



# Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 05:45 pm BST

PDB ID : 6G32  
Title : Crystal structure of human geranylgeranyl diphosphate synthase mutant D188Y  
Authors : Lisnyansky, M.; Kapelushnik, N.; Ben-Bassat, A.; Marom, M.; Loewenstein, A.; Khananshvili, D.; Giladi, M.; Haitin, Y.  
Deposited on : 2018-03-24  
Resolution : 3.28 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

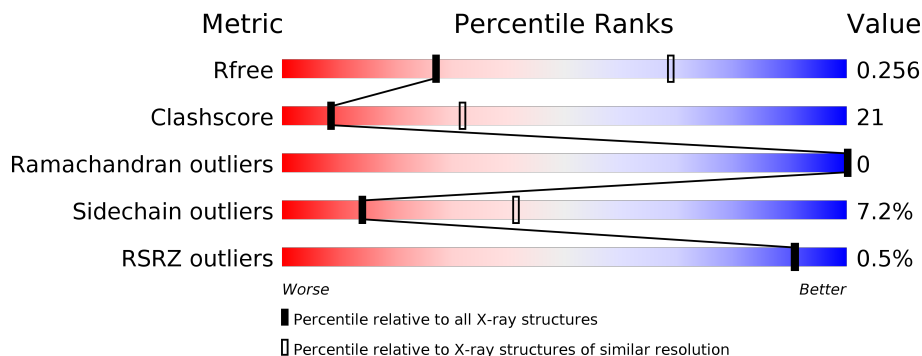
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.28 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	1177 (3.32-3.24)
Clashscore	141614	1044 (3.30-3.26)
Ramachandran outliers	138981	1026 (3.30-3.26)
Sidechain outliers	138945	1025 (3.30-3.26)
RSRZ outliers	127900	1141 (3.32-3.24)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	307	
1	B	307	
1	C	307	
1	D	307	
1	E	307	
1	F	307	

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 12999 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 protein called Geranylgeranyl pyrophosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	286	Total 2191	C 1426	N 359	O 399	S 7	0	0	0
1	B	285	Total 2224	C 1446	N 366	O 405	S 7	0	0	0
1	C	285	Total 2156	C 1401	N 355	O 393	S 7	0	0	0
1	D	284	Total 2191	C 1426	N 362	O 396	S 7	0	0	0
1	E	284	Total 2124	C 1370	N 352	O 395	S 7	0	0	0
1	F	285	Total 2076	C 1341	N 349	O 379	S 7	0	0	0

There are 54 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	GLY	-	expression tag	UNP O95749
A	-5	SER	-	expression tag	UNP O95749
A	-4	GLY	-	expression tag	UNP O95749
A	-3	SER	-	expression tag	UNP O95749
A	-2	GLY	-	expression tag	UNP O95749
A	-1	SER	-	expression tag	UNP O95749
A	0	GLY	-	expression tag	UNP O95749
A	109	GLN	PRO	conflict	UNP O95749
A	188	TYR	ASP	engineered mutation	UNP O95749
B	-6	GLY	-	expression tag	UNP O95749
B	-5	SER	-	expression tag	UNP O95749
B	-4	GLY	-	expression tag	UNP O95749
B	-3	SER	-	expression tag	UNP O95749
B	-2	GLY	-	expression tag	UNP O95749
B	-1	SER	-	expression tag	UNP O95749
B	0	GLY	-	expression tag	UNP O95749
B	109	GLN	PRO	conflict	UNP O95749

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Chain	Residue	Modelled	Actual	Comment	Reference
B	188	TYR	ASP	engineered mutation	UNP 095749
C	-6	GLY	-	expression tag	UNP 095749
C	-5	SER	-	expression tag	UNP 095749
C	-4	GLY	-	expression tag	UNP 095749
C	-3	SER	-	expression tag	UNP 095749
C	-2	GLY	-	expression tag	UNP 095749
C	-1	SER	-	expression tag	UNP 095749
C	0	GLY	-	expression tag	UNP 095749
C	109	GLN	PRO	conflict	UNP 095749
C	188	TYR	ASP	engineered mutation	UNP 095749
D	-6	GLY	-	expression tag	UNP 095749
D	-5	SER	-	expression tag	UNP 095749
D	-4	GLY	-	expression tag	UNP 095749
D	-3	SER	-	expression tag	UNP 095749
D	-2	GLY	-	expression tag	UNP 095749
D	-1	SER	-	expression tag	UNP 095749
D	0	GLY	-	expression tag	UNP 095749
D	109	GLN	PRO	conflict	UNP 095749
D	188	TYR	ASP	engineered mutation	UNP 095749
E	-6	GLY	-	expression tag	UNP 095749
E	-5	SER	-	expression tag	UNP 095749
E	-4	GLY	-	expression tag	UNP 095749
E	-3	SER	-	expression tag	UNP 095749
E	-2	GLY	-	expression tag	UNP 095749
E	-1	SER	-	expression tag	UNP 095749
E	0	GLY	-	expression tag	UNP 095749
E	109	GLN	PRO	conflict	UNP 095749
E	188	TYR	ASP	engineered mutation	UNP 095749
F	-6	GLY	-	expression tag	UNP 095749
F	-5	SER	-	expression tag	UNP 095749
F	-4	GLY	-	expression tag	UNP 095749
F	-3	SER	-	expression tag	UNP 095749
F	-2	GLY	-	expression tag	UNP 095749
F	-1	SER	-	expression tag	UNP 095749
F	0	GLY	-	expression tag	UNP 095749
F	109	GLN	PRO	conflict	UNP 095749
F	188	TYR	ASP	engineered mutation	UNP 095749

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 6 3 3	0	0
2	B	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	E	1	Total C O 6 3 3	0	0
2	F	1	Total C O 6 3 3	0	0

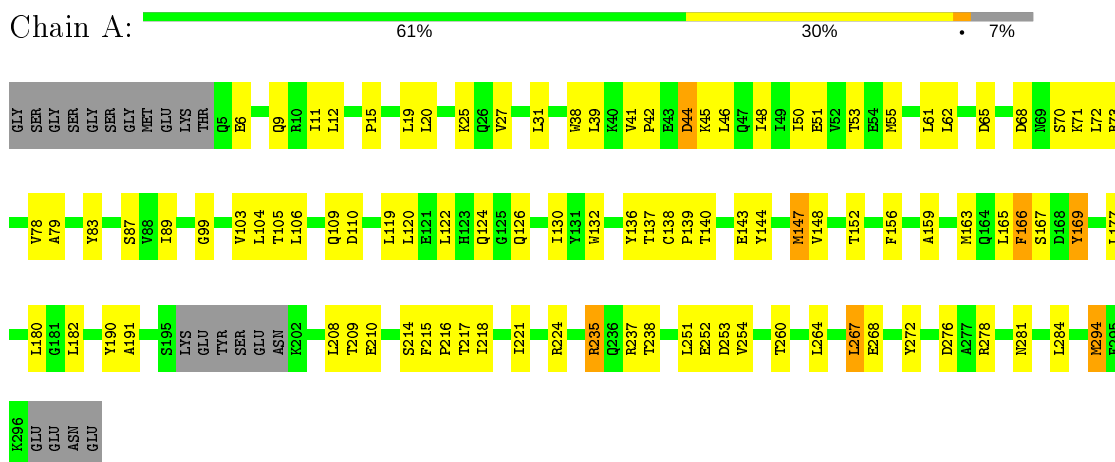
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total O 1 1	0	0

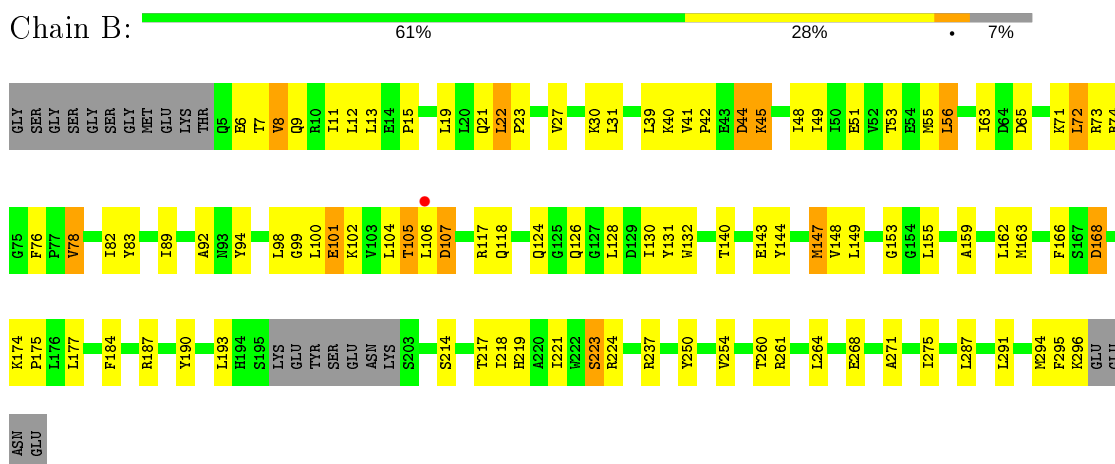
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

