



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 16, 2023 – 12:19 AM EDT

PDB ID : 1U76  
Title : Crystal structure of hPCNA bound to residues 452-466 of the DNA polymerase-delta-p66 subunit  
Authors : Bruning, J.B.; Shamoo, Y.  
Deposited on : 2004-08-02  
Resolution : 2.60 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.

The types of validation reports are described at

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

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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  
Xtriage (Phenix) : 1.13  
EDS : 2.35  
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) : 2.35

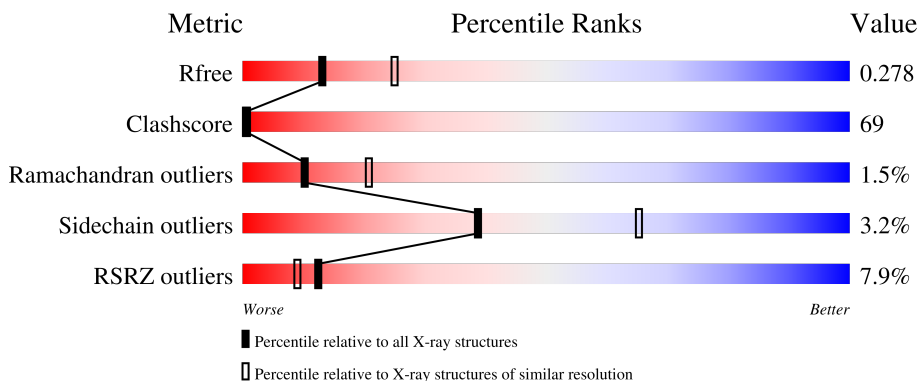
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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.



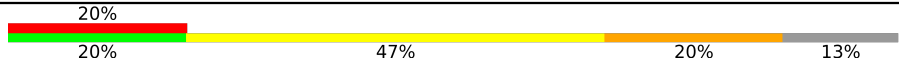
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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  $\geq 3$ , 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	261	
1	C	261	
1	E	261	
2	B	15	
2	D	15	

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Mol	Chain	Length	Quality of chain
2	F	15	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into five segments: a red segment (20%), a green segment (20%), a yellow segment (47%), an orange segment (20%), and a grey segment (13%).</p>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6701 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 Proliferating cell nuclear antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	253	Total 1934	C 1217	N 316	O 385	S 16	0	0	0
1	C	251	Total 1923	C 1211	N 314	O 382	S 16	0	0	0
1	E	253	Total 1933	C 1216	N 318	O 383	S 16	0	0	0

- Molecule 2 is a protein called KANRQVSITGFFQRK peptide from DNA polymerase delta subunit 3.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	12	Total 96	C 61	N 18	O 17	0	0	0
2	D	13	Total 107	C 67	N 22	O 18	0	0	0
2	F	13	Total 107	C 67	N 22	O 18	0	0	0

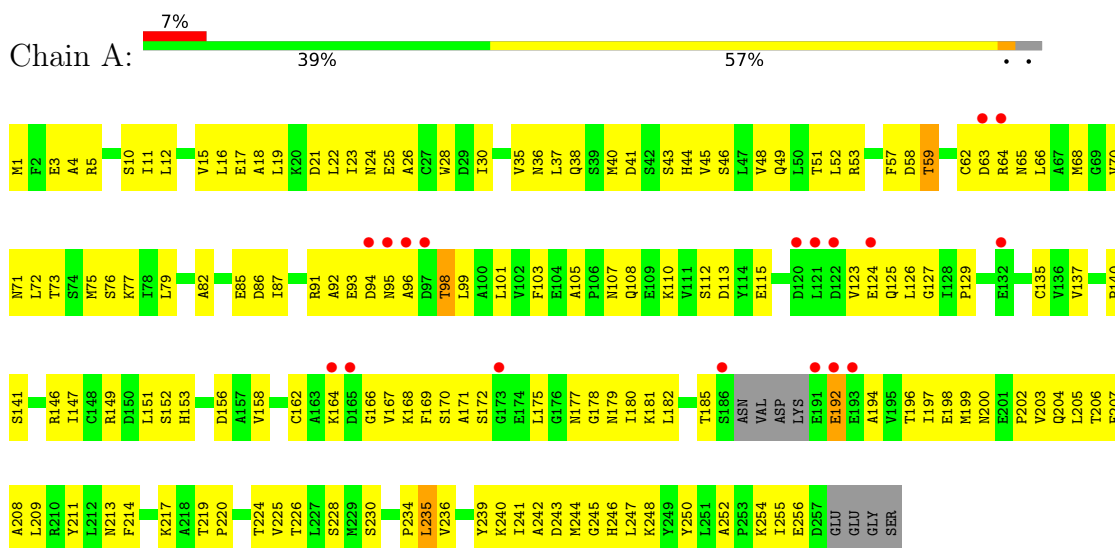
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	181	Total 181	O 181	0	0
3	B	10	Total 10	O 10	0	0
3	C	197	Total 197	O 197	0	0
3	D	11	Total 11	O 11	0	0
3	E	187	Total 187	O 187	0	0
3	F	15	Total 15	O 15	0	0

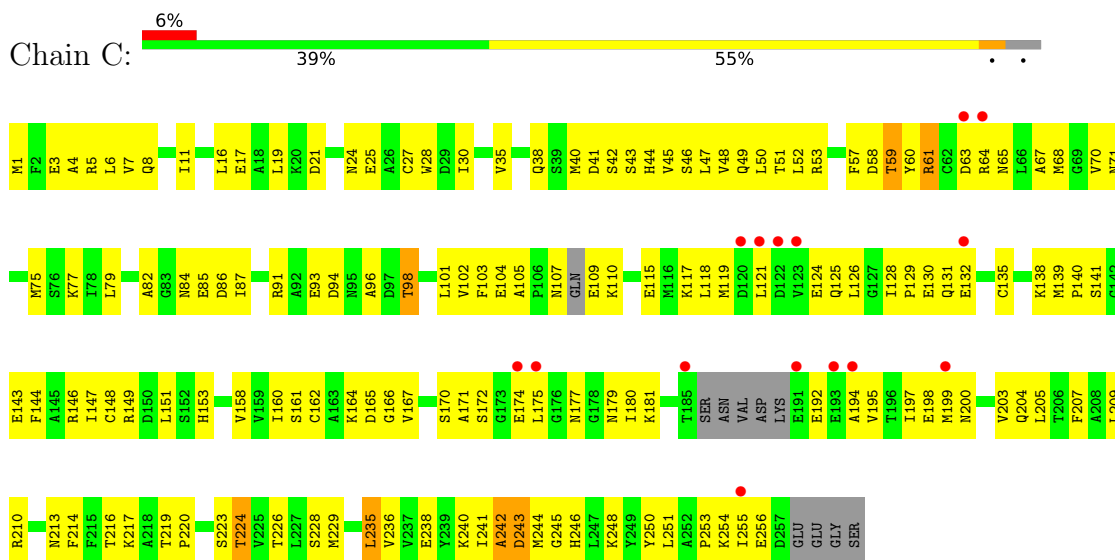
### 3 Residue-property plots

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: Proliferating cell nuclear antigen

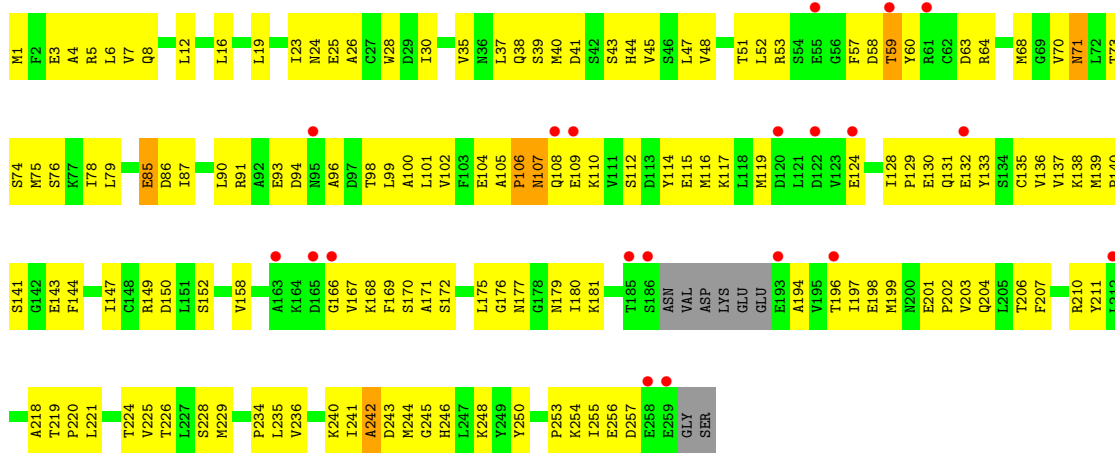


- Molecule 1: Proliferating cell nuclear antigen

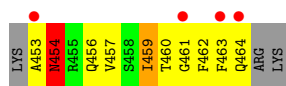
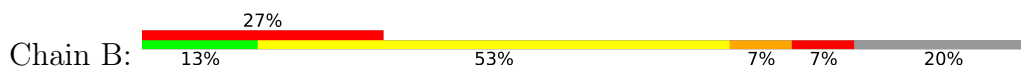


- Molecule 1: Proliferating cell nuclear antigen

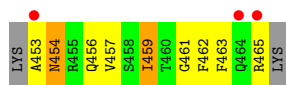




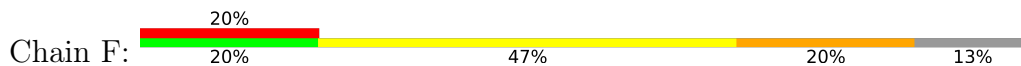
● Molecule 2: KANRQVSITGFFQRK peptide from DNA polymerase delta subunit 3



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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	82.48Å 82.48Å 203.80Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.60 29.25 – 2.60	Depositor EDS
% Data completeness (in resolution range)	94.0 (10.00-2.60) 95.5 (29.25-2.60)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.50 (at 2.61Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.240 , 0.278 0.241 , 0.278	Depositor DCC
$R_{free}$ test set	1150 reflections (4.70%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	58.0	Xtrriage
Anisotropy	0.391	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 71.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	0.000 for -h,-k,l	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6701	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	65.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 25.58 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.0861e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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](#)

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	0.38	0/1959	0.63	0/2646
1	C	0.39	0/1947	0.62	0/2628
1	E	0.38	0/1958	0.64	0/2645
2	B	0.51	0/97	0.61	0/129
2	D	0.47	0/108	0.58	0/143
2	F	0.50	0/108	0.62	0/143
All	All	0.39	0/6177	0.63	0/8334

There are no bond length outliers.

There are no bond angle outliers.

