:1180:A:O2'	2.50	0.43
1:A:1152:A:OP1	10:J:68:HIS:ND1	2.50	0.43
1:A:1223:C:OP1	1:A:1224:G:H3'	2.18	0.43
1:A:1296:C:H5'	1:A:1297:C:OP2	2.18	0.43
1:A:1316:G:H4'	14:N:18:VAL:HG11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:9:GLU:CD	2:B:217:ARG:NH2	2.65	0.43
2:B:130:ARG:CB	2:B:131:PRO:HD2	2.41	0.43
3:C:9:GLY:O	3:C:12:LEU:HG	2.18	0.43
3:C:85:ARG:HH11	3:C:85:ARG:HG3	1.83	0.43
3:C:93:LYS:HE2	3:C:93:LYS:CA	2.40	0.43
3:C:130:VAL:HB	3:C:131:ARG:H	1.63	0.43
3:C:153:VAL:HG12	3:C:154:SER:N	2.33	0.43
4:D:28:SER:O	4:D:30:LYS:N	2.51	0.43
4:D:91:SER:O	4:D:94:LEU:N	2.51	0.43
4:D:117:ALA:C	4:D:121:VAL:HG23	2.38	0.43
4:D:156:GLU:HG2	4:D:157:LEU:N	2.24	0.43
7:G:103:TRP:CD1	7:G:137:LYS:HD3	2.53	0.43
8:H:5:PRO:HB3	8:H:32:LYS:HE3	2.00	0.43
8:H:14:ARG:NH1	8:H:14:ARG:CB	2.80	0.43
8:H:108:GLY:HA3	8:H:138:TRP:HB3	1.99	0.43
9:I:14:VAL:O	9:I:15:ALA:HB2	2.17	0.43
9:I:89:ASN:C	9:I:91:ASP:N	2.69	0.43
10:J:53:PRO:HA	14:N:41:ARG:HH21	1.83	0.43
11:K:14:VAL:HG12	11:K:16:SER:H	1.82	0.43
11:K:27:ASN:OD1	11:K:28:THR:N	2.43	0.43
13:M:39:ILE:O	13:M:41:PRO:HD3	2.18	0.43
13:M:81:LEU:CD1	13:M:88:ARG:HD3	2.48	0.43
18:R:72:ARG:O	18:R:75:ILE:HB	2.17	0.43
19:S:40:ILE:CD1	19:S:62:ILE:HD11	2.46	0.43
20:T:44:ALA:O	20:T:45:GLN:C	2.56	0.43
20:T:73:HIS:O	20:T:76:ALA:HB3	2.18	0.43
21:V:10:ARG:HH11	21:V:10:ARG:CG	2.29	0.43
1:A:112:G:C4'	1:A:389:A:H5''	2.44	0.43
1:A:129:U:O3'	1:A:129(A):G:H3'	2.18	0.43
1:A:289:G:H5'	1:A:289:G:H8	1.83	0.43
1:A:290:C:C6	1:A:290:C:H3'	2.53	0.43
1:A:322:C:H2'	1:A:323:U:H5'	1.99	0.43
1:A:533:A:H2'	1:A:534:U:H5''	1.99	0.43
1:A:880:C:H5	12:L:9:GLN:HE22	1.65	0.43
1:A:930:C:H2'	1:A:931:C:C5'	2.46	0.43
1:A:942:G:C2	1:A:943:U:C6	3.06	0.43
1:A:1347:G:H3'	9:I:109:VAL:HA	1.99	0.43
1:A:1477:C:H2'	1:A:1478:C:C5	2.53	0.43
2:B:19:HIS:HD2	2:B:189:ASP:OD1	2.01	0.43
3:C:61:ALA:O	3:C:63:ASN:ND2	2.50	0.43
3:C:114:PRO:HG3	3:C:185:GLY:CA	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:126:ARG:O	3:C:127:ARG:HB2	2.18	0.43
4:D:22:LYS:O	4:D:23:GLY:C	2.57	0.43
5:E:98:THR:HB	5:E:117:ASP:HB3	1.99	0.43
6:F:18:GLN:O	6:F:21:LEU:CB	2.63	0.43
7:G:31:MET:HG2	7:G:36:LYS:CB	2.48	0.43
7:G:72:ARG:HA	7:G:96:GLN:OE1	2.18	0.43
9:I:65:VAL:HG22	9:I:66:ARG:N	2.32	0.43
11:K:94:ALA:O	11:K:95:ILE:C	2.55	0.43
12:L:59:ARG:HH12	12:L:65:GLU:HG2	1.83	0.43
12:L:87:GLY:O	12:L:99:HIS:N	2.52	0.43
13:M:74:VAL:C	13:M:76:ALA:H	2.21	0.43
15:O:10:LYS:O	15:O:14:GLU:HB2	2.18	0.43
16:P:75:ARG:CA	16:P:80:PHE:HE1	2.31	0.43
19:S:46:GLY:O	19:S:47:HIS:O	2.36	0.43
20:T:65:LYS:O	20:T:66:ALA:C	2.55	0.43
20:T:93:GLU:HA	20:T:93:GLU:OE2	2.18	0.43
1:A:122:G:H8	1:A:122:G:OP1	2.02	0.43
1:A:124:G:C5	1:A:125:U:C5	3.06	0.43
1:A:129(A):G:H4'	1:A:130:A:OP2	2.18	0.43
1:A:381:C:C2	1:A:382:A:C8	3.07	0.43
1:A:448:A:C2	1:A:449:C:C4	3.06	0.43
1:A:642:A:C6	1:A:643:C:C4	3.06	0.43
1:A:642:A:C5	1:A:643:C:C5	3.07	0.43
1:A:688:G:H2'	1:A:689:C:H6	1.84	0.43
1:A:836:G:H1	1:A:850:U:H3	1.66	0.43
1:A:858:G:C6	1:A:869:G:C8	3.06	0.43
1:A:877:C:H1'	8:H:3:THR:HG22	2.00	0.43
1:A:891:U:C2'	1:A:892:A:H5'	2.48	0.43
1:A:959:A:H5''	1:A:960:U:OP2	2.19	0.43
1:A:1142:G:C3'	1:A:1143:G:H8	2.26	0.43
1:A:1360:A:H3'	1:A:1361:G:C8	2.53	0.43
2:B:21:ARG:NH1	2:B:23:ARG:CZ	2.81	0.43
2:B:44:LEU:O	2:B:47:THR:N	2.51	0.43
2:B:80:ILE:HG22	2:B:81:VAL:N	2.32	0.43
3:C:11:ARG:NH1	3:C:179:ARG:H	1.95	0.43
3:C:48:TYR:CD1	3:C:52:LEU:HD13	2.53	0.43
3:C:85:ARG:O	3:C:87:LEU:N	2.51	0.43
4:D:17:VAL:CG1	4:D:18:LYS:N	2.63	0.43
4:D:151:LYS:HD2	4:D:151:LYS:N	2.30	0.43
5:E:151:LEU:CD1	8:H:77:GLU:OE2	2.67	0.43
6:F:1:MET:HE2	6:F:36:ARG:HH21	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:49:ILE:HG22	7:G:49:ILE:O	2.18	0.43
11:K:91:ARG:O	11:K:94:ALA:N	2.50	0.43
12:L:106:ASP:N	12:L:106:ASP:OD2	2.51	0.43
15:O:68:ARG:HH11	15:O:68:ARG:HB3	1.83	0.43
16:P:60:LEU:HA	16:P:60:LEU:HD23	1.73	0.43
16:P:67:THR:HG22	16:P:68:ASP:N	2.32	0.43
17:Q:9:VAL:CG1	17:Q:10:VAL:H	2.30	0.43
17:Q:75:ARG:HG3	17:Q:76:LEU:H	1.83	0.43
20:T:50:GLU:HG2	20:T:100:ILE:CG1	2.39	0.43
20:T:77:ALA:O	20:T:78:ALA:C	2.56	0.43
1:A:11:G:O2'	1:A:12:U:H5'	2.18	0.43
1:A:23:C:H2'	1:A:24:U:C6	2.49	0.43
1:A:149:A:C2	1:A:150:C:C6	3.07	0.43
1:A:186:C:H2'	1:A:187:C:C5	2.53	0.43
1:A:259:G:H2'	1:A:260:G:C8	2.54	0.43
1:A:299:G:C5	1:A:300:A:C6	3.07	0.43
1:A:338:A:H2'	1:A:339:C:O4'	2.19	0.43
1:A:351:G:O2'	1:A:352:C:O5'	2.32	0.43
1:A:575:G:C6	1:A:821:G:C8	3.07	0.43
1:A:602:A:C6	1:A:637:G:C6	3.06	0.43
1:A:1117:G:N2	1:A:1180:A:C1'	2.72	0.43
1:A:1241:G:C2	1:A:1242:C:C4	3.06	0.43
1:A:1251:A:H2'	1:A:1252:A:C8	2.53	0.43
1:A:1298:C:H4'	1:A:1299:A:C4	2.54	0.43
1:A:1372:U:C2'	1:A:1373:G:H5'	2.48	0.43
1:A:1431:C:C2	1:A:1470:G:C2	3.06	0.43
2:B:42:ILE:HD12	2:B:203:GLY:HA2	1.99	0.43
3:C:133:ALA:HA	3:C:136:GLN:NE2	2.32	0.43
3:C:137:ALA:CA	3:C:140:ARG:NH1	2.81	0.43
4:D:6:GLY:O	4:D:8:VAL:N	2.51	0.43
4:D:31:CYS:O	4:D:33:MET:N	2.36	0.43
4:D:117:ALA:O	4:D:118:ARG:C	2.56	0.43
4:D:126:ILE:O	4:D:132:ARG:HB2	2.18	0.43
5:E:100:VAL:HG23	5:E:100:VAL:O	2.19	0.43
6:F:23:LYS:O	6:F:27:GLN:HG3	2.19	0.43
6:F:32:ASN:HD22	6:F:32:ASN:C	2.19	0.43
7:G:139:GLU:C	7:G:141:VAL:N	2.72	0.43
8:H:87:SER:HB3	8:H:133:LEU:O	2.19	0.43
10:J:48:THR:HG23	10:J:62:HIS:HA	2.00	0.43
12:L:115:LYS:C	12:L:117:ARG:H	2.20	0.43
13:M:78:ILE:O	13:M:80:ARG:N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:T:14:LYS:NZ	20:T:17:ARG:NH2	2.67	0.43
1:A:97:G:H2'	1:A:98:U:O4'	2.19	0.43
1:A:116:A:O2'	1:A:117:G:H5'	2.19	0.43
1:A:655:A:C2	1:A:754:C:N4	2.87	0.43
1:A:733:A:H4'	1:A:734:G:OP1	2.19	0.43
1:A:949:A:C6	1:A:950:U:N3	2.86	0.43
1:A:965:A:C2'	1:A:966:G:OP2	2.67	0.43
2:B:216:SER:O	2:B:219:VAL:HG23	2.19	0.43
3:C:50:ALA:O	3:C:72:LYS:HG3	2.18	0.43
3:C:119:ARG:C	3:C:121:ALA:N	2.71	0.43
3:C:191:THR:CG2	3:C:194:GLY:C	2.86	0.43
3:C:203:PHE:C	3:C:204:LEU:HG	2.38	0.43
4:D:3:ARG:N	4:D:3:ARG:NE	2.67	0.43
4:D:93:PHE:O	4:D:97:LEU:HB2	2.18	0.43
5:E:13:ILE:HG22	5:E:30:ALA:CA	2.43	0.43
7:G:101:LEU:C	7:G:103:TRP:H	2.21	0.43
7:G:113:GLU:HB3	7:G:118:VAL:CG1	2.49	0.43
9:I:4:TYR:CD2	9:I:88:TYR:CB	3.02	0.43
9:I:28:VAL:HG13	9:I:65:VAL:HG12	2.01	0.43
9:I:54:ASP:C	9:I:58:ARG:HH21	2.22	0.43
9:I:113:LYS:HD2	9:I:113:LYS:H	1.84	0.43
9:I:123:PRO:O	9:I:124:GLN:HB3	2.18	0.43
10:J:49:VAL:O	10:J:60:ARG:O	2.36	0.43
13:M:39:ILE:HD12	13:M:56:LEU:HD21	2.01	0.43
13:M:74:VAL:O	13:M:75:ALA:C	2.57	0.43
13:M:80:ARG:HG2	13:M:81:LEU:HD23	2.00	0.43
14:N:23:ARG:HG2	14:N:23:ARG:HH11	1.83	0.43
14:N:24:CYS:O	14:N:25:VAL:C	2.56	0.43
14:N:53:LEU:O	14:N:56:VAL:HB	2.19	0.43
17:Q:79:SER:OG	17:Q:80:GLY:N	2.52	0.43
18:R:44:LEU:CD1	18:R:48:GLY:O	2.67	0.43
1:A:51:A:H4'	1:A:52:G:H5'	2.01	0.43
1:A:130:A:H1'	1:A:263:A:O2'	2.18	0.43
1:A:254:G:C2	1:A:273:A:C2	3.05	0.43
1:A:324:G:P	20:T:22:ARG:HH12	2.42	0.43
1:A:381:C:C4	1:A:382:A:C5	3.07	0.43
1:A:603:U:H2'	1:A:604:G:C8	2.45	0.43
1:A:666:G:C2	1:A:741:G:C4	3.06	0.43
1:A:717:C:O2'	1:A:734:G:O4'	2.37	0.43
1:A:877:C:O2	8:H:3:THR:CG2	2.67	0.43
1:A:1262:C:N4	1:A:1273:G:H1	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1367:C:C2	1:A:1368:G:C8	3.07	0.43
1:A:1480:G:N3	1:A:1480:G:H2'	2.34	0.43
1:A:1502:A:H2	1:A:1505:G:H1	1.66	0.43
1:A:1503:A:H5'	1:A:1531:A:H1'	2.01	0.43
2:B:60:ASP:O	2:B:61:LEU:C	2.56	0.43
2:B:73:THR:HG22	2:B:169:LYS:HE3	1.99	0.43
2:B:160:ASP:O	2:B:161:ALA:CB	2.66	0.43
3:C:95:THR:HG23	3:C:98:ASN:CA	2.48	0.43
3:C:157:ILE:C	3:C:159:GLY:N	2.72	0.43
7:G:139:GLU:O	7:G:142:GLU:N	2.51	0.43
11:K:34:ASP:OD2	11:K:38:ASN:N	2.52	0.43
11:K:69:ALA:O	11:K:73:MET:HG2	2.19	0.43
14:N:23:ARG:HG2	14:N:23:ARG:NH1	2.32	0.43
16:P:28:ARG:HG3	16:P:29:ASP:H	1.82	0.43
16:P:38:TYR:OH	16:P:50:LYS:HD3	2.18	0.43
17:Q:85:VAL:O	17:Q:86:GLU:C	2.56	0.43
17:Q:95:TYR:O	17:Q:98:LEU:N	2.52	0.43
19:S:33:THR:O	19:S:34:TRP:C	2.57	0.43
20:T:53:LEU:C	20:T:53:LEU:HD13	2.39	0.43
1:A:181:G:C2	1:A:195:A:C8	3.06	0.43
1:A:253:U:OP1	17:Q:67:LYS:HE3	2.18	0.43
1:A:287:U:H2'	1:A:288:A:C5'	2.49	0.43
1:A:317:G:O5'	1:A:317:G:H8	2.01	0.43
1:A:483:C:H3'	1:A:484:G:H2'	2.01	0.43
1:A:518:C:H5''	1:A:530:G:O4'	2.19	0.43
1:A:644:G:C6	1:A:645:C:C4	3.06	0.43
1:A:679:C:O5'	1:A:679:C:H6	2.02	0.43
1:A:964:A:O2'	10:J:55:LYS:HD3	2.19	0.43
1:A:1050:G:O2'	1:A:1051:C:H5'	2.18	0.43
1:A:1118:C:C1'	1:A:1179:A:H1'	2.49	0.43
1:A:1127:G:N1	1:A:1144:G:N2	2.55	0.43
2:B:55:PHE:CE1	2:B:218:ALA:HA	2.54	0.43
2:B:97:TRP:HB2	2:B:170:GLU:OE2	2.18	0.43
3:C:44:GLU:HG2	3:C:55:VAL:HG22	2.01	0.43
3:C:123:GLN:HE21	3:C:126:ARG:NH2	2.16	0.43
4:D:128:VAL:HG11	4:D:138:TYR:CE2	2.53	0.43
5:E:41:VAL:O	5:E:66:MET:HA	2.18	0.43
6:F:63:TYR:CD1	6:F:63:TYR:N	2.86	0.43
6:F:91:VAL:HG12	6:F:92:LYS:O	2.19	0.43
7:G:38:LEU:HG	7:G:42:ILE:CD1	2.31	0.43
7:G:120:ILE:O	7:G:124:LEU:HG	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:54:ASP:O	8:H:56:LYS:HD3	2.19	0.43
10:J:8:LEU:HD21	10:J:20:ALA:HB2	2.00	0.43
10:J:8:LEU:O	10:J:16:LEU:HD11	2.19	0.43
13:M:33:ALA:O	13:M:37:THR:CB	2.65	0.43
13:M:106:ASN:O	13:M:107:ALA:CB	2.62	0.43
14:N:11:LYS:O	14:N:12:ARG:CB	2.49	0.43
18:R:73:ALA:HB1	18:R:79:LEU:HD13	2.01	0.43
19:S:10:PHE:CZ	19:S:12:ASP:HA	2.53	0.43
19:S:64:GLU:OE1	19:S:64:GLU:HA	2.19	0.43
1:A:116:A:H2'	1:A:117:G:O4'	2.18	0.43
1:A:246:A:N1	1:A:279:A:N7	2.67	0.43
1:A:321:A:O2'	1:A:322:C:O4'	2.34	0.43
1:A:665:A:N1	1:A:732:C:C4	2.86	0.43
1:A:1029:C:H2'	1:A:1030:C:H5'	1.99	0.43
1:A:1074:G:O2'	2:B:103:THR:CG2	2.67	0.43
1:A:1076:C:O2'	1:A:1077:G:H5'	2.18	0.43
1:A:1298:C:C3'	1:A:1299:A:H5'	2.47	0.43
1:A:1345:U:OP1	1:A:1345:U:H3'	2.18	0.43
1:A:1424:C:H2'	1:A:1425:U:C5'	2.49	0.43
1:A:1470:G:H2'	1:A:1471:G:H8	1.84	0.43
1:A:1503:A:H5'	1:A:1531:A:O4'	2.18	0.43
2:B:174:VAL:O	2:B:175:ARG:C	2.56	0.43
3:C:22:TRP:CD1	3:C:59:ARG:NE	2.87	0.43
4:D:152:SER:O	4:D:153:ARG:C	2.57	0.43
7:G:43:PHE:HD2	7:G:44:TYR:CD2	2.37	0.43
8:H:95:VAL:HB	8:H:99:GLU:CB	2.48	0.43
8:H:96:GLY:O	8:H:97:VAL:HG13	2.19	0.43
8:H:116:LYS:HZ3	8:H:127:LEU:HD12	1.83	0.43
9:I:8:GLY:HA2	9:I:79:LEU:CD1	2.49	0.43
9:I:17:VAL:HG13	9:I:63:ILE:CG1	2.46	0.43
10:J:18:ALA:C	10:J:20:ALA:H	2.22	0.43
12:L:86:ARG:HH11	12:L:86:ARG:HG3	1.83	0.43
13:M:19:LEU:HB3	13:M:25:ILE:HG21	1.99	0.43
14:N:39:LEU:HB2	14:N:40:CYS:H	1.62	0.43
15:O:69:TYR:CE1	15:O:73:GLU:OE1	2.71	0.43
17:Q:5:VAL:CG1	17:Q:6:LEU:N	2.82	0.43
20:T:57:ARG:NH2	20:T:100:ILE:HG22	2.31	0.43
1:A:39:G:O2'	1:A:40:C:H5'	2.18	0.43
1:A:57:G:C6	1:A:58:C:N3	2.86	0.43
1:A:226:G:H2'	1:A:227:G:H8	1.84	0.43
1:A:327:A:N6	1:A:329:A:C6	2.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:435:C:C2'	1:A:436:C:O5'	2.67	0.43
1:A:885:G:H1	1:A:912:C:N4	2.16	0.43
1:A:979:C:H3'	1:A:980:C:H6	1.84	0.43
1:A:1002:G:H2'	1:A:1002:G:N3	2.34	0.43
1:A:1036:G:H2'	1:A:1037:C:C5'	2.49	0.43
1:A:1088:G:O5'	1:A:1088:G:C8	2.68	0.43
1:A:1129:C:H4'	1:A:1130:A:C8	2.46	0.43
1:A:1298:C:H5''	1:A:1299:A:H5''	2.00	0.43
2:B:24:TRP:CB	2:B:190:THR:HG23	2.49	0.43
2:B:55:PHE:O	2:B:56:ARG:C	2.57	0.43
2:B:73:THR:CG2	2:B:73:THR:O	2.66	0.43
2:B:92:TYR:CE1	2:B:151:GLY:N	2.87	0.43
3:C:48:TYR:C	3:C:48:TYR:CD1	2.90	0.43
4:D:94:LEU:HD22	4:D:196:LEU:HD12	2.01	0.43
4:D:98:GLU:HG2	4:D:189:PRO:HG3	2.01	0.43
4:D:110:PHE:HA	4:D:162:LEU:HD11	2.01	0.43
7:G:61:VAL:O	7:G:62:PHE:C	2.57	0.43
8:H:96:GLY:C	8:H:97:VAL:CG1	2.86	0.43
13:M:29:ARG:HB3	13:M:64:TRP:CH2	2.53	0.43
15:O:45:VAL:HB	15:O:46:HIS:H	1.58	0.43
17:Q:95:TYR:HB3	17:Q:96:GLN:H	1.44	0.43
20:T:79:ARG:O	20:T:83:ARG:HB2	2.19	0.43
21:V:18:TYR:HA	21:V:22:ARG:HB3	2.01	0.43
1:A:37:U:H2'	1:A:38:G:H5'	2.00	0.43
1:A:102:G:H2'	1:A:103:C:C6	2.48	0.43
1:A:127:G:C2	1:A:128:G:C8	3.07	0.43
1:A:132:C:C4	1:A:133:U:C5	3.07	0.43
1:A:234:C:H2'	1:A:235:C:H6	1.84	0.43
1:A:854:G:C6	1:A:855:G:N7	2.87	0.43
1:A:881:G:P	12:L:12:ARG:HH22	2.41	0.43
1:A:915:A:O2'	1:A:916:G:H5'	2.18	0.43
1:A:1127:G:N2	1:A:1147:C:H42	2.16	0.43
1:A:1257:U:H5''	1:A:1258:G:O5'	2.18	0.43
1:A:1455:G:O5'	1:A:1455:G:C8	2.71	0.43
2:B:9:GLU:CD	2:B:217:ARG:HH12	2.22	0.43
2:B:91:PRO:HG3	2:B:154:LEU:HD12	2.01	0.43
2:B:108:ILE:O	2:B:108:ILE:CG1	2.67	0.43
2:B:126:GLU:HG2	2:B:129:GLU:HB2	2.00	0.43
3:C:102:ASN:HD22	3:C:102:ASN:H	1.67	0.43
3:C:111:LEU:HD21	3:C:146:ALA:H	1.84	0.43
3:C:122:GLU:C	3:C:124:ILE:N	2.71	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:137:ALA:O	3:C:139:GLN:N	2.51	0.43
3:C:155:GLY:O	3:C:156:ARG:O	2.36	0.43
4:D:156:GLU:CD	4:D:157:LEU:H	2.22	0.43
9:I:23:ASN:HB3	9:I:60:ASP:OD1	2.19	0.43
9:I:86:VAL:HA	9:I:89:ASN:C	2.39	0.43
10:J:18:ALA:O	10:J:20:ALA:N	2.52	0.43
10:J:19:SER:OG	10:J:91:PRO:CB	2.67	0.43
10:J:38:ILE:HA	10:J:39:PRO:HD3	1.90	0.43
11:K:62:GLN:O	11:K:63:LEU:C	2.57	0.43
12:L:83:VAL:CG2	12:L:100:ILE:HD13	2.45	0.43
17:Q:90:ILE:HA	17:Q:93:GLN:HB3	2.01	0.43
19:S:22:LEU:CD2	19:S:28:LYS:HB2	2.49	0.43
19:S:34:TRP:HE3	19:S:34:TRP:N	2.17	0.43
20:T:40:ALA:HA	20:T:55:ILE:CD1	2.30	0.43
1:A:593:G:C6	1:A:594:G:C5	3.06	0.42
1:A:680:C:O2'	1:A:681:C:H5'	2.19	0.42
1:A:1003(A):G:C2	1:A:1004:A:C4	3.07	0.42
1:A:1406:U:H5'	1:A:1518:A:O2'	2.19	0.42
1:A:1475:G:H2'	1:A:1476:G:O4'	2.19	0.42
1:A:1503:A:C4	1:A:1531:A:C2	3.07	0.42
2:B:17:PHE:CD1	2:B:17:PHE:C	2.92	0.42
2:B:58:ILE:O	2:B:60:ASP:N	2.52	0.42
3:C:140:ARG:O	3:C:141:VAL:C	2.57	0.42
3:C:179:ARG:HD2	3:C:180:ALA:HA	2.01	0.42
4:D:49:ARG:HG2	4:D:49:ARG:NH1	2.34	0.42
4:D:57:ARG:NH2	4:D:205:GLU:OE2	2.44	0.42
4:D:62:GLN:O	4:D:63:LYS:C	2.58	0.42
5:E:12:LEU:C	5:E:12:LEU:HD22	2.39	0.42
5:E:28:PHE:CD1	5:E:50:GLU:HA	2.53	0.42
6:F:30:LEU:C	6:F:32:ASN:H	2.22	0.42
10:J:7:LYS:HB2	10:J:97:GLU:HB2	2.00	0.42
10:J:47:PHE:HB2	14:N:44:LEU:HD21	1.99	0.42
11:K:48:ILE:N	11:K:48:ILE:HD13	2.34	0.42
16:P:39:TYR:CD2	16:P:39:TYR:C	2.92	0.42
17:Q:82:MET:O	17:Q:86:GLU:N	2.35	0.42
19:S:12:ASP:CB	19:S:38:SER:HB2	2.33	0.42
19:S:40:ILE:HB	19:S:67:VAL:O	2.19	0.42
20:T:14:LYS:O	20:T:15:ARG:C	2.57	0.42
20:T:51:GLU:CA	20:T:54:LYS:HB3	2.49	0.42
1:A:36:C:O5'	1:A:36:C:H6	2.03	0.42
1:A:363:A:N6	1:A:364:A:N1	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:664:G:N2	1:A:742:G:C2	2.87	0.42
1:A:789:U:H3	1:A:792:A:P	2.42	0.42
1:A:945:G:H2'	1:A:945:G:N3	2.34	0.42
1:A:988:G:O2'	1:A:989:C:H5'	2.19	0.42
1:A:1152:A:C4	1:A:1153:C:C5	3.07	0.42
1:A:1249:C:C6	1:A:1249:C:C3'	3.02	0.42
1:A:1432:G:H1'	1:A:1468:A:N6	2.33	0.42
1:A:1459:C:C2	1:A:1460:A:C8	3.06	0.42
4:D:28:SER:O	4:D:30:LYS:HB2	2.18	0.42
4:D:148:VAL:HG21	4:D:181:MET:HB3	2.00	0.42
4:D:201:GLN:HE22	5:E:99:GLY:HA3	1.83	0.42
7:G:37:ASN:O	7:G:40:ALA:HB3	2.19	0.42
7:G:38:LEU:C	7:G:42:ILE:HG13	2.37	0.42
8:H:104:ARG:HD3	8:H:104:ARG:HA	1.75	0.42
10:J:3:LYS:CG	10:J:76:ASN:HA	2.49	0.42
10:J:31:GLY:CA	10:J:81:THR:OG1	2.67	0.42
12:L:103:GLY:H	12:L:108:ALA:N	2.17	0.42
14:N:20:ALA:O	14:N:22:THR:N	2.52	0.42
15:O:36:ILE:O	15:O:37:ASN:C	2.58	0.42
16:P:56:ALA:O	16:P:60:LEU:N	2.52	0.42
17:Q:90:ILE:HA	17:Q:93:GLN:HB2	2.00	0.42
1:A:123:C:H4'	1:A:290:C:O2	2.19	0.42
1:A:321:A:C2'	1:A:322:C:H6	2.32	0.42
1:A:498:U:O2	1:A:498:U:C2'	2.67	0.42
1:A:597:G:N2	8:H:94:TYR:HE2	2.17	0.42
1:A:818:G:N3	1:A:820:U:C5	2.87	0.42
1:A:868:C:C2'	1:A:869:G:H5'	2.49	0.42
1:A:908:A:O2'	1:A:909:A:H5'	2.19	0.42
1:A:944:G:H3'	1:A:945:G:C5'	2.49	0.42
1:A:1022:G:C2	1:A:1023:G:N7	2.87	0.42
1:A:1059:C:H42	1:A:1198:G:H1	1.65	0.42
1:A:1200:C:O2	1:A:1200:C:C2'	2.60	0.42
1:A:1287:A:N7	1:A:1288:A:N6	2.67	0.42
1:A:1299:A:C8	1:A:1301:U:O2	2.72	0.42
1:A:1313:U:H5	19:S:4:SER:OG	2.01	0.42
1:A:1358:U:OP1	14:N:35:ARG:HB2	2.19	0.42
1:A:1520:G:C2	1:A:1521:G:C5	3.07	0.42
2:B:21:ARG:HA	2:B:39:ILE:HG12	2.01	0.42
2:B:55:PHE:O	2:B:57:PHE:N	2.52	0.42
2:B:82:ARG:O	2:B:92:TYR:HE2	2.02	0.42
2:B:208:ILE:CA	2:B:211:ILE:HG12	2.48	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:26:LYS:O	3:C:30:ARG:NH1	2.53	0.42
3:C:83:ARG:CA	3:C:86:VAL:HG23	2.33	0.42
4:D:57:ARG:CD	4:D:202:LEU:HD22	2.49	0.42
4:D:200:GLU:HG2	4:D:201:GLN:N	2.34	0.42
5:E:109:ILE:HD13	5:E:109:ILE:N	2.34	0.42
7:G:135:VAL:CG1	7:G:138:LYS:HZ2	2.10	0.42
9:I:97:LYS:HZ1	9:I:102:LEU:HD11	1.84	0.42
11:K:53:SER:C	11:K:55:LYS:N	2.72	0.42
12:L:24:VAL:C	12:L:26:ALA:N	2.70	0.42
12:L:58:VAL:CG1	12:L:59:ARG:N	2.81	0.42
13:M:14:ARG:HH12	13:M:16:ASP:CG	2.23	0.42
13:M:32:GLU:HG2	13:M:32:GLU:O	2.19	0.42
15:O:4:THR:N	15:O:7:GLU:OE1	2.46	0.42
16:P:39:TYR:HD1	16:P:73:LEU:HD22	1.83	0.42
17:Q:45:HIS:NE2	17:Q:47:PRO:HG3	2.33	0.42
17:Q:46:ASP:C	17:Q:46:ASP:OD2	2.58	0.42
17:Q:76:LEU:C	17:Q:76:LEU:CD1	2.83	0.42
18:R:38:GLU:CA	18:R:41:LYS:HE3	2.49	0.42
19:S:22:LEU:HB3	19:S:28:LYS:N	2.33	0.42
20:T:84:LEU:HD22	20:T:88:VAL:CG2	2.50	0.42
1:A:37:U:O2'	1:A:38:G:H5'	2.19	0.42
1:A:255:G:N2	1:A:272:C:C1'	2.81	0.42
1:A:433:C:C2	1:A:434:U:C5	3.07	0.42
1:A:451:A:H1'	1:A:452:A:N7	2.33	0.42
1:A:525:C:H2'	1:A:526:C:H6	1.81	0.42
1:A:581:G:OP1	15:O:65:ARG:NH2	2.27	0.42
1:A:582:U:C2	1:A:760:G:C6	3.07	0.42
1:A:951:G:C5	1:A:952:U:C5	3.07	0.42
1:A:981:U:O4	1:A:1222:G:O6	2.37	0.42
1:A:993:G:H3'	1:A:993:G:N3	2.34	0.42
1:A:1013:G:O2'	1:A:1014:A:C8	2.72	0.42
1:A:1120:G:N1	1:A:1121:U:C4	2.87	0.42
1:A:1128:C:N3	1:A:1144:G:N2	2.67	0.42
1:A:1286:A:H2'	1:A:1287:A:C4'	2.35	0.42
1:A:1310:G:C6	19:S:2:PRO:HD3	2.55	0.42
1:A:1343:G:H2'	1:A:1344:C:C5	2.48	0.42
1:A:1507:A:C5	1:A:1508:G:N7	2.87	0.42
2:B:10:LEU:C	2:B:12:GLU:N	2.73	0.42
2:B:164:VAL:HG12	2:B:165:VAL:H	1.84	0.42
3:C:10:PHE:CE2	3:C:178:LEU:HD13	2.54	0.42
3:C:23:TYR:HD1	10:J:11:PHE:CZ	2.37	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:184:TYR:OH	3:C:186:PHE:HB2	2.18	0.42
5:E:41:VAL:HG12	5:E:42:GLY:O	2.20	0.42
5:E:74:GLY:O	5:E:116:THR:HG22	2.19	0.42
5:E:90:VAL:HG22	5:E:121:LYS:O	2.19	0.42