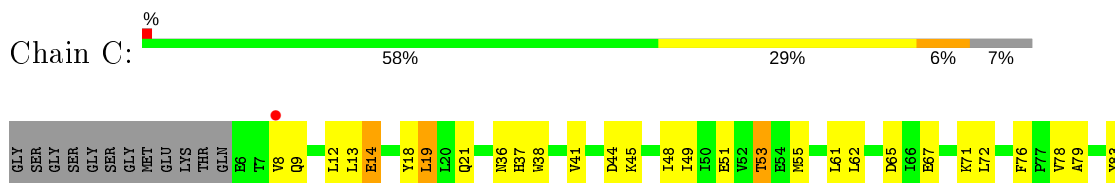
- Molecule 1: Geranylgeranyl pyrophosphate synthase

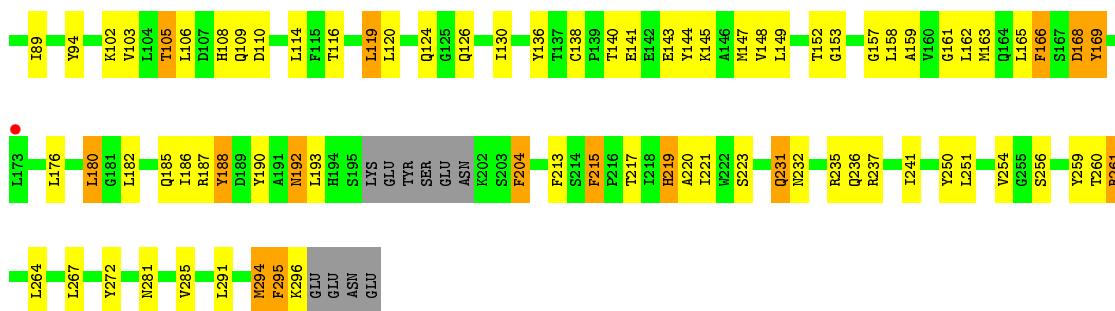


- Molecule 1: Geranylgeranyl pyrophosphate synthase



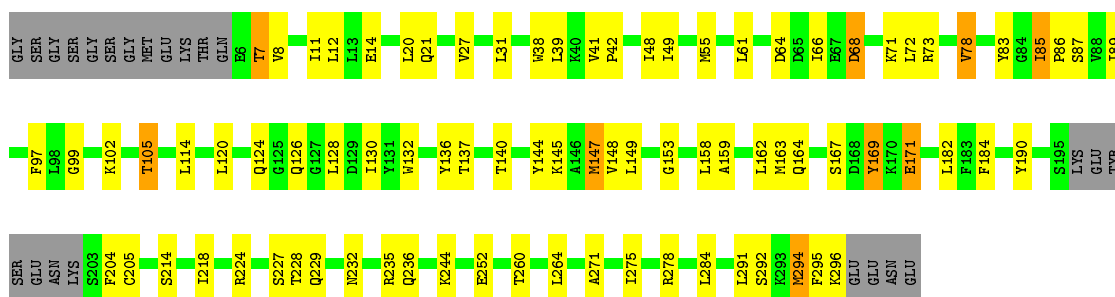
- Molecule 1: Geranylgeranyl pyrophosphate synthase





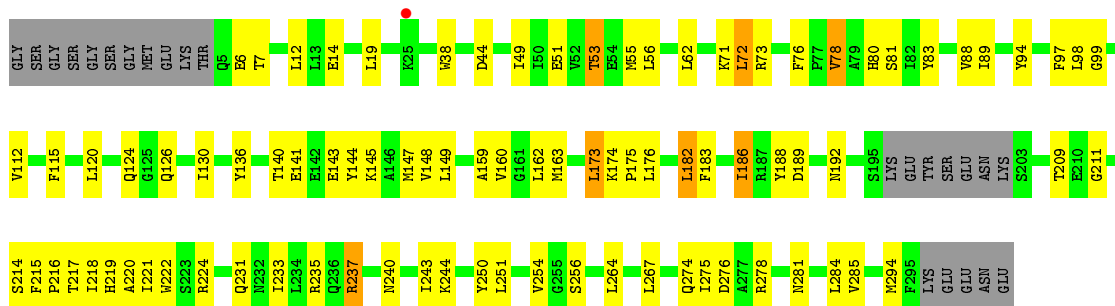
- Molecule 1: Geranylgeranyl pyrophosphate synthase

Chain D: 65% 24% 7%



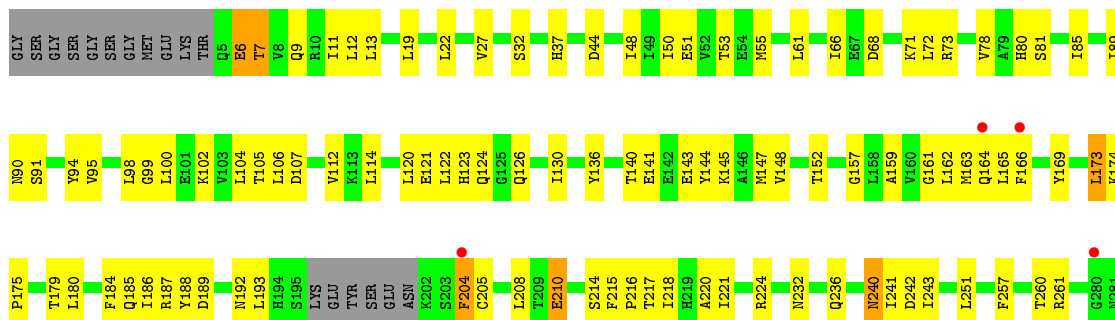
- Molecule 1: Geranylgeranyl pyrophosphate synthase

Chain E: 64% 27% 7%



- Molecule 1: Geranylgeranyl pyrophosphate synthase

Chain F: 59% 32% 7%



P282	
M294	
F295	
LYS	
GLU	
GLU	
ASN	
GLU	



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	148.54Å 148.54Å 268.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.90 – 3.28 48.90 – 3.28	Depositor EDS
% Data completeness (in resolution range)	84.3 (48.90-3.28) 84.4 (48.90-3.28)	Depositor EDS
$R_{merge}$	0.20	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.08 (at 3.25Å)	Xtrriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, $R_{free}$	0.206 , 0.258 0.210 , 0.256	Depositor DCC
$R_{free}$ test set	1978 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	89.9	Xtrriage
Anisotropy	0.030	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 86.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	12999	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	76.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.97% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: GOL

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	0.40	0/2242	0.55	0/3062
1	B	0.44	0/2275	0.55	1/3099 (0.0%)
1	C	0.36	0/2207	0.53	0/3020
1	D	0.40	0/2242	0.59	2/3059 (0.1%)
1	E	0.37	0/2175	0.53	0/2979
1	F	0.33	0/2126	0.50	0/2918
All	All	0.38	0/13267	0.54	3/18137 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	68	ASP	N-CA-C	-6.97	92.18	111.00
1	D	21	GLN	C-N-CA	-5.99	106.74	121.70
1	B	21	GLN	C-N-CA	-5.89	106.98	121.70

There are no chirality outliers.

There are no planarity outliers.

### 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	2191	0	2047	87	0
1	B	2224	0	2120	86	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2156	0	1975	101	0
1	D	2191	0	2076	83	0
1	E	2124	0	1889	84	0
1	F	2076	0	1809	110	0
2	A	6	0	8	1	0
2	B	6	0	8	0	0
2	C	6	0	8	0	0
2	D	6	0	8	0	0
2	E	6	0	8	0	0
2	F	6	0	8	1	0
3	B	1	0	0	0	0
All	All	12999	0	11964	515	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 21.