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	1934	0	1934	265	0
1	C	1923	0	1926	270	0
1	E	1933	0	1936	269	1
2	B	96	0	92	23	0
2	D	107	0	105	29	0
2	F	107	0	105	24	0
3	A	181	0	0	166	0
3	B	10	0	0	11	0
3	C	197	0	0	204	0
3	D	11	0	0	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	187	0	0	199	1
3	F	15	0	0	10	0
All	All	6701	0	6098	844	2

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 69.

All (844) 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:E:64:ARG:HA	3:E:444:HOH:O	1.33	1.27
1:E:236:VAL:HA	3:E:420:HOH:O	1.35	1.26
1:A:40:MET:HB3	3:A:351:HOH:O	1.37	1.24
1:E:73:THR:HB	3:E:395:HOH:O	1.30	1.23
2:F:456:GLN:HG3	3:F:540:HOH:O	1.38	1.21
1:C:194:ALA:HB1	3:C:343:HOH:O	1.45	1.14
2:B:460:THR:HA	3:B:525:HOH:O	1.48	1.13
1:A:126:LEU:HB3	3:A:366:HOH:O	1.47	1.12
1:E:152:SER:HA	3:E:443:HOH:O	1.49	1.11
1:E:133:TYR:HA	3:E:448:HOH:O	1.47	1.11
1:E:171:ALA:HB2	3:E:349:HOH:O	1.48	1.11
1:A:23:ILE:HA	3:A:343:HOH:O	1.49	1.11
1:C:102:VAL:HG23	3:C:368:HOH:O	1.50	1.10
1:A:217:LYS:HG3	3:A:339:HOH:O	1.51	1.09
1:E:45:VAL:HB	3:E:376:HOH:O	1.53	1.09
1:A:73:THR:HA	3:A:402:HOH:O	1.53	1.08
1:A:203:VAL:HA	3:A:422:HOH:O	1.51	1.08
1:A:49:GLN:HB2	3:A:439:HOH:O	1.54	1.07
1:E:105:ALA:HB3	3:E:348:HOH:O	1.53	1.07
3:C:385:HOH:O	2:D:454:ASN:HB3	1.53	1.06
1:A:36:ASN:HB2	3:A:334:HOH:O	1.55	1.06
1:C:131:GLN:HB2	3:C:426:HOH:O	1.56	1.04
1:A:16:LEU:HD22	1:A:79:LEU:HD12	1.38	1.04
1:E:108:GLN:HA	3:E:383:HOH:O	1.57	1.04
1:A:203:VAL:HB	3:A:348:HOH:O	1.57	1.04
1:A:58:ASP:HB3	3:A:338:HOH:O	1.57	1.03
1:A:248:LYS:HD3	3:A:409:HOH:O	1.57	1.03
1:E:16:LEU:HD22	1:E:79:LEU:HD12	1.41	1.01
1:E:242:ALA:HB2	3:E:417:HOH:O	1.59	1.01
1:C:24:ASN:HB3	3:C:374:HOH:O	1.58	1.01
1:E:35:VAL:HB	3:E:398:HOH:O	1.56	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:LEU:HA	3:A:357:HOH:O	1.61	1.00
1:E:114:TYR:HA	3:E:340:HOH:O	1.60	0.99
1:C:243:ASP:HB2	3:C:401:HOH:O	1.64	0.97
1:E:73:THR:HA	3:E:406:HOH:O	1.63	0.97
1:A:239:TYR:HB3	3:A:427:HOH:O	1.64	0.97
1:A:252:ALA:HB3	3:A:346:HOH:O	1.63	0.97
1:C:131:GLN:HB3	3:C:373:HOH:O	1.64	0.97
1:C:194:ALA:HB3	3:C:377:HOH:O	1.65	0.97
1:A:169:PHE:HB2	3:A:425:HOH:O	1.62	0.96
1:A:101:LEU:HG	3:A:405:HOH:O	1.63	0.96
1:E:23:ILE:HA	3:E:425:HOH:O	1.65	0.96
1:C:124:GLU:HG3	2:D:465:ARG:HH21	1.28	0.96
1:A:247:LEU:HD12	3:A:404:HOH:O	1.65	0.95
1:E:85:GLU:HG2	3:E:360:HOH:O	1.66	0.95
1:E:206:THR:OG1	2:F:453:ALA:HB2	1.65	0.95
1:E:48:VAL:HG23	3:E:401:HOH:O	1.66	0.95
1:A:103:PHE:HE2	3:A:377:HOH:O	1.50	0.94
2:D:457:VAL:HA	3:D:272:HOH:O	1.65	0.93
1:E:102:VAL:HG12	3:E:346:HOH:O	1.66	0.93
1:C:170:SER:HB3	1:C:179:ASN:HB3	1.51	0.93
1:A:44:HIS:HA	3:A:351:HOH:O	1.67	0.93
1:A:103:PHE:CE2	3:A:377:HOH:O	2.21	0.93
1:E:96:ALA:HB3	3:E:338:HOH:O	1.67	0.93
1:C:70:VAL:HG11	3:C:405:HOH:O	1.68	0.92
1:C:130:GLU:HG3	3:C:389:HOH:O	1.69	0.92
1:A:153:HIS:HE1	3:A:431:HOH:O	1.54	0.91
1:E:225:VAL:HG22	3:E:385:HOH:O	1.67	0.91
1:A:76:SER:HB3	3:A:402:HOH:O	1.72	0.90
1:E:170:SER:HB3	1:E:179:ASN:HB3	1.54	0.88
1:C:240:LYS:HE2	3:C:439:HOH:O	1.71	0.88
1:E:25:GLU:HG2	3:E:370:HOH:O	1.74	0.88
1:A:94:ASP:HB3	3:A:365:HOH:O	1.73	0.87
1:E:52:LEU:HB2	3:E:398:HOH:O	1.72	0.87
1:C:117:LYS:HB3	3:C:421:HOH:O	1.73	0.87
1:A:170:SER:HB3	1:A:179:ASN:HB3	1.55	0.87
1:E:76:SER:HB3	3:E:406:HOH:O	1.74	0.87
1:C:153:HIS:CB	3:C:347:HOH:O	2.21	0.86
1:A:48:VAL:HG12	3:A:404:HOH:O	1.75	0.86
1:A:151:LEU:HB2	3:A:437:HOH:O	1.75	0.85
1:E:167:VAL:HG22	3:E:367:HOH:O	1.76	0.85
1:C:16:LEU:HD22	1:C:79:LEU:HD12	1.55	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:197:ILE:HG22	3:A:396:HOH:O	1.76	0.84
1:A:167:VAL:HG22	3:A:378:HOH:O	1.78	0.84
1:C:160:ILE:HG22	3:C:408:HOH:O	1.76	0.84
1:A:18:ALA:C	3:A:435:HOH:O	2.16	0.84
1:C:91:ARG:HB2	3:C:397:HOH:O	1.76	0.83
1:C:224:THR:HB	3:C:386:HOH:O	1.76	0.83
1:E:139:MET:HG2	3:E:426:HOH:O	1.78	0.83
1:A:199:MET:HB2	3:A:396:HOH:O	1.77	0.82
1:E:150:ASP:N	3:E:433:HOH:O	2.11	0.82
1:A:203:VAL:CA	3:A:422:HOH:O	2.17	0.82
1:A:219:THR:HA	3:A:354:HOH:O	1.79	0.82
1:C:181:LYS:HG3	3:C:348:HOH:O	1.78	0.82
1:A:169:PHE:HD1	3:A:425:HOH:O	1.62	0.82
1:C:75:MET:SD	3:C:405:HOH:O	2.36	0.81
1:C:214:PHE:HA	3:C:330:HOH:O	1.79	0.81
1:E:107:ASN:HB2	3:E:342:HOH:O	1.79	0.81
1:E:225:VAL:HG13	3:E:385:HOH:O	1.81	0.81
1:E:5:ARG:HB3	1:E:59:THR:HB	1.63	0.81
1:E:149:ARG:HB3	3:E:433:HOH:O	1.81	0.81
1:E:43:SER:HB3	3:E:364:HOH:O	1.80	0.80
1:E:48:VAL:N	3:E:401:HOH:O	2.14	0.80
1:E:94:ASP:N	3:E:439:HOH:O	2.15	0.80
1:A:103:PHE:CB	3:A:412:HOH:O	2.29	0.80
1:C:91:ARG:HD3	3:C:397:HOH:O	1.81	0.80
1:C:148:CYS:SG	3:C:444:HOH:O	2.40	0.80
1:A:254:LYS:HE2	3:B:271:HOH:O	1.82	0.80
1:C:57:PHE:HB3	3:C:455:HOH:O	1.82	0.79
1:C:61:ARG:HG2	3:C:396:HOH:O	1.81	0.79
1:E:201:GLU:HA	3:E:388:HOH:O	1.83	0.79
1:A:169:PHE:CD1	3:A:425:HOH:O	2.33	0.79
1:A:156:ASP:HB3	3:A:423:HOH:O	1.82	0.78
1:C:58:ASP:HA	3:C:356:HOH:O	1.84	0.78
1:A:239:TYR:C	3:A:427:HOH:O	2.21	0.78
1:A:213:ASN:C	3:A:434:HOH:O	2.21	0.78
2:B:464:GLN:NE2	2:B:464:GLN:H	1.80	0.78
1:E:106:PRO:HB2	3:E:350:HOH:O	1.84	0.78
1:A:182:LEU:N	3:A:393:HOH:O	2.16	0.78
1:E:211:TYR:CE1	3:F:449:HOH:O	2.35	0.78
1:A:5:ARG:HB3	1:A:59:THR:HB	1.64	0.77
1:C:194:ALA:HA	3:C:403:HOH:O	1.83	0.77
1:C:5:ARG:HB3	1:C:59:THR:HB	1.66	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:194:ALA:C	3:C:403:HOH:O	2.23	0.77
1:E:254:LYS:HG3	3:E:358:HOH:O	1.84	0.77
1:E:201:GLU:CA	3:E:388:HOH:O	2.33	0.77
1:A:110:LYS:HE2	3:A:420:HOH:O	1.84	0.77
1:E:90:LEU:HA	3:E:405:HOH:O	1.83	0.76
1:A:203:VAL:CB	3:A:348:HOH:O	2.21	0.76
1:E:196:THR:HG23	3:E:434:HOH:O	1.85	0.76
1:E:256:GLU:HA	3:E:374:HOH:O	1.85	0.76
1:C:50:LEU:HA	3:C:431:HOH:O	1.85	0.76
1:C:5:ARG:N	3:C:455:HOH:O	2.17	0.76
1:C:42:SER:HA	3:C:355:HOH:O	1.87	0.75
1:C:192:GLU:HB2	3:C:338:HOH:O	1.85	0.75
1:C:94:ASP:N	3:C:415:HOH:O	2.19	0.75
1:C:242:ALA:O	3:C:401:HOH:O	2.04	0.75
1:A:156:ASP:N	3:A:423:HOH:O	2.20	0.75
1:C:146:ARG:CZ	3:C:400:HOH:O	2.35	0.75
1:A:225:VAL:CA	3:A:358:HOH:O	2.34	0.74
1:A:167:VAL:C	3:A:378:HOH:O	2.26	0.74
1:C:6:LEU:C	3:C:422:HOH:O	2.25	0.74
1:A:147:ILE:HG23	1:A:180:ILE:HD12	1.69	0.74
1:A:207:PHE:HA	3:A:428:HOH:O	1.88	0.74
1:A:92:ALA:HA	3:A:357:HOH:O	1.87	0.74
1:E:139:MET:N	3:E:426:HOH:O	2.19	0.74
1:E:5:ARG:HB2	3:E:422:HOH:O	1.87	0.74
1:A:203:VAL:HG13	3:A:422:HOH:O	1.88	0.73
1:E:114:TYR:CA	3:E:340:HOH:O	2.24	0.73
1:C:124:GLU:HG3	2:D:465:ARG:NH2	2.02	0.73
1:C:175:LEU:HG	3:C:360:HOH:O	1.89	0.73
3:C:372:HOH:O	2:D:456:GLN:NE2	2.20	0.73
1:E:137:VAL:HG23	3:E:347:HOH:O	1.88	0.73
1:E:59:THR:C	3:E:381:HOH:O	2.27	0.73
1:E:255:ILE:HD13	2:F:457:VAL:HG21	1.69	0.73
1:A:254:LYS:HB3	3:B:271:HOH:O	1.89	0.73
1:E:5:ARG:CG	3:E:422:HOH:O	2.36	0.72
2:D:454:ASN:N	3:D:393:HOH:O	2.20	0.72
1:A:123:VAL:HG12	3:A:370:HOH:O	1.88	0.72
1:E:199:MET:HG2	3:E:396:HOH:O	1.89	0.72
1:A:49:GLN:HG2	3:A:386:HOH:O	1.87	0.72
1:E:202:PRO:HD2	3:E:388:HOH:O	1.88	0.72
2:F:456:GLN:C	3:F:540:HOH:O	2.26	0.72
1:C:64:ARG:HG2	3:C:351:HOH:O	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:58:ASP:HB2	3:C:328:HOH:O	1.89	0.72
1:E:94:ASP:HB3	3:E:375:HOH:O	1.90	0.72
1:A:203:VAL:N	3:A:348:HOH:O	2.22	0.71
1:A:225:VAL:C	3:A:406:HOH:O	2.28	0.71
1:A:180:ILE:HD11	3:E:421:HOH:O	1.89	0.71
1:E:3:GLU:HG2	3:E:368:HOH:O	1.90	0.71