6:F:16:GLN:NE2	6:F:16:GLN:C	2.73	0.42
7:G:17:VAL:HG12	7:G:18:TYR:HD1	1.84	0.42
8:H:1:MET:HG2	8:H:2:LEU:O	2.18	0.42
11:K:31:THR:OG1	11:K:42:TRP:HB3	2.19	0.42
12:L:109:GLY:CA	12:L:121:GLY:O	2.67	0.42
13:M:3:ARG:HD3	13:M:7:VAL:HA	2.01	0.42
13:M:11:ARG:O	13:M:12:ASN:CB	2.66	0.42
13:M:59:TYR:O	13:M:63:THR:HB	2.18	0.42
13:M:95:GLY:C	13:M:96:LEU:HD12	2.39	0.42
16:P:10:GLY:CA	16:P:16:HIS:H	2.32	0.42
17:Q:13:ASP:O	17:Q:14:LYS:HB2	2.18	0.42
17:Q:76:LEU:HD11	17:Q:78:GLU:C	2.39	0.42
1:A:187:C:H1'	20:T:104:LEU:CD2	2.49	0.42
1:A:440:A:C6	1:A:495:U:N3	2.88	0.42
1:A:485:G:C2'	1:A:486:U:OP2	2.68	0.42
1:A:500:G:C6	1:A:546:G:C2	3.07	0.42
1:A:951:G:C6	1:A:1231:G:C6	3.08	0.42
1:A:988:G:H1'	1:A:1015:A:N1	2.35	0.42
1:A:994:A:N3	1:A:994:A:C2'	2.79	0.42
1:A:1151:A:C5'	10:J:42:THR:H	2.31	0.42
1:A:1239:A:H61	1:A:1296:C:H2'	1.84	0.42
1:A:1241:G:C4	1:A:1242:C:C5	3.08	0.42
1:A:1261:A:H5'	1:A:1284:C:OP1	2.19	0.42
1:A:1371:G:C5	1:A:1372:U:C5	3.07	0.42
1:A:1428:A:C6	1:A:1473:A:N6	2.88	0.42
2:B:25:ASN:ND2	2:B:27:LYS:HG3	2.35	0.42
2:B:131:PRO:O	2:B:134:GLU:HG3	2.19	0.42
4:D:54:TYR:O	4:D:55:ALA:C	2.58	0.42
4:D:124:GLY:HA3	4:D:132:ARG:HE	1.81	0.42
4:D:162:LEU:HD12	4:D:162:LEU:N	2.35	0.42
5:E:110:LEU:HA	5:E:110:LEU:HD23	1.81	0.42
5:E:146:ALA:O	5:E:149:GLU:N	2.52	0.42
7:G:118:VAL:O	7:G:119:ARG:C	2.58	0.42
7:G:153:HIS:CD2	7:G:154:TYR:CZ	3.07	0.42
8:H:1:MET:HG2	8:H:2:LEU:N	2.35	0.42
8:H:105:ARG:HG3	8:H:105:ARG:O	2.20	0.42
9:I:43:ALA:HB2	9:I:74:ILE:HD13	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:74:ILE:O	9:I:74:ILE:HG22	2.19	0.42
11:K:63:LEU:O	11:K:64:ALA:C	2.57	0.42
12:L:42:THR:HG23	12:L:53:ARG:C	2.40	0.42
13:M:60:VAL:HG13	13:M:64:TRP:CE3	2.54	0.42
13:M:96:LEU:O	13:M:110:ARG:CZ	2.68	0.42
15:O:56:LEU:O	15:O:59:MET:CB	2.68	0.42
15:O:61:GLY:O	15:O:62:GLN:C	2.57	0.42
15:O:78:TYR:OH	15:O:82:ILE:HD12	2.19	0.42
16:P:72:ARG:O	16:P:75:ARG:N	2.50	0.42
17:Q:94:ASN:N	17:Q:94:ASN:ND2	2.67	0.42
19:S:53:ASN:ND2	19:S:53:ASN:O	2.52	0.42
20:T:52:ALA:O	20:T:55:ILE:HG12	2.19	0.42
1:A:79:G:C2	1:A:91:C:N3	2.88	0.42
1:A:117:G:O2'	1:A:118:U:H5'	2.18	0.42
1:A:120:A:C6	1:A:122:G:C2	3.07	0.42
1:A:401:C:P	4:D:73:ARG:NH2	2.93	0.42
1:A:508:C:P	4:D:209:ARG:NH2	2.93	0.42
1:A:635:G:C4	1:A:636:U:C5	3.07	0.42
1:A:649:G:N2	1:A:650:G:C4	2.88	0.42
1:A:781:A:H4'	1:A:1522:U:O2'	2.20	0.42
1:A:956:U:H2'	1:A:957:U:C5'	2.49	0.42
1:A:1114:C:H2'	1:A:1115:C:C5'	2.50	0.42
1:A:1419:G:H3'	1:A:1420:C:C6	2.54	0.42
1:A:1437:C:H2'	1:A:1438:G:H8	1.83	0.42
2:B:19:HIS:CD2	2:B:189:ASP:OD1	2.73	0.42
2:B:136:VAL:HA	2:B:139:LYS:CB	2.50	0.42
2:B:203:GLY:O	2:B:204:ASN:C	2.57	0.42
2:B:214:ILE:HG22	2:B:215:LEU:N	2.35	0.42
2:B:235:SER:C	2:B:237:ALA:H	2.21	0.42
3:C:45:LYS:HD3	3:C:46:GLU:OE2	2.19	0.42
3:C:47:LEU:HD11	3:C:76:VAL:HG13	1.97	0.42
3:C:116:VAL:O	3:C:119:ARG:N	2.52	0.42
3:C:118:GLN:O	3:C:122:GLU:N	2.42	0.42
3:C:120:VAL:O	3:C:124:ILE:CB	2.61	0.42
4:D:59:ARG:O	4:D:60:GLU:C	2.58	0.42
4:D:63:LYS:C	4:D:65:ARG:N	2.73	0.42
4:D:136:PRO:O	4:D:138:TYR:N	2.45	0.42
5:E:12:LEU:HD13	5:E:12:LEU:N	2.35	0.42
5:E:135:THR:C	5:E:137:GLU:N	2.72	0.42
5:E:141:GLN:O	5:E:142:LEU:O	2.37	0.42
6:F:45:LEU:O	6:F:46:ARG:CG	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:58:PRO:HA	7:G:61:VAL:HG23	2.01	0.42
9:I:43:ALA:HA	9:I:74:ILE:CG2	2.46	0.42
9:I:97:LYS:C	9:I:99:LEU:N	2.72	0.42
10:J:49:VAL:HG12	10:J:50:ILE:H	1.83	0.42
10:J:56:HIS:O	10:J:58:ASP:N	2.53	0.42
12:L:74:GLY:H	12:L:110:VAL:CG1	2.32	0.42
13:M:59:TYR:O	13:M:63:THR:CG2	2.67	0.42
15:O:34:LEU:O	15:O:35:ARG:C	2.58	0.42
18:R:38:GLU:HA	18:R:41:LYS:CE	2.49	0.42
18:R:50:ILE:CG2	18:R:51:LEU:N	2.83	0.42
20:T:67:ALA:HB1	20:T:73:HIS:HA	2.02	0.42
21:V:5:ASP:O	21:V:6:ARG:C	2.57	0.42
1:A:154:C:H2'	1:A:155:C:H6	1.84	0.42
1:A:185:A:C4	1:A:186:C:C5	3.07	0.42
1:A:302:G:N3	1:A:556:C:H4'	2.35	0.42
1:A:310:G:N2	1:A:311:C:C2	2.87	0.42
1:A:356:A:C4	1:A:357:G:C8	3.08	0.42
1:A:675:A:H1'	11:K:116:HIS:CG	2.55	0.42
1:A:701:C:H4'	1:A:702:A:OP2	2.18	0.42
1:A:803:G:C6	1:A:804:U:C4	3.07	0.42
1:A:821:G:H2'	1:A:822:C:H6	1.84	0.42
1:A:850:U:O2'	1:A:851:G:H5'	2.19	0.42
1:A:1013:G:HO2'	1:A:1014:A:H8	1.66	0.42
1:A:1074:G:N2	1:A:1102:A:C8	2.88	0.42
1:A:1086:U:H2'	1:A:1087:G:O4'	2.20	0.42
1:A:1327:C:O2'	1:A:1328:C:H5'	2.19	0.42
1:A:1343:G:C6	1:A:1344:C:C4	3.08	0.42
1:A:1372:U:H2'	1:A:1373:G:C5'	2.49	0.42
1:A:1403:C:O2	1:A:1499:A:N1	2.53	0.42
2:B:140:HIS:O	2:B:141:GLU:C	2.57	0.42
2:B:239:VAL:CG1	2:B:240:GLN:H	2.21	0.42
3:C:5:ILE:O	3:C:6:HIS:HB2	2.20	0.42
3:C:21:ARG:NH1	3:C:56:ASP:HB3	2.34	0.42
3:C:82:GLU:HA	3:C:82:GLU:OE1	2.19	0.42
3:C:131:ARG:O	3:C:135:LYS:N	2.45	0.42
4:D:6:GLY:H	4:D:115:ARG:NH2	2.15	0.42
4:D:61:LYS:CD	4:D:62:GLN:N	2.73	0.42
4:D:200:GLU:OE1	4:D:200:GLU:N	2.46	0.42
5:E:82:VAL:HG11	5:E:134:ALA:O	2.20	0.42
7:G:92:SER:CB	7:G:93:PRO:HD2	2.46	0.42
9:I:10:ARG:NH2	9:I:11:LYS:HB2	2.33	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:99:LEU:HB3	9:I:101:PHE:CE1	2.54	0.42
10:J:8:LEU:CD2	10:J:20:ALA:HB2	2.50	0.42
11:K:69:ALA:HB1	11:K:101:SER:HB2	2.02	0.42
12:L:97:ARG:C	12:L:98:TYR:CG	2.93	0.42
13:M:17:VAL:O	13:M:19:LEU:N	2.52	0.42
13:M:102:ARG:NH1	13:M:106:ASN:HD22	2.17	0.42
13:M:110:ARG:HG3	13:M:110:ARG:O	2.19	0.42
15:O:54:ARG:O	15:O:55:GLY:C	2.57	0.42
16:P:55:ARG:O	16:P:56:ALA:C	2.58	0.42
19:S:15:LEU:HD11	19:S:38:SER:HG	1.84	0.42
20:T:63:ILE:O	20:T:66:ALA:HB3	2.20	0.42
1:A:191:G:O2'	20:T:102:GLY:O	2.33	0.42
1:A:355:C:H1'	1:A:388:G:H1'	2.02	0.42
1:A:391:G:C2'	1:A:392:G:O5'	2.68	0.42
1:A:508:C:O2	1:A:508:C:O4'	2.36	0.42
1:A:525:C:OP1	12:L:91:LYS:HD3	2.19	0.42
1:A:528:C:O2'	1:A:529:G:H5'	2.19	0.42
1:A:686:U:O2	1:A:687:A:C5	2.73	0.42
1:A:718:G:C4'	11:K:117:ASN:HD21	2.32	0.42
1:A:737:A:OP1	6:F:91:VAL:HG13	2.19	0.42
1:A:865:A:C6	1:A:866:C:N4	2.88	0.42
1:A:1074:G:N2	1:A:1102:A:C5	2.87	0.42
1:A:1095:U:H5''	1:A:1109:C:O2	2.20	0.42
1:A:1173:G:C5	1:A:1174:G:C5	3.08	0.42
1:A:1178:G:H8	1:A:1178:G:O5'	2.03	0.42
1:A:1399:C:C4	1:A:1401:G:C2	3.07	0.42
1:A:1502:A:H2	1:A:1505:G:N1	2.18	0.42
1:A:1507:A:H2'	1:A:1508:G:H8	1.84	0.42
1:A:1518:A:C8	1:A:1518:A:C3'	3.02	0.42
2:B:57:PHE:CZ	2:B:61:LEU:CD1	3.03	0.42
2:B:134:GLU:C	2:B:136:VAL:N	2.72	0.42
3:C:19:GLU:O	3:C:56:ASP:HA	2.20	0.42
3:C:175:LEU:O	3:C:176:HIS:C	2.58	0.42
3:C:179:ARG:HH11	3:C:179:ARG:HG3	1.85	0.42
4:D:165:MET:O	4:D:166:LYS:HG3	2.20	0.42
5:E:105:VAL:HG11	5:E:132:ALA:N	2.34	0.42
7:G:54:THR:HB	7:G:56:GLN:CD	2.40	0.42
8:H:4:ASP:CG	8:H:85:ARG:NH1	2.73	0.42
9:I:40:LEU:HD23	9:I:40:LEU:H	1.85	0.42
10:J:9:ARG:H	10:J:9:ARG:CD	2.26	0.42
10:J:70:ARG:HH11	10:J:70:ARG:CG	2.32	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:J:79:ARG:HH11	10:J:79:ARG:CB	2.32	0.42
12:L:38:THR:HG22	12:L:39:VAL:N	2.35	0.42
13:M:56:LEU:O	13:M:60:VAL:N	2.38	0.42
14:N:28:GLY:C	14:N:29:ARG:O	2.58	0.42
17:Q:98:LEU:HA	17:Q:102:GLY:N	2.34	0.42
19:S:34:TRP:CD1	19:S:52:TYR:HB2	2.55	0.42
1:A:51:A:H1'	1:A:52:G:OP2	2.20	0.42
1:A:170:U:O2'	1:A:171:A:C5'	2.64	0.42
1:A:184:G:C6	1:A:185:A:N7	2.88	0.42
1:A:186:C:C5'	20:T:78:ALA:HB1	2.49	0.42
1:A:410:G:OP1	4:D:30:LYS:HE3	2.19	0.42
1:A:666:G:N2	1:A:741:G:H1'	2.35	0.42
1:A:746:A:C2'	1:A:747:C:H5'	2.50	0.42
1:A:916:G:H2'	1:A:917:G:H8	1.84	0.42
1:A:1177:G:C5	1:A:1178:G:N7	2.88	0.42
1:A:1224:G:H4'	1:A:1225:A:OP1	2.20	0.42
3:C:112:SER:OG	3:C:115:LEU:HG	2.20	0.42
4:D:77:ASN:O	4:D:81:GLU:HG3	2.19	0.42
4:D:130:GLY:C	4:D:132:ARG:N	2.73	0.42
4:D:134:ASP:OD1	4:D:135:LEU:HD23	2.20	0.42
5:E:19:MET:HE3	5:E:20:GLN:N	2.31	0.42
5:E:48:ALA:C	5:E:50:GLU:H	2.23	0.42
5:E:145:LYS:O	5:E:149:GLU:HG3	2.20	0.42
7:G:26:PHE:C	7:G:28:ASN:N	2.74	0.42
7:G:59:LEU:O	7:G:63:LYS:CG	2.68	0.42
8:H:32:LYS:O	8:H:36:LEU:HG	2.19	0.42
8:H:100:ILE:HG22	8:H:125:ARG:HH21	1.83	0.42
9:I:38:GLN:O	9:I:40:LEU:N	2.53	0.42
10:J:8:LEU:CD1	10:J:8:LEU:O	2.68	0.42
10:J:27:ALA:HB3	10:J:74:ILE:HG22	2.01	0.42
11:K:109:VAL:CG1	18:R:84:LYS:HB3	2.50	0.42
11:K:114:VAL:HA	11:K:115:PRO:HD3	1.87	0.42
13:M:94:ARG:HG2	13:M:94:ARG:NH1	2.35	0.42
19:S:31:ILE:CG2	19:S:32:LYS:N	2.83	0.42
20:T:62:LEU:CA	20:T:65:LYS:HB3	2.50	0.42
1:A:29:G:H8	1:A:29:G:O5'	2.03	0.42
1:A:38:G:O2'	1:A:39:G:H5''	2.20	0.42
1:A:67:C:H4'	1:A:172:A:O4'	2.19	0.42
1:A:145:G:H2'	1:A:146:G:C8	2.53	0.42
1:A:313:A:N6	1:A:314:C:N4	2.68	0.42
1:A:520:A:N6	1:A:529:G:H1'	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:545:C:C2'	1:A:546:G:H5'	2.50	0.42
1:A:715:A:H2'	1:A:716:A:H8	1.84	0.42
1:A:810:C:O2	1:A:810:C:H2'	2.20	0.42
1:A:830:G:H2'	1:A:831:U:O4'	2.20	0.42
1:A:882:C:H2'	1:A:883:C:H6	1.85	0.42
1:A:941:G:C6	1:A:942:G:N7	2.88	0.42
1:A:1075:C:H2'	1:A:1076:C:H6	1.85	0.42
1:A:1104:G:H4'	2:B:111:ARG:HH22	1.76	0.42
1:A:1240:U:C2	7:G:32:ARG:HD2	2.54	0.42
1:A:1305:G:O2'	1:A:1306:A:C8	2.61	0.42
1:A:1368:G:O2'	1:A:1369:C:H5'	2.20	0.42
2:B:112:VAL:HG22	2:B:149:LEU:HD13	2.02	0.42
4:D:13:ARG:HB3	4:D:40:PRO:HD3	2.02	0.42
4:D:61:LYS:NZ	4:D:207:TYR:OH	2.50	0.42
5:E:127:ASN:OD1	5:E:129:ILE:HB	2.20	0.42
7:G:75:VAL:HA	7:G:87:VAL:O	2.20	0.42
7:G:121:ALA:HA	7:G:124:LEU:HD12	2.02	0.42
9:I:23:ASN:H	9:I:60:ASP:N	2.16	0.42
9:I:31:GLN:HB3	9:I:35:GLU:CB	2.48	0.42
9:I:46:ALA:HB1	9:I:81:ILE:HD12	2.02	0.42
11:K:128:ALA:O	11:K:129:SER:HB2	2.20	0.42
16:P:8:ARG:HE	16:P:15:PRO:HB3	1.85	0.42
16:P:58:TYR:HA	16:P:61:SER:CB	2.50	0.42
18:R:59:SER:O	18:R:63:GLN:N	2.43	0.42
19:S:61:TYR:C	19:S:61:TYR:CD2	2.93	0.42
20:T:23:ARG:O	20:T:25:ARG:N	2.53	0.42
1:A:17:U:N3	1:A:919:A:H2	2.18	0.41
1:A:102:G:N3	1:A:151:A:H2	2.18	0.41
1:A:223:U:C5'	20:T:68:LYS:HZ1	2.33	0.41
1:A:268:C:C2'	1:A:269:C:O5'	2.68	0.41
1:A:308:C:O2	1:A:309:G:C8	2.73	0.41
1:A:345:C:O4'	1:A:346:G:N2	2.52	0.41
1:A:391:G:H2'	1:A:392:G:O5'	2.20	0.41
1:A:792:A:C6	1:A:794:A:C6	3.08	0.41
1:A:1117:G:O2'	9:I:104:ARG:HD3	2.20	0.41
1:A:1271:G:H5'	1:A:1314:C:OP1	2.19	0.41
1:A:1281:U:HO2'	1:A:1282:C:P	2.42	0.41
1:A:1346:A:H4'	1:A:1347:G:O5'	2.20	0.41
2:B:91:PRO:HG2	2:B:155:LEU:HD23	1.99	0.41
2:B:97:TRP:HH2	2:B:102:LEU:H	1.67	0.41
2:B:119:GLU:OE1	2:B:153:ARG:NH2	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:91:LEU:HD21	3:C:98:ASN:HD22	1.85	0.41
3:C:175:LEU:CD2	3:C:175:LEU:H	2.33	0.41
3:C:201:TYR:C	3:C:202:ILE:HG13	2.40	0.41
6:F:9:VAL:CG2	6:F:87:ARG:HB2	2.50	0.41
7:G:14:PRO:O	7:G:15:ASP:O	2.38	0.41
7:G:141:VAL:O	7:G:145:ALA:N	2.49	0.41
8:H:18:ARG:HA	8:H:18:ARG:HD3	1.83	0.41
14:N:6:LEU:HD22	14:N:23:ARG:CZ	2.49	0.41
18:R:22:VAL:HG13	18:R:26:LEU:CD1	2.50	0.41
20:T:10:LEU:O	20:T:11:SER:C	2.58	0.41
20:T:43:LEU:O	20:T:44:ALA:C	2.58	0.41
20:T:82:SER:O	20:T:84:LEU:N	2.52	0.41
1:A:202:U:O2	1:A:202:U:H2'	2.20	0.41
1:A:377:G:O2'	1:A:378:G:H5'	2.20	0.41
1:A:409:G:O2'	1:A:410:G:H5'	2.20	0.41
1:A:451:A:C6	1:A:481:G:C4	3.08	0.41
1:A:455:C:H42	1:A:477:G:H1	1.66	0.41
1:A:491:G:C2	1:A:492:G:C5	3.08	0.41
1:A:507:C:C4	1:A:508:C:C5	3.09	0.41
1:A:575:G:C4	1:A:881:G:C2	3.08	0.41
1:A:707:C:C2	1:A:708:C:C5	3.08	0.41
1:A:832:C:C2'	1:A:833:U:H5'	2.50	0.41
1:A:967:C:O2'	9:I:128:ARG:NE	2.52	0.41
1:A:982:U:H1'	1:A:983:A:C5	2.55	0.41
1:A:1319:A:N6	1:A:1361:G:N2	2.60	0.41
1:A:1495:U:C4	1:A:1496:C:N4	2.88	0.41
2:B:24:TRP:CA	2:B:190:THR:HG23	2.50	0.41
2:B:104:ASN:CG	2:B:104:ASN:O	2.59	0.41
2:B:129:GLU:CD	2:B:130:ARG:H	2.23	0.41
2:B:182:ILE:O	2:B:183:PRO:O	2.37	0.41
2:B:193:ASP:C	2:B:193:ASP:OD1	2.59	0.41
3:C:22:TRP:CB	3:C:59:ARG:HG2	2.50	0.41
3:C:42:LEU:O	3:C:45:LYS:N	2.53	0.41
3:C:206:GLU:O	3:C:207:VAL:CB	2.68	0.41
4:D:22:LYS:HB3	4:D:22:LYS:HE2	1.56	0.41
4:D:66:ARG:O	4:D:67:ILE:C	2.56	0.41
4:D:91:SER:O	4:D:93:PHE:N	2.53	0.41
4:D:134:ASP:CG	4:D:135:LEU:N	2.73	0.41
4:D:192:GLU:O	4:D:194:LEU:N	2.53	0.41
5:E:71:LEU:CD1	5:E:114:GLY:O	2.67	0.41
6:F:63:TYR:HD1	6:F:63:TYR:N	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:17:VAL:CG1	7:G:18:TYR:N	2.61	0.41
9:I:118:LYS:CB	9:I:118:LYS:HZ2	2.31	0.41
10:J:26:ALA:O	10:J:29:ARG:HG2	2.19	0.41
11:K:33:THR:OG1	11:K:38:ASN:O	2.38	0.41
11:K:86:GLY:CA	11:K:112:THR:HG23	2.49	0.41
13:M:78:ILE:C	13:M:80:ARG:H	2.23	0.41
13:M:109:THR:CG2	13:M:110:ARG:N	2.83	0.41
14:N:27:CYS:SG	14:N:29:ARG:HB3	2.60	0.41
16:P:8:ARG:HH21	16:P:15:PRO:HD3	1.84	0.41
16:P:22:THR:CG2	16:P:23:ASP:H	2.12	0.41
17:Q:97:SER:CB	17:Q:103:GLY:CA	2.89	0.41
18:R:31:LEU:HD21	18:R:66:LEU:H	1.85	0.41
20:T:51:GLU:C	20:T:54:LYS:HB3	2.40	0.41
20:T:54:LYS:N	20:T:100:ILE:HD12	2.35	0.41
21:V:22:ARG:HG2	21:V:22:ARG:HH11	1.85	0.41
1:A:29:G:N2	1:A:555:C:C4	2.88	0.41
1:A:110:C:H2'	1:A:111:G:H5'	2.02	0.41
1:A:250:A:C4'	1:A:251:G:O5'	2.61	0.41
1:A:345:C:H5'	1:A:346:G:C4	2.55	0.41
1:A:355:C:H5'	1:A:389:A:OP2	2.21	0.41
1:A:451:A:N6	1:A:481:G:H1'	2.36	0.41
1:A:584:G:C2	1:A:585:G:C4	3.09	0.41
1:A:740:U:O4'	15:O:42:HIS:ND1	2.53	0.41
1:A:837:G:C2	1:A:850:U:O2	2.74	0.41
1:A:932:C:C6	7:G:3:ARG:HD3	2.54	0.41
1:A:938:A:C8	1:A:938:A:O5'	2.73	0.41
1:A:941:G:C5	1:A:942:G:N7	2.89	0.41
1:A:949:A:C6	1:A:1233:G:C2	3.08	0.41
1:A:1015:A:H1'	1:A:1218:C:O2'	2.20	0.41
1:A:1038:C:H2'	1:A:1039:C:C5	2.54	0.41
1:A:1061:G:C6	1:A:1062:U:C4	3.09	0.41
1:A:1243:C:OP2	21:V:10:ARG:CZ	2.67	0.41
1:A:1276:G:N2	1:A:1282:C:O2	2.54	0.41
1:A:1369:C:H2'	1:A:1370:G:H8	1.83	0.41
1:A:1436:U:C2'	1:A:1437:C:H5'	2.49	0.41
3:C:95:THR:HG23	3:C:98:ASN:H	1.84	0.41
3:C:201:TYR:N	3:C:201:TYR:HD1	2.17	0.41
4:D:125:HIS:HA	4:D:149:ALA:HB3	2.01	0.41
4:D:127:THR:HG23	4:D:127:THR:O	2.20	0.41
4:D:157:LEU:O	4:D:160:GLN:N	2.54	0.41
5:E:63:ARG:O	5:E:66:MET:HG2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:143:ARG:NH1	8:H:77:GLU:CD	2.73	0.41
9:I:118:LYS:HB3	9:I:118:LYS:HZ3	1.84	0.41
12:L:7:ILE:O	12:L:10:LEU:CB	2.68	0.41
13:M:34:LEU:HD11	13:M:41:PRO:HA	2.02	0.41
14:N:9:LYS:O	14:N:9:LYS:CD	2.68	0.41
15:O:74:ASP:O	15:O:74:ASP:OD1	2.38	0.41
18:R:30:ASP:C	18:R:32:ARG:N	2.73	0.41
20:T:67:ALA:O	20:T:73:HIS:ND1	2.53	0.41
1:A:51:A:C2	1:A:116:A:H1'	2.55	0.41
1:A:259:G:H2'	1:A:260:G:H8	1.86	0.41
1:A:273:A:H2'	1:A:274:A:C5'	2.50	0.41
1:A:322:C:H2'	1:A:323:U:C5'	2.50	0.41
1:A:325:A:N6	1:A:326:G:C2	2.88	0.41
1:A:362:G:H2'	1:A:364:A:OP2	2.20	0.41
1:A:416:G:O5'	1:A:416:G:H8	2.03	0.41
1:A:593:G:N1	1:A:594:G:C4	2.88	0.41
1:A:602:A:N3	1:A:637:G:C2	2.88	0.41
1:A:718:G:C5	11:K:116:HIS:CD2	3.08	0.41
1:A:922:G:C2	1:A:1396:A:C2	3.07	0.41
1:A:1030(D):A:H3'	1:A:1031:G:C5'	2.50	0.41
1:A:1047:G:H1	1:A:1210:C:H42	1.68	0.41
1:A:1070:U:H2'	1:A:1071:C:C6	2.56	0.41
1:A:1116:C:H2'	1:A:1117:G:C5'	2.50	0.41
1:A:1185:G:C2	1:A:1186:G:C8	3.08	0.41
1:A:1191:A:OP2	3:C:3:ASN:OD1	2.39	0.41
1:A:1202:G:O2'	14:N:2:ALA:HB1	2.21	0.41
1:A:1222:G:H5''	19:S:78:ARG:HH11	1.81	0.41
2:B:19:HIS:H	2:B:39:ILE:CG2	2.34	0.41
2:B:21:ARG:HA	2:B:39:ILE:N	2.35	0.41
2:B:29:ALA:C	2:B:31:TYR:N	2.73	0.41
3:C:40:ARG:HG3	3:C:40:ARG:HH11	1.86	0.41
3:C:94:LEU:C	3:C:94:LEU:CD2	2.88	0.41
3:C:112:SER:OG	3:C:115:LEU:CG	2.68	0.41
4:D:150:GLU:C	4:D:152:SER:N	2.74	0.41
5:E:12:LEU:HD12	5:E:12:LEU:N	2.35	0.41
5:E:131:ILE:HD13	5:E:131:ILE:HA	1.82	0.41
8:H:126:LYS:HG2	8:H:127:LEU:HD22	2.02	0.41
8:H:127:LEU:HD22	8:H:127:LEU:N	2.35	0.41
9:I:86:VAL:O	9:I:87:GLN:C	2.56	0.41
9:I:88:TYR:CE1	9:I:89:ASN:HB2	2.56	0.41
11:K:78:GLN:O	11:K:79:SER:O	2.39	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:K:100:ALA:O	11:K:102:GLY:N	2.53	0.41
11:K:127:LYS:CE	11:K:127:LYS:CA	2.93	0.41
12:L:33:ARG:HB3	12:L:60:LEU:HD12	2.01	0.41
12:L:69:TYR:CE1	12:L:70:ILE:O	2.73	0.41
13:M:76:ALA:O	13:M:79:LYS:N	2.53	0.41
15:O:16:ALA:CB	15:O:21:ASP:O	2.68	0.41
17:Q:71:PHE:C	17:Q:72:ARG:HG2	2.40	0.41
20:T:57:ARG:HE	20:T:100:ILE:HG21	1.85	0.41
20:T:76:ALA:O	20:T:80:ARG:CG	2.60	0.41
21:V:15:ARG:HH11	21:V:15:ARG:HB2	1.86	0.41
1:A:22:G:O2'	1:A:23:C:H5'	2.19	0.41
1:A:59:A:H1'	1:A:354:G:N2	2.36	0.41
1:A:118:U:O4	1:A:288:A:H2'	2.19	0.41
1:A:393:A:N3	1:A:394:G:C8	2.88	0.41
1:A:411:A:C2	1:A:431:A:N6	2.88	0.41
1:A:427:U:H4'	1:A:541:G:H5''	2.01	0.41
1:A:437:U:O2'	1:A:438:G:H5'	2.21	0.41
1:A:451:A:N6	1:A:481:G:N9	2.69	0.41
1:A:451:A:H61	1:A:481:G:H1'	1.85	0.41
1:A:513:C:H2'	1:A:514:C:C6	2.55	0.41
1:A:521:G:OP1	12:L:73:GLU:O	2.38	0.41
1:A:533:A:HO2'	1:A:534:U:P	2.43	0.41
1:A:622:A:C2	1:A:623:C:H1'	2.55	0.41
1:A:633:G:C2	1:A:634:C:O2	2.72	0.41
1:A:862:C:O4'	1:A:874:G:H4'	2.20	0.41
1:A:1047:G:C3'	1:A:1048:G:C5'	2.95	0.41
1:A:1148:U:O2'	9:I:14:VAL:HG11	2.20	0.41
1:A:1157:A:O4'	1:A:1158:C:C2	2.73	0.41
1:A:1177:G:H2'	1:A:1178:G:O5'	2.20	0.41
1:A:1427:U:H3	1:A:1473:A:H61	1.68	0.41
2:B:82:ARG:HA	2:B:92:TYR:CD2	2.51	0.41
2:B:163:PHE:HZ	2:B:201:ILE:HD12	1.85	0.41
2:B:178:ARG:O	8:H:71:GLY:HA2	2.20	0.41
3:C:16:ARG:NH2	3:C:183:ASP:CB	2.84	0.41
3:C:120:VAL:O	3:C:120:VAL:HG12	2.20	0.41
4:D:31:CYS:C	4:D:33:MET:N	2.74	0.41
4:D:57:ARG:NE	4:D:202:LEU:HD22	2.35	0.41
7:G:47:CYS:O	7:G:49:ILE:N	2.49	0.41
7:G:100:ALA:HB3	7:G:101:LEU:CD2	2.51	0.41
8:H:60:ARG:HG3	8:H:60:ARG:NH1	2.35	0.41
9:I:13:ALA:HB3	9:I:76:ALA:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:93:ARG:HG2	9:I:96:LEU:HD23	2.03	0.41
10:J:18:ALA:HA	10:J:21:GLN:CB	2.51	0.41
14:N:47:LEU:O	14:N:50:LYS:HG2	2.21	0.41
15:O:63:ARG:C	15:O:65:ARG:H	2.24	0.41
15:O:79:ARG:O	15:O:82:ILE:HG22	2.21	0.41
1:A:16:A:H2'	1:A:17:U:H5'	2.02	0.41
1:A:103:C:P	20:T:17:ARG:HH12	2.43	0.41
1:A:322:C:C2'	1:A:323:U:C5'	2.98	0.41
1:A:413:G:H2'	1:A:428:G:N2	2.35	0.41
1:A:438:G:C4'	1:A:439:A:OP1	2.68	0.41
1:A:564:C:C6	17:Q:31:LEU:HD11	2.56	0.41
1:A:621:A:C4	1:A:622:A:C8	3.09	0.41
1:A:622:A:C4	1:A:623:C:C6	3.08	0.41
1:A:658:G:C2	1:A:749:C:N3	2.88	0.41
1:A:880:C:OP2	12:L:9:GLN:OE1	2.38	0.41