All (515) 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)
1:F:19:LEU:HD11	1:F:73:ARG:HD3	1.37	1.07
1:A:89:ILE:HD11	1:B:89:ILE:HD11	1.40	1.03
1:E:72:LEU:H	1:E:78:VAL:HG22	1.24	1.03
1:B:31:LEU:HD21	1:B:184:PHE:CE2	1.97	0.99
1:A:48:ILE:HG21	1:A:106:LEU:HD11	1.41	0.98
1:A:53:THR:HG21	1:A:159:ALA:HB2	1.45	0.97
1:B:22:LEU:HG	1:B:76:PHE:HD2	1.30	0.95
1:B:140:THR:OG1	1:B:143:GLU:HG3	1.68	0.93
1:E:140:THR:OG1	1:E:143:GLU:OE1	1.85	0.92
1:C:102:LYS:O	1:C:105:THR:OG1	1.90	0.90
1:D:55:MET:HE1	1:D:99:GLY:HA2	1.55	0.89
1:C:148:VAL:HG13	1:C:185:GLN:HG3	1.55	0.88
1:A:140:THR:OG1	1:A:143:GLU:HG3	1.74	0.86
1:C:219:HIS:CD2	1:C:256:SER:HA	2.11	0.85
1:B:31:LEU:CD2	1:B:184:PHE:HE2	1.88	0.85
1:D:102:LYS:O	1:D:105:THR:OG1	1.95	0.85
1:F:73:ARG:HH21	1:F:73:ARG:HG3	1.40	0.84
1:D:38:TRP:HB3	1:D:169:TYR:CD2	2.13	0.83
1:C:250:TYR:O	1:C:254:VAL:HG22	1.79	0.83
1:B:102:LYS:O	1:B:105:THR:OG1	1.96	0.82
1:B:31:LEU:HD21	1:B:184:PHE:HE2	1.38	0.81
1:B:22:LEU:HG	1:B:76:PHE:CD2	2.13	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:12:LEU:HD21	1:D:124:GLN:HA	1.62	0.81
1:B:71:LYS:O	1:B:78:VAL:HG22	1.81	0.80
1:C:38:TRP:HB3	1:C:169:TYR:CD2	2.18	0.79
1:C:53:THR:HG21	1:C:159:ALA:HB2	1.65	0.79
1:E:72:LEU:HD23	1:E:72:LEU:O	1.82	0.79
1:E:94:TYR:HE1	1:F:120:LEU:HD22	1.45	0.79
1:E:112:VAL:HG11	1:F:112:VAL:HG11	1.65	0.79
1:F:94:TYR:O	1:F:98:LEU:HD23	1.82	0.79
1:C:71:LYS:O	1:C:78:VAL:HG23	1.83	0.78
1:C:162:LEU:HA	1:C:165:LEU:HD13	1.63	0.78
1:B:159:ALA:O	1:B:163:MET:HG3	1.84	0.78
1:E:89:ILE:HD11	1:F:89:ILE:CD1	2.14	0.78
1:F:11:ILE:HD12	1:F:11:ILE:H	1.49	0.78
1:B:31:LEU:CD2	1:B:184:PHE:CE2	2.63	0.77
1:D:271:ALA:O	1:D:275:ILE:HG13	1.85	0.77
1:D:64:ASP:OD2	1:D:68:ASP:OD2	2.02	0.77
1:E:211:GLY:HA3	1:E:237:ARG:HH11	1.51	0.76
1:C:140:THR:OG1	1:C:143:GLU:HG3	1.86	0.76
1:E:214:SER:O	1:E:218:ILE:HG13	1.85	0.76
1:D:145:LYS:HG2	1:D:182:LEU:CD2	2.16	0.76
1:E:12:LEU:HD21	1:F:124:GLN:HA	1.68	0.75
1:E:120:LEU:HD21	1:F:98:LEU:HD22	1.67	0.75
1:F:73:ARG:NH2	1:F:73:ARG:HG3	1.98	0.74
1:E:89:ILE:HD11	1:F:89:ILE:HD11	1.70	0.74
1:F:19:LEU:HD11	1:F:73:ARG:CD	2.18	0.74
1:B:72:LEU:HD12	1:B:72:LEU:N	2.01	0.73
1:F:71:LYS:O	1:F:72:LEU:HD23	1.87	0.73
1:A:38:TRP:HB3	1:A:169:TYR:CD2	2.24	0.73
1:D:8:VAL:HA	1:D:11:ILE:HD13	1.69	0.73
1:A:109:GLN:HG3	1:A:110:ASP:OD1	1.89	0.72
1:B:31:LEU:HD21	1:B:184:PHE:CZ	2.23	0.72
1:D:71:LYS:O	1:D:72:LEU:HD12	1.90	0.72
1:E:19:LEU:HD11	1:E:73:ARG:HD2	1.71	0.72
1:E:19:LEU:HD22	1:E:62:LEU:HD23	1.72	0.71
1:A:12:LEU:HD21	1:B:124:GLN:HA	1.71	0.71
1:A:224:ARG:NH1	1:A:254:VAL:HG13	2.05	0.71
1:E:250:TYR:O	1:E:254:VAL:HG23	1.90	0.71
1:F:27:VAL:HG11	1:F:184:PHE:HE1	1.56	0.71
1:D:184:PHE:HE1	1:D:291:LEU:HD13	1.56	0.70
1:B:65:ASP:OD2	1:B:73:ARG:NE	2.24	0.70
1:F:152:THR:HG21	1:F:185:GLN:HG2	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:271:ALA:O	1:B:275:ILE:HG13	1.92	0.69
1:C:72:LEU:HB2	1:C:76:PHE:O	1.93	0.69
1:A:120:LEU:HD11	1:B:101:GLU:HG3	1.74	0.68
1:C:219:HIS:HD2	1:C:256:SER:HA	1.59	0.68
1:D:144:TYR:O	1:D:148:VAL:HG23	1.93	0.68
1:F:240:ASN:HD22	1:F:241:ILE:N	1.90	0.68
1:B:187:ARG:HE	1:B:294:MET:HG2	1.55	0.68
1:A:126:GLN:O	1:A:130:ILE:HG13	1.93	0.68
1:E:126:GLN:O	1:E:130:ILE:HG13	1.93	0.68
1:A:38:TRP:HB3	1:A:169:TYR:CE2	2.29	0.68
1:A:281:ASN:ND2	1:A:284:LEU:HB2	2.09	0.67
1:E:144:TYR:OH	1:E:189:ASP:OD2	2.12	0.67
1:C:19:LEU:HD13	1:C:19:LEU:O	1.94	0.67
1:B:250:TYR:O	1:B:254:VAL:HG23	1.93	0.67
1:B:72:LEU:HD12	1:B:72:LEU:H	1.57	0.67
1:A:89:ILE:CD1	1:B:89:ILE:HD11	2.23	0.67
1:D:136:TYR:HE1	1:D:235:ARG:HG2	1.60	0.67
1:D:55:MET:HE1	1:D:99:GLY:CA	2.23	0.67
1:C:38:TRP:HB3	1:C:169:TYR:CE2	2.29	0.67
1:E:281:ASN:ND2	1:E:284:LEU:HB2	2.10	0.66
1:A:224:ARG:CZ	1:A:254:VAL:HG13	2.26	0.66
1:C:161:GLY:O	1:C:165:LEU:HD12	1.95	0.66
1:E:94:TYR:CE1	1:F:120:LEU:HD22	2.30	0.66
1:D:145:LYS:HG2	1:D:182:LEU:HD21	1.76	0.66
1:E:55:MET:HE1	1:E:99:GLY:HA2	1.77	0.66
1:C:176:LEU:O	1:C:180:LEU:HD22	1.96	0.66
1:F:144:TYR:O	1:F:148:VAL:HG23	1.96	0.66
1:C:152:THR:HG21	1:C:185:GLN:HG2	1.78	0.65
1:E:233:ILE:HD13	1:E:243:ILE:HG23	1.77	0.65
1:D:7:THR:O	1:D:11:ILE:HD12	1.97	0.65
1:A:48:ILE:HG21	1:A:106:LEU:CD1	2.21	0.65
1:C:126:GLN:O	1:C:130:ILE:HG13	1.96	0.65
1:F:173:LEU:HD12	1:F:173:LEU:H	1.60	0.65
1:A:39:LEU:O	1:A:167:SER:OG	2.14	0.64
1:E:44:ASP:OD1	1:E:44:ASP:N	2.27	0.64
1:A:214:SER:O	1:A:218:ILE:HG13	1.98	0.64
1:A:48:ILE:CG2	1:A:106:LEU:HD11	2.21	0.64
1:C:8:VAL:HG11	1:D:128:LEU:HD21	1.79	0.64
1:F:11:ILE:N	1:F:11:ILE:HD12	2.12	0.64
1:F:94:TYR:CE2	1:F:98:LEU:HD21	2.31	0.64
1:F:126:GLN:O	1:F:130:ILE:HG13	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:GLN:HA	1:B:12:LEU:HD11	1.79	0.63
1:F:162:LEU:HA	1:F:165:LEU:HD13	1.80	0.63
1:C:187:ARG:NH2	1:C:291:LEU:HD22	2.14	0.63
1:C:51:GLU:O	1:C:55:MET:HG3	1.98	0.63
1:D:214:SER:O	1:D:218:ILE:HG13	1.99	0.63
1:A:217:THR:O	1:A:221:ILE:HG13	1.97	0.63
1:F:148:VAL:HG13	1:F:185:GLN:HG3	1.81	0.62
1:F:78:VAL:HG21	1:F:80:HIS:CE1	2.33	0.62
1:B:41:VAL:HG12	1:B:42:PRO:O	1.99	0.62
1:B:65:ASP:CG	1:B:73:ARG:HE	2.03	0.62
1:F:78:VAL:HG22	1:F:81:SER:HB3	1.81	0.62
1:A:51:GLU:O	1:A:55:MET:HG3	1.99	0.62
1:D:85:ILE:HD11	1:D:89:ILE:CD1	2.30	0.62
1:B:224:ARG:CZ	1:B:254:VAL:HG13	2.29	0.62
1:A:144:TYR:O	1:A:148:VAL:HG23	2.00	0.61
1:D:8:VAL:HA	1:D:11:ILE:CD1	2.30	0.61
1:C:217:THR:O	1:C:221:ILE:HG13	2.00	0.61
1:F:217:THR:O	1:F:221:ILE:HG13	2.00	0.61
1:A:25:LYS:CB	1:A:27:VAL:HG23	2.30	0.61
1:B:184:PHE:HE1	1:B:291:LEU:HD13	1.66	0.61
1:C:124:GLN:HA	1:D:12:LEU:HD11	1.82	0.61
1:F:140:THR:HB	1:F:143:GLU:OE1	2.01	0.61
1:F:257:PHE:O	1:F:261:ARG:HD3	2.00	0.61
1:A:103:VAL:O	1:A:105:THR:O	2.18	0.61
1:D:190:TYR:HB2	1:D:260:THR:HG21	1.81	0.61
1:F:48:ILE:HG21	1:F:106:LEU:CD2	2.31	0.60