1:A:18:ALA:HB3	3:A:435:HOH:O	1.90	0.71
1:A:22:LEU:HD21	3:A:356:HOH:O	1.90	0.71
1:C:6:LEU:N	3:C:422:HOH:O	2.22	0.71
2:F:464:GLN:HG2	3:F:303:HOH:O	1.90	0.71
1:A:112:SER:HB3	3:C:458:HOH:O	1.91	0.71
1:C:118:LEU:HB3	3:C:345:HOH:O	1.90	0.71
1:C:194:ALA:CA	3:C:403:HOH:O	2.38	0.71
1:C:224:THR:CB	3:C:386:HOH:O	2.34	0.71
1:E:130:GLU:HA	3:E:399:HOH:O	1.89	0.71
1:E:175:LEU:HG	3:E:294:HOH:O	1.91	0.70
1:E:58:ASP:HB2	3:E:341:HOH:O	1.90	0.70
1:A:18:ALA:CB	3:A:435:HOH:O	2.39	0.70
1:A:200:ASN:N	3:A:436:HOH:O	2.23	0.70
1:E:35:VAL:CB	3:E:398:HOH:O	2.25	0.70
1:E:101:LEU:HD12	3:E:392:HOH:O	1.90	0.70
1:A:214:PHE:N	3:A:434:HOH:O	2.24	0.70
1:C:151:LEU:HD21	3:C:362:HOH:O	1.91	0.70
1:E:78:ILE:CG2	3:E:442:HOH:O	2.39	0.70
1:E:119:MET:SD	3:E:370:HOH:O	2.49	0.70
1:E:199:MET:CG	3:E:396:HOH:O	2.39	0.70
1:C:153:HIS:ND1	3:C:347:HOH:O	2.23	0.70
1:E:3:GLU:OE2	1:E:91:ARG:HD3	1.92	0.70
1:A:66:LEU:HD23	3:A:400:HOH:O	1.92	0.69
1:A:214:PHE:HA	3:A:434:HOH:O	1.91	0.69
1:A:244:MET:C	3:A:382:HOH:O	2.30	0.69
1:C:253:PRO:C	3:C:372:HOH:O	2.31	0.69
1:A:225:VAL:HA	3:A:358:HOH:O	1.90	0.69
1:C:119:MET:HE1	3:C:337:HOH:O	1.91	0.69
1:A:63:ASP:C	3:A:359:HOH:O	2.31	0.69
1:C:50:LEU:CA	3:C:431:HOH:O	2.39	0.69
1:A:180:ILE:HG13	3:E:421:HOH:O	1.92	0.69
1:C:70:VAL:C	3:C:381:HOH:O	2.30	0.69
1:E:255:ILE:HD12	3:E:296:HOH:O	1.90	0.69
1:A:126:LEU:HD21	3:B:525:HOH:O	1.93	0.69
1:E:60:TYR:C	3:E:381:HOH:O	2.31	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:LYS:N	3:A:378:HOH:O	2.25	0.69
1:E:53:ARG:NE	3:E:365:HOH:O	2.25	0.69
1:E:135:CYS:SG	3:E:396:HOH:O	2.50	0.69
1:E:138:LYS:C	3:E:426:HOH:O	2.31	0.69
2:F:465:ARG:HG2	2:F:465:ARG:HH11	1.56	0.69
1:C:255:ILE:HD13	2:D:457:VAL:HG21	1.75	0.69
1:E:254:LYS:HG2	2:F:456:GLN:HA	1.74	0.69
1:C:138:LYS:HG3	3:C:369:HOH:O	1.92	0.68
1:C:144:PHE:CD1	3:C:444:HOH:O	2.46	0.68
1:E:37:LEU:HD23	3:E:404:HOH:O	1.93	0.68
1:E:58:ASP:C	3:E:341:HOH:O	2.32	0.68
1:A:66:LEU:HA	3:A:400:HOH:O	1.93	0.68
3:C:385:HOH:O	2:D:454:ASN:CB	2.26	0.68
1:E:3:GLU:CG	3:E:368:HOH:O	2.41	0.68
1:C:41:ASP:O	3:C:355:HOH:O	2.12	0.68
1:E:60:TYR:CA	3:E:381:HOH:O	2.42	0.67
1:E:133:TYR:CD1	3:E:371:HOH:O	2.46	0.67
1:E:158:VAL:HA	3:E:349:HOH:O	1.95	0.67
1:A:3:GLU:OE2	1:A:91:ARG:HD3	1.95	0.67
1:A:180:ILE:CG1	3:E:421:HOH:O	2.42	0.67
1:A:64:ARG:HD2	3:A:365:HOH:O	1.93	0.67
1:A:103:PHE:HB3	3:A:412:HOH:O	1.92	0.67
1:A:180:ILE:HB	3:A:425:HOH:O	1.94	0.67
1:A:181:LYS:C	3:A:393:HOH:O	2.33	0.67
1:C:30:ILE:HD12	1:C:35:VAL:HG22	1.76	0.67
1:E:202:PRO:HG3	3:E:366:HOH:O	1.95	0.67
2:B:461:GLY:N	3:B:471:HOH:O	2.19	0.67
1:E:5:ARG:CB	3:E:422:HOH:O	2.42	0.67
1:E:7:VAL:O	3:E:354:HOH:O	2.13	0.67
1:E:30:ILE:HD12	1:E:35:VAL:HG22	1.76	0.67
1:E:253:PRO:HB2	3:E:296:HOH:O	1.94	0.67
1:C:254:LYS:HE2	3:D:393:HOH:O	1.92	0.67
1:A:225:VAL:HG22	3:A:406:HOH:O	1.95	0.67
1:A:87:ILE:CG1	3:A:417:HOH:O	2.43	0.67
1:A:169:PHE:CB	3:A:425:HOH:O	2.30	0.67
1:E:139:MET:SD	3:E:427:HOH:O	2.53	0.67
1:C:254:LYS:HG2	2:D:456:GLN:HA	1.77	0.66
1:E:5:ARG:HG3	3:E:422:HOH:O	1.94	0.66
1:C:195:VAL:HG23	3:C:403:HOH:O	1.93	0.66
3:C:359:HOH:O	2:D:463:PHE:HA	1.95	0.66
1:E:117:LYS:NZ	3:E:384:HOH:O	2.27	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:172:SER:HB3	1:A:177:ASN:HB3	1.78	0.66
1:E:94:ASP:CB	3:E:439:HOH:O	2.42	0.66
1:E:168:LYS:C	3:E:367:HOH:O	2.33	0.66
1:E:207:PHE:HA	3:E:358:HOH:O	1.95	0.66
1:A:182:LEU:HG	3:A:393:HOH:O	1.96	0.66
1:C:238:GLU:CD	3:C:439:HOH:O	2.34	0.66
1:E:138:LYS:HE3	3:E:434:HOH:O	1.95	0.66
1:A:180:ILE:CD1	3:E:421:HOH:O	2.43	0.66
1:C:103:PHE:O	3:C:380:HOH:O	2.14	0.66
1:C:144:PHE:CE1	3:C:444:HOH:O	2.48	0.66
1:E:59:THR:HG22	3:E:381:HOH:O	1.95	0.66
1:C:40:MET:HE3	2:D:459:ILE:HD13	1.78	0.66
1:A:10:SER:N	3:A:362:HOH:O	2.28	0.65
1:E:201:GLU:CB	3:E:388:HOH:O	2.44	0.65
1:A:108:GLN:CB	3:A:341:HOH:O	2.44	0.65
1:C:6:LEU:HG	3:C:414:HOH:O	1.96	0.65
1:C:11:ILE:HG13	3:C:388:HOH:O	1.97	0.65
1:E:41:ASP:OD2	1:E:43:SER:HB2	1.97	0.65
1:A:226:THR:N	3:A:406:HOH:O	2.29	0.65
1:C:165:ASP:HA	3:C:402:HOH:O	1.96	0.65
1:E:5:ARG:HB3	3:E:341:HOH:O	1.96	0.65
1:E:133:TYR:CE1	3:E:371:HOH:O	2.48	0.65
1:A:24:ASN:N	3:A:343:HOH:O	2.29	0.65
1:A:214:PHE:CA	3:A:434:HOH:O	2.45	0.65
1:C:250:TYR:N	3:C:429:HOH:O	2.23	0.65
1:E:229:MET:N	3:E:409:HOH:O	2.23	0.65
1:A:252:ALA:N	3:A:346:HOH:O	2.29	0.65
1:C:102:VAL:N	3:C:368:HOH:O	2.29	0.65
1:C:121:LEU:HA	3:C:437:HOH:O	1.95	0.65
1:C:214:PHE:CD1	3:C:330:HOH:O	2.50	0.65
1:C:229:MET:HA	3:C:315:HOH:O	1.97	0.64
1:A:26:ALA:HA	3:A:394:HOH:O	1.97	0.64
1:C:153:HIS:CG	3:C:347:HOH:O	2.47	0.64
1:C:51:THR:N	3:C:431:HOH:O	2.31	0.64
1:C:5:ARG:HD3	3:C:328:HOH:O	1.96	0.64
1:E:39:SER:N	3:E:404:HOH:O	2.30	0.64
1:A:53:ARG:HB2	3:A:408:HOH:O	1.97	0.64
1:A:243:ASP:C	3:A:382:HOH:O	2.35	0.64
1:C:8:GLN:N	3:C:414:HOH:O	2.30	0.64
1:C:11:ILE:CG1	3:C:388:HOH:O	2.45	0.64
1:E:47:LEU:C	3:E:401:HOH:O	2.36	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:170:SER:CB	1:E:179:ASN:HB3	2.28	0.64
1:E:172:SER:HB3	1:E:177:ASN:HB3	1.78	0.64
1:A:19:LEU:HG	3:A:435:HOH:O	1.96	0.64
1:A:64:ARG:N	3:A:359:HOH:O	2.30	0.64
1:A:151:LEU:N	3:A:437:HOH:O	2.31	0.64
1:A:252:ALA:CB	3:A:346:HOH:O	2.33	0.64
1:E:43:SER:CB	3:E:364:HOH:O	2.44	0.64
1:C:216:THR:HB	3:C:413:HOH:O	1.96	0.63
1:E:94:ASP:CA	3:E:375:HOH:O	2.46	0.63
1:A:170:SER:CB	1:A:179:ASN:HB3	2.27	0.63
1:C:47:LEU:O	3:C:429:HOH:O	2.16	0.63
1:C:77:LYS:HE3	3:C:353:HOH:O	1.98	0.63
1:C:126:LEU:CA	3:C:412:HOH:O	2.45	0.63
1:E:140:PRO:HB3	3:E:429:HOH:O	1.98	0.63
1:A:44:HIS:ND1	3:A:345:HOH:O	2.31	0.63
1:A:124:GLU:HB3	3:A:332:HOH:O	1.97	0.63
1:C:172:SER:HB3	1:C:177:ASN:HB3	1.80	0.62
1:C:110:LYS:CE	3:C:354:HOH:O	2.47	0.62
1:A:30:ILE:HD12	1:A:35:VAL:HG22	1.80	0.62
1:C:126:LEU:C	3:C:412:HOH:O	2.37	0.62
1:C:149:ARG:NH1	3:C:371:HOH:O	2.32	0.62
1:E:87:ILE:HG12	3:E:354:HOH:O	1.99	0.62
1:E:210:ARG:NH1	3:E:382:HOH:O	2.32	0.62
1:E:175:LEU:C	3:E:379:HOH:O	2.38	0.62
1:E:168:LYS:CD	3:E:430:HOH:O	2.47	0.62
1:A:204:GLN:N	3:A:422:HOH:O	2.30	0.62
1:C:164:LYS:O	3:C:402:HOH:O	2.16	0.62
1:C:216:THR:CB	3:C:413:HOH:O	2.46	0.62
1:E:78:ILE:HG22	3:E:442:HOH:O	1.96	0.62
1:E:257:ASP:N	3:E:374:HOH:O	2.33	0.62
1:A:40:MET:HG3	3:A:383:HOH:O	1.98	0.62
2:D:461:GLY:C	3:D:254:HOH:O	2.38	0.61
1:A:87:ILE:HG12	3:A:417:HOH:O	1.99	0.61
1:E:35:VAL:C	3:E:398:HOH:O	2.38	0.61
1:C:207:PHE:HA	3:C:288:HOH:O	1.99	0.61
1:E:181:LYS:HE2	3:E:430:HOH:O	2.00	0.61
2:D:465:ARG:N	3:D:246:HOH:O	2.33	0.61
2:D:459:ILE:N	2:D:459:ILE:HD12	2.15	0.61
1:E:171:ALA:O	1:E:177:ASN:HB2	2.01	0.61
1:C:77:LYS:HD3	3:C:441:HOH:O	2.00	0.61
2:D:456:GLN:O	3:D:272:HOH:O	2.16	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:202:PRO:CD	3:E:403:HOH:O	2.49	0.61
1:C:103:PHE:CE2	3:C:432:HOH:O	2.51	0.61
1:E:229:MET:CA	3:E:409:HOH:O	2.49	0.61
1:A:234:PRO:HD3	2:B:462:PHE:CD2	2.36	0.61
1:C:180:ILE:HG13	3:C:362:HOH:O	2.01	0.61
1:A:246:HIS:NE2	3:A:407:HOH:O	2.31	0.61
1:A:168:LYS:CA	3:A:378:HOH:O	2.49	0.60
1:A:171:ALA:O	1:A:177:ASN:HB2	2.02	0.60
1:E:131:GLN:HB3	3:E:264:HOH:O	2.00	0.60
1:C:174:GLU:N	3:C:433:HOH:O	2.33	0.60
1:E:52:LEU:HB3	1:E:244:MET:HE1	1.83	0.60
1:C:57:PHE:HD2	3:C:455:HOH:O	1.84	0.60
1:E:94:ASP:CA	3:E:439:HOH:O	2.50	0.60
1:A:245:GLY:N	3:A:382:HOH:O	2.34	0.60
1:C:58:ASP:CA	3:C:356:HOH:O	2.46	0.60