1:A:1240:U:P	7:G:119:ARG:HH22	2.42	0.41
1:A:1351:U:C4'	7:G:33:ASP:OD1	2.69	0.41
2:B:175:ARG:HB2	2:B:175:ARG:CZ	2.51	0.41
4:D:25:ARG:O	4:D:27:TYR:N	2.52	0.41
4:D:59:ARG:H	4:D:59:ARG:HG2	1.75	0.41
6:F:12:PRO:HB2	6:F:45:LEU:HD21	2.01	0.41
6:F:73:ASN:O	6:F:76:ALA:N	2.54	0.41
7:G:44:TYR:CE1	9:I:41:VAL:HG21	2.55	0.41
7:G:72:ARG:HG2	7:G:142:GLU:HG2	2.01	0.41
7:G:109:ASN:OD1	7:G:119:ARG:NH1	2.53	0.41
9:I:3:GLN:O	9:I:4:TYR:CD2	2.74	0.41
9:I:17:VAL:HG11	9:I:81:ILE:HG23	2.01	0.41
9:I:26:VAL:CG1	9:I:28:VAL:HG23	2.50	0.41
9:I:55:ALA:O	9:I:56:LEU:C	2.59	0.41
11:K:124:LYS:HB3	11:K:125:PHE:CE1	2.56	0.41
12:L:120:TYR:CD2	12:L:120:TYR:N	2.84	0.41
13:M:6:GLY:O	13:M:7:VAL:HG23	2.20	0.41
15:O:9:GLN:O	15:O:10:LYS:C	2.58	0.41
15:O:18:PHE:CE2	15:O:21:ASP:HB2	2.56	0.41
16:P:21:VAL:CG2	16:P:59:TRP:HE1	2.31	0.41
16:P:42:ARG:O	16:P:43:LYS:C	2.59	0.41
16:P:43:LYS:HA	16:P:48:TRP:HB3	2.02	0.41
17:Q:18:THR:HG22	17:Q:19:VAL:N	2.36	0.41
1:A:83:U:O2'	1:A:84:U:H5'	2.20	0.41
1:A:147:G:C2	1:A:176:C:N3	2.89	0.41
1:A:232:G:N3	1:A:263:A:H2	2.18	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:U:O2'	1:A:265:G:H5'	2.20	0.41
1:A:491:G:C2	1:A:492:G:C4	3.09	0.41
1:A:514:C:H2'	1:A:515:G:O4'	2.20	0.41
1:A:673:G:O3'	6:F:87:ARG:NH2	2.54	0.41
1:A:792:A:H1'	1:A:793:U:C5'	2.50	0.41
1:A:865:A:C6	1:A:866:C:C4	3.09	0.41
1:A:1162:C:N3	1:A:1175:G:C2	2.89	0.41
1:A:1262:C:O2'	1:A:1263:C:H5'	2.20	0.41
1:A:1285:A:H5'	1:A:1286:A:C2	2.55	0.41
1:A:1333:A:H2'	1:A:1334:G:C8	2.56	0.41
2:B:73:THR:CB	2:B:96:ARG:HH21	2.34	0.41
2:B:178:ARG:CZ	2:B:196:LEU:O	2.69	0.41
3:C:88:ARG:C	3:C:90:GLU:H	2.22	0.41
3:C:118:GLN:CA	3:C:121:ALA:HB3	2.49	0.41
4:D:102:ASP:CG	4:D:103:ASN:H	2.23	0.41
4:D:133:VAL:O	4:D:134:ASP:HB3	2.21	0.41
5:E:144:THR:O	5:E:146:ALA:N	2.53	0.41
7:G:17:VAL:C	7:G:19:GLY:N	2.74	0.41
7:G:23:VAL:O	7:G:24:THR:C	2.58	0.41
7:G:120:ILE:O	7:G:121:ALA:C	2.59	0.41
8:H:104:ARG:CZ	8:H:138:TRP:CZ2	3.04	0.41
9:I:16:ARG:HD3	9:I:16:ARG:N	2.20	0.41
9:I:56:LEU:CD2	9:I:57:GLY:N	2.83	0.41
11:K:59:TYR:CE1	11:K:63:LEU:HD21	2.56	0.41
11:K:61:ALA:O	11:K:64:ALA:HB3	2.20	0.41
11:K:115:PRO:C	11:K:117:ASN:N	2.70	0.41
12:L:27:LEU:HD22	12:L:64:TYR:CE1	2.55	0.41
13:M:14:ARG:HH12	13:M:16:ASP:CB	2.34	0.41
17:Q:75:ARG:HH12	17:Q:77:VAL:HG22	1.85	0.41
17:Q:88:TYR:CD2	17:Q:89:LEU:HD23	2.55	0.41
19:S:69:HIS:CB	19:S:73:GLU:OE1	2.67	0.41
20:T:57:ARG:HE	20:T:100:ILE:CG2	2.33	0.41
20:T:60:GLU:CA	20:T:63:ILE:HD12	2.44	0.41
20:T:66:ALA:O	20:T:68:LYS:N	2.53	0.41
1:A:102:G:C5	1:A:103:C:C5	3.08	0.41
1:A:175:C:H2'	1:A:176:C:O4'	2.20	0.41
1:A:376:G:O2'	1:A:377:G:C5'	2.64	0.41
1:A:453:A:N6	1:A:480:U:H3	2.17	0.41
1:A:758:G:H8	1:A:758:G:O5'	2.04	0.41
1:A:879:C:C5	12:L:6:THR:HG21	2.55	0.41
1:A:879:C:H3'	1:A:879:C:H6	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:939:G:N2	1:A:1344:C:O2	2.53	0.41
1:A:1060:C:C5	3:C:2:GLY:N	2.89	0.41
1:A:1118:C:H1'	1:A:1179:A:N3	2.34	0.41
1:A:1302:U:O4	13:M:14:ARG:HD3	2.21	0.41
1:A:1520:G:N1	1:A:1521:G:C6	2.89	0.41
2:B:15:VAL:HG11	2:B:209:ARG:C	2.41	0.41
2:B:28:PHE:CZ	2:B:189:ASP:CA	2.86	0.41
3:C:18:TRP:CD1	14:N:54:PRO:HA	2.55	0.41
3:C:58:GLU:CG	3:C:67:THR:OG1	2.69	0.41
3:C:109:PRO:CB	3:C:115:LEU:HD13	2.51	0.41
3:C:153:VAL:HG12	3:C:154:SER:H	1.85	0.41
6:F:19:LEU:C	6:F:21:LEU:N	2.73	0.41
8:H:33:GLU:HA	8:H:36:LEU:HD12	2.01	0.41
9:I:50:LEU:HD11	9:I:81:ILE:CB	2.50	0.41
10:J:20:ALA:O	10:J:23:ILE:HB	2.20	0.41
10:J:72:VAL:O	10:J:73:ASP:O	2.38	0.41
12:L:86:ARG:HG3	12:L:86:ARG:NH1	2.36	0.41
13:M:20:THR:CA	13:M:25:ILE:HG22	2.43	0.41
15:O:83:GLU:O	15:O:83:GLU:HG2	2.20	0.41
19:S:17:GLU:O	19:S:19:VAL:N	2.43	0.41
1:A:35:G:OP1	12:L:104:VAL:HG22	2.20	0.41
1:A:37:U:C2'	1:A:38:G:H5'	2.51	0.41
1:A:101:A:C2	1:A:102:G:N9	2.89	0.41
1:A:124:G:C5	1:A:125:U:C4	3.09	0.41
1:A:156:G:C2	1:A:166:G:C6	3.09	0.41
1:A:191:G:C5	1:A:192:U:C5	3.09	0.41
1:A:195:A:C6	1:A:196:A:N1	2.89	0.41
1:A:203:U:H5''	1:A:204:U:O5'	2.21	0.41
1:A:256:U:C5'	17:Q:17:LYS:NZ	2.82	0.41
1:A:451:A:N7	1:A:481:G:C2	2.89	0.41
1:A:547:A:H4'	1:A:548:G:O5'	2.19	0.41
1:A:588:G:C4	1:A:753:A:C6	3.09	0.41
1:A:590:C:C2	1:A:650:G:C2	3.09	0.41
1:A:596:C:OP2	1:A:596:C:H3'	2.20	0.41
1:A:622:A:C2	1:A:623:C:C1'	3.03	0.41
1:A:658:G:C2'	1:A:659:U:H5'	2.50	0.41
1:A:763:G:N2	1:A:764:C:C2	2.89	0.41
1:A:830:G:C2'	1:A:831:U:H5'	2.50	0.41
1:A:923:A:C2	1:A:1395:C:N3	2.89	0.41
1:A:944:G:C3'	1:A:945:G:C5'	2.99	0.41
1:A:953:G:O6	1:A:1228:C:N4	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1004:A:N7	1:A:1037:C:O2	2.53	0.41
1:A:1025:U:C2'	1:A:1026:G:H8	2.34	0.41
1:A:1046:A:C2	1:A:1047:G:H1'	2.56	0.41
1:A:1078:U:H2'	1:A:1079:G:O4'	2.21	0.41
1:A:1101:A:C4'	1:A:1102:A:O5'	2.68	0.41
1:A:1213:A:C5	1:A:1215:G:N7	2.88	0.41
1:A:1253:G:N2	1:A:1254:C:H1'	2.36	0.41
1:A:1326:C:H5''	21:V:18:TYR:O	2.21	0.41
1:A:1402:C:O2'	1:A:1403:C:H5'	2.20	0.41
1:A:1405:G:O2'	1:A:1406:U:H5'	2.21	0.41
1:A:1408:A:H2'	1:A:1409:C:H6	1.86	0.41
1:A:1465:C:H2'	1:A:1466:C:H6	1.84	0.41
1:A:1500:A:OP1	1:A:1508:G:OP1	2.39	0.41
2:B:53:ARG:CG	2:B:54:THR:N	2.83	0.41
2:B:105:PHE:O	2:B:107:THR:N	2.53	0.41
2:B:122:PHE:O	2:B:127:ILE:HG13	2.21	0.41
2:B:142:LEU:HD23	2:B:146:GLN:HG3	2.01	0.41
3:C:10:PHE:O	3:C:14:ILE:HD11	2.20	0.41
3:C:18:TRP:CZ2	14:N:56:VAL:O	2.73	0.41
3:C:59:ARG:O	3:C:60:ALA:HB2	2.20	0.41
3:C:61:ALA:C	3:C:63:ASN:N	2.73	0.41
3:C:157:ILE:C	3:C:159:GLY:H	2.24	0.41
4:D:2:GLY:C	4:D:3:ARG:HE	2.23	0.41
4:D:100:ARG:HH12	4:D:137:SER:HA	1.85	0.41
4:D:163:GLU:C	4:D:165:MET:H	2.24	0.41
5:E:63:ARG:HA	5:E:66:MET:SD	2.61	0.41
5:E:93:PRO:HG2	5:E:93:PRO:O	2.21	0.41
6:F:1:MET:SD	6:F:66:GLU:OE2	2.78	0.41
6:F:16:GLN:C	6:F:16:GLN:CD	2.80	0.41
6:F:18:GLN:CA	6:F:21:LEU:HB3	2.50	0.41
6:F:32:ASN:ND2	6:F:32:ASN:C	2.75	0.41
6:F:80:ARG:NH1	6:F:88:VAL:HB	2.35	0.41
6:F:91:VAL:CG1	18:R:72:ARG:NH2	2.84	0.41
6:F:100:ASN:HD22	18:R:23:LYS:HG2	1.86	0.41
7:G:21:VAL:O	7:G:24:THR:CB	2.69	0.41
7:G:143:ARG:O	7:G:145:ALA:O	2.39	0.41
8:H:7:ALA:HA	8:H:10:LEU:HD12	2.03	0.41
8:H:18:ARG:HB2	8:H:18:ARG:NH1	2.30	0.41
8:H:40:ALA:O	8:H:42:GLU:N	2.54	0.41
8:H:100:ILE:CG2	8:H:125:ARG:HE	2.34	0.41
9:I:18:PHE:O	9:I:61:ALA:CB	2.66	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:97:LYS:O	9:I:98:PRO:C	2.57	0.41
10:J:3:LYS:CG	10:J:76:ASN:HD22	2.30	0.41
10:J:18:ALA:C	10:J:20:ALA:N	2.74	0.41
10:J:82:ILE:O	10:J:86:MET:HB2	2.21	0.41
10:J:90:LEU:HD23	10:J:90:LEU:O	2.20	0.41
12:L:53:ARG:NH1	12:L:93:LEU:HD21	2.36	0.41
12:L:115:LYS:C	12:L:117:ARG:N	2.74	0.41
13:M:10:PRO:CB	13:M:18:ALA:CB	2.89	0.41
13:M:116:THR:HG22	13:M:117:VAL:H	1.86	0.41
14:N:9:LYS:O	14:N:9:LYS:NZ	2.37	0.41
15:O:68:ARG:O	15:O:72:ARG:HB2	2.21	0.41
16:P:1:MET:CE	16:P:65:GLN:HB2	2.50	0.41
16:P:81:ARG:NE	16:P:83:GLU:OE2	2.51	0.41
17:Q:5:VAL:HA	17:Q:59:ILE:O	2.20	0.41
17:Q:56:VAL:O	17:Q:57:VAL:CG1	2.69	0.41
17:Q:75:ARG:CG	17:Q:76:LEU:H	2.34	0.41
18:R:20:ALA:C	18:R:55:ARG:HH12	2.24	0.41
19:S:17:GLU:C	19:S:19:VAL:N	2.73	0.41
19:S:40:ILE:HD13	19:S:62:ILE:CD1	2.47	0.41
20:T:30:LYS:CD	20:T:72:LEU:HD21	2.41	0.41
20:T:100:ILE:HG13	20:T:100:ILE:H	1.73	0.41
1:A:51:A:C1'	1:A:52:G:OP2	2.69	0.41
1:A:60:A:H1'	1:A:61:G:O4'	2.21	0.41
1:A:76:C:H2'	1:A:77:G:C8	2.50	0.41
1:A:109:A:C2	1:A:327:A:N1	2.89	0.41
1:A:397:A:C6	1:A:548:G:N7	2.89	0.41
1:A:448:A:OP2	1:A:485:G:N2	2.44	0.41
1:A:501:C:H1'	1:A:549:C:H1'	2.02	0.41
1:A:505:G:C6	1:A:535:A:C2	3.09	0.41
1:A:552:U:O2'	12:L:31:PRO:HB3	2.21	0.41
1:A:578:C:O2'	1:A:579:G:H5'	2.21	0.41
1:A:588:G:C6	1:A:753:A:C8	3.09	0.41
1:A:594:G:N2	1:A:646:U:O2	2.53	0.41
1:A:761:G:H4'	17:Q:102:GLY:CA	2.46	0.41
1:A:872:A:C8	1:A:874:G:C8	3.09	0.41
1:A:892:A:C6	1:A:893:C:N4	2.89	0.41
1:A:1057:G:C5	1:A:1058:G:C5	3.09	0.41
1:A:1151:A:H5'	10:J:41:PRO:HA	2.03	0.41
1:A:1261:A:H62	1:A:1274:G:N2	2.05	0.41
1:A:1361:G:C8	1:A:1361:G:P	3.14	0.41
2:B:21:ARG:O	2:B:22:LYS:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:73:THR:CG2	2:B:96:ARG:HH21	2.33	0.41
2:B:115:LEU:O	2:B:118:LEU:HB2	2.21	0.41
2:B:196:LEU:HA	2:B:196:LEU:HD23	1.80	0.41
3:C:95:THR:CG2	3:C:95:THR:O	2.69	0.41
3:C:138:VAL:HG12	3:C:170:GLN:HB2	2.03	0.41
4:D:91:SER:OG	4:D:92:VAL:N	2.54	0.41
4:D:205:GLU:O	4:D:206:PHE:C	2.59	0.41
5:E:14:ARG:NH1	5:E:129:ILE:CD1	2.84	0.41
6:F:12:PRO:CB	6:F:45:LEU:HD21	2.51	0.41
9:I:36:TYR:CZ	9:I:70:LYS:HE2	2.55	0.41
9:I:65:VAL:CG2	9:I:66:ARG:N	2.84	0.41
10:J:8:LEU:HD21	10:J:20:ALA:CB	2.50	0.41
11:K:14:VAL:HG21	11:K:40:ILE:HD11	2.03	0.41
11:K:44:SER:O	11:K:45:GLY:C	2.58	0.41
11:K:93:GLN:O	11:K:96:ARG:HB3	2.21	0.41
11:K:93:GLN:CD	11:K:96:ARG:NH2	2.74	0.41
13:M:8:GLU:OE2	13:M:22:ILE:HG23	2.21	0.41
15:O:68:ARG:CB	15:O:68:ARG:NH1	2.84	0.41
16:P:74:LEU:HD22	16:P:79:VAL:CG2	2.50	0.41
21:V:18:TYR:CE2	21:V:24:ARG:HD3	2.56	0.41
1:A:106:C:H2'	1:A:107:G:C5'	2.50	0.40
1:A:260:G:C4	1:A:261:U:C5	3.09	0.40
1:A:302:G:H21	1:A:556:C:C4'	2.33	0.40
1:A:370:C:C2	1:A:371:G:C8	3.09	0.40
1:A:376:G:N3	1:A:389:A:C2	2.88	0.40
1:A:597:G:N7	1:A:598:U:C5	2.89	0.40
1:A:623:C:O2'	1:A:624:C:H5'	2.21	0.40
1:A:792:A:HO2'	1:A:793:U:P	2.37	0.40
1:A:872:A:C2	1:A:874:G:C6	3.09	0.40
1:A:1118:C:O4'	1:A:1179:A:H1'	2.21	0.40
1:A:1125:U:OP2	1:A:1145:C:N4	2.54	0.40
1:A:1299:A:O2'	1:A:1300:G:H4'	2.21	0.40
1:A:1367:C:N3	1:A:1368:G:N7	2.69	0.40
1:A:1414:U:H2'	1:A:1415:G:H8	1.86	0.40
1:A:1520:G:C6	1:A:1521:G:C6	3.09	0.40
2:B:90:MET:HA	2:B:91:PRO:HD3	1.81	0.40
2:B:136:VAL:HA	2:B:139:LYS:HB2	2.03	0.40
2:B:187:LEU:HD23	2:B:201:ILE:HB	2.03	0.40
3:C:66:VAL:HG12	3:C:66:VAL:O	2.20	0.40
3:C:131:ARG:O	3:C:134:ILE:N	2.54	0.40
4:D:49:ARG:O	4:D:51:PRO:N	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:119:GLN:C	4:D:121:VAL:N	2.74	0.40
5:E:36:ASP:OD1	5:E:38:GLN:N	2.48	0.40
7:G:28:ASN:CA	7:G:31:MET:HB2	2.37	0.40
8:H:42:GLU:O	8:H:42:GLU:OE1	2.39	0.40
8:H:126:LYS:O	8:H:128:GLY:N	2.54	0.40
9:I:19:LEU:CD2	9:I:59:PHE:HB3	2.33	0.40
9:I:96:LEU:CG	9:I:102:LEU:HD13	2.45	0.40
10:J:76:ASN:N	10:J:76:ASN:ND2	2.68	0.40
11:K:83:ILE:HA	11:K:109:VAL:O	2.21	0.40
12:L:24:VAL:HG12	12:L:26:ALA:HB3	2.01	0.40
13:M:108:ARG:NE	13:M:108:ARG:HA	2.36	0.40
14:N:11:LYS:NZ	14:N:13:THR:HB	2.36	0.40
14:N:47:LEU:O	14:N:48:ALA:C	2.60	0.40
15:O:78:TYR:C	15:O:82:ILE:HG22	2.41	0.40
18:R:68:LYS:HE3	18:R:72:ARG:NH2	2.36	0.40
21:V:14:TRP:O	21:V:15:ARG:C	2.59	0.40
1:A:44:G:N3	1:A:399:G:C2	2.89	0.40
1:A:320:C:O2'	1:A:321:A:H5'	2.22	0.40
1:A:363:A:N1	1:A:364:A:C2	2.89	0.40
1:A:377:G:C2	1:A:378:G:C5	3.09	0.40
1:A:394:G:O2'	1:A:395:C:H5'	2.21	0.40
1:A:420:U:O2	1:A:424:G:N2	2.54	0.40
1:A:455:C:C2	1:A:456:C:C5	3.09	0.40
1:A:475:G:C4	1:A:476:G:C8	3.10	0.40
1:A:528:C:H5'	1:A:535:A:N6	2.36	0.40
1:A:620:C:H2'	1:A:621:A:O4'	2.21	0.40
1:A:714:G:H2'	1:A:715:A:C8	2.56	0.40
1:A:773:G:H2'	1:A:774:G:H8	1.86	0.40
1:A:939:G:H2'	1:A:940:C:C6	2.57	0.40
1:A:969:A:H2'	1:A:970:C:O4'	2.21	0.40
1:A:1059:C:N4	1:A:1198:G:H1	2.20	0.40
1:A:1073:U:P	5:E:57:LYS:NZ	2.94	0.40
1:A:1230:C:O2'	1:A:1231:G:H5'	2.21	0.40
1:A:1255:G:N1	1:A:1283:G:N2	2.70	0.40
1:A:1305:G:C8	1:A:1305:G:OP2	2.74	0.40
1:A:1321:C:H2'	1:A:1322:C:H5	1.86	0.40
1:A:1382:C:O2'	1:A:1383:C:H5'	2.21	0.40
1:A:1494:G:H2'	1:A:1495:U:C5'	2.51	0.40
2:B:30:ARG:C	2:B:31:TYR:CD1	2.94	0.40
2:B:123:ALA:O	2:B:124:SER:O	2.39	0.40
2:B:223:ILE:HD13	2:B:230:VAL:CG2	2.46	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:7:PRO:HA	3:C:10:PHE:CB	2.41	0.40
3:C:114:PRO:HG3	3:C:185:GLY:HA2	2.03	0.40
4:D:57:ARG:HD2	4:D:202:LEU:HD22	2.03	0.40
4:D:127:THR:HG1	4:D:130:GLY:CA	2.34	0.40
5:E:39:GLY:HA2	5:E:71:LEU:CG	2.51	0.40
5:E:41:VAL:CG2	5:E:113:ALA:CB	2.99	0.40
5:E:135:THR:O	5:E:138:ALA:N	2.55	0.40
8:H:60:ARG:CB	8:H:62:TYR:HE2	2.34	0.40
8:H:94:TYR:CD1	8:H:94:TYR:N	2.89	0.40
9:I:23:ASN:HB3	9:I:24:GLY:H	1.67	0.40
9:I:42:ARG:HE	9:I:42:ARG:HB3	1.51	0.40
10:J:18:ALA:HA	10:J:21:GLN:HB2	2.04	0.40
11:K:15:ALA:HB1	11:K:78:GLN:CB	2.49	0.40
13:M:44:ARG:HB2	13:M:46:LYS:HG2	2.03	0.40
18:R:45:SER:C	18:R:47:THR:N	2.69	0.40
19:S:5:LEU:CD2	19:S:9:VAL:HG13	2.50	0.40
20:T:75:ASN:O	20:T:78:ALA:HB3	2.21	0.40
1:A:14:U:H3'	1:A:14:U:H6	1.87	0.40
1:A:74:C:C4	1:A:75:G:C5	3.10	0.40
1:A:89:C:N1	1:A:90:U:H5	2.19	0.40
1:A:106:C:HO2'	1:A:107:G:H5'	1.83	0.40
1:A:119:A:N6	1:A:240:C:C2	2.90	0.40
1:A:255:G:H1'	17:Q:16:GLN:CD	2.42	0.40
1:A:301:G:H2'	1:A:302:G:C8	2.56	0.40
1:A:596:C:O2	1:A:596:C:C2'	2.70	0.40
1:A:684:A:H2'	1:A:685:G:C8	2.55	0.40
1:A:701:C:O2'	1:A:702:A:P	2.79	0.40
1:A:781:A:N7	1:A:802:A:C2	2.90	0.40
1:A:1015:A:C2'	1:A:1016:A:C8	2.93	0.40
1:A:1152:A:O2'	1:A:1153:C:O5'	2.37	0.40
1:A:1255:G:N1	1:A:1283:G:N3	2.70	0.40
1:A:1424:C:C2'	1:A:1425:U:C5'	2.99	0.40
2:B:57:PHE:CZ	2:B:61:LEU:HD12	2.56	0.40
2:B:69:LEU:CD2	2:B:102:LEU:HD11	2.52	0.40
2:B:86:GLU:H	2:B:86:GLU:HG2	1.70	0.40
3:C:179:ARG:HD3	3:C:207:VAL:HA	2.02	0.40
4:D:64:LEU:HD12	4:D:75:PHE:CZ	2.56	0.40
4:D:97:LEU:HD23	4:D:97:LEU:O	2.21	0.40
5:E:72:GLN:O	5:E:73:ASN:HB3	2.22	0.40
6:F:15:ASP:OD1	6:F:15:ASP:C	2.60	0.40
7:G:77:SER:O	7:G:78:ARG:HB2	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:H:11:THR:C	8:H:13:ILE:N	2.70	0.40
10:J:26:ALA:HB3	10:J:85:LEU:HD11	2.03	0.40
11:K:31:THR:O	11:K:31:THR:HG22	2.22	0.40
13:M:17:VAL:O	13:M:20:THR:N	2.54	0.40
13:M:103:THR:OG1	13:M:103:THR:O	2.32	0.40
17:Q:26:GLN:O	17:Q:27:PHE:HB3	2.22	0.40
18:R:76:LEU:HD23	18:R:76:LEU:HA	1.72	0.40
20:T:11:SER:O	20:T:14:LYS:HE2	2.21	0.40
20:T:39:LYS:HG2	20:T:55:ILE:CD1	2.51	0.40
1:A:279:A:C2	17:Q:98:LEU:HD12	2.57	0.40
1:A:283:C:C2	1:A:284:G:C8	3.09	0.40
1:A:287:U:C2'	1:A:288:A:C5'	3.00	0.40
1:A:382:A:C4	1:A:383:A:C8	3.10	0.40
1:A:408:A:C6	1:A:409:G:C6	3.10	0.40
1:A:533:A:O2'	1:A:534:U:P	2.79	0.40
1:A:541:G:C2	1:A:542:G:C4	3.10	0.40
1:A:613:C:C2	1:A:628:G:N2	2.89	0.40
1:A:686:U:O2	1:A:687:A:N9	2.54	0.40
1:A:877:C:C2'	1:A:878:G:C5'	2.99	0.40
1:A:883:C:N3	1:A:884:U:C4	2.89	0.40
1:A:942:G:N3	1:A:943:U:C6	2.90	0.40
1:A:1003(A):G:H8	1:A:1003(A):G:O5'	2.04	0.40
1:A:1222:G:H5''	19:S:78:ARG:HH12	1.84	0.40
1:A:1269:A:N3	1:A:1313:U:H1'	2.36	0.40
1:A:1299:A:C4	1:A:1301:U:C2	3.10	0.40
2:B:17:PHE:C	2:B:17:PHE:HD1	2.24	0.40
2:B:39:ILE:HG22	2:B:40:HIS:N	2.37	0.40
3:C:140:ARG:HB2	3:C:141:VAL:H	1.71	0.40
3:C:157:ILE:H	3:C:160:ALA:HB3	1.86	0.40
4:D:5:ILE:HA	4:D:115:ARG:HH22	1.85	0.40
4:D:47:ARG:HG3	4:D:47:ARG:NH1	2.36	0.40
4:D:48:ALA:O	4:D:49:ARG:C	2.60	0.40
4:D:125:HIS:C	4:D:126:ILE:HD13	2.42	0.40
4:D:125:HIS:C	4:D:149:ALA:HB2	2.41	0.40
7:G:136:LYS:C	7:G:140:ASP:HB2	2.41	0.40
7:G:152:ALA:HA	7:G:155:ARG:NH2	2.36	0.40
9:I:29:ASN:OD1	9:I:64:THR:HG23	2.20	0.40
11:K:33:THR:HG1	11:K:37:GLY:C	2.24	0.40
11:K:50:TYR:HB2	11:K:55:LYS:HG2	2.04	0.40
14:N:23:ARG:NH1	14:N:29:ARG:C	2.74	0.40
15:O:4:THR:OG1	15:O:7:GLU:HB3	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:Q:6:LEU:HD23	17:Q:6:LEU:HA	1.91	0.40
17:Q:52:LYS:O	17:Q:55:ASP:OD2	2.38	0.40
17:Q:71:PHE:O	17:Q:72:ARG:HG2	2.22	0.40
1:A:59:A:H2'	1:A:331:G:H22	1.85	0.40
1:A:186:C:O2'	20:T:82:SER:HA	2.21	0.40
1:A:187:C:O2'	20:T:89:ARG:NE	2.55	0.40
1:A:357:G:N3	1:A:358:U:C6	2.90	0.40
1:A:560:U:O4'	1:A:566:G:N2	2.54	0.40
1:A:568:G:N2	1:A:883:C:C5	2.90	0.40
1:A:604:G:C5	1:A:605:U:C5	3.09	0.40
1:A:694:A:H5'	11:K:53:SER:HB2	2.03	0.40
1:A:706:A:C8	1:A:707:C:C5	3.10	0.40
1:A:710:G:H5''	6:F:54:LYS:HZ1	1.77	0.40
1:A:777:A:N3	1:A:777:A:H2'	2.35	0.40
1:A:894:G:N3	1:A:895:G:C8	2.90	0.40
1:A:949:A:H61	1:A:1232:U:H3	1.68	0.40
1:A:990:C:H4'	1:A:1018:C:P	2.61	0.40
1:A:1011:G:H2'	1:A:1012:U:O4'	2.22	0.40
1:A:1186:G:N2	1:A:1187:G:H1'	2.36	0.40
1:A:1215:G:C2'	1:A:1216:G:C5'	2.99	0.40
1:A:1238:A:H5'	1:A:1336:C:H41	1.86	0.40
1:A:1481:U:H2'	1:A:1482:G:O4'	2.21	0.40
1:A:1486:G:H2'	1:A:1487:G:H8	1.87	0.40
2:B:101:MET:O	2:B:105:PHE:HA	2.22	0.40
2:B:168:THR:O	2:B:168:THR:HG22	2.21	0.40
2:B:181:PHE:HD2	8:H:70:GLN:HB3	1.82	0.40
3:C:59:ARG:CG	3:C:60:ALA:H	2.12	0.40
4:D:72:GLU:C	4:D:74:GLN:N	2.75	0.40
4:D:87:GLY:O	4:D:89:THR:N	2.55	0.40
4:D:89:THR:O	4:D:91:SER:N	2.55	0.40
5:E:148:VAL:HG21	8:H:107:LEU:CD1	2.49	0.40
6:F:35:ALA:HA	6:F:67:MET:HB3	2.03	0.40
6:F:97:PHE:O	6:F:98:LEU:HD23	2.20	0.40
7:G:30:ILE:O	7:G:31:MET:C	2.59	0.40
11:K:82:VAL:CG1	11:K:83:ILE:H	2.35	0.40
12:L:46:LYS:HG2	12:L:47:LYS:CG	2.51	0.40
12:L:89:ARG:HB2	12:L:90:VAL:H	1.60	0.40
13:M:81:LEU:HD13	13:M:88:ARG:HD3	2.04	0.40
16:P:23:ASP:C	16:P:25:ARG:N	2.75	0.40
17:Q:79:SER:O	17:Q:80:GLY:O	2.39	0.40
19:S:28:LYS:CG	19:S:29:ARG:N	2.84	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	232/256 (91%)	104 (45%)	70 (30%)	58 (25%)	0	0
3	C	204/239 (85%)	95 (47%)	59 (29%)	50 (24%)	0	0
4	D	206/208 (99%)	94 (46%)	62 (30%)	50 (24%)	0	0
5	E	148/161 (92%)	86 (58%)	38 (26%)	24 (16%)	0	3
6	F	99/101 (98%)	72 (73%)	15 (15%)	12 (12%)	0	5
7	G	153/155 (99%)	78 (51%)	45 (29%)	30 (20%)	0	1
8	H	136/138 (99%)	81 (60%)	32 (24%)	23 (17%)	0	2
9	I	125/128 (98%)	66 (53%)	38 (30%)	21 (17%)	0	2
10	J	96/104 (92%)	54 (56%)	19 (20%)	23 (24%)	0	0
11	K	117/129 (91%)	53 (45%)	33 (28%)	31 (26%)	0	0
12	L	122/135 (90%)	65 (53%)	28 (23%)	29 (24%)	0	0
13	M	116/126 (92%)	67 (58%)	28 (24%)	21 (18%)	0	1
14	N	58/60 (97%)	25 (43%)	19 (33%)	14 (24%)	0	0
15	O	86/88 (98%)	44 (51%)	23 (27%)	19 (22%)	0	1
16	P	81/88 (92%)	41 (51%)	18 (22%)	22 (27%)	0	0
17	Q	102/104 (98%)	66 (65%)	25 (24%)	11 (11%)	0	6
18	R	71/88 (81%)	38 (54%)	16 (22%)	17 (24%)	0	0
19	S	78/92 (85%)	43 (55%)	24 (31%)	11 (14%)	0	3
20	T	97/106 (92%)	31 (32%)	36 (37%)	30 (31%)	0	0
21	V	22/26 (85%)	14 (64%)	7 (32%)	1 (4%)	2	23
All	All	2349/2532 (93%)	1217 (52%)	635 (27%)	497 (21%)	0	1

