



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 02:24 PM GMT

PDB ID : 1E7B
Title : CRYSTAL STRUCTURE OF HUMAN SERUM ALBUMIN COMPLEXED
WITH THE GENERAL ANESTHETIC HALOTHANE
Authors : Bhattacharya, A.A.; Curry, S.; Franks, N.P.
Deposited on : 2000-08-26
Resolution : 2.38 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

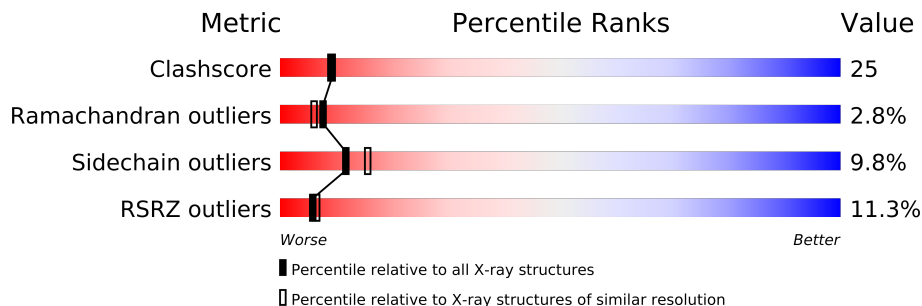
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance



The reported resolution of this entry is 2.38 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))
Clashscore	79885	3668 (2.40-2.36)
Ramachandran outliers	78287	3600 (2.40-2.36)
Sidechain outliers	78261	3602 (2.40-2.36)
RSRZ outliers	66119	2966 (2.40-2.36)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	585	
1	B	585	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	HLT	A	4001	-	X
2	HLT	A	4003	-	X
2	HLT	B	4001	-	X
2	HLT	B	4002	-	X
2	HLT	B	4003	-	X

2 Entry composition i

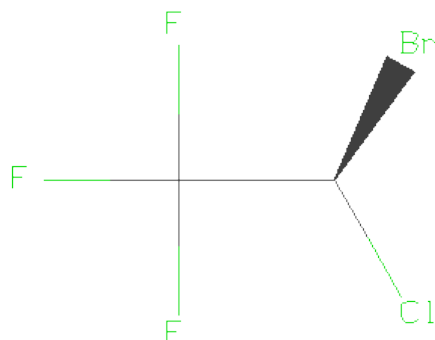
There are 3 unique types of molecules in this entry. The entry contains 8650 atoms, of which 0 are hydrogen and 0 are deuterium.

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 SERUM ALBUMIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	578	Total	C	N	O	S	0	0	0
			4303	2732	728	803	40			
1	B	576	Total	C	N	O	S	0	0	0
			4248	2684	717	806	41			

- Molecule 2 is 2-BROMO-2-CHLORO-1,1,1-TRIFLUOROETHANE (three-letter code: HLT) (formula: $C_2HBrClF_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	Br	C	Cl	F	0	0
			7	1	2	1	3		
2	A	1	Total	Br	C	Cl	F	0	0
			7	1	2	1	3		
2	A	1	Total	Br	C	Cl	F	0	0
			7	1	2	1	3		
2	B	1	Total	Br	C	Cl	F	0	0
			7	1	2	1	3		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total 7	Br 1	C 2	Cl 1	F 3	0	0
2	B	1	Total 7	Br 1	C 2	Cl 1	F 3	0	0

- Molecule 3 is water.