1:A:53:ARG:NH2	3:A:270:HOH:O	2.29	0.60
1:A:178:GLY:HA2	3:E:340:HOH:O	2.01	0.60
1:C:214:PHE:CA	3:C:330:HOH:O	2.42	0.60
1:C:170:SER:CB	1:C:179:ASN:HB3	2.28	0.60
1:C:226:THR:HG22	3:C:369:HOH:O	2.01	0.60
1:C:42:SER:CA	3:C:355:HOH:O	2.49	0.59
1:C:195:VAL:N	3:C:403:HOH:O	2.33	0.59
1:E:202:PRO:N	3:E:403:HOH:O	2.34	0.59
1:C:161:SER:N	3:C:408:HOH:O	2.35	0.59
1:E:250:TYR:HA	3:E:420:HOH:O	2.01	0.59
1:C:254:LYS:HG3	3:C:288:HOH:O	2.01	0.59
1:E:64:ARG:HD2	1:E:94:ASP:HB3	1.84	0.59
1:C:135:CYS:SG	1:C:199:MET:HG2	2.42	0.59
1:A:252:ALA:CA	3:A:346:HOH:O	2.50	0.59
1:E:115:GLU:HG3	3:E:304:HOH:O	2.02	0.59
1:E:115:GLU:N	3:E:340:HOH:O	2.36	0.59
2:B:464:GLN:H	2:B:464:GLN:CD	2.05	0.59
1:C:224:THR:CG2	3:C:386:HOH:O	2.51	0.59
1:A:52:LEU:HB3	1:A:244:MET:HE1	1.84	0.58
1:C:87:ILE:HG23	3:C:422:HOH:O	2.02	0.58
1:A:70:VAL:HB	3:A:399:HOH:O	2.02	0.58
1:A:226:THR:C	3:A:406:HOH:O	2.40	0.58
1:E:202:PRO:HD2	3:E:403:HOH:O	2.02	0.58
1:E:225:VAL:C	3:E:385:HOH:O	2.40	0.58
1:A:255:ILE:HD13	2:B:457:VAL:HG21	1.85	0.58
1:E:70:VAL:HB	3:E:353:HOH:O	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:30:ILE:HG13	1:C:68:MET:CE	2.33	0.58
1:C:58:ASP:C	3:C:328:HOH:O	2.40	0.58
2:D:462:PHE:N	3:D:254:HOH:O	2.36	0.58
1:E:12:LEU:CA	3:E:344:HOH:O	2.52	0.58
1:E:254:LYS:HE3	3:F:132:HOH:O	2.04	0.58
1:C:3:GLU:OE2	1:C:91:ARG:HD3	2.03	0.58
1:E:4:ALA:HB1	1:E:57:PHE:CD2	2.38	0.58
1:E:78:ILE:HG23	3:E:442:HOH:O	2.03	0.58
1:C:105:ALA:N	3:C:380:HOH:O	2.37	0.58
1:C:101:LEU:C	3:C:368:HOH:O	2.41	0.58
1:C:149:ARG:HG2	1:C:149:ARG:HH11	1.68	0.57
1:A:141:SER:HB2	1:A:219:THR:HG23	1.86	0.57
1:A:175:LEU:HG	3:A:316:HOH:O	2.04	0.57
1:C:61:ARG:NH2	3:C:451:HOH:O	2.37	0.57
1:E:202:PRO:CD	3:E:388:HOH:O	2.50	0.57
1:E:176:GLY:HA3	3:E:379:HOH:O	2.04	0.57
1:A:113:ASP:N	3:C:458:HOH:O	2.37	0.57
1:A:244:MET:N	3:A:382:HOH:O	2.37	0.57
1:C:51:THR:O	1:C:245:GLY:HA3	2.03	0.57
1:C:124:GLU:HB2	3:C:435:HOH:O	2.03	0.57
1:E:28:TRP:HE3	1:E:35:VAL:HG11	1.70	0.57
1:E:86:ASP:OD1	1:E:105:ALA:HA	2.04	0.57
1:A:49:GLN:N	3:A:439:HOH:O	2.37	0.57
2:B:459:ILE:HD12	2:B:459:ILE:N	2.20	0.57
1:C:153:HIS:HB2	3:C:347:HOH:O	1.93	0.57
1:E:40:MET:HE1	1:E:44:HIS:ND1	2.19	0.57
1:E:94:ASP:HB2	3:E:439:HOH:O	2.03	0.57
1:A:40:MET:HE3	2:B:459:ILE:HD13	1.87	0.57
1:A:48:VAL:C	3:A:439:HOH:O	2.43	0.57
1:E:51:THR:O	1:E:245:GLY:HA3	2.05	0.57
1:A:226:THR:CA	3:A:406:HOH:O	2.53	0.56
1:C:144:PHE:CG	3:C:444:HOH:O	2.57	0.56
1:A:156:ASP:CB	3:A:423:HOH:O	2.46	0.56
1:A:37:LEU:O	3:A:386:HOH:O	2.18	0.56
1:A:91:ARG:NH1	3:A:432:HOH:O	2.37	0.56
1:C:144:PHE:CZ	3:C:444:HOH:O	2.58	0.56
1:E:53:ARG:NH2	3:E:278:HOH:O	2.36	0.56
2:F:464:GLN:N	3:F:303:HOH:O	2.38	0.56
2:B:460:THR:HG22	3:B:525:HOH:O	2.05	0.56
1:C:16:LEU:HD13	1:C:79:LEU:CD1	2.35	0.56
1:E:60:TYR:N	3:E:381:HOH:O	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:107:ASN:O	1:E:109:GLU:OE1	2.23	0.56
1:E:168:LYS:CA	3:E:367:HOH:O	2.53	0.56
1:C:210:ARG:HB3	3:C:367:HOH:O	2.05	0.56
1:E:43:SER:CA	3:E:364:HOH:O	2.52	0.56
1:E:107:ASN:C	3:E:342:HOH:O	2.42	0.56
1:C:129:PRO:HD3	3:C:359:HOH:O	2.05	0.56
1:A:208:ALA:HB2	3:B:50:HOH:O	2.05	0.56
1:C:64:ARG:HD2	1:C:94:ASP:HB3	1.88	0.56
1:C:171:ALA:O	1:C:177:ASN:HB2	2.05	0.56
2:F:457:VAL:N	3:F:540:HOH:O	2.37	0.56
1:C:194:ALA:CB	3:C:377:HOH:O	2.39	0.56
1:E:87:ILE:HG22	3:E:292:HOH:O	2.04	0.56
1:A:211:TYR:HB3	3:A:289:HOH:O	2.05	0.56
1:E:90:LEU:HD22	3:E:405:HOH:O	2.07	0.55
1:C:52:LEU:HB3	1:C:244:MET:HE1	1.88	0.55
1:A:16:LEU:HD21	1:A:75:MET:CG	2.37	0.55
1:A:49:GLN:CA	3:A:439:HOH:O	2.54	0.55
1:A:240:LYS:N	3:A:427:HOH:O	2.38	0.55
1:C:8:GLN:CA	3:C:414:HOH:O	2.55	0.55
1:C:19:LEU:CD2	1:C:48:VAL:HG11	2.37	0.55
1:E:100:ALA:HA	3:E:392:HOH:O	2.06	0.55
1:C:141:SER:HB2	1:C:219:THR:HG23	1.89	0.55
1:E:3:GLU:CD	3:E:368:HOH:O	2.45	0.55
1:E:132:GLU:O	3:E:448:HOH:O	2.18	0.55
1:A:86:ASP:OD1	1:A:105:ALA:HA	2.07	0.55
1:A:112:SER:CA	3:C:458:HOH:O	2.54	0.55
2:B:464:GLN:CD	2:B:464:GLN:N	2.61	0.55
1:C:53:ARG:NH2	3:C:266:HOH:O	2.37	0.55
1:C:93:GLU:HB2	1:C:96:ALA:HB3	1.89	0.55
1:C:167:VAL:HG23	3:C:408:HOH:O	2.07	0.55
1:A:93:GLU:HB2	1:A:96:ALA:HB3	1.89	0.55
1:A:135:CYS:SG	1:A:199:MET:HG2	2.46	0.55
1:A:151:LEU:HD21	3:E:421:HOH:O	2.06	0.55
1:A:254:LYS:HG2	2:B:456:GLN:HA	1.89	0.55
1:C:174:GLU:CG	3:C:433:HOH:O	2.55	0.55
1:A:135:CYS:HG	1:A:162:CYS:HG	1.55	0.54
1:E:139:MET:HG3	3:E:427:HOH:O	2.05	0.54
1:C:194:ALA:CA	3:C:377:HOH:O	2.54	0.54
1:E:19:LEU:CD2	1:E:48:VAL:HG11	2.38	0.54
1:E:135:CYS:HB3	3:E:347:HOH:O	2.05	0.54
1:E:99:LEU:O	3:E:392:HOH:O	2.18	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:THR:CA	3:A:354:HOH:O	2.44	0.54
1:C:30:ILE:CD1	1:C:35:VAL:HG22	2.37	0.54
1:C:132:GLU:N	3:C:426:HOH:O	2.40	0.54
1:A:51:THR:O	1:A:245:GLY:HA3	2.07	0.54
1:A:26:ALA:CA	3:A:394:HOH:O	2.54	0.54
1:A:49:GLN:CB	3:A:439:HOH:O	2.30	0.54
2:B:461:GLY:CA	3:B:471:HOH:O	2.56	0.54
1:E:6:LEU:HD12	3:E:378:HOH:O	2.07	0.54
2:D:463:PHE:N	3:D:254:HOH:O	2.40	0.54
1:E:139:MET:CG	3:E:427:HOH:O	2.56	0.54
1:A:28:TRP:HE3	1:A:35:VAL:HG11	1.73	0.54
1:E:229:MET:C	3:E:409:HOH:O	2.46	0.54
1:A:40:MET:HE1	1:A:44:HIS:ND1	2.23	0.54
1:C:135:CYS:HG	1:C:162:CYS:HG	1.56	0.53
1:E:229:MET:O	3:E:409:HOH:O	2.18	0.53
1:E:236:VAL:HG22	3:E:420:HOH:O	2.07	0.53
1:C:65:ASN:N	3:C:351:HOH:O	2.38	0.53
1:C:224:THR:HG21	3:C:386:HOH:O	2.08	0.53
1:E:40:MET:HE3	2:F:459:ILE:HD13	1.90	0.53
1:A:1:MET:H3	1:A:63:ASP:HB2	1.72	0.53
1:A:25:GLU:O	3:A:394:HOH:O	2.19	0.53
1:A:203:VAL:CG1	3:A:422:HOH:O	2.52	0.53
1:A:239:TYR:CB	3:A:427:HOH:O	2.37	0.53
1:C:82:ALA:HA	3:C:430:HOH:O	2.07	0.53
1:E:37:LEU:CD2	3:E:404:HOH:O	2.54	0.53
2:B:454:ASN:N	3:B:271:HOH:O	2.41	0.53
1:C:77:LYS:CD	3:C:441:HOH:O	2.55	0.53
1:E:147:ILE:HG23	1:E:180:ILE:HD12	1.91	0.53
1:A:30:ILE:HG13	1:A:68:MET:CE	2.39	0.53
2:B:461:GLY:HA3	3:B:471:HOH:O	2.09	0.53
1:E:199:MET:HG3	3:E:396:HOH:O	2.06	0.53
2:F:459:ILE:N	2:F:459:ILE:HD12	2.23	0.53
1:A:45:VAL:HG13	3:A:346:HOH:O	2.08	0.53
1:C:98:THR:HG22	3:C:276:HOH:O	2.09	0.53
1:E:1:MET:H3	1:E:63:ASP:HB2	1.73	0.53
1:E:143:GLU:HB3	3:E:427:HOH:O	2.08	0.53
1:A:94:ASP:CA	3:A:365:HOH:O	2.56	0.53
1:A:151:LEU:CA	3:A:437:HOH:O	2.57	0.53
1:C:40:MET:HE3	2:D:459:ILE:CD1	2.39	0.53
1:C:40:MET:HE1	1:C:44:HIS:ND1	2.24	0.53
1:C:179:ASN:HA	3:C:362:HOH:O	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:454:ASN:CA	3:D:393:HOH:O	2.57	0.53
1:E:108:GLN:N	3:E:342:HOH:O	2.42	0.53
1:A:246:HIS:CD2	3:A:407:HOH:O	2.61	0.53
1:C:8:GLN:C	3:C:414:HOH:O	2.46	0.53
1:C:205:LEU:HD23	3:C:268:HOH:O	2.09	0.52
1:E:30:ILE:HG13	1:E:68:MET:CE	2.39	0.52
1:C:204:GLN:HG2	3:C:285:HOH:O	2.09	0.52
1:C:140:PRO:HD3	3:C:457:HOH:O	2.10	0.52
1:C:174:GLU:HG2	3:C:433:HOH:O	2.10	0.52
1:E:16:LEU:HD13	1:E:79:LEU:CD1	2.40	0.52
1:A:16:LEU:HD21	1:A:75:MET:HG2	1.90	0.52
1:A:170:SER:HB3	1:A:179:ASN:CB	2.34	0.52
1:C:254:LYS:HG3	3:C:372:HOH:O	2.08	0.52
1:C:50:LEU:C	3:C:431:HOH:O	2.48	0.52
1:E:135:CYS:HA	1:E:198:GLU:O	2.09	0.52
1:E:194:ALA:HA	3:E:318:HOH:O	2.09	0.52
1:C:217:LYS:HG3	3:C:387:HOH:O	2.09	0.52
1:E:168:LYS:NZ	3:E:430:HOH:O	2.43	0.52
1:C:248:LYS:HE3	3:C:341:HOH:O	2.08	0.52
1:E:94:ASP:CB	3:E:375:HOH:O	2.54	0.52
1:A:196:THR:C	3:A:397:HOH:O	2.48	0.52
1:E:256:GLU:CA	3:E:374:HOH:O	2.51	0.52
1:E:104:GLU:HG3	3:E:346:HOH:O	2.09	0.52