All (497) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	15	VAL
2	B	17	PHE
2	B	39	ILE
2	B	78	GLN
2	B	82	ARG
2	B	96	ARG
2	B	183	PRO
2	B	190	THR
2	B	204	ASN
2	B	211	ILE
3	C	15	THR
3	C	17	ASP
3	C	18	TRP
3	C	60	ALA
3	C	72	LYS
3	C	77	ILE
3	C	111	LEU
3	C	156	ARG
3	C	168	ALA
3	C	171	GLY
3	C	176	HIS
3	C	179	ARG
3	C	180	ALA
3	C	183	ASP
3	C	189	ALA
4	D	29	PRO
4	D	36	ARG
4	D	40	PRO
4	D	44	GLY
4	D	53	ASP
4	D	91	SER
4	D	92	VAL
4	D	132	ARG
4	D	156	GLU
4	D	177	ASP
4	D	191	ARG
5	E	21	ALA
5	E	71	LEU
5	E	104	ALA
5	E	112	LEU
5	E	142	LEU
5	E	146	ALA
6	F	37	VAL

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Mol	Chain	Res	Type
6	F	39	LYS
6	F	83	ASP
7	G	50	ILE
7	G	63	LYS
7	G	97	GLN
7	G	114	ARG
7	G	125	MET
7	G	135	VAL
8	H	31	PHE
8	H	32	LYS
8	H	40	ALA
8	H	45	ILE
8	H	97	VAL
8	H	105	ARG
8	H	134	ILE
9	I	21	PRO
9	I	31	GLN
9	I	38	GLN
9	I	41	VAL
9	I	56	LEU
9	I	85	LEU
9	I	92	TYR
9	I	120	ARG
9	I	123	PRO
10	J	32	ALA
10	J	39	PRO
10	J	48	THR
10	J	54	PHE
10	J	55	LYS
10	J	58	ASP
10	J	59	SER
10	J	60	ARG
10	J	65	LEU
10	J	73	ASP
10	J	78	ASN
10	J	85	LEU
10	J	86	MET
10	J	90	LEU
11	K	12	ARG
11	K	47	VAL
11	K	48	ILE
11	K	50	TYR