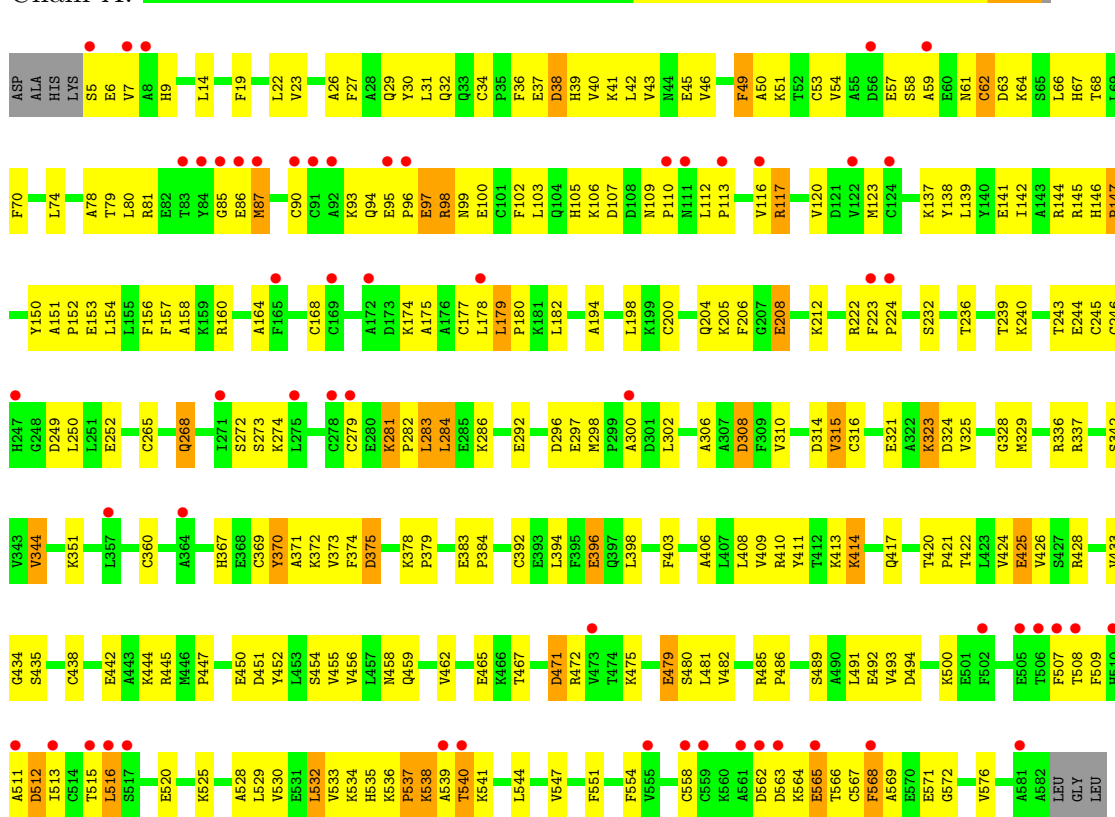
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	30	Total 30	O 30	0	0
3	B	27	Total 27	O 27	0	0

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 errors 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.