1:A:4:ALA:HB1	1:A:57:PHE:CD2	2.44	0.51
1:C:147:ILE:HG23	1:C:180:ILE:HD12	1.92	0.51
1:C:16:LEU:HD21	1:C:75:MET:CG	2.40	0.51
1:C:21:ASP:HA	3:C:322:HOH:O	2.10	0.51
1:C:126:LEU:N	3:C:412:HOH:O	2.44	0.51
1:E:45:VAL:HG13	2:F:456:GLN:NE2	2.25	0.51
1:E:107:ASN:CB	3:E:342:HOH:O	2.50	0.51
1:A:16:LEU:HD13	1:A:79:LEU:CD1	2.41	0.51
1:A:92:ALA:HA	3:A:323:HOH:O	2.10	0.51
1:A:254:LYS:NZ	3:A:315:HOH:O	2.41	0.51
1:C:28:TRP:HE3	1:C:35:VAL:HG11	1.75	0.51
1:E:85:GLU:CG	3:E:360:HOH:O	2.38	0.51
1:C:1:MET:H3	1:C:63:ASP:HB2	1.75	0.51
1:C:214:PHE:N	3:C:330:HOH:O	2.43	0.51
3:C:350:HOH:O	2:D:462:PHE:CE2	2.64	0.51
1:E:16:LEU:HD21	1:E:75:MET:CG	2.41	0.51
1:E:59:THR:O	3:E:381:HOH:O	2.18	0.51
1:E:107:ASN:CA	3:E:342:HOH:O	2.58	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:374:HOH:O	2:F:454:ASN:CB	2.59	0.51
2:D:462:PHE:C	3:D:254:HOH:O	2.49	0.51
1:A:19:LEU:CD2	1:A:48:VAL:HG11	2.41	0.51
1:A:200:ASN:CG	3:A:436:HOH:O	2.49	0.51
1:E:110:LYS:HB3	3:E:348:HOH:O	2.10	0.51
1:C:255:ILE:HD11	3:C:350:HOH:O	2.11	0.51
1:E:93:GLU:HB2	1:E:96:ALA:HB3	1.92	0.51
1:C:110:LYS:HE3	3:C:354:HOH:O	2.10	0.50
1:C:172:SER:HB3	1:C:177:ASN:CB	2.41	0.50
1:A:95:ASN:OD1	3:A:387:HOH:O	2.19	0.50
1:A:112:SER:CB	3:C:458:HOH:O	2.56	0.50
1:C:146:ARG:NE	3:C:400:HOH:O	2.44	0.50
1:C:174:GLU:CA	3:C:433:HOH:O	2.58	0.50
1:E:24:ASN:OD1	3:E:397:HOH:O	2.19	0.50
1:E:136:VAL:N	3:E:347:HOH:O	2.44	0.50
1:A:107:ASN:CB	3:A:372:HOH:O	2.59	0.50
1:A:151:LEU:CD2	3:E:421:HOH:O	2.59	0.50
1:C:135:CYS:HA	1:C:198:GLU:O	2.10	0.50
1:C:194:ALA:O	3:C:457:HOH:O	2.20	0.50
1:E:47:LEU:CA	3:E:401:HOH:O	2.58	0.50
1:A:254:LYS:HG3	3:A:428:HOH:O	2.11	0.50
1:A:72:LEU:HD21	3:A:399:HOH:O	2.10	0.50
1:C:226:THR:CG2	3:C:369:HOH:O	2.58	0.50
1:C:180:ILE:N	3:C:362:HOH:O	2.43	0.50
1:A:40:MET:HE2	1:A:44:HIS:CG	2.47	0.50
1:C:86:ASP:OD1	1:C:105:ALA:HA	2.11	0.50
1:C:161:SER:C	3:C:408:HOH:O	2.49	0.50
1:E:168:LYS:N	3:E:367:HOH:O	2.45	0.50
1:A:149:ARG:HG2	1:A:149:ARG:HH11	1.76	0.50
1:E:170:SER:HB3	1:E:179:ASN:CB	2.33	0.50
1:A:45:VAL:CG1	3:A:346:HOH:O	2.59	0.49
1:C:8:GLN:HB2	3:C:414:HOH:O	2.11	0.49
1:C:45:VAL:HG13	2:D:456:GLN:NE2	2.26	0.49
1:C:119:MET:CE	3:C:337:HOH:O	2.56	0.49
1:A:49:GLN:HA	3:A:386:HOH:O	2.12	0.49
1:E:143:GLU:O	1:E:147:ILE:HG13	2.12	0.49
2:F:462:PHE:C	2:F:463:PHE:HD2	2.16	0.49
1:C:70:VAL:CG1	3:C:405:HOH:O	2.41	0.49
1:E:52:LEU:HB3	1:E:244:MET:CE	2.42	0.49
1:E:228:SER:HB2	1:E:236:VAL:HB	1.94	0.49
1:A:28:TRP:CE3	1:A:35:VAL:HG11	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:207:PHE:CA	3:A:428:HOH:O	2.54	0.49
1:E:141:SER:HB2	1:E:219:THR:HG23	1.93	0.49
1:E:225:VAL:CG2	3:E:385:HOH:O	2.43	0.49
1:C:143:GLU:O	1:C:147:ILE:HG13	2.12	0.49
1:A:40:MET:CG	3:A:383:HOH:O	2.59	0.49
1:C:139:MET:HA	3:C:457:HOH:O	2.11	0.49
1:E:131:GLN:N	3:E:399:HOH:O	2.44	0.49
1:C:5:ARG:HB3	3:C:328:HOH:O	2.11	0.49
1:C:40:MET:HE2	1:C:44:HIS:CG	2.48	0.49
1:E:12:LEU:HA	3:E:344:HOH:O	2.13	0.49
1:E:94:ASP:HA	3:E:375:HOH:O	2.10	0.49
1:A:77:LYS:NZ	3:A:350:HOH:O	2.45	0.48
3:A:366:HOH:O	2:B:463:PHE:HD1	1.96	0.48
1:E:166:GLY:HA2	1:E:197:ILE:CD1	2.43	0.48
1:A:40:MET:HE3	2:B:459:ILE:CD1	2.43	0.48
1:A:103:PHE:CG	3:A:412:HOH:O	2.61	0.48
1:C:107:ASN:O	1:C:109:GLU:OE1	2.32	0.48
1:E:172:SER:HB3	1:E:177:ASN:CB	2.43	0.48
1:C:38:GLN:HA	1:C:48:VAL:O	2.14	0.48
1:E:30:ILE:CD1	1:E:35:VAL:HG22	2.43	0.48
1:C:255:ILE:HG23	3:C:286:HOH:O	2.13	0.48
1:E:149:ARG:HG2	1:E:149:ARG:HH11	1.79	0.48
1:A:172:SER:HB3	1:A:177:ASN:CB	2.42	0.48
1:E:28:TRP:CE3	1:E:35:VAL:HG11	2.46	0.48
1:E:225:VAL:CG1	3:E:385:HOH:O	2.52	0.48
1:A:87:ILE:HG13	3:A:417:HOH:O	2.13	0.48
1:A:30:ILE:CD1	1:A:35:VAL:HG22	2.43	0.48
1:C:153:HIS:CD2	3:C:298:HOH:O	2.66	0.48
1:C:7:VAL:HG13	3:C:420:HOH:O	2.14	0.48
1:C:21:ASP:HB2	3:C:277:HOH:O	2.13	0.48
1:C:166:GLY:HA2	1:C:197:ILE:HD12	1.96	0.48
2:B:457:VAL:HG11	2:B:462:PHE:HE1	1.79	0.47
1:C:28:TRP:CE3	1:C:35:VAL:HG11	2.49	0.47
1:C:144:PHE:CD2	3:C:444:HOH:O	2.66	0.47
1:A:68:MET:O	1:A:70:VAL:HG23	2.14	0.47
1:C:124:GLU:N	3:C:435:HOH:O	2.33	0.47
1:E:168:LYS:HD3	3:E:430:HOH:O	2.11	0.47
1:E:203:VAL:HG12	1:E:204:GLN:N	2.28	0.47
1:A:23:ILE:HG21	3:A:394:HOH:O	2.15	0.47
1:A:225:VAL:C	3:A:358:HOH:O	2.51	0.47
1:C:16:LEU:HD21	1:C:75:MET:HG2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:63:ASP:O	3:E:444:HOH:O	2.20	0.47
1:E:254:LYS:HG2	2:F:456:GLN:CA	2.44	0.47
1:A:135:CYS:HA	1:A:198:GLU:O	2.14	0.47
1:E:207:PHE:CZ	1:E:235:LEU:HB2	2.49	0.47
1:C:64:ARG:HB3	1:C:94:ASP:OD1	2.15	0.47
1:C:241:ILE:O	1:C:244:MET:HB2	2.15	0.47
1:A:73:THR:CA	3:A:402:HOH:O	2.34	0.47
1:C:172:SER:CB	1:C:177:ASN:HB3	2.44	0.47
1:E:16:LEU:HD21	1:E:75:MET:HG2	1.97	0.47
1:E:124:GLU:OE2	2:F:465:ARG:NH1	2.48	0.47
1:E:196:THR:HA	3:E:355:HOH:O	2.14	0.47
1:E:234:PRO:HD3	2:F:462:PHE:CD2	2.50	0.47
1:A:129:PRO:HD3	2:B:463:PHE:CG	2.49	0.47
1:C:4:ALA:HB1	1:C:57:PHE:CD2	2.50	0.47
1:C:24:ASN:CA	3:C:374:HOH:O	2.61	0.47
1:C:24:ASN:C	3:C:374:HOH:O	2.53	0.47
1:E:240:LYS:NZ	3:E:278:HOH:O	2.45	0.47
1:A:95:ASN:HB2	3:A:297:HOH:O	2.13	0.47
1:E:236:VAL:CA	3:E:420:HOH:O	2.20	0.47
1:E:175:LEU:HD12	1:E:175:LEU:O	2.15	0.47
1:C:44:HIS:C	2:D:459:ILE:HD11	2.35	0.46
1:C:110:LYS:HE3	1:E:180:ILE:HG21	1.96	0.46
1:C:144:PHE:CE2	3:C:444:HOH:O	2.67	0.46
2:F:457:VAL:C	3:F:540:HOH:O	2.54	0.46
1:A:64:ARG:C	3:A:359:HOH:O	2.54	0.46
1:A:152:SER:N	3:A:437:HOH:O	2.44	0.46
1:A:166:GLY:HA2	1:A:197:ILE:CD1	2.45	0.46
1:E:166:GLY:HA2	1:E:197:ILE:HD12	1.97	0.46
1:A:26:ALA:HB2	3:A:394:HOH:O	2.15	0.46
1:C:24:ASN:CB	3:C:374:HOH:O	2.36	0.46
1:C:27:CYS:SG	3:C:449:HOH:O	2.61	0.46
1:E:5:ARG:HD3	3:E:341:HOH:O	2.15	0.46
1:E:110:LYS:HG2	3:E:314:HOH:O	2.15	0.46
1:A:211:TYR:HD2	3:A:289:HOH:O	1.97	0.46
1:C:41:ASP:OD2	1:C:43:SER:HB2	2.15	0.46
1:C:153:HIS:HB3	3:C:347:HOH:O	2.03	0.46
1:C:104:GLU:CA	3:C:380:HOH:O	2.64	0.46
1:E:248:LYS:NZ	3:E:408:HOH:O	2.46	0.46
1:C:200:ASN:HB2	3:C:274:HOH:O	2.16	0.46
1:E:74:SER:N	3:E:395:HOH:O	2.48	0.46
1:C:207:PHE:CZ	1:C:235:LEU:HB2	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:254:LYS:NZ	3:C:335:HOH:O	2.44	0.46
1:C:170:SER:HB3	1:C:179:ASN:CB	2.34	0.46
1:A:110:LYS:HE3	1:C:180:ILE:HG21	1.97	0.45
1:A:166:GLY:HA2	1:A:197:ILE:HD12	1.98	0.45
1:E:40:MET:HE3	2:F:459:ILE:CD1	2.46	0.45
2:F:465:ARG:HG2	2:F:465:ARG:NH1	2.28	0.45
1:E:40:MET:HE2	1:E:44:HIS:CG	2.51	0.45
1:A:19:LEU:N	3:A:435:HOH:O	2.45	0.45
1:A:82:ALA:HA	3:A:412:HOH:O	2.17	0.45
1:A:206:THR:OG1	2:B:453:ALA:HB2	2.16	0.45
1:C:128:ILE:HA	1:C:129:PRO:HD3	1.85	0.45
1:E:38:GLN:HA	1:E:48:VAL:O	2.17	0.45
1:E:149:ARG:O	1:E:152:SER:HB2	2.16	0.45
1:E:170:SER:CB	3:E:387:HOH:O	2.64	0.45
1:A:52:LEU:HB3	1:A:244:MET:CE	2.46	0.45
1:A:127:GLY:N	3:A:366:HOH:O	2.49	0.45
1:C:125:GLN:C	3:C:412:HOH:O	2.55	0.45
1:C:166:GLY:HA2	1:C:197:ILE:CD1	2.46	0.45
1:E:241:ILE:HG22	1:E:244:MET:HB2	1.99	0.45
1:C:149:ARG:NH1	3:C:456:HOH:O	2.48	0.45
1:E:246:HIS:HD2	1:E:248:LYS:HG3	1.81	0.45
1:A:64:ARG:HD2	1:A:94:ASP:HB3	1.99	0.45
1:C:135:CYS:HB3	1:C:162:CYS:HG	1.81	0.45
1:A:45:VAL:HG13	2:B:456:GLN:NE2	2.32	0.45
1:C:61:ARG:NH1	3:C:396:HOH:O	2.49	0.45
1:C:244:MET:HE2	1:C:244:MET:HB3	1.52	0.45
1:E:138:LYS:CA	3:E:426:HOH:O	2.63	0.45
1:C:67:ALA:HA	3:C:449:HOH:O	2.16	0.45
1:C:217:LYS:CG	3:C:387:HOH:O	2.64	0.45
1:C:255:ILE:HD13	2:D:457:VAL:CG2	2.45	0.45