1:A:105:THR:O	1:A:106:LEU:HD13	2.02	0.60
1:D:38:TRP:HB3	1:D:169:TYR:HD2	1.66	0.60
1:C:186:ILE:HD13	1:C:215:PHE:CE1	2.36	0.60
1:E:38:TRP:CZ2	1:E:281:ASN:HB2	2.36	0.60
1:B:31:LEU:CD1	1:B:287:LEU:HD21	2.31	0.59
1:F:180:LEU:O	1:F:184:PHE:HD2	1.85	0.59
1:A:12:LEU:CD2	1:B:124:GLN:HA	2.32	0.59
1:B:44:ASP:N	1:B:44:ASP:OD1	2.29	0.59
1:C:148:VAL:CG1	1:C:185:GLN:HG3	2.30	0.59
1:B:9:GLN:OE1	1:B:94:TYR:OH	2.16	0.59
1:C:149:LEU:O	1:C:153:GLY:HA3	2.03	0.59
1:F:11:ILE:CD1	1:F:11:ILE:H	2.15	0.59
1:A:44:ASP:O	1:A:48:ILE:HD13	2.03	0.59
1:C:219:HIS:O	1:C:223:SER:OG	2.21	0.59
1:A:106:LEU:HB3	1:A:166:PHE:HZ	1.68	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:126:GLN:O	1:D:130:ILE:HG13	2.03	0.59
1:C:219:HIS:ND1	1:C:223:SER:OG	2.35	0.58
1:C:109:GLN:HG3	1:C:110:ASP:OD1	2.03	0.58
1:E:140:THR:HG22	1:E:222:TRP:CZ2	2.38	0.58
1:B:55:MET:HE1	1:B:99:GLY:N	2.18	0.58
1:D:171:GLU:OE1	1:D:278:ARG:NE	2.25	0.58
1:F:51:GLU:O	1:F:55:MET:HG3	2.03	0.58
1:D:11:ILE:H	1:D:11:ILE:HD12	1.68	0.58
1:E:78:VAL:HG23	1:E:81:SER:HB2	1.85	0.58
1:E:51:GLU:O	1:E:55:MET:HG3	2.04	0.58
1:D:38:TRP:HB3	1:D:169:TYR:CE2	2.39	0.57
1:E:294:MET:HG3	1:E:294:MET:O	2.03	0.57
1:F:161:GLY:O	1:F:165:LEU:HD12	2.03	0.57
1:A:6:GLU:HA	1:A:6:GLU:OE2	2.04	0.57
1:A:252:GLU:HA	1:A:252:GLU:OE2	2.04	0.57
1:D:136:TYR:CE1	1:D:235:ARG:HG2	2.38	0.57
1:E:120:LEU:CD2	1:F:98:LEU:HD22	2.32	0.57
1:E:173:LEU:H	1:E:173:LEU:HD12	1.69	0.57
1:E:211:GLY:HA3	1:E:237:ARG:NH1	2.19	0.57
1:F:204:PHE:HE1	1:F:205:CYS:HG	1.52	0.56
1:F:48:ILE:HG21	1:F:106:LEU:HD21	1.87	0.56
1:B:8:VAL:HA	1:B:11:ILE:HD12	1.87	0.56
1:D:39:LEU:HD13	1:D:163:MET:HB2	1.87	0.56
1:D:66:ILE:HD13	1:D:85:ILE:CD1	2.35	0.56
1:D:149:LEU:O	1:D:153:GLY:HA3	2.05	0.56
1:D:164:GLN:HA	1:D:167:SER:HB3	1.87	0.56
1:B:49:ILE:CD1	1:B:162:LEU:HB3	2.35	0.56
1:D:275:ILE:HD13	1:D:284:LEU:HD23	1.87	0.56
1:F:164:GLN:HA	1:F:164:GLN:OE1	2.06	0.56
1:C:219:HIS:CD2	1:C:259:TYR:HD2	2.24	0.56
1:A:235:ARG:HG2	1:A:235:ARG:NH1	2.21	0.56
1:B:264:LEU:HD13	1:B:295:PHE:CD1	2.41	0.56
1:C:182:LEU:HB3	1:C:267:LEU:HD21	1.87	0.56
1:C:152:THR:HG21	1:C:185:GLN:CG	2.36	0.55
1:A:105:THR:HG22	1:A:106:LEU:N	2.21	0.55
1:C:188:TYR:HD1	1:C:188:TYR:O	1.88	0.55
1:E:53:THR:HG21	1:E:159:ALA:HB2	1.88	0.55
1:C:141:GLU:O	1:C:145:LYS:HG3	2.07	0.55
1:E:186:ILE:HD13	1:E:215:PHE:CD2	2.41	0.55
1:F:164:GLN:NE2	1:F:173:LEU:HD11	2.22	0.55
1:A:209:THR:OG1	1:A:238:THR:O	2.18	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:189:ASP:O	1:F:193:LEU:HD13	2.06	0.55
1:E:281:ASN:O	1:E:285:VAL:HG23	2.07	0.55
1:F:215:PHE:HB3	1:F:216:PRO:HD3	1.87	0.55
1:A:191:ALA:HB2	1:A:294:MET:SD	2.47	0.55
1:C:204:PHE:HE2	1:C:241:ILE:HG23	1.72	0.55
1:F:240:ASN:HD22	1:F:240:ASN:C	2.09	0.55
1:F:73:ARG:HH21	1:F:73:ARG:CG	2.14	0.54
1:A:159:ALA:O	1:A:163:MET:HG3	2.07	0.54
1:A:89:ILE:HD11	1:B:89:ILE:CD1	2.28	0.54
1:E:124:GLN:HA	1:F:12:LEU:HD21	1.90	0.54
1:B:15:PRO:HA	1:B:83:TYR:CE2	2.41	0.54
1:C:18:TYR:O	1:C:21:GLN:HG2	2.07	0.54
1:D:71:LYS:C	1:D:72:LEU:HD12	2.28	0.54
1:E:174:LYS:N	1:E:175:PRO:HD2	2.22	0.54
1:F:141:GLU:O	1:F:145:LYS:HG3	2.06	0.54
1:B:168:ASP:N	1:B:168:ASP:OD1	2.39	0.54
1:F:243:ILE:HD12	1:F:243:ILE:N	2.22	0.54
1:E:72:LEU:HA	1:E:76:PHE:O	2.07	0.54
1:F:9:GLN:O	1:F:13:LEU:HD13	2.08	0.54
1:D:85:ILE:HD12	1:D:89:ILE:HG13	1.90	0.54
1:D:204:PHE:HE2	1:D:244:LYS:HB3	1.73	0.54
1:C:67:GLU:HA	1:D:86:PRO:HB3	1.90	0.54
1:F:61:LEU:HD22	1:F:73:ARG:NH1	2.23	0.54
1:F:48:ILE:CG2	1:F:106:LEU:HD21	2.39	0.54
1:B:190:TYR:HB2	1:B:260:THR:HG21	1.89	0.53
1:B:23:PRO:O	1:B:74:ARG:HD3	2.08	0.53
1:C:186:ILE:HD13	1:C:215:PHE:HE1	1.72	0.53
1:C:45:LYS:O	1:C:49:ILE:HG13	2.07	0.53
1:B:144:TYR:O	1:B:148:VAL:HG23	2.09	0.53
1:F:214:SER:O	1:F:218:ILE:HG13	2.08	0.53
1:A:105:THR:O	1:A:106:LEU:HB2	2.08	0.53
1:B:214:SER:O	1:B:218:ILE:HG13	2.09	0.53
1:B:82:ILE:HG22	1:B:83:TYR:CD1	2.44	0.53
1:D:159:ALA:O	1:D:163:MET:HG3	2.09	0.53
1:E:220:ALA:HB2	1:E:251:LEU:CD2	2.39	0.53
1:B:174:LYS:HA	1:B:177:LEU:HD12	1.89	0.53
1:C:162:LEU:HA	1:C:165:LEU:CD1	2.35	0.53
1:C:192:ASN:HB3	1:C:193:LEU:HD12	1.91	0.53
1:C:144:TYR:O	1:C:148:VAL:HG23	2.08	0.53
1:F:159:ALA:O	1:F:163:MET:HG3	2.09	0.53
1:A:235:ARG:HH11	1:A:235:ARG:HG2	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:LEU:HD11	1:B:287:LEU:HD21	1.90	0.52
1:C:38:TRP:HB3	1:C:169:TYR:HD2	1.71	0.52
1:E:219:HIS:HD1	1:E:256:SER:HG	0.61	0.52
1:A:214:SER:HB3	1:A:216:PRO:HD2	1.92	0.52
1:D:85:ILE:HD11	1:D:89:ILE:HD12	1.90	0.52
1:F:114:LEU:HD23	1:F:114:LEU:O	2.10	0.52
1:A:140:THR:HG1	1:A:143:GLU:HG3	1.74	0.52
1:C:295:PHE:O	1:C:296:LYS:CB	2.57	0.52
1:D:66:ILE:HD13	1:D:85:ILE:HD12	1.91	0.52
1:A:61:LEU:HD22	1:A:73:ARG:NH2	2.25	0.52
1:C:231:GLN:O	1:C:235:ARG:HG3	2.09	0.51
1:A:166:PHE:CD2	1:A:166:PHE:N	2.79	0.51
1:F:140:THR:CB	1:F:143:GLU:OE1	2.58	0.51
1:B:126:GLN:O	1:B:130:ILE:HG13	2.10	0.51
1:C:219:HIS:CE1	1:C:223:SER:OG	2.63	0.51
1:E:160:VAL:HA	1:E:163:MET:CE	2.41	0.51
1:E:186:ILE:HB	1:E:264:LEU:HD21	1.92	0.51
1:E:224:ARG:NE	1:E:254:VAL:HG13	2.26	0.51
1:E:55:MET:CE	1:E:99:GLY:HA2	2.40	0.50
1:B:219:HIS:O	1:B:223:SER:OG	2.23	0.50
1:C:159:ALA:O	1:C:163:MET:HG3	2.11	0.50
1:C:219:HIS:HD1	1:C:223:SER:HG	1.58	0.50
1:F:13:LEU:HD12	1:F:13:LEU:N	2.26	0.50
1:F:9:GLN:OE1	1:F:94:TYR:OH	2.27	0.50
1:C:294:MET:CE	1:C:294:MET:HA	2.42	0.50
1:C:71:LYS:C	1:C:78:VAL:CG2	2.80	0.50
1:C:108:HIS:CE1	1:C:165:LEU:HD23	2.47	0.50
1:E:49:ILE:HD13	1:E:162:LEU:HB3	1.94	0.50
1:E:176:LEU:HA	1:E:274:GLN:OE1	2.12	0.50
1:F:173:LEU:HD12	1:F:173:LEU:N	2.26	0.49
1:C:193:LEU:N	1:C:193:LEU:HD12	2.27	0.49
1:E:243:ILE:HD12	1:E:243:ILE:N	2.27	0.49
1:A:235:ARG:HH11	1:A:235:ARG:CG	2.26	0.49
1:A:45:LYS:HD2	1:A:166:PHE:CE1	2.47	0.49