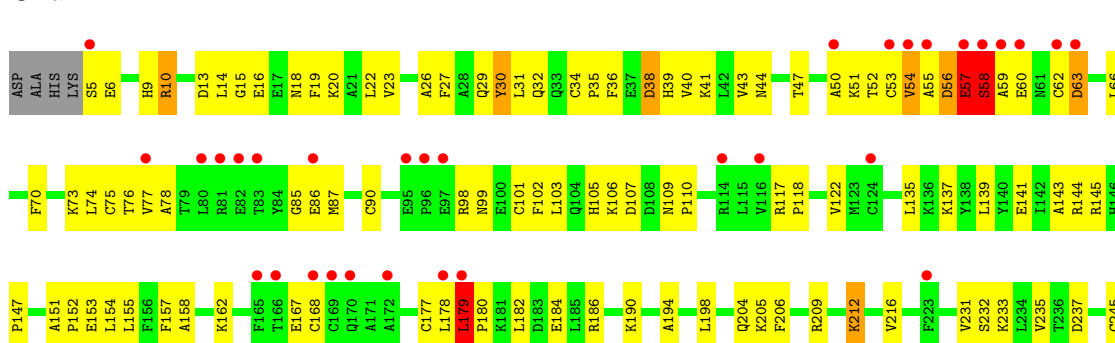
• Molecule 1: SERUM ALBUMIN

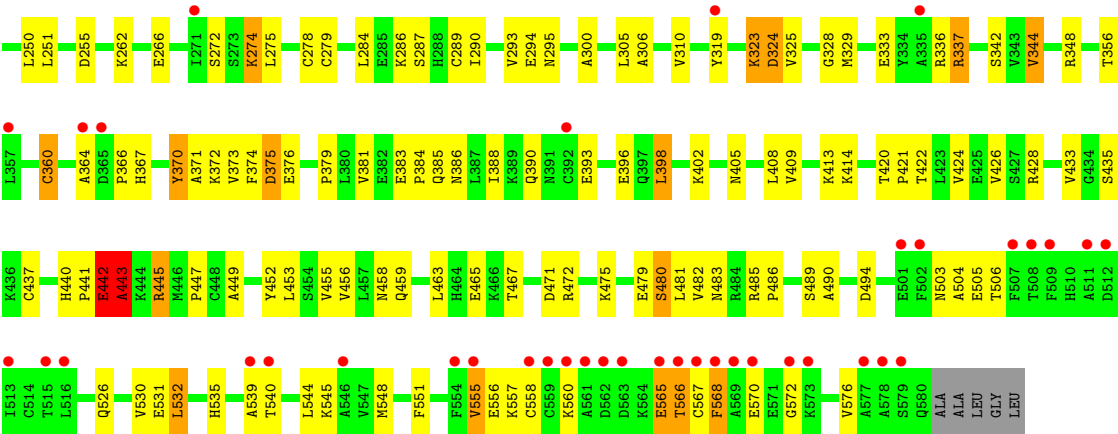
Chain A:



• Molecule 1: SERUM ALBUMIN

Chain B:





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	54.58Å 54.96Å 120.00Å 81.39° 90.79° 65.55°	Depositor
Resolution (Å)	17.00 – 2.38 14.97 – 2.38	Depositor EDS
% Data completeness (in resolution range)	96.0 (17.00-2.38) 96.2 (14.97-2.38)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 2.37Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.270 , 0.303 0.252 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	53.4	Xtriage
Anisotropy	0.430	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 81.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 47994 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8650	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.50% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0.34	0/4388	0.58	1/5965 (0.0%)
1	B	0.34	1/4329 (0.0%)	0.54	2/5892 (0.0%)
All	All	0.34	1/8717 (0.0%)	0.56	3/11857 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	442	GLU	CB-CG	7.29	1.66	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	443	ALA	N-CA-C	-6.56	93.28	111.00
1	B	539	ALA	N-CA-C	5.88	126.89	111.00
1	A	541	LYS	N-CA-C	-5.30	96.68	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts

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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4303	0	3992	233	0
1	B	4248	0	3880	175	0
2	A	21	0	0	2	0
2	B	21	0	0	4	0
3	A	30	0	0	2	0
3	B	27	0	0	1	0
All	All	8650	0	7872	409	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 25.

All (409) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:67:HIS:HB3	1:A:98:ARG:HH21	1.27	0.99
1:A:540:THR:HG23	1:A:544:LEU:HG	1.42	0.98
1:A:98:ARG:HH11	1:A:98:ARG:H	1.05	0.95
1:B:556:GLU:HG3	1:B:557:LYS:H	1.38	0.89
1:A:383:GLU:HB3	1:A:384:PRO:HD3	1.55	0.88
1:A:53:CYS:O	1:A:57:GLU:HG2	1.74	0.87
1:A:328:GLY:HA2	2:A:4001:HLT:BR	2.30	0.87
1:B:151:ALA:HB3	1:B:152:PRO:HD3	1.57	0.87
1:B:106:LYS:HD3	1:B:147:PRO:HB2	1.55	0.86
1:A:511:ALA:HB2	1:A:565:GLU:HB3	1.56	0.86
1:A:485:ARG:HB3	1:A:486:PRO:HD3	1.57	0.85
1:B:409:VAL:HG12	1:B:413:LYS:HE3	1.59	0.84
1:A:367:HIS:O	1:A:371:ALA:HB2	1.78	0.83
1:B:441:PRO:O	1:B:443:ALA:N	2.11	0.83
1:A:98:ARG:NH1	1:A:99:ASN:H	1.76	0.83
1:B:383:GLU:HB3	1:B:384:PRO:HD3	1.59	0.82
1:A:281:LYS:HB2	1:A:282:PRO:HD2	1.62	0.81
1:B:39:HIS:O	1:B:43:VAL:HG23	1.80	0.81
1:A:297:GLU:O	1:A:297:GLU:CA	2.29	0.81
1:A:472:ARG:HH12	1:A:494:ASP:HA	1.47	0.79
1:A:94:GLN:O	1:A:98:ARG:HB3	1.84	0.78
1:A:424:VAL:O	1:A:428:ARG:HG3	1.81	0.78
1:A:98:ARG:CZ	1:A:99:ASN:HB2	2.15	0.77
1:A:151:ALA:HB3	1:A:152:PRO:HD3	1.67	0.76
1:A:511:ALA:HA	1:A:568:PHE:CE2	2.20	0.75
1:A:14:LEU:HD13	1:A:22:LEU:HD12	1.68	0.75
1:A:306:ALA:HA	1:A:310:VAL:HG22	1.69	0.75
1:A:378:LYS:HB3	1:A:379:PRO:HD3	1.67	0.75
1:A:425:GLU:OE1	1:A:425:GLU:HA	1.85	0.75
1:B:22:LEU:HD21	1:B:155:LEU:HD11	1.68	0.75