1:E:71:ASN:ND2	3:E:395:HOH:O	2.50	0.45
1:E:73:THR:CA	3:E:406:HOH:O	2.39	0.45
1:C:91:ARG:CD	3:C:397:HOH:O	2.52	0.44
1:C:213:ASN:C	3:C:330:HOH:O	2.55	0.44
1:C:235:LEU:HA	3:C:315:HOH:O	2.17	0.44
1:E:26:ALA:HB3	3:E:353:HOH:O	2.16	0.44
1:C:52:LEU:HD22	1:C:244:MET:HE1	1.99	0.44
1:C:217:LYS:CD	3:C:387:HOH:O	2.65	0.44
1:E:12:LEU:CB	3:E:344:HOH:O	2.66	0.44
1:E:176:GLY:CA	3:E:379:HOH:O	2.64	0.44
1:A:41:ASP:OD2	1:A:43:SER:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:HIS:HD2	1:A:248:LYS:HG3	1.82	0.44
1:A:205:LEU:HD11	1:A:230:SER:O	2.18	0.44
1:C:158:VAL:HB	1:C:209:LEU:HD21	1.99	0.44
1:E:105:ALA:O	3:E:383:HOH:O	2.21	0.44
1:C:64:ARG:CG	3:C:351:HOH:O	2.54	0.44
1:E:255:ILE:HD13	2:F:457:VAL:CG2	2.44	0.44
1:A:16:LEU:CD2	1:A:75:MET:HG2	2.48	0.44
1:A:226:THR:HG21	3:A:410:HOH:O	2.17	0.44
1:E:7:VAL:HG13	3:E:354:HOH:O	2.16	0.44
1:A:140:PRO:HG3	1:A:192:GLU:O	2.17	0.44
1:C:254:LYS:HB3	3:D:393:HOH:O	2.17	0.44
1:E:246:HIS:CD2	1:E:248:LYS:HG3	2.53	0.44
1:A:46:SER:HB2	1:A:250:TYR:O	2.18	0.44
1:C:45:VAL:HG12	1:C:251:LEU:HD12	2.00	0.44
1:C:135:CYS:CB	1:C:162:CYS:HG	2.30	0.44
1:A:26:ALA:CB	3:A:394:HOH:O	2.66	0.44
1:A:228:SER:HB2	1:A:236:VAL:HB	1.98	0.44
1:C:84:ASN:ND2	3:C:269:HOH:O	2.50	0.44
1:C:140:PRO:CD	3:C:457:HOH:O	2.65	0.44
1:A:21:ASP:HB2	3:A:388:HOH:O	2.17	0.43
1:A:72:LEU:CD2	3:A:399:HOH:O	2.65	0.43
1:A:158:VAL:HB	1:A:209:LEU:HD21	1.99	0.43
1:C:25:GLU:HG3	3:C:361:HOH:O	2.17	0.43
1:E:176:GLY:N	3:E:379:HOH:O	2.49	0.43
1:A:65:ASN:HB3	3:A:426:HOH:O	2.18	0.43
1:C:63:ASP:CG	3:C:396:HOH:O	2.56	0.43
1:C:115:GLU:HG2	3:C:376:HOH:O	2.18	0.43
1:A:16:LEU:HD21	1:A:75:MET:SD	2.59	0.43
1:A:38:GLN:HA	1:A:48:VAL:O	2.18	0.43
1:A:197:ILE:CG2	3:A:396:HOH:O	2.47	0.43
1:C:128:ILE:HG22	2:D:463:PHE:CE2	2.53	0.43
2:D:453:ALA:C	3:D:393:HOH:O	2.53	0.43
1:A:48:VAL:CG1	3:A:404:HOH:O	2.50	0.43
1:A:146:ARG:HD3	1:A:149:ARG:NH2	2.33	0.43
1:A:172:SER:CB	1:A:177:ASN:HB3	2.48	0.43
1:A:246:HIS:CD2	1:A:248:LYS:HG3	2.53	0.43
1:C:16:LEU:HD21	1:C:75:MET:SD	2.59	0.43
1:C:52:LEU:HB3	1:C:244:MET:CE	2.48	0.43
1:E:99:LEU:HD12	1:E:100:ALA:H	1.83	0.43
1:E:108:GLN:O	1:E:108:GLN:HG3	2.18	0.43
1:E:203:VAL:CG1	1:E:204:GLN:N	2.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:LEU:CD2	1:A:79:LEU:HD12	2.28	0.43
1:A:23:ILE:CA	3:A:343:HOH:O	2.28	0.43
1:A:225:VAL:HG13	3:A:406:HOH:O	2.19	0.43
1:C:60:TYR:O	1:C:61:ARG:HB2	2.19	0.43
1:E:105:ALA:O	1:E:106:PRO:C	2.57	0.43
1:E:244:MET:HB3	1:E:244:MET:HE2	1.48	0.43
1:A:110:LYS:HA	1:C:181:LYS:O	2.18	0.43
1:A:115:GLU:HG3	3:C:390:HOH:O	2.19	0.43
1:C:11:ILE:HG12	3:C:388:HOH:O	2.12	0.43
1:C:131:GLN:CB	3:C:426:HOH:O	2.35	0.43
1:E:229:MET:HB2	3:E:409:HOH:O	2.18	0.43
1:A:207:PHE:CZ	1:A:235:LEU:HB2	2.54	0.43
1:A:219:THR:N	1:A:220:PRO:HD2	2.33	0.43
1:C:47:LEU:N	3:C:429:HOH:O	2.42	0.43
1:C:160:ILE:CG2	3:C:408:HOH:O	2.51	0.43
1:E:1:MET:N	1:E:94:ASP:OD2	2.50	0.43
1:A:181:LYS:CA	3:A:393:HOH:O	2.67	0.42
1:C:210:ARG:CB	3:C:367:HOH:O	2.65	0.42
1:A:112:SER:HA	3:C:458:HOH:O	2.19	0.42
1:C:219:THR:N	1:C:220:PRO:HD2	2.34	0.42
1:C:253:PRO:O	3:C:372:HOH:O	2.20	0.42
1:E:64:ARG:HD2	3:E:375:HOH:O	2.18	0.42
1:E:135:CYS:SG	1:E:199:MET:HG2	2.59	0.42
1:A:5:ARG:CB	1:A:59:THR:HB	2.42	0.42
1:A:107:ASN:O	1:A:108:GLN:CB	2.66	0.42
1:A:137:VAL:O	1:A:226:THR:HA	2.18	0.42
1:A:241:ILE:O	1:A:244:MET:HB2	2.19	0.42
1:C:129:PRO:HD2	3:C:309:HOH:O	2.18	0.42
1:C:223:SER:HB3	3:C:338:HOH:O	2.19	0.42
1:C:228:SER:HB2	1:C:236:VAL:HB	2.00	0.42
1:E:64:ARG:HB3	1:E:94:ASP:OD1	2.19	0.42
1:A:208:ALA:N	3:A:428:HOH:O	2.53	0.42
1:C:246:HIS:CD2	1:C:248:LYS:HG3	2.55	0.42
1:E:5:ARG:CB	1:E:59:THR:HB	2.42	0.42
1:A:125:GLN:NE2	3:A:331:HOH:O	2.53	0.42
1:A:199:MET:CE	1:A:202:PRO:HG3	2.50	0.42
1:C:24:ASN:HB3	3:C:361:HOH:O	2.19	0.42
1:C:203:VAL:HG12	1:C:204:GLN:N	2.33	0.42
1:E:116:MET:HG3	3:E:290:HOH:O	2.18	0.42
1:C:103:PHE:CB	3:C:430:HOH:O	2.67	0.42
1:C:110:LYS:HA	1:E:181:LYS:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:217:LYS:HD2	3:C:387:HOH:O	2.20	0.42
1:E:172:SER:CB	1:E:177:ASN:HB3	2.47	0.42
1:A:12:LEU:HD12	1:A:12:LEU:HA	1.83	0.42
1:A:25:GLU:HG3	3:A:352:HOH:O	2.19	0.42
1:E:201:GLU:HG2	3:E:388:HOH:O	2.18	0.42
1:A:64:ARG:HB3	1:A:94:ASP:OD1	2.19	0.42
1:C:160:ILE:C	3:C:408:HOH:O	2.58	0.42
1:A:1:MET:N	1:A:94:ASP:OD2	2.49	0.42
1:A:96:ALA:HB1	1:A:98:THR:HG23	2.02	0.42
1:A:203:VAL:CG1	1:A:204:GLN:N	2.83	0.42
1:A:248:LYS:HB3	3:A:409:HOH:O	2.20	0.42
1:C:149:ARG:HD3	3:C:427:HOH:O	2.19	0.42
2:D:454:ASN:C	2:D:454:ASN:HD22	2.22	0.42
1:E:106:PRO:HD2	3:E:350:HOH:O	2.19	0.42
1:A:17:GLU:HA	1:A:17:GLU:OE1	2.19	0.41
1:E:106:PRO:CB	3:E:350:HOH:O	2.57	0.41
1:E:112:SER:HB3	1:E:114:TYR:HE1	1.86	0.41
1:E:168:LYS:HD2	3:E:430:HOH:O	2.17	0.41
1:A:11:ILE:O	1:A:15:VAL:HG23	2.21	0.41
1:A:87:ILE:HG22	3:A:330:HOH:O	2.19	0.41
1:C:194:ALA:C	3:C:377:HOH:O	2.59	0.41
1:E:169:PHE:CE1	3:E:367:HOH:O	2.73	0.41
2:F:458:SER:HB3	3:F:517:HOH:O	2.20	0.41
1:A:40:MET:N	3:A:383:HOH:O	2.38	0.41
1:A:219:THR:C	3:A:354:HOH:O	2.59	0.41
1:C:46:SER:HB2	3:C:429:HOH:O	2.20	0.41
1:E:12:LEU:HD13	3:E:344:HOH:O	2.19	0.41
1:E:114:TYR:C	3:E:340:HOH:O	2.54	0.41
1:A:164:LYS:HG2	3:A:328:HOH:O	2.21	0.41
1:A:254:LYS:CB	3:B:271:HOH:O	2.61	0.41
1:C:103:PHE:CD2	3:C:430:HOH:O	2.57	0.41
1:E:68:MET:O	1:E:70:VAL:HG23	2.21	0.41
1:E:128:ILE:HA	1:E:129:PRO:HD3	1.83	0.41
1:E:144:PHE:N	3:E:427:HOH:O	2.53	0.41
1:A:203:VAL:HG12	1:A:204:GLN:N	2.36	0.41
1:C:16:LEU:CD2	1:C:75:MET:HG2	2.50	0.41
1:C:17:GLU:OE1	1:C:17:GLU:HA	2.20	0.41
1:C:131:GLN:CG	3:C:426:HOH:O	2.68	0.41
1:A:64:ARG:NH1	3:A:365:HOH:O	2.53	0.41
2:D:459:ILE:HD12	2:D:459:ILE:H	1.85	0.41
1:E:1:MET:HB3	1:E:63:ASP:OD2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:ILE:CG2	3:A:394:HOH:O	2.69	0.41
1:A:44:HIS:C	2:B:459:ILE:HD11	2.40	0.41
1:A:62:CYS:C	3:A:359:HOH:O	2.58	0.41
1:A:244:MET:HB3	1:A:244:MET:HE2	1.53	0.41
1:C:3:GLU:HB2	3:C:399:HOH:O	2.21	0.41
1:C:52:LEU:HD22	1:C:244:MET:CE	2.51	0.41
1:C:131:GLN:CB	3:C:373:HOH:O	2.43	0.41
1:C:248:LYS:CE	3:C:341:HOH:O	2.68	0.41
1:E:115:GLU:HG2	3:E:340:HOH:O	2.21	0.41
1:E:169:PHE:CD1	3:E:367:HOH:O	2.57	0.41
1:C:203:VAL:HA	3:C:384:HOH:O	2.21	0.41
1:E:218:ALA:O	1:E:221:LEU:HB2	2.21	0.41
1:C:45:VAL:C	2:D:459:ILE:HG13	2.41	0.40
1:C:205:LEU:CD2	3:C:391:HOH:O	2.68	0.40
1:C:256:GLU:HA	3:C:385:HOH:O	2.20	0.40
1:E:196:THR:N	3:E:434:HOH:O	2.54	0.40
1:E:226:THR:N	3:E:385:HOH:O	2.54	0.40
2:F:463:PHE:C	3:F:303:HOH:O	2.60	0.40
1:A:129:PRO:CG	2:B:463:PHE:HB3	2.51	0.40
1:E:12:LEU:HB2	3:E:344:HOH:O	2.21	0.40
1:E:219:THR:N	1:E:220:PRO:HD2	2.35	0.40
1:A:126:LEU:CD2	2:B:464:GLN:HA	2.52	0.40
1:A:185:THR:OG1	1:A:194:ALA:HB1	2.21	0.40
1:A:255:ILE:CG2	1:A:256:GLU:N	2.84	0.40
1:E:112:SER:HB3	1:E:114:TYR:CE1	2.56	0.40
1:E:149:ARG:C	3:E:433:HOH:O	2.56	0.40
1:E:202:PRO:CG	3:E:366:HOH:O	2.63	0.40
1:C:49:GLN:O	3:C:431:HOH:O	2.22	0.40
1:E:23:ILE:C	3:E:397:HOH:O	2.60	0.40
1:A:164:LYS:CG	3:A:328:HOH:O	2.69	0.40
1:A:207:PHE:C	3:A:428:HOH:O	2.59	0.40
1:C:254:LYS:N	3:C:372:HOH:O	2.53	0.40
1:E:71:ASN:C	1:E:71:ASN:HD22	2.23	0.40
2:F:465:ARG:NH1	2:F:465:ARG:CG	2.85	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:8:GLN:NE2	1:E:8:GLN:NE2[5_674]	1.61	0.59
3:E:315:HOH:O	3:E:360:HOH:O[5_674]	2.17	0.03