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Mol	Chain	Res	Type
11	K	53	SER
11	K	54	ARG
11	K	79	SER
11	K	80	VAL
11	K	90	GLY
11	K	101	SER
11	K	126	ARG
12	L	7	ILE
12	L	17	LYS
12	L	19	ARG
12	L	27	LEU
12	L	41	ARG
12	L	56	ALA
12	L	90	VAL
12	L	108	ALA
12	L	121	GLY
13	M	42	ALA
13	M	43	THR
13	M	44	ARG
13	M	63	THR
13	M	97	PRO
13	M	118	ALA
14	N	14	PRO
14	N	15	LYS
14	N	22	THR
14	N	23	ARG
14	N	25	VAL
14	N	29	ARG
14	N	33	VAL
14	N	49	HIS
16	P	10	GLY
16	P	12	LYS
16	P	20	VAL
16	P	21	VAL
16	P	23	ASP
16	P	31	LYS
16	P	34	GLU
16	P	36	ILE
16	P	48	TRP
16	P	81	ARG
17	Q	14	LYS
17	Q	69	LYS

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Mol	Chain	Res	Type
17	Q	80	GLY
17	Q	81	ARG
17	Q	95	TYR
17	Q	96	GLN
17	Q	103	GLY
18	R	19	LYS
18	R	22	VAL
18	R	23	LYS
18	R	36	ASN
18	R	37	VAL
18	R	70	ILE
19	S	6	LYS
19	S	24	ALA
19	S	35	SER
19	S	45	VAL
19	S	47	HIS
19	S	77	THR
19	S	78	ARG
20	T	11	SER
20	T	14	LYS
20	T	15	ARG
20	T	49	ALA
20	T	50	GLU
20	T	51	GLU
20	T	54	LYS
20	T	70	SER
20	T	74	LYS
20	T	89	ARG
21	V	3	LYS
2	B	18	GLY
2	B	22	LYS
2	B	37	ASN
2	B	61	LEU
2	B	80	ILE
2	B	84	GLU
2	B	88	ALA
2	B	106	LYS
2	B	117	GLU
2	B	130	ARG
2	B	132	LYS
2	B	135	GLN
2	B	150	SER