1:D:31:LEU:N	1:D:31:LEU:HD12	2.28	0.49
1:E:149:LEU:HD21	1:E:182:LEU:CD1	2.42	0.49
1:E:250:TYR:CE2	1:E:254:VAL:HG21	2.47	0.49
1:F:102:LYS:O	1:F:105:THR:HG23	2.12	0.49
1:F:104:LEU:HD23	1:F:112:VAL:HG22	1.95	0.49
1:F:186:ILE:HD13	1:F:215:PHE:CG	2.48	0.49
1:C:37:HIS:HB3	1:C:281:ASN:ND2	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:228:THR:N	1:E:14:GLU:OE2	2.43	0.49
1:D:232:ASN:O	1:D:236:GLN:HG3	2.12	0.49
1:D:49:ILE:HD13	1:D:162:LEU:HB3	1.94	0.49
1:F:162:LEU:HA	1:F:165:LEU:CD1	2.41	0.49
1:A:46:LEU:O	1:A:50:ILE:HG13	2.13	0.49
1:F:44:ASP:O	1:F:48:ILE:HD13	2.13	0.49
1:C:136:TYR:CD1	1:C:235:ARG:HG2	2.48	0.49
1:A:106:LEU:HB3	1:A:166:PHE:CZ	2.45	0.48
1:E:144:TYR:O	1:E:148:VAL:HG23	2.13	0.48
1:E:217:THR:O	1:E:221:ILE:HG13	2.13	0.48
1:E:160:VAL:HA	1:E:163:MET:HE2	1.94	0.48
1:E:89:ILE:HD11	1:F:89:ILE:HD13	1.91	0.48
1:D:41:VAL:HG13	1:D:42:PRO:HD2	1.95	0.48
1:D:55:MET:CE	1:D:99:GLY:HA2	2.37	0.48
1:F:215:PHE:HD2	1:F:260:THR:HG1	1.59	0.48
1:A:31:LEU:HD12	1:A:31:LEU:N	2.29	0.48
1:A:41:VAL:CG1	1:A:42:PRO:HD2	2.44	0.48
1:D:227:SER:HA	1:E:14:GLU:OE2	2.14	0.48
1:C:65:ASP:HB3	1:C:78:VAL:HG11	1.96	0.48
1:C:71:LYS:C	1:C:78:VAL:HG23	2.33	0.48
1:E:173:LEU:C	1:E:175:PRO:HD2	2.34	0.48
1:F:37:HIS:CD2	1:F:282:PRO:HD2	2.48	0.48
1:B:295:PHE:HD2	1:B:295:PHE:O	1.97	0.48
1:C:295:PHE:N	1:C:295:PHE:CD1	2.82	0.48
1:A:55:MET:HE1	1:A:99:GLY:N	2.29	0.48
1:C:190:TYR:HB2	1:C:260:THR:HG21	1.96	0.48
1:D:295:PHE:O	1:D:296:LYS:HG2	2.13	0.48
1:A:20:LEU:HA	1:A:20:LEU:HD23	1.47	0.48
1:C:72:LEU:C	1:C:72:LEU:HD12	2.34	0.48
1:F:243:ILE:HD12	1:F:243:ILE:H	1.79	0.48
1:B:31:LEU:HD12	1:B:287:LEU:CD2	2.44	0.48
1:D:132:TRP:HZ3	1:D:147:MET:HG2	1.79	0.48
1:B:117:ARG:HG3	1:B:118:GLN:N	2.29	0.47
1:E:97:PHE:HD2	1:F:123:HIS:HD2	1.60	0.47
1:A:190:TYR:HB2	1:A:260:THR:HG21	1.95	0.47
1:F:6:GLU:OE2	1:F:6:GLU:HA	2.14	0.47
1:C:72:LEU:HA	1:C:78:VAL:HG22	1.96	0.47
1:F:106:LEU:O	1:F:107:ASP:HB3	2.15	0.47
1:F:68:ASP:O	1:F:210:GLU:HG2	2.13	0.47
1:A:65:ASP:OD1	1:A:73:ARG:NH1	2.47	0.47
1:A:105:THR:HG22	1:A:106:LEU:H	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:157:GLY:O	1:C:161:GLY:N	2.42	0.47
1:F:53:THR:HG21	1:F:159:ALA:HB2	1.97	0.47
1:B:27:VAL:O	1:B:31:LEU:HD13	2.15	0.47
1:B:149:LEU:O	1:B:153:GLY:HA3	2.13	0.47
1:C:215:PHE:HD2	1:C:215:PHE:O	1.97	0.47
1:D:145:LYS:HG2	1:D:182:LEU:HD22	1.95	0.47
1:F:66:ILE:HD13	1:F:85:ILE:HG23	1.97	0.47
1:A:104:LEU:HA	1:A:104:LEU:HD23	1.76	0.47
1:A:165:LEU:HB2	1:A:166:PHE:CD2	2.50	0.47
1:C:72:LEU:HA	1:C:78:VAL:CG2	2.45	0.47
1:B:184:PHE:CE1	1:B:291:LEU:HD13	2.48	0.47
1:C:36:ASN:OD1	1:C:41:VAL:HG23	2.15	0.47
1:F:7:THR:O	1:F:11:ILE:CD1	2.63	0.47
1:A:53:THR:HG21	1:A:159:ALA:CB	2.31	0.46
1:E:233:ILE:HG23	1:E:243:ILE:CG2	2.45	0.46
1:E:56:LEU:HD11	1:E:115:PHE:HE1	1.80	0.46
1:F:187:ARG:HE	1:F:294:MET:HG2	1.80	0.46
1:B:19:LEU:O	1:B:19:LEU:HD23	2.15	0.46
1:B:107:ASP:CG	1:B:107:ASP:O	2.53	0.46
1:B:39:LEU:HD13	1:B:163:MET:HB2	1.97	0.46
1:B:71:LYS:O	1:B:78:VAL:CG2	2.59	0.46
1:C:37:HIS:HB3	1:C:281:ASN:HD21	1.81	0.46
1:E:278:ARG:HA	1:E:278:ARG:HE	1.79	0.46
1:E:6:GLU:O	1:E:7:THR:OG1	2.26	0.46
1:F:122:LEU:HD12	1:F:122:LEU:O	2.15	0.46
1:D:235:ARG:HD3	1:F:136:TYR:CE2	2.51	0.46
1:F:208:LEU:HA	1:F:208:LEU:HD23	1.70	0.46
1:F:48:ILE:HG21	1:F:106:LEU:HD23	1.97	0.46
1:A:65:ASP:CG	1:A:73:ARG:HH11	2.18	0.46
1:B:31:LEU:CD1	1:B:287:LEU:CD2	2.93	0.46
1:C:103:VAL:HA	1:C:106:LEU:HD23	1.98	0.46
1:C:44:ASP:O	1:C:48:ILE:HD13	2.16	0.46
1:E:141:GLU:HG2	1:E:145:LYS:HE3	1.96	0.46
1:E:224:ARG:CD	1:E:254:VAL:HG13	2.45	0.46
1:A:62:LEU:HD22	1:A:79:ALA:CB	2.45	0.46
1:E:55:MET:HE1	1:E:99:GLY:CA	2.44	0.46
1:A:180:LEU:HA	1:A:180:LEU:HD23	1.80	0.46
1:B:49:ILE:HD11	1:B:162:LEU:HB3	1.97	0.46
1:D:49:ILE:CD1	1:D:162:LEU:HB3	2.46	0.46
1:F:94:TYR:CE2	1:F:98:LEU:CD2	2.99	0.46
1:F:94:TYR:CZ	1:F:98:LEU:HD21	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:ASP:O	1:A:210:GLU:HG2	2.15	0.45
1:A:46:LEU:HD11	1:A:163:MET:HE1	1.98	0.45
1:F:121:GLU:OE1	1:F:121:GLU:HA	2.17	0.45
1:F:165:LEU:H	1:F:165:LEU:HD12	1.80	0.45
1:C:232:ASN:O	1:C:236:GLN:HG3	2.15	0.45
1:C:261:ARG:HG3	1:C:295:PHE:HB3	1.99	0.45
1:C:62:LEU:HD22	1:C:79:ALA:CB	2.46	0.45
1:D:275:ILE:CD1	1:D:284:LEU:HD23	2.47	0.45
1:D:264:LEU:HD13	1:D:295:PHE:CE1	2.51	0.45
1:D:64:ASP:CG	1:D:68:ASP:OD2	2.54	0.45
1:B:100:LEU:O	1:B:104:LEU:HG	2.16	0.45
1:B:44:ASP:O	1:B:48:ILE:HD13	2.16	0.45
1:B:56:LEU:H	1:B:56:LEU:HD23	1.81	0.45
1:D:85:ILE:CD1	1:D:89:ILE:CD1	2.94	0.45
1:C:166:PHE:CD2	1:C:166:PHE:N	2.85	0.45
1:D:61:LEU:HD22	1:D:73:ARG:NH1	2.31	0.45
1:F:220:ALA:HB2	1:F:251:LEU:CD2	2.47	0.45
1:C:138:CYS:HB2	1:C:213:PHE:CE1	2.52	0.45
1:C:89:ILE:HD12	1:D:89:ILE:CD1	2.47	0.45
1:C:8:VAL:O	1:C:8:VAL:HG12	2.16	0.45
1:E:218:ILE:O	1:E:222:TRP:HD1	2.00	0.45
1:E:72:LEU:C	1:E:72:LEU:HD23	2.37	0.45
1:A:208:LEU:HD23	1:A:208:LEU:HA	1.81	0.45
1:A:19:LEU:HD22	1:A:62:LEU:CD2	2.47	0.45
1:B:56:LEU:HD23	1:B:56:LEU:N	2.32	0.45
1:C:94:TYR:HE1	1:D:120:LEU:CD2	2.29	0.45
1:A:105:THR:C	1:A:106:LEU:CD1	2.85	0.45
1:B:132:TRP:HZ3	1:B:147:MET:HG2	1.82	0.45
1:D:14:GLU:HG2	1:D:83:TYR:CE1	2.51	0.45
1:D:264:LEU:HA	1:D:264:LEU:HD23	1.61	0.45
1:F:55:MET:HE1	1:F:99:GLY:N	2.32	0.45
1:C:8:VAL:HG11	1:D:128:LEU:CD2	2.47	0.45
1:D:204:PHE:CD2	1:D:205:CYS:HB2	2.52	0.45
1:E:221:ILE:HD13	1:E:231:GLN:HG3	1.99	0.45
1:F:224:ARG:NH1	1:F:224:ARG:HG3	2.32	0.45
1:F:48:ILE:HD12	1:F:48:ILE:N	2.32	0.45
1:F:104:LEU:CD2	1:F:112:VAL:HG22	2.47	0.44
1:F:251:LEU:HD23	1:F:251:LEU:HA	1.74	0.44
1:B:31:LEU:HD12	1:B:287:LEU:HD21	1.99	0.44
1:A:68:ASP:OD1	2:A:1001:GOL:H2	2.17	0.44
1:B:190:TYR:OH	1:B:261:ARG:NH1	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:13:LEU:HD11	1:B:94:TYR:HE2	1.83	0.44