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:556:GLU:HG3	1:B:557:LYS:N	2.01	0.75
1:A:558:CYS:HA	1:A:567:CYS:SG	2.27	0.74
1:A:198:LEU:HA	1:A:458:ASN:ND2	2.02	0.74
1:A:98:ARG:HH11	1:A:98:ARG:N	1.85	0.73
1:B:433:VAL:HG22	1:B:452:TYR:CD2	2.24	0.73
1:B:531:GLU:O	1:B:535:HIS:HD2	1.72	0.72
1:A:98:ARG:NH2	1:A:99:ASN:HB2	2.03	0.72
1:A:323:LYS:HG3	1:A:324:ASP:N	2.04	0.72
1:A:110:PRO:HG2	1:A:145:ARG:HA	1.72	0.71
1:A:26:ALA:HB2	1:A:250:LEU:HD12	1.70	0.71
1:A:564:LYS:O	1:A:566:THR:N	2.19	0.71
1:B:373:VAL:HG13	1:B:374:PHE:HD1	1.56	0.71
1:A:511:ALA:HB2	1:A:565:GLU:CB	2.20	0.71
1:A:540:THR:CG2	1:A:544:LEU:HG	2.20	0.70
1:A:120:VAL:HG21	1:A:175:ALA:HA	1.73	0.70
1:B:262:LYS:O	1:B:266:GLU:HG3	1.92	0.70
1:A:516:LEU:HD22	1:A:520:GLU:OE1	1.91	0.70
1:B:279:CYS:HA	1:B:286:LYS:HD2	1.74	0.70
1:A:141:GLU:OE1	1:A:144:ARG:HD3	1.92	0.69
1:B:16:GLU:O	1:B:20:LYS:HG2	1.92	0.69
1:A:34:CYS:HB3	1:A:39:HIS:NE2	2.08	0.69
1:B:52:THR:HA	1:B:56:ASP:OD2	1.92	0.69
1:A:179:LEU:HB2	1:A:180:PRO:HD3	1.75	0.69
1:A:279:CYS:HA	1:A:286:LYS:HD2	1.74	0.69
1:B:306:ALA:HA	1:B:310:VAL:HG22	1.75	0.68
1:B:34:CYS:HB3	1:B:39:HIS:NE2	2.09	0.68
1:A:98:ARG:NH1	1:A:98:ARG:H	1.87	0.68
1:A:110:PRO:C	1:A:112:LEU:H	1.96	0.68
1:A:107:ASP:O	1:A:147:PRO:HG2	1.94	0.67
1:B:449:ALA:O	1:B:453:LEU:HG	1.95	0.66
1:A:87:MET:HE3	1:A:105:HIS:HB3	1.76	0.66
1:A:141:GLU:O	1:A:145:ARG:HG3	1.95	0.65
1:A:208:GLU:OE2	1:A:212:LYS:HE3	1.96	0.65
1:B:367:HIS:O	1:B:371:ALA:HB2	1.96	0.65
1:B:420:THR:HB	1:B:421:PRO:HD3	1.78	0.65
1:B:373:VAL:HG13	1:B:374:PHE:CD1	2.31	0.65
1:B:153:GLU:O	1:B:157:PHE:HD1	1.80	0.65
1:A:23:VAL:O	1:A:27:PHE:HD1	1.78	0.65
1:B:424:VAL:O	1:B:428:ARG:HG3	1.97	0.64
1:B:26:ALA:HB2	1:B:250:LEU:HD12	1.80	0.64
1:A:472:ARG:NH1	1:A:494:ASP:HA	2.13	0.64
1:A:106:LYS:HD3	1:A:147:PRO:HB3	1.78	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:310:VAL:HG11	1:B:374:PHE:CE1	2.32	0.64
1:A:471:ASP:N	1:A:471:ASP:OD1	2.26	0.64
1:B:15:GLY:O	1:B:19:PHE:HB3	1.98	0.64