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Mol	Chain	Res	Type
2	B	152	PHE
2	B	161	ALA
2	B	174	VAL
2	B	177	ALA
2	B	207	ALA
2	B	214	ILE
2	B	224	GLN
2	B	239	VAL
3	C	29	TYR
3	C	43	LEU
3	C	53	ALA
3	C	61	ALA
3	C	67	THR
3	C	78	GLY
3	C	84	ILE
3	C	101	LEU
3	C	119	ARG
3	C	129	ALA
3	C	130	VAL
3	C	155	GLY
3	C	178	LEU
3	C	182	ILE
4	D	7	PRO
4	D	23	GLY
4	D	30	LYS
4	D	33	MET
4	D	41	GLY
4	D	58	LEU
4	D	63	LYS
4	D	101	LEU
4	D	104	VAL
4	D	131	ARG
4	D	133	VAL
4	D	134	ASP
4	D	151	LYS
4	D	157	LEU
4	D	158	ILE
4	D	172	PRO
4	D	175	SER
4	D	206	PHE
5	E	49	PRO
5	E	79	GLU

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Mol	Chain	Res	Type
5	E	108	ALA
5	E	111	GLU
5	E	129	ILE
6	F	19	LEU
6	F	23	LYS
6	F	70	ASP
6	F	72	VAL
6	F	74	ASP
6	F	84	ASN
7	G	6	ARG
7	G	15	ASP
7	G	18	TYR
7	G	33	ASP
7	G	42	ILE
7	G	46	ALA
7	G	80	VAL
7	G	93	PRO
7	G	118	VAL
7	G	119	ARG
7	G	134	ALA
7	G	140	ASP
8	H	6	ILE
8	H	39	LEU
8	H	41	ARG
8	H	42	GLU
8	H	90	GLY
8	H	91	ARG
8	H	126	LYS
9	I	12	GLU
9	I	29	ASN
9	I	33	PHE
9	I	39	GLY
9	I	114	TYR
10	J	24	VAL
10	J	52	GLY
10	J	68	HIS
11	K	64	ALA
11	K	68	ALA
11	K	83	ILE
11	K	91	ARG
11	K	102	GLY
11	K	106	LYS

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Mol	Chain	Res	Type
11	K	117	ASN
11	K	118	GLY
12	L	6	THR
12	L	14	GLY
12	L	18	VAL
12	L	51	ALA
12	L	55	VAL
12	L	63	GLY
12	L	73	GLU
12	L	79	GLU
12	L	96	VAL
12	L	109	GLY
12	L	116	SER
12	L	123	LYS
13	M	5	ALA
13	M	14	ARG
13	M	20	THR
13	M	27	LYS
13	M	67	GLU
13	M	75	ALA
13	M	81	LEU
13	M	100	GLY
13	M	117	VAL
14	N	12	ARG
14	N	50	LYS
14	N	57	ARG
15	O	9	GLN
15	O	34	LEU
15	O	36	ILE
15	O	48	LYS
15	O	61	GLY
15	O	62	GLN
15	O	79	ARG
15	O	87	ILE
16	P	27	LYS
16	P	51	VAL
16	P	63	GLY
16	P	67	THR
16	P	78	GLY
17	Q	33	GLY
17	Q	47	PRO
18	R	17	SER

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Mol	Chain	Res	Type
18	R	18	ARG
18	R	31	LEU
18	R	57	GLY
18	R	69	THR
18	R	87	ARG
19	S	28	LYS
19	S	68	GLY
20	T	12	ALA
20	T	60	GLU
20	T	67	ALA
20	T	84	LEU
20	T	95	ALA
20	T	101	GLY
20	T	102	GLY
20	T	103	GLY
2	B	8	LYS
2	B	11	LEU
2	B	20	GLU
2	B	26	PRO
2	B	110	GLN
2	B	113	HIS
2	B	133	LYS
2	B	147	LYS
2	B	179	LYS
2	B	209	ARG
2	B	229	VAL
2	B	232	PRO
2	B	234	PRO
3	C	6	HIS
3	C	74	GLY
3	C	98	ASN
3	C	144	SER
3	C	146	ALA
3	C	177	THR
4	D	3	ARG
4	D	47	ARG
4	D	64	LEU
4	D	137	SER
4	D	196	LEU
5	E	16	THR
5	E	22	GLY
5	E	52	PRO