1:E:215:PHE:HB3	1:E:216:PRO:HD3	2.00	0.44
1:E:98:LEU:HD23	1:E:98:LEU:HA	1.87	0.44
1:B:56:LEU:CD2	1:B:56:LEU:N	2.81	0.44
1:C:220:ALA:HB2	1:C:251:LEU:CD2	2.48	0.44
1:A:215:PHE:HB3	1:A:216:PRO:HD3	2.00	0.44
1:D:27:VAL:O	1:D:31:LEU:HD13	2.18	0.44
1:A:138:CYS:SG	1:A:139:PRO:HD2	2.58	0.44
1:B:128:LEU:HA	1:B:128:LEU:HD23	1.73	0.44
1:C:116:THR:HG22	1:C:120:LEU:HD12	1.98	0.44
1:C:168:ASP:N	1:C:168:ASP:OD1	2.36	0.44
1:C:219:HIS:CG	1:C:259:TYR:HD2	2.36	0.44
1:C:187:ARG:CZ	1:C:291:LEU:HD22	2.47	0.44
1:D:49:ILE:HD13	1:D:162:LEU:CB	2.48	0.44
1:D:85:ILE:HD11	1:D:89:ILE:HD11	2.00	0.44
1:F:188:TYR:CE1	2:F:1001:GOL:O2	2.69	0.43
1:A:138:CYS:SG	1:A:221:ILE:HD12	2.58	0.43
1:C:14:GLU:OE1	1:C:83:TYR:CE1	2.72	0.43
1:E:186:ILE:CG2	1:E:264:LEU:HD21	2.48	0.43
1:E:186:ILE:HD13	1:E:215:PHE:CG	2.53	0.43
1:E:71:LYS:O	1:E:72:LEU:HD22	2.18	0.43
1:E:141:GLU:O	1:E:145:LYS:HG3	2.18	0.43
1:F:224:ARG:HG3	1:F:224:ARG:HH11	1.83	0.43
1:A:46:LEU:HD12	1:A:163:MET:HE2	2.00	0.43
1:C:149:LEU:O	1:C:153:GLY:CA	2.67	0.43
1:E:149:LEU:HD21	1:E:182:LEU:HD12	2.01	0.43
1:E:240:ASN:OD1	1:E:243:ILE:HD13	2.17	0.43
1:A:182:LEU:HB3	1:A:267:LEU:HD11	2.00	0.43
1:F:157:GLY:O	1:F:161:GLY:N	2.47	0.43
1:F:106:LEU:O	1:F:107:ASP:CB	2.67	0.43
1:A:122:LEU:C	1:A:122:LEU:HD23	2.39	0.43
1:A:251:LEU:HA	1:A:251:LEU:HD23	1.80	0.43
1:B:63:ILE:HD11	1:B:92:ALA:HB3	2.01	0.43
1:D:264:LEU:HB3	1:D:295:PHE:CZ	2.54	0.43
1:C:119:LEU:HD12	1:D:97:PHE:CE2	2.54	0.43
1:B:41:VAL:HG13	1:B:42:PRO:HD2	2.01	0.43
1:B:55:MET:HE3	1:B:55:MET:HB3	1.78	0.43
1:D:114:LEU:CD2	1:D:158:LEU:HD12	2.49	0.43
1:A:264:LEU:HD23	1:A:264:LEU:HA	1.70	0.42
1:C:41:VAL:HG21	1:C:163:MET:HB3	2.01	0.42
1:E:209:THR:OG1	1:E:244:LYS:HE2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:152:THR:HG21	1:F:185:GLN:CG	2.46	0.42
1:F:32:SER:CB	1:F:50:ILE:CD1	2.97	0.42
1:D:41:VAL:HG21	1:D:163:MET:HB3	2.01	0.42
1:B:45:LYS:HD3	1:B:166:PHE:CZ	2.54	0.42
1:E:173:LEU:H	1:E:173:LEU:CD1	2.29	0.42
1:A:124:GLN:HG2	1:B:12:LEU:CD1	2.49	0.42
1:C:294:MET:O	1:C:294:MET:HG3	2.19	0.42
1:D:291:LEU:O	1:D:294:MET:HB3	2.19	0.42
1:F:166:PHE:CD2	1:F:166:PHE:N	2.88	0.42
1:A:177:LEU:HA	1:A:177:LEU:HD23	1.90	0.42
1:C:158:LEU:HD23	1:C:158:LEU:O	2.20	0.42
1:C:186:ILE:HG23	1:C:215:PHE:CD1	2.54	0.42
1:F:140:THR:HG22	1:F:141:GLU:N	2.35	0.42
1:B:217:THR:O	1:B:221:ILE:HG13	2.19	0.42
1:C:188:TYR:C	1:C:188:TYR:CD1	2.93	0.42
1:E:183:PHE:HA	1:E:267:LEU:HD23	2.01	0.42
1:F:91:SER:O	1:F:95:VAL:HG23	2.19	0.42
1:A:235:ARG:NH1	1:A:235:ARG:CG	2.83	0.42
1:B:155:LEU:N	1:B:155:LEU:HD22	2.35	0.42
1:C:94:TYR:HE1	1:D:120:LEU:HD22	1.85	0.42
1:D:292:SER:C	1:D:294:MET:H	2.23	0.42
1:D:85:ILE:HG13	1:D:86:PRO:N	2.34	0.42
1:E:120:LEU:CD2	1:F:98:LEU:CD2	2.98	0.42
1:F:100:LEU:O	1:F:104:LEU:HG	2.19	0.42
1:A:268:GLU:HG2	1:A:272:TYR:CE2	2.55	0.42
1:A:41:VAL:HG13	1:A:42:PRO:HD2	2.02	0.42
1:A:71:LYS:O	1:A:72:LEU:HG	2.20	0.42
1:A:38:TRP:CZ2	1:A:278:ARG:HB2	2.55	0.42
1:B:6:GLU:O	1:B:6:GLU:HG3	2.20	0.42
1:C:9:GLN:O	1:C:13:LEU:HD13	2.20	0.42
1:D:169:TYR:CE1	1:D:171:GLU:HG3	2.55	0.42
1:E:136:TYR:CE2	1:E:235:ARG:HG3	2.55	0.42
1:A:15:PRO:HA	1:A:83:TYR:CE2	2.55	0.41
1:B:224:ARG:NH1	1:B:254:VAL:HG13	2.35	0.41
1:C:48:ILE:HD12	1:C:48:ILE:N	2.35	0.41
1:E:6:GLU:C	1:E:7:THR:HG23	2.39	0.41
1:F:32:SER:CB	1:F:50:ILE:HD11	2.50	0.41
1:F:19:LEU:CD1	1:F:73:ARG:CD	2.94	0.41
1:C:264:LEU:HB3	1:C:295:PHE:CZ	2.55	0.41
1:C:272:TYR:CE1	1:C:285:VAL:HG13	2.54	0.41
1:F:105:THR:O	1:F:106:LEU:C	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:268:GLU:OE2	1:B:296:LYS:NZ	2.42	0.41
1:D:66:ILE:HD13	1:D:85:ILE:HD13	2.02	0.41
1:A:105:THR:O	1:A:106:LEU:CB	2.69	0.41
1:F:232:ASN:O	1:F:236:GLN:HG3	2.20	0.41
1:A:132:TRP:CZ3	1:A:147:MET:HG2	2.56	0.41
1:C:165:LEU:H	1:C:165:LEU:HD12	1.86	0.41
1:A:11:ILE:HG22	1:B:131:TYR:CE1	2.56	0.41
1:B:174:LYS:N	1:B:175:PRO:CD	2.84	0.41
1:E:188:TYR:HE2	1:E:192:ASN:HD22	1.67	0.41
1:E:55:MET:HE3	1:E:55:MET:HB3	1.85	0.41
1:D:169:TYR:CZ	1:D:171:GLU:CD	2.94	0.41
1:E:80:HIS:HA	1:E:88:VAL:HG21	2.03	0.41
1:F:174:LYS:N	1:F:175:PRO:CD	2.83	0.41
1:A:55:MET:HE1	1:A:99:GLY:CA	2.50	0.41
1:C:188:TYR:HD1	1:C:188:TYR:C	2.24	0.41
1:F:192:ASN:HB3	1:F:193:LEU:HD12	2.03	0.41
1:A:152:THR:O	1:A:156:PHE:HD2	2.03	0.41
1:B:193:LEU:N	1:B:193:LEU:HD12	2.36	0.41
1:B:19:LEU:C	1:B:19:LEU:HD23	2.41	0.41
1:B:98:LEU:HD23	1:B:98:LEU:HA	1.82	0.41
1:C:18:TYR:O	1:C:21:GLN:CG	2.69	0.41
1:C:61:LEU:HA	1:C:61:LEU:HD23	1.87	0.41
1:D:39:LEU:HD13	1:D:163:MET:CB	2.49	0.41
1:D:55:MET:CE	1:D:99:GLY:CA	2.97	0.41
1:E:97:PHE:HE2	1:F:123:HIS:NE2	2.19	0.41
1:A:136:TYR:CE2	1:A:235:ARG:CZ	3.04	0.40
1:D:85:ILE:N	1:D:86:PRO:HD2	2.36	0.40
1:E:112:VAL:HG11	1:F:112:VAL:CG1	2.42	0.40
1:B:51:GLU:O	1:B:55:MET:HG3	2.20	0.40
1:C:13:LEU:N	1:C:13:LEU:HD12	2.35	0.40
1:D:169:TYR:CE1	1:D:171:GLU:CG	3.05	0.40
1:F:22:LEU:HA	1:F:22:LEU:HD12	1.87	0.40
1:C:264:LEU:HD23	1:C:264:LEU:HA	1.84	0.40
1:D:229:GLN:HG2	1:E:83:TYR:OH	2.22	0.40
1:D:71:LYS:O	1:D:78:VAL:HG22	2.21	0.40
1:F:193:LEU:N	1:F:193:LEU:HD12	2.36	0.40
1:B:106:LEU:HA	1:B:106:LEU:HD23	1.85	0.40
1:B:55:MET:HE1	1:B:99:GLY:CA	2.51	0.40
1:C:158:LEU:HD23	1:C:162:LEU:HG	2.02	0.40
1:D:48:ILE:HD12	1:D:48:ILE:N	2.37	0.40
1:F:68:ASP:HB3	1:F:210:GLU:HG2	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:240:ASN:ND2	1:F:242:ASP:H	2.19	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	
1	A	282/307 (92%)	278 (99%)	4 (1%)	0	100	100
1	B	281/307 (92%)	278 (99%)	3 (1%)	0	100	100
1	C	281/307 (92%)	279 (99%)	2 (1%)	0	100	100
1	D	280/307 (91%)	277 (99%)	3 (1%)	0	100	100
1	E	280/307 (91%)	277 (99%)	3 (1%)	0	100	100
1	F	281/307 (92%)	277 (99%)	4 (1%)	0	100	100
All	All	1685/1842 (92%)	1666 (99%)	19 (1%)	0	100	100