1:A:511:ALA:CB	1:A:565:GLU:HB3	2.28	0.63
1:A:7:VAL:HG22	1:A:66:LEU:HD23	1.81	0.63
1:A:394:LEU:HD11	1:A:398:LEU:HD11	1.80	0.63
1:A:67:HIS:CB	1:A:98:ARG:HH21	2.08	0.63
1:A:138:TYR:CE1	1:A:142:ILE:HD11	2.34	0.63
1:B:23:VAL:O	1:B:27:PHE:HD1	1.82	0.62
1:A:507:PHE:HD1	1:A:509:PHE:CE2	2.17	0.62
1:A:383:GLU:HB3	1:A:384:PRO:CD	2.29	0.62
1:A:281:LYS:HB2	1:A:282:PRO:CD	2.29	0.62
1:A:283:LEU:HG	1:A:284:LEU:HD23	1.81	0.62
1:B:141:GLU:O	1:B:145:ARG:HG3	1.98	0.62
1:A:420:THR:HB	1:A:421:PRO:HD3	1.82	0.62
1:A:406:ALA:O	1:A:409:VAL:HG12	2.00	0.61
1:B:22:LEU:CD2	1:B:155:LEU:HD11	2.29	0.61
1:A:472:ARG:HH21	1:A:491:LEU:HD22	1.65	0.61
1:A:297:GLU:CA	1:A:298:MET:N	2.63	0.61
1:A:205:LYS:HE3	1:A:465:GLU:OE2	2.00	0.61
1:B:90:CYS:O	1:B:98:ARG:HG3	2.01	0.60
1:A:433:VAL:HG22	1:A:452:TYR:CD2	2.36	0.60
1:B:290:ILE:O	1:B:293:VAL:HG12	2.01	0.60
1:A:507:PHE:HZ	1:A:576:VAL:HG22	1.67	0.60
1:A:39:HIS:O	1:A:43:VAL:HG23	2.02	0.60
1:A:31:LEU:HG	1:A:74:LEU:HD22	1.82	0.60
1:B:233:LYS:HE3	1:B:237:ASP:OD2	2.01	0.59
1:A:572:GLY:O	1:A:576:VAL:HG23	2.02	0.59
1:B:14:LEU:HD13	1:B:22:LEU:HD12	1.84	0.59
1:B:556:GLU:O	1:B:560:LYS:HG2	2.02	0.59
1:B:59:ALA:HB3	1:B:62:CYS:SG	2.43	0.59
1:A:49:PHE:HE1	1:A:62:CYS:SG	2.25	0.59
1:A:281:LYS:CB	1:A:282:PRO:HD2	2.33	0.59
1:A:141:GLU:OE1	1:A:141:GLU:HA	2.03	0.59
1:B:485:ARG:HB3	1:B:486:PRO:HD3	1.84	0.58
1:B:437:CYS:O	1:B:440:HIS:HB2	2.02	0.58
1:A:34:CYS:HB3	1:A:39:HIS:HE2	1.67	0.58
1:A:61:ASN:O	1:A:64:LYS:HB2	2.03	0.58
1:B:323:LYS:HG3	1:B:324:ASP:N	2.19	0.58
1:B:29:GLN:HG2	1:B:143:ALA:O	2.03	0.58
1:A:328:GLY:CA	2:A:4001:HLT:BR	3.06	0.58
1:A:66:LEU:O	1:A:70:PHE:HD2	1.84	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:212:LYS:O	1:B:216:VAL:HG23	2.03	0.57
1:A:485:ARG:HB3	1:A:486:PRO:CD	2.31	0.57
1:B:475:LYS:O	1:B:479:GLU:HB2	2.03	0.57
1:B:5:SER:HA	1:B:62:CYS:O	2.05	0.57
1:A:297:GLU:O	1:A:298:MET:N	2.38	0.57
1:A:49:PHE:CE1	1:A:53:CYS:SG	2.98	0.56
1:A:507:PHE:HD1	1:A:509:PHE:CD2	2.22	0.56