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Mol	Chain	Res	Type
5	E	55	VAL
5	E	137	GLU
5	E	138	ALA
6	F	38	GLU
7	G	17	VAL
8	H	89	PRO
9	I	15	ALA
10	J	19	SER
10	J	40	LEU
10	J	76	ASN
11	K	13	GLN
11	K	42	TRP
11	K	44	SER
11	K	46	GLY
11	K	55	LYS
11	K	62	GLN
11	K	89	ALA
11	K	128	ALA
14	N	31	ARG
15	O	17	ARG
15	O	78	TYR
16	P	25	ARG
17	Q	64	PRO
18	R	21	LYS
18	R	25	THR
18	R	76	LEU
19	S	61	TYR
20	T	34	LYS
20	T	55	ILE
20	T	76	ALA
20	T	79	ARG
2	B	52	GLU
2	B	119	GLU
2	B	165	VAL
2	B	208	ILE
2	B	228	GLY
3	C	49	SER
3	C	52	LEU
3	C	76	VAL
3	C	137	ALA
3	C	140	ARG
4	D	5	ILE

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Mol	Chain	Res	Type
4	D	12	CYS
4	D	26	CYS
4	D	51	PRO
4	D	150	GLU
4	D	164	ALA
4	D	180	GLY
4	D	193	ASP
5	E	72	GLN
5	E	109	ILE
6	F	82	ARG
7	G	78	ARG
7	G	91	VAL
7	G	104	LEU
7	G	105	VAL
7	G	120	ILE
7	G	129	GLU
8	H	63	LEU
8	H	73	ASP
9	I	23	ASN
9	I	67	GLY
10	J	61	GLU
11	K	73	MET
11	K	92	GLU
12	L	29	GLY
12	L	57	LYS
13	M	24	GLY
13	M	46	LYS
14	N	45	ARG
15	O	16	ALA
15	O	59	MET
15	O	60	VAL
15	O	74	ASP
16	P	13	HIS
16	P	72	ARG
17	Q	66	SER
18	R	71	LYS
19	S	60	VAL
20	T	59	ALA
20	T	97	ALA
2	B	56	ARG
3	C	68	VAL
3	C	113	ALA

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Mol	Chain	Res	Type
3	C	192	THR
4	D	4	TYR
4	D	9	CYS
4	D	105	VAL
5	E	78	HIS
5	E	136	MET
6	F	27	GLN
7	G	35	LYS
7	G	48	LYS
8	H	5	PRO
8	H	86	ILE
8	H	127	LEU
9	I	24	GLY
12	L	47	LYS
12	L	71	PRO
12	L	72	GLY
12	L	89	ARG
13	M	49	THR
13	M	82	MET
15	O	38	ARG
15	O	80	ALA
15	O	88	ARG
16	P	43	LYS
16	P	52	ASP
20	T	44	ALA
20	T	77	ALA
20	T	78	ALA
2	B	54	THR
2	B	59	GLU
2	B	121	LEU
2	B	124	SER
2	B	194	PRO
3	C	172	ARG
3	C	206	GLU
4	D	135	LEU
4	D	184	LYS
5	E	34	VAL
11	K	14	VAL
13	M	106	ASN
15	O	75	PRO
16	P	41	PRO
18	R	39	VAL

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Mol	Chain	Res	Type
20	T	52	ALA
20	T	96	GLY
2	B	81	VAL
3	C	51	GLY
3	C	81	GLY
4	D	17	VAL
7	G	61	VAL
8	H	4	ASP
14	N	56	VAL
15	O	45	VAL
2	B	112	VAL
2	B	125	PRO
3	C	138	VAL
3	C	202	ILE
8	H	103	VAL
9	I	97	LYS
13	M	4	ILE
3	C	5	ILE
4	D	88	VAL
5	E	106	PRO
7	G	49	ILE
9	I	6	GLY
10	J	72	VAL
20	T	41	ILE
5	E	80	ILE
12	L	48	PRO
12	L	101	VAL
16	P	2	VAL
4	D	178	VAL
7	G	133	GLY
8	H	101	PRO
9	I	98	PRO
10	J	77	PRO
11	K	29	ILE

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	202/220 (92%)	163 (81%)	39 (19%)	1	9
3	C	160/188 (85%)	127 (79%)	33 (21%)	1	7
4	D	180/180 (100%)	147 (82%)	33 (18%)	1	10
5	E	115/122 (94%)	98 (85%)	17 (15%)	3	18
6	F	90/90 (100%)	79 (88%)	11 (12%)	5	24
7	G	126/126 (100%)	105 (83%)	21 (17%)	2	14
8	H	119/119 (100%)	103 (87%)	16 (13%)	4	21
9	I	98/99 (99%)	71 (72%)	27 (28%)	0	3
10	J	87/91 (96%)	63 (72%)	24 (28%)	0	3
11	K	90/99 (91%)	73 (81%)	17 (19%)	1	9
12	L	104/111 (94%)	90 (86%)	14 (14%)	4	21
13	M	94/101 (93%)	80 (85%)	14 (15%)	3	17
14	N	49/49 (100%)	31 (63%)	18 (37%)	0	0
15	O	79/79 (100%)	68 (86%)	11 (14%)	3	20
16	P	72/74 (97%)	59 (82%)	13 (18%)	1	10
17	Q	96/96 (100%)	82 (85%)	14 (15%)	3	18
18	R	64/77 (83%)	55 (86%)	9 (14%)	3	19
19	S	71/79 (90%)	62 (87%)	9 (13%)	4	23
20	T	76/82 (93%)	68 (90%)	8 (10%)	7	30
21	V	19/21 (90%)	13 (68%)	6 (32%)	0	1
All	All	1991/2103 (95%)	1637 (82%)	354 (18%)	2	11

All (354) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	8	LYS
2	B	15	VAL
2	B	17	PHE
2	B	21	ARG
2	B	25	ASN
2	B	30	ARG
2	B	44	LEU
2	B	55	PHE
2	B	59	GLU
2	B	60	ASP
2	B	64	ARG

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Mol	Chain	Res	Type
2	B	71	VAL
2	B	76	GLN
2	B	82	ARG
2	B	92	TYR
2	B	96	ARG
2	B	97	TRP
2	B	105	PHE
2	B	107	THR
2	B	108	ILE
2	B	111	ARG
2	B	113	HIS
2	B	129	GLU
2	B	130	ARG
2	B	139	LYS
2	B	144	ARG
2	B	162	ILE
2	B	165	VAL
2	B	172	ILE
2	B	178	ARG
2	B	183	PRO
2	B	195	ASP
2	B	200	ILE
2	B	204	ASN
2	B	210	SER
2	B	224	GLN
2	B	231	GLU
2	B	232	PRO
2	B	236	TYR
3	C	3	ASN
3	C	14	ILE
3	C	15	THR
3	C	21	ARG
3	C	34	LEU
3	C	35	GLU
3	C	37	GLN
3	C	48	TYR
3	C	52	LEU
3	C	56	ASP
3	C	68	VAL
3	C	79	ARG
3	C	82	GLU
3	C	86	VAL

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Mol	Chain	Res	Type
3	C	91	LEU
3	C	97	LYS
3	C	98	ASN
3	C	102	ASN
3	C	104	GLN
3	C	108	ASN
3	C	116	VAL
3	C	118	GLN
3	C	119	ARG
3	C	127	ARG
3	C	128	PHE
3	C	139	GLN
3	C	143	GLU
3	C	167	TRP
3	C	175	LEU
3	C	181	ASN
3	C	191	THR
3	C	196	LEU
3	C	201	TYR
4	D	3	ARG
4	D	9	CYS
4	D	10	ARG
4	D	26	CYS
4	D	29	PRO
4	D	47	ARG
4	D	59	ARG
4	D	61	LYS
4	D	62	GLN
4	D	76	ARG
4	D	77	ASN
4	D	78	LEU
4	D	79	PHE
4	D	80	GLU
4	D	99	SER
4	D	112	VAL
4	D	115	ARG
4	D	122	ARG
4	D	134	ASP
4	D	135	LEU
4	D	137	SER
4	D	139	ARG
4	D	156	GLU

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Mol	Chain	Res	Type
4	D	157	LEU
4	D	158	ILE
4	D	160	GLN
4	D	170	VAL
4	D	179	GLU
4	D	188	LEU
4	D	191	ARG
4	D	194	LEU
4	D	199	ASN
4	D	209	ARG
5	E	10	MET
5	E	12	LEU
5	E	15	ARG
5	E	20	GLN
5	E	24	ARG
5	E	31	LEU
5	E	33	VAL
5	E	51	VAL
5	E	53	LEU
5	E	56	GLN
5	E	67	VAL
5	E	68	GLU
5	E	79	GLU
5	E	80	ILE
5	E	89	ILE
5	E	126	ARG
5	E	147	ASP
6	F	7	ASN
6	F	16	GLN
6	F	21	LEU
6	F	24	GLU
6	F	31	GLU
6	F	32	ASN
6	F	47	ARG
6	F	57	GLN
6	F	63	TYR
6	F	82	ARG
6	F	86	ARG
7	G	4	ARG
7	G	12	LEU
7	G	16	LEU
7	G	37	ASN

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Mol	Chain	Res	Type
7	G	38	LEU
7	G	50	ILE
7	G	56	GLN
7	G	57	GLU
7	G	62	PHE
7	G	64	GLN
7	G	73	MET
7	G	84	ASN
7	G	101	LEU
7	G	115	ARG
7	G	126	ASP
7	G	138	LYS
7	G	140	ASP
7	G	144	MET
7	G	153	HIS
7	G	154	TYR
7	G	155	ARG
8	H	18	ARG
8	H	25	ASP
8	H	30	ARG
8	H	31	PHE
8	H	39	LEU
8	H	56	LYS
8	H	65	TYR
8	H	85	ARG
8	H	91	ARG
8	H	92	ARG
8	H	97	VAL
8	H	104	ARG
8	H	115	SER
8	H	119	LEU
8	H	120	THR
8	H	133	LEU
9	I	3	GLN
9	I	10	ARG
9	I	11	LYS
9	I	16	ARG
9	I	19	LEU
9	I	21	PRO
9	I	23	ASN
9	I	25	LYS
9	I	27	THR

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Mol	Chain	Res	Type
9	I	34	ASN
9	I	37	PHE
9	I	40	LEU
9	I	42	ARG
9	I	66	ARG
9	I	81	ILE
9	I	86	VAL
9	I	88	TYR
9	I	91	ASP
9	I	92	TYR
9	I	97	LYS
9	I	102	LEU
9	I	104	ARG
9	I	111	ARG
9	I	114	TYR
9	I	121	ARG
9	I	125	TYR
9	I	127	LYS
10	J	6	ILE
10	J	8	LEU
10	J	9	ARG
10	J	13	HIS
10	J	23	ILE
10	J	33	GLN
10	J	34	VAL
10	J	40	LEU
10	J	45	ARG
10	J	48	THR
10	J	49	VAL
10	J	50	ILE
10	J	60	ARG
10	J	68	HIS
10	J	70	ARG
10	J	71	LEU
10	J	74	ILE
10	J	75	ILE
10	J	79	ARG
10	J	81	THR
10	J	83	GLU
10	J	87	THR
10	J	95	GLU
10	J	98	ILE

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Mol	Chain	Res	Type
11	K	18	ARG
11	K	28	THR
11	K	29	ILE
11	K	31	THR
11	K	33	THR
11	K	57	THR
11	K	66	LEU
11	K	75	TYR
11	K	77	MET
11	K	81	ASP
11	K	91	ARG
11	K	92	GLU
11	K	98	LEU
11	K	109	VAL
11	K	116	HIS
11	K	126	ARG
11	K	127	LYS
12	L	9	GLN
12	L	37	CYS
12	L	41	ARG
12	L	49	ASN
12	L	54	LYS
12	L	65	GLU
12	L	66	VAL
12	L	81	SER
12	L	99	HIS
12	L	106	ASP
12	L	110	VAL
12	L	112	ASP
12	L	113	ARG
12	L	126	LYS
13	M	9	ILE
13	M	11	ARG
13	M	19	LEU
13	M	35	GLU
13	M	40	ASN
13	M	48	LEU
13	M	56	LEU
13	M	70	LEU
13	M	79	LYS
13	M	82	MET
13	M	83	ASP

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Mol	Chain	Res	Type
13	M	87	TYR
13	M	102	ARG
13	M	103	THR
14	N	3	ARG
14	N	6	LEU
14	N	8	GLU
14	N	11	LYS
14	N	12	ARG
14	N	21	TYR
14	N	23	ARG
14	N	24	CYS
14	N	27	CYS
14	N	31	ARG
14	N	37	PHE
14	N	39	LEU
14	N	41	ARG
14	N	44	LEU
14	N	46	GLU
14	N	47	LEU
14	N	53	LEU
14	N	58	LYS
15	O	4	THR
15	O	13	GLN
15	O	22	THR
15	O	24	SER
15	O	39	LEU
15	O	46	HIS
15	O	48	LYS
15	O	57	LEU
15	O	81	LEU
15	O	82	ILE
15	O	83	GLU
16	P	5	ARG
16	P	17	TYR
16	P	20	VAL
16	P	21	VAL
16	P	29	ASP
16	P	43	LYS
16	P	45	THR
16	P	59	TRP
16	P	62	VAL
16	P	69	THR

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Mol	Chain	Res	Type
16	P	71	ARG
16	P	80	PHE
16	P	82	GLN
17	Q	20	THR
17	Q	34	LYS
17	Q	36	ILE
17	Q	38	ARG
17	Q	53	LEU
17	Q	60	ILE
17	Q	68	ARG
17	Q	73	VAL
17	Q	74	LEU
17	Q	76	LEU
17	Q	82	MET
17	Q	96	GLN
17	Q	98	LEU
17	Q	104	LYS
18	R	18	ARG
18	R	28	GLU
18	R	36	ASN
18	R	39	VAL
18	R	47	THR
18	R	53	ARG
18	R	56	THR
18	R	86	VAL
18	R	87	ARG
19	S	12	ASP
19	S	19	VAL
19	S	25	LYS
19	S	32	LYS
19	S	34	TRP
19	S	39	THR
19	S	52	TYR
19	S	58	VAL
19	S	61	TYR
20	T	10	LEU
20	T	14	LYS
20	T	23	ARG
20	T	75	ASN
20	T	80	ARG
20	T	83	ARG
20	T	84	LEU

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Mol	Chain	Res	Type
20	T	89	ARG
21	V	5	ASP
21	V	10	ARG
21	V	13	ILE
21	V	15	ARG
21	V	21	TYR
21	V	23	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (66) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	19	HIS
2	B	25	ASN
2	B	37	ASN
2	B	78	GLN
2	B	204	ASN
3	C	3	ASN
3	C	28	GLN
3	C	37	GLN
3	C	98	ASN
3	C	102	ASN
3	C	104	GLN
3	C	123	GLN
3	C	136	GLN
3	C	139	GLN
4	D	42	GLN
4	D	62	GLN
4	D	74	GLN
4	D	77	ASN
4	D	123	HIS
4	D	129	ASN
4	D	160	GLN
4	D	199	ASN
4	D	201	GLN
5	E	73	ASN
5	E	130	ASN
6	F	7	ASN
6	F	13	ASN
6	F	16	GLN
6	F	18	GLN
6	F	32	ASN
6	F	57	GLN

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Mol	Chain	Res	Type
6	F	73	ASN
6	F	100	ASN
7	G	11	GLN
7	G	37	ASN
7	G	56	GLN
7	G	148	ASN
9	I	23	ASN
9	I	34	ASN
10	J	33	GLN
10	J	76	ASN
10	J	84	GLN
11	K	62	GLN
11	K	78	GLN
11	K	93	GLN
11	K	99	GLN
11	K	117	ASN
12	L	49	ASN
12	L	75	HIS
12	L	78	GLN
13	M	12	ASN
13	M	40	ASN
13	M	62	ASN
13	M	77	ASN
13	M	106	ASN
15	O	13	GLN
15	O	37	ASN
15	O	50	HIS
16	P	82	GLN
17	Q	16	GLN
17	Q	94	ASN
17	Q	96	GLN
18	R	36	ASN
19	S	53	ASN
19	S	56	GLN
20	T	75	ASN

5.3.3 RNA

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	1511/1522 (99%)	253 (16%)	62 (4%)

All (253) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	8	A
1	A	9	G
1	A	31	G
1	A	32	A
1	A	39	G
1	A	41	G
1	A	42	G
1	A	47	C
1	A	48	C
1	A	49	U
1	A	50	A
1	A	51	A
1	A	52	G
1	A	61	G
1	A	64	G
1	A	65	U
1	A	73	C
1	A	81	U
1	A	82	U
1	A	89	C
1	A	91	C
1	A	116	A
1	A	120	A
1	A	121	C
1	A	129(A)	G
1	A	130	A
1	A	131	C
1	A	163	C
1	A	182	U
1	A	183	G
1	A	190(E)	U
1	A	190(F)	G
1	A	190(G)	G
1	A	195	A
1	A	197	A
1	A	199	G
1	A	201	C
1	A	202	U
1	A	216	G
1	A	217	C
1	A	244	U
1	A	247	G
1	A	251	G

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Mol	Chain	Res	Type
1	A	252	U
1	A	266	G
1	A	267	C
1	A	280	C
1	A	282	A
1	A	283	C
1	A	289	G
1	A	326	G
1	A	328	C
1	A	329	A
1	A	330	C
1	A	345	C
1	A	352	C
1	A	353	A
1	A	354	G
1	A	366	C
1	A	367	U
1	A	372	C
1	A	373	A
1	A	384	G
1	A	392	G
1	A	397	A
1	A	398	C
1	A	412	A
1	A	413	G
1	A	414	A
1	A	421	U
1	A	423	G
1	A	424	G
1	A	428	G
1	A	429	U
1	A	430	A
1	A	436	C
1	A	439	A
1	A	452	A
1	A	453	A
1	A	460	A
1	A	461	C
1	A	481	G
1	A	482	A
1	A	484	G
1	A	485	G

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Mol	Chain	Res	Type
1	A	497	A
1	A	498	U
1	A	500	G
1	A	509	A
1	A	510	A
1	A	511	C
1	A	512	U
1	A	517	G
1	A	518	C
1	A	519	C
1	A	527	G
1	A	531	U
1	A	533	A
1	A	534	U
1	A	545	C
1	A	547	A
1	A	559	A
1	A	560	U
1	A	561	U
1	A	562	C
1	A	572	A
1	A	573	A
1	A	575	G
1	A	576	G
1	A	577	G
1	A	596	C
1	A	598	U
1	A	632	A
1	A	653	A
1	A	665	A
1	A	687	A
1	A	688	G
1	A	695	A
1	A	702	A
1	A	703	G
1	A	718	G
1	A	723	U
1	A	724	G
1	A	731	G
1	A	734	G
1	A	749	C
1	A	755	G