There are no Ramachandran outliers to report.

#### 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	
1	A	214/275 (78%)	198 (92%)	16 (8%)	13	39
1	B	224/275 (82%)	206 (92%)	18 (8%)	12	37

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	205/275 (74%)	184 (90%)	21 (10%)	7	27
1	D	218/275 (79%)	204 (94%)	14 (6%)	17	47
1	E	198/275 (72%)	188 (95%)	10 (5%)	24	54
1	F	183/275 (66%)	172 (94%)	11 (6%)	19	49
All	All	1242/1650 (75%)	1152 (93%)	90 (7%)	14	41

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	44	ASP
1	A	70	SER
1	A	78	VAL
1	A	87	SER
1	A	119	LEU
1	A	137	THR
1	A	147	MET
1	A	166	PHE
1	A	169	TYR
1	A	235	ARG
1	A	237	ARG
1	A	253	ASP
1	A	267	LEU
1	A	276	ASP
1	A	294	MET
1	B	7	THR
1	B	8	VAL
1	B	22	LEU
1	B	30	LYS
1	B	40	LYS
1	B	44	ASP
1	B	45	LYS
1	B	53	THR
1	B	56	LEU
1	B	72	LEU
1	B	78	VAL
1	B	101	GLU
1	B	105	THR
1	B	107	ASP
1	B	147	MET
1	B	168	ASP