## 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	249/261 (95%)	226 (91%)	21 (8%)	2 (1%)	19	39
1	C	245/261 (94%)	226 (92%)	15 (6%)	4 (2%)	9	19
1	E	249/261 (95%)	228 (92%)	17 (7%)	4 (2%)	9	19
2	B	10/15 (67%)	9 (90%)	0	1 (10%)	0	0
2	D	11/15 (73%)	10 (91%)	1 (9%)	0	100	100
2	F	11/15 (73%)	10 (91%)	0	1 (9%)	1	0
All	All	775/828 (94%)	709 (92%)	54 (7%)	12 (2%)	10	21

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	242	ALA
1	C	242	ALA
1	E	242	ALA
2	F	454	ASN
1	A	59	THR
1	C	59	THR
1	C	61	ARG
2	B	454	ASN
1	E	106	PRO
1	C	243	ASP
1	E	59	THR
1	E	243	ASP

### 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	217/228 (95%)	211 (97%)	6 (3%)	43	69
1	C	216/228 (95%)	211 (98%)	5 (2%)	50	75
1	E	217/228 (95%)	212 (98%)	5 (2%)	50	75
2	B	10/13 (77%)	8 (80%)	2 (20%)	1	2
2	D	11/13 (85%)	9 (82%)	2 (18%)	1	2
2	F	11/13 (85%)	9 (82%)	2 (18%)	1	2
All	All	682/723 (94%)	660 (97%)	22 (3%)	39	65