1:B:325:VAL:HG12	1:B:329:MET:CE	2.35	0.56
1:A:206:PHE:CE2	1:A:481:LEU:HD13	2.40	0.56
1:B:433:VAL:HG22	1:B:452:TYR:HD2	1.71	0.56
1:A:178:LEU:O	1:A:179:LEU:C	2.43	0.56
1:A:511:ALA:HA	1:A:568:PHE:CD2	2.40	0.56
1:A:138:TYR:CZ	1:A:142:ILE:HD11	2.41	0.56
1:A:99:ASN:HA	1:A:102:PHE:HD2	1.71	0.55
1:A:38:ASP:OD1	1:A:38:ASP:N	2.38	0.55
1:A:297:GLU:O	1:A:298:MET:HA	2.06	0.55
1:A:156:PHE:CE1	1:A:160:ARG:HD2	2.42	0.55
1:A:87:MET:CE	1:A:105:HIS:HB3	2.37	0.54
1:B:408:LEU:HD11	1:B:526:GLN:HB3	1.89	0.54
1:B:405:ASN:O	1:B:409:VAL:HG23	2.07	0.54
1:B:306:ALA:CA	1:B:310:VAL:HG22	2.37	0.54
1:A:38:ASP:O	1:A:42:LEU:HG	2.08	0.54
1:A:408:LEU:HD23	1:A:529:LEU:HD23	1.88	0.54
1:B:198:LEU:HA	1:B:458:ASN:ND2	2.22	0.54
1:A:49:PHE:HE1	1:A:53:CYS:SG	2.31	0.54
1:A:511:ALA:CB	1:A:565:GLU:CB	2.85	0.54
1:A:567:CYS:O	1:A:571:GLU:N	2.36	0.54
1:A:145:ARG:O	1:A:146:HIS:HD2	1.89	0.54
1:A:408:LEU:HD21	1:A:530:VAL:HG23	1.90	0.54
1:A:249:ASP:HB3	1:A:252:GLU:CD	2.28	0.54
1:B:383:GLU:HB3	1:B:384:PRO:CD	2.35	0.54
1:A:564:LYS:C	1:A:566:THR:H	2.09	0.54
1:A:32:GLN:NE2	1:A:147:PRO:HG3	2.23	0.54
1:B:98:ARG:O	1:B:101:CYS:HB3	2.08	0.54
1:A:452:TYR:O	1:A:456:VAL:HG23	2.07	0.54
1:B:274:LYS:CE	1:B:294:GLU:HG3	2.38	0.54
1:A:178:LEU:HG	1:A:182:LEU:HG	1.88	0.53
1:B:186:ARG:O	1:B:190:LYS:HG3	2.08	0.53
1:A:281:LYS:CB	1:A:282:PRO:CD	2.85	0.53
1:B:54:VAL:HG12	1:B:55:ALA:N	2.23	0.53
1:A:533:VAL:HG12	1:A:533:VAL:O	2.09	0.53
1:A:139:LEU:HD21	1:A:158:ALA:HB2	1.91	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:283:LEU:HG	1:A:284:LEU:N	2.22	0.53
1:A:49:PHE:O	1:A:49:PHE:HD1	1.91	0.53
1:B:370:TYR:C	1:B:370:TYR:CD1	2.81	0.53
1:A:372:LYS:O	1:A:375:ASP:HB2	2.09	0.53
1:A:558:CYS:CB	1:A:567:CYS:SG	2.97	0.53
1:A:370:TYR:CD1	1:A:370:TYR:C	2.83	0.52
1:A:23:VAL:HG13	1:A:70:PHE:HE1	1.73	0.52
1:B:565:GLU:HG3	1:B:565:GLU:O	2.08	0.52
1:B:348:ARG:HG3	1:B:482:VAL:HG12	1.90	0.52
1:A:351:LYS:HD3	3:A:2016:HOH:O	2.08	0.52
1:A:94:GLN:O	1:A:98:ARG:HD3	2.09	0.52
1:A:562:ASP:OD1	1:A:562:ASP:O	2.26	0.52