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Mol	Chain	Res	Type
1	A	777	A
1	A	792	A
1	A	793	U
1	A	794	A
1	A	817	C
1	A	820	U
1	A	828	A
1	A	839	U
1	A	840	C
1	A	841	U
1	A	848	C
1	A	859	A
1	A	867	G
1	A	902	G
1	A	913	A
1	A	914	A
1	A	926	G
1	A	927	G
1	A	934	C
1	A	935	A
1	A	945	G
1	A	960	U
1	A	961	U
1	A	966	G
1	A	968	A
1	A	969	A
1	A	971	G
1	A	974	A
1	A	975	A
1	A	976	G
1	A	977	A
1	A	983	A
1	A	984	C
1	A	992	U
1	A	993	G
1	A	994	A
1	A	1000	U
1	A	1024	G
1	A	1026	G
1	A	1027	C
1	A	1029	C
1	A	1031	G

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Mol	Chain	Res	Type
1	A	1038	C
1	A	1048	G
1	A	1050	G
1	A	1054	C
1	A	1066	C
1	A	1068	G
1	A	1086	U
1	A	1094	G
1	A	1095	U
1	A	1101	A
1	A	1102	A
1	A	1104	G
1	A	1108	G
1	A	1116	C
1	A	1117	G
1	A	1125	U
1	A	1126	U
1	A	1129	C
1	A	1131	G
1	A	1136	U
1	A	1137	C
1	A	1138	G
1	A	1139	G
1	A	1142	G
1	A	1146	A
1	A	1150	U
1	A	1152	A
1	A	1159	U
1	A	1183	A
1	A	1184	G
1	A	1196	U
1	A	1197	G
1	A	1201	A
1	A	1202	G
1	A	1211	U
1	A	1212	U
1	A	1215	G
1	A	1225	A
1	A	1226	C
1	A	1227	A
1	A	1241	G
1	A	1250	A

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Mol	Chain	Res	Type
1	A	1256	A
1	A	1257	U
1	A	1258	G
1	A	1279	A
1	A	1280	A
1	A	1281	U
1	A	1282	C
1	A	1285	A
1	A	1286	A
1	A	1287	A
1	A	1297	C
1	A	1299	A
1	A	1301	U
1	A	1303	C
1	A	1305	G
1	A	1306	A
1	A	1323	G
1	A	1331	G
1	A	1338	G
1	A	1347	G
1	A	1348	U
1	A	1359	C
1	A	1362	C
1	A	1363	A
1	A	1398	A
1	A	1427	U
1	A	1428	A
1	A	1442	G
1	A	1446	A
1	A	1452	C
1	A	1476	G
1	A	1492	A
1	A	1493	A
1	A	1499	A
1	A	1504	G
1	A	1506	U
1	A	1517	G
1	A	1518	A
1	A	1520	G
1	A	1529	G
1	A	1530	G
1	A	1533	C

All (62) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	7	G
1	A	30	U
1	A	51	A
1	A	60	A
1	A	64	G
1	A	115	G
1	A	119	A
1	A	129(A)	G
1	A	204	U
1	A	243	A
1	A	250	A
1	A	251	G
1	A	266	G
1	A	279	A
1	A	281	G
1	A	328	C
1	A	344	A
1	A	351	G
1	A	353	A
1	A	366	C
1	A	372	C
1	A	428	G
1	A	429	U
1	A	438	G
1	A	484	G
1	A	496	A
1	A	509	A
1	A	533	A
1	A	559	A
1	A	560	U
1	A	575	G
1	A	687	A
1	A	701	C
1	A	748	C
1	A	792	A
1	A	840	C
1	A	913	A
1	A	960	U
1	A	975	A
1	A	992	U
1	A	993	G
1	A	1049	U

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Mol	Chain	Res	Type
1	A	1065	U
1	A	1067	A
1	A	1085	U
1	A	1101	A
1	A	1182	G
1	A	1196	U
1	A	1201	A
1	A	1214	C
1	A	1224	G
1	A	1226	C
1	A	1257	U
1	A	1279	A
1	A	1281	U
1	A	1285	A
1	A	1305	G
1	A	1346	A
1	A	1347	G
1	A	1397	C
1	A	1451	A
1	A	1498	U

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	1512/1522 (99%)	0.83	169 (11%) 5 3	12, 87, 171, 195	0
2	B	234/256 (91%)	-0.14	3 (1%) 77 65	21, 87, 149, 181	0
3	C	206/239 (86%)	0.38	20 (9%) 7 5	31, 115, 164, 182	0
4	D	208/208 (100%)	0.04	6 (2%) 51 37	15, 83, 140, 189	0
5	E	150/161 (93%)	0.01	1 (0%) 87 80	4, 55, 110, 132	0
6	F	101/101 (100%)	0.12	5 (4%) 28 20	32, 111, 151, 174	0
7	G	155/155 (100%)	-0.16	6 (3%) 39 27	27, 116, 156, 189	0
8	H	138/138 (100%)	-0.07	1 (0%) 87 80	0, 45, 102, 156	0
9	I	127/128 (99%)	0.66	21 (16%) 1 1	30, 120, 164, 186	0
10	J	98/104 (94%)	1.32	33 (33%) 0 0	43, 122, 168, 183	0
11	K	119/129 (92%)	0.17	4 (3%) 45 33	21, 87, 147, 195	0
12	L	124/135 (91%)	0.13	6 (4%) 30 21	2, 77, 127, 154	0
13	M	118/126 (93%)	0.47	13 (11%) 5 4	45, 118, 163, 195	0
14	N	60/60 (100%)	1.98	24 (40%) 0 0	50, 118, 171, 186	0
15	O	88/88 (100%)	-0.10	3 (3%) 45 33	23, 79, 143, 195	0
16	P	83/88 (94%)	0.06	2 (2%) 59 45	0, 60, 119, 138	0
17	Q	104/104 (100%)	-0.06	3 (2%) 51 37	8, 59, 132, 195	0
18	R	73/88 (82%)	-0.04	3 (4%) 37 26	3, 81, 147, 182	0
19	S	80/92 (86%)	1.33	22 (27%) 0 0	49, 124, 178, 181	0
20	T	99/106 (93%)	-0.33	1 (1%) 82 72	10, 66, 124, 141	0
21	V	24/26 (92%)	2.00	11 (45%) 0 0	84, 114, 145, 180	0
All	All	3901/4054 (96%)	0.46	357 (9%) 9 6	0, 90, 162, 195	0

All (357) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
14	N	4	LYS	10.9
19	S	3	ARG	8.4
17	Q	105	ALA	7.9
4	D	42	GLN	7.8
21	V	24	ARG	7.3
11	K	129	SER	7.2
14	N	17	LYS	7.1
19	S	31	ILE	7.1
9	I	9	ARG	6.8
19	S	77	THR	6.5
1	A	1220	G	6.3
3	C	2	GLY	6.2
1	A	1005	A	6.2
9	I	128	ARG	6.1
13	M	98	VAL	6.0
10	J	55	LYS	5.9
3	C	67	THR	5.9
10	J	46	ARG	5.8
9	I	37	PHE	5.8
1	A	1053	G	5.8
1	A	1200	C	5.7
13	M	99	ARG	5.5
7	G	2	ALA	5.5
14	N	30	ALA	5.5
1	A	975	A	5.5
1	A	1361(A)	C	5.4
14	N	3	ARG	5.4
14	N	31	ARG	5.3
1	A	1117	G	5.3
1	A	985	C	5.2
14	N	18	VAL	5.1
3	C	196	LEU	5.0
10	J	54	PHE	5.0
1	A	1030(B)	C	5.0
10	J	47	PHE	4.9
4	D	41	GLY	4.8
14	N	2	ALA	4.7
1	A	1031	G	4.6
1	A	984	C	4.6
21	V	17	THR	4.6
1	A	1004	A	4.6
1	A	1129	C	4.5
8	H	1	MET	4.5

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Mol	Chain	Res	Type	RSRZ
1	A	1320	C	4.5
14	N	22	THR	4.5
13	M	88	ARG	4.5
13	M	119	GLY	4.5
1	A	978	A	4.4
3	C	195	VAL	4.4
10	J	53	PRO	4.4
1	A	388	G	4.3
1	A	1191	A	4.3
21	V	4	GLY	4.3
19	S	36	ARG	4.2
1	A	979	C	4.2
1	A	993	G	4.2
10	J	43	ARG	4.1
9	I	17	VAL	4.1
1	A	723	U	4.1
19	S	37	ARG	4.1
1	A	1219	U	4.0
1	A	1362	C	4.0
1	A	1270	C	4.0
3	C	43	LEU	4.0
14	N	34	TYR	3.9
14	N	44	LEU	3.9
1	A	976	G	3.9
10	J	34	VAL	3.9
1	A	1260	C	3.9
14	N	32	SER	3.9
1	A	1317	C	3.9
1	A	547	A	3.9
14	N	61	TRP	3.9
10	J	64	GLU	3.8
14	N	29	ARG	3.7
19	S	78	ARG	3.7
1	A	964	A	3.7
3	C	47	LEU	3.6
13	M	10	PRO	3.6
1	A	1283	G	3.6
14	N	20	ALA	3.6
1	A	1319	A	3.5
19	S	35	SER	3.5
10	J	24	VAL	3.5
3	C	57	ILE	3.5

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Mol	Chain	Res	Type	RSRZ
14	N	21	TYR	3.5
1	A	1540	U	3.5
13	M	104	ARG	3.5
21	V	9	ARG	3.5
1	A	1221	G	3.5
11	K	51	LYS	3.5
1	A	1144	G	3.5
1	A	1202	G	3.5
19	S	62	ILE	3.4
17	Q	104	LYS	3.4
7	G	74	GLU	3.4
9	I	63	ILE	3.4
1	A	1250	A	3.4
10	J	28	ARG	3.4
4	D	2	GLY	3.4
13	M	116	THR	3.4
14	N	19	ARG	3.4
1	A	1003(A)	G	3.4
1	A	1201	A	3.3
10	J	62	HIS	3.3
10	J	96	ILE	3.3
1	A	1017	G	3.3
1	A	1041	A	3.3
1	A	1363	A	3.3
10	J	72	VAL	3.3
13	M	4	ILE	3.3
19	S	28	LYS	3.3
3	C	65	ALA	3.2
9	I	65	VAL	3.2
10	J	38	ILE	3.2
19	S	30	LEU	3.2
1	A	1226	C	3.2
1	A	1030(C)	G	3.2
13	M	97	PRO	3.2
9	I	126	SER	3.2
1	A	1135	U	3.2
1	A	1236	A	3.1
1	A	934	C	3.1
9	I	70	LYS	3.1
19	S	34	TRP	3.1
10	J	63	PHE	3.1
21	V	7	ARG	3.1

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Mol	Chain	Res	Type	RSRZ
10	J	44	VAL	3.1
7	G	5	ARG	3.1
1	A	536	C	3.1
10	J	71	LEU	3.1
1	A	977	A	3.1
12	L	33	ARG	3.1
1	A	1124	G	3.1
1	A	1334	G	3.1
19	S	47	HIS	3.1
1	A	1159	U	3.1
1	A	958	A	3.1
14	N	23	ARG	3.1
1	A	971	G	3.1
1	A	983	A	3.0
14	N	5	ALA	3.0
1	A	1360	A	3.0
1	A	81	U	3.0
1	A	1361	G	3.0
1	A	1214	C	3.0
12	L	115	LYS	3.0
3	C	66	VAL	3.0
19	S	2	PRO	3.0
1	A	427	U	3.0
19	S	53	ASN	3.0
1	A	973	G	3.0
1	A	353	A	3.0
19	S	48	THR	3.0
1	A	1321	C	2.9
11	K	19	ALA	2.9
10	J	74	ILE	2.9
6	F	67	MET	2.9
1	A	1224	G	2.9
1	A	1030(D)	A	2.9
21	V	5	ASP	2.9
1	A	1064	G	2.9
10	J	73	ASP	2.9
1	A	548	G	2.9
1	A	1049	U	2.9
3	C	26	LYS	2.9
1	A	1356	G	2.9
9	I	119	ALA	2.8
1	A	80	G	2.8

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Mol	Chain	Res	Type	RSRZ
5	E	5	ASP	2.8
6	F	63	TYR	2.8
9	I	77	ILE	2.8
1	A	1357	A	2.8
1	A	67	C	2.8
1	A	1434	A	2.8
1	A	982	U	2.8
19	S	79	THR	2.8
1	A	992	U	2.8
1	A	1036	G	2.8
10	J	27	ALA	2.8
1	A	499	A	2.8
13	M	105	THR	2.8
1	A	1139	G	2.7
1	A	1277	C	2.7
4	D	3	ARG	2.7
10	J	33	GLN	2.7
6	F	68	PRO	2.7
1	A	998	G	2.7
1	A	1251	A	2.7
3	C	189	ALA	2.7
1	A	1216	G	2.7
4	D	25	ARG	2.7
17	Q	101	ARG	2.7
1	A	1190	G	2.7
7	G	35	LYS	2.7
1	A	1014	A	2.7
1	A	1047	G	2.7
9	I	66	ARG	2.6
1	A	1128	C	2.6
9	I	85	LEU	2.6
1	A	1030	C	2.6
1	A	521	G	2.6
1	A	1068	G	2.6
1	A	1371	G	2.6
1	A	1094	G	2.6
1	A	1302	U	2.6
6	F	69	GLU	2.6
1	A	1050	G	2.6
1	A	1132	C	2.6
9	I	110	GLU	2.6
14	N	37	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
20	T	104	LEU	2.5
1	A	1084	G	2.5
13	M	64	TRP	2.5
19	S	32	LYS	2.5
3	C	103	VAL	2.5
1	A	1198	G	2.5
12	L	31	PRO	2.5
1	A	1052	U	2.5
3	C	68	VAL	2.5
1	A	1060	C	2.5
4	D	18	LYS	2.5
9	I	4	TYR	2.5
21	V	6	ARG	2.5
1	A	1002	G	2.5
1	A	952	U	2.5
1	A	1215	G	2.5
1	A	1467	G	2.5
1	A	373	A	2.5
1	A	1123	A	2.5
1	A	1287	A	2.5
21	V	18	TYR	2.5
14	N	39	LEU	2.5
15	O	54	ARG	2.5
1	A	1127	G	2.4
13	M	24	GLY	2.4
1	A	1042	G	2.4
1	A	1134	G	2.4
2	B	72	GLY	2.4
2	B	40	HIS	2.4
3	C	197	GLY	2.4
1	A	1267	C	2.4
2	B	16	HIS	2.4
1	A	904	C	2.4
14	N	38	GLY	2.4
1	A	410	G	2.4
1	A	1048	G	2.4
21	V	23	PRO	2.4
1	A	537	G	2.4
1	A	718	G	2.4
1	A	365	U	2.4
1	A	1199	U	2.4
1	A	161	A	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	995	C	2.4
15	O	46	HIS	2.4
18	R	80	PRO	2.4
1	A	1255	G	2.4
6	F	89	MET	2.4
3	C	44	GLU	2.4
1	A	776	G	2.4
1	A	1323	G	2.4
14	N	35	ARG	2.4
9	I	106	ALA	2.4
1	A	1531	A	2.4
1	A	970	C	2.4
14	N	16	PHE	2.4
10	J	45	ARG	2.3
10	J	66	ARG	2.3
10	J	70	ARG	2.3
3	C	17	ASP	2.3
16	P	83	GLU	2.3
10	J	48	THR	2.3
10	J	78	ASN	2.3
1	A	1130	A	2.3
1	A	1138	G	2.3
9	I	41	VAL	2.3
9	I	127	LYS	2.3
21	V	20	LYS	2.3
14	N	41	ARG	2.3
3	C	56	ASP	2.3
16	P	54	GLU	2.3
19	S	15	LEU	2.3
9	I	7	THR	2.3
1	A	1541	U	2.3
1	A	1544	U	2.3
11	K	39	PRO	2.3
19	S	12	ASP	2.3
1	A	1268	A	2.3
1	A	267	C	2.3
18	R	71	LYS	2.2
19	S	44	MET	2.2
1	A	327	A	2.2
1	A	675	A	2.2
1	A	696	A	2.2
1	A	1364	U	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	47	C	2.2
9	I	83	ARG	2.2
1	A	878	G	2.2
7	G	34	GLY	2.2
1	A	697	U	2.2
3	C	178	LEU	2.2
1	A	816	A	2.2
12	L	72	GLY	2.2
1	A	374	A	2.2
1	A	793	U	2.2
1	A	950	U	2.2
1	A	963	G	2.2
10	J	56	HIS	2.2
12	L	96	VAL	2.2
1	A	674	G	2.2
1	A	1037	C	2.2
1	A	1314	C	2.2
1	A	797	C	2.1
10	J	37	PRO	2.1
3	C	111	LEU	2.1
1	A	389	A	2.1
1	A	965	A	2.1
1	A	1333	A	2.1
1	A	726	C	2.1
1	A	1359	C	2.1
12	L	32	PHE	2.1
1	A	1365	G	2.1
7	G	109	ASN	2.1
1	A	1539	C	2.1
1	A	1266	G	2.1
18	R	74	ARG	2.1
1	A	1067	A	2.1
1	A	1280	A	2.1
19	S	41	VAL	2.1
1	A	1345	U	2.1
1	A	1149	C	2.1
10	J	17	ASP	2.1
1	A	1235	U	2.1
3	C	76	VAL	2.1
1	A	980	C	2.1
1	A	1006	C	2.1
9	I	18	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	1519	A	2.1
1	A	1218	C	2.1
21	V	25	LYS	2.1
13	M	117	VAL	2.1
15	O	51	HIS	2.1
10	J	60	ARG	2.1
1	A	1339	A	2.0
1	A	951	G	2.0
10	J	39	PRO	2.0
19	S	54	GLY	2.0
1	A	1056	U	2.0
1	A	39	G	2.0
1	A	1228	C	2.0
1	A	1530	G	2.0
10	J	35	SER	2.0
1	A	363	A	2.0
9	I	111	ARG	2.0
10	J	99	LYS	2.0
1	A	962	C	2.0
1	A	1355	G	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
22	ZN	D	306	1/1	0.97	0.45	74,74,74,74	0
22	ZN	N	307	1/1	0.97	0.10	74,74,74,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.