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	B	223	SER
1	B	237	ARG
1	C	14	GLU
1	C	19	LEU
1	C	53	THR
1	C	105	THR
1	C	114	LEU
1	C	119	LEU
1	C	147	MET
1	C	166	PHE
1	C	168	ASP
1	C	169	TYR
1	C	180	LEU
1	C	188	TYR
1	C	192	ASN
1	C	204	PHE
1	C	215	PHE
1	C	219	HIS
1	C	231	GLN
1	C	237	ARG
1	C	261	ARG
1	C	294	MET
1	C	295	PHE
1	D	7	THR
1	D	20	LEU
1	D	78	VAL
1	D	85	ILE
1	D	87	SER
1	D	105	THR
1	D	137	THR
1	D	140	THR
1	D	147	MET
1	D	169	TYR
1	D	171	GLU
1	D	224	ARG
1	D	252	GLU
1	D	294	MET
1	E	53	THR
1	E	72	LEU
1	E	78	VAL
1	E	147	MET
1	E	173	LEU

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Mol	Chain	Res	Type
1	E	182	LEU
1	E	186	ILE
1	E	237	ARG
1	E	275	ILE
1	E	276	ASP
1	F	6	GLU
1	F	7	THR
1	F	90	ASN
1	F	147	MET
1	F	169	TYR
1	F	173	LEU
1	F	179	THR
1	F	204	PHE
1	F	210	GLU
1	F	240	ASN
1	F	294	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	240	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

### 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 carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	GOL	F	1001	-	5,5,5	0.34	0	5,5,5	0.13	0
2	GOL	B	1001	-	5,5,5	0.46	0	5,5,5	0.55	0
2	GOL	E	1001	-	5,5,5	0.39	0	5,5,5	0.26	0
2	GOL	A	1001	-	5,5,5	0.70	0	5,5,5	0.70	0
2	GOL	C	1001	-	5,5,5	0.40	0	5,5,5	0.18	0
2	GOL	D	1001	-	5,5,5	0.26	0	5,5,5	0.63	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	F	1001	-	-	3/4/4/4	-
2	GOL	B	1001	-	-	4/4/4/4	-
2	GOL	E	1001	-	-	4/4/4/4	-
2	GOL	A	1001	-	-	0/4/4/4	-
2	GOL	C	1001	-	-	0/4/4/4	-
2	GOL	D	1001	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1001	GOL	O1-C1-C2-C3
2	E	1001	GOL	O1-C1-C2-C3
2	E	1001	GOL	C1-C2-C3-O3
2	D	1001	GOL	O1-C1-C2-C3
2	B	1001	GOL	O1-C1-C2-O2
2	F	1001	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
2	E	1001	GOL	O1-C1-C2-O2
2	E	1001	GOL	O2-C2-C3-O3
2	D	1001	GOL	O1-C1-C2-O2
2	F	1001	GOL	O1-C1-C2-O2
2	B	1001	GOL	C1-C2-C3-O3
2	B	1001	GOL	O2-C2-C3-O3
2	F	1001	GOL	O1-C1-C2-C3

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	1001	GOL	1	0
2	A	1001	GOL	1	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	286/307 (93%)	-0.06	0 <a href="#">100</a>   <a href="#">100</a>	53, 68, 90, 107	0
1	B	285/307 (92%)	-0.14	1 (0%) <a href="#">92</a>   <a href="#">93</a>	54, 64, 77, 92	0
1	C	285/307 (92%)	0.02	2 (0%) <a href="#">87</a>   <a href="#">88</a>	28, 76, 97, 113	0
1	D	284/307 (92%)	-0.07	0 <a href="#">100</a>   <a href="#">100</a>	53, 70, 84, 96	0
1	E	284/307 (92%)	-0.05	1 (0%) <a href="#">92</a>   <a href="#">93</a>	63, 87, 114, 126	0
1	F	285/307 (92%)	0.01	4 (1%) <a href="#">75</a>   <a href="#">74</a>	61, 90, 116, 128	0
All	All	1709/1842 (92%)	-0.05	8 (0%) <a href="#">91</a>   <a href="#">91</a>	28, 74, 108, 128	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	25	LYS	2.9
1	F	204	PHE	2.6
1	C	173	LEU	2.2
1	F	280	GLY	2.2
1	F	164	GLN	2.2
1	F	166	PHE	2.0
1	B	106	LEU	2.0
1	C	8	VAL	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 carbohydrates 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
2	GOL	A	1001	6/6	0.68	0.29	76,76,76,76	0
2	GOL	C	1001	6/6	0.82	0.30	76,76,76,76	0
2	GOL	B	1001	6/6	0.83	0.70	62,62,62,62	0
2	GOL	F	1001	6/6	0.84	0.36	92,92,92,92	0
2	GOL	E	1001	6/6	0.88	0.34	90,90,90,90	0
2	GOL	D	1001	6/6	0.93	0.47	61,61,61,61	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.