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	ASN
1	A	85	GLU
1	A	98	THR
1	A	192	GLU
1	A	224	THR
1	A	235	LEU
2	B	454	ASN
2	B	459	ILE
1	C	71	ASN
1	C	85	GLU
1	C	98	THR
1	C	224	THR
1	C	235	LEU
2	D	454	ASN
2	D	459	ILE
1	E	71	ASN
1	E	85	GLU
1	E	98	THR
1	E	107	ASN
1	E	224	THR
2	F	459	ILE
2	F	465	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	GLN

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Mol	Chain	Res	Type
1	A	65	ASN
1	A	71	ASN
2	B	454	ASN
2	B	464	GLN
1	C	38	GLN
1	C	65	ASN
1	C	71	ASN
1	C	184	GLN
2	D	454	ASN
1	E	38	GLN
1	E	65	ASN
1	E	71	ASN
1	E	107	ASN
1	E	184	GLN
2	F	464	GLN

### 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 monosaccharides in this entry.

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

There are no ligands in this entry.

### 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

### 6.1 Protein, DNA and RNA chains

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	253/261 (96%)	0.15	18 (7%) 16 11	33, 59, 100, 122	0
1	C	251/261 (96%)	0.15	15 (5%) 21 16	35, 58, 99, 124	0
1	E	253/261 (96%)	0.18	20 (7%) 12 9	35, 59, 100, 122	0
2	B	12/15 (80%)	1.06	4 (33%) 0 0	41, 66, 90, 99	0
2	D	13/15 (86%)	0.94	3 (23%) 0 0	40, 60, 101, 109	0
2	F	13/15 (86%)	1.02	3 (23%) 0 0	41, 58, 109, 118	0
All	All	795/828 (96%)	0.20	63 (7%) 12 9	33, 59, 100, 124	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	185	THR	7.7
2	F	464	GLN	5.7
1	A	191	GLU	5.6
2	B	453	ALA	5.3
1	C	191	GLU	5.3
2	F	453	ALA	5.0
2	D	465	ARG	4.9
1	A	95	ASN	4.8
1	E	186	SER	4.7
1	A	164	LYS	4.6
1	C	120	ASP	4.6
1	A	192	GLU	4.1
1	C	121	LEU	3.9
2	D	453	ALA	3.8
1	C	185	THR	3.7
1	C	194	ALA	3.5
2	F	465	ARG	3.5
2	B	464	GLN	3.4
1	E	259	GLU	3.4

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	E	95	ASN	3.3
1	E	122	ASP	3.1
1	C	174	GLU	3.0
1	C	123	VAL	3.0
1	E	196	THR	3.0
1	A	124	GLU	2.9
1	A	64	ARG	2.9
1	C	122	ASP	2.9
1	E	163	ALA	2.8
1	E	166	GLY	2.8
1	A	96	ALA	2.8
2	D	464	GLN	2.7
1	C	199	MET	2.7
1	E	193	GLU	2.7
1	A	97	ASP	2.7
1	E	120	ASP	2.7
1	A	165	ASP	2.6
1	A	186	SER	2.6
1	E	109	GLU	2.6
1	E	55	GLU	2.5
2	B	461	GLY	2.5
1	C	193	GLU	2.5
1	C	63	ASP	2.5
1	A	122	ASP	2.5
1	E	258	GLU	2.4
1	C	64	ARG	2.4
1	E	108	GLN	2.3
1	A	132	GLU	2.3
1	C	175	LEU	2.3
1	E	165	ASP	2.3
1	C	132	GLU	2.2
1	A	63	ASP	2.2
2	B	463	PHE	2.2
1	E	59	THR	2.2
1	E	132	GLU	2.2
1	A	173	GLY	2.2
1	A	121	LEU	2.1
1	E	124	GLU	2.1
1	A	120	ASP	2.1
1	A	193	GLU	2.1
1	E	212	LEU	2.1
1	C	255	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	94	ASP	2.1
1	E	61	ARG	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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.