1:A:99:ASN:HA	1:A:102:PHE:CD2	2.45	0.52
1:A:530:VAL:O	1:A:534:LYS:HG3	2.09	0.52
1:B:99:ASN:HA	1:B:102:PHE:HD2	1.74	0.52
1:B:422:THR:O	1:B:426:VAL:HG23	2.09	0.52
1:A:30:TYR:HE1	1:A:103:LEU:HD23	1.75	0.52
1:A:5:SER:HA	1:A:62:CYS:O	2.10	0.52
1:A:507:PHE:CD1	1:A:509:PHE:CE2	2.98	0.52
1:A:41:LYS:O	1:A:45:GLU:HG3	2.09	0.52
1:B:205:LYS:HE2	1:B:465:GLU:OE1	2.10	0.52
1:A:509:PHE:CZ	1:A:551:PHE:CZ	2.98	0.51
1:B:14:LEU:HD13	1:B:22:LEU:CD1	2.39	0.51
1:A:116:VAL:HG22	1:A:117:ARG:N	2.24	0.51
1:B:342:SER:OG	1:B:344:VAL:HG23	2.10	0.51
1:B:36:PHE:O	1:B:40:VAL:HG23	2.10	0.51
1:A:558:CYS:CA	1:A:567:CYS:SG	2.97	0.51
1:B:558:CYS:SG	1:B:568:PHE:N	2.83	0.51
1:B:356:THR:O	1:B:360:CYS:HB2	2.11	0.51
1:A:239:THR:O	1:A:243:THR:OG1	2.27	0.51
1:B:422:THR:HG23	1:B:463:LEU:HD13	1.93	0.51
1:B:551:PHE:O	1:B:555:VAL:HG23	2.10	0.51
1:B:504:ALA:C	1:B:506:THR:H	2.13	0.51
1:A:279:CYS:HA	1:A:286:LYS:CD	2.38	0.51
1:A:373:VAL:HG13	1:A:374:PHE:CD1	2.46	0.51
1:A:342:SER:HA	1:A:447:PRO:HA	1.93	0.51
1:B:10:ARG:NH1	1:B:255:ASP:OD2	2.44	0.51
1:B:19:PHE:CD1	1:B:19:PHE:C	2.84	0.51
1:B:23:VAL:O	1:B:27:PHE:CD1	2.64	0.51
1:A:106:LYS:HD3	1:A:147:PRO:CB	2.41	0.51
1:B:480:SER:OG	1:B:483:ASN:HB2	2.11	0.51
1:A:66:LEU:HB3	1:A:70:PHE:CE2	2.45	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:107:ASP:HB3	1:B:110:PRO:HG3	1.92	0.50
1:B:398:LEU:O	1:B:402:LYS:HB2	2.11	0.50
1:B:118:PRO:HB2	1:B:122:VAL:HB	1.92	0.50
1:A:344:VAL:HG12	1:A:482:VAL:HG13	1.93	0.50
1:A:6:GLU:O	1:A:9:HIS:HB3	2.10	0.50
1:B:384:PRO:O	1:B:388:ILE:HG12	2.11	0.50
1:B:57:GLU:OE1	1:B:57:GLU:HA	2.10	0.50
1:A:67:HIS:CE1	1:A:99:ASN:ND2	2.80	0.50
1:A:297:GLU:O	1:A:298:MET:CA	2.60	0.50
1:B:566:THR:O	1:B:570:GLU:N	2.45	0.49
1:A:49:PHE:CD1	1:A:49:PHE:C	2.85	0.49
1:B:274:LYS:HE3	1:B:294:GLU:HG3	1.93	0.49
1:B:490:ALA:HB3	3:B:2024:HOH:O	2.10	0.49
1:B:370:TYR:C	1:B:370:TYR:HD1	2.15	0.49
1:A:51:LYS:HA	1:A:54:VAL:HG23	1.95	0.49
1:A:392:CYS:O	1:A:396:GLU:HG3	2.12	0.49
1:A:422:THR:O	1:A:426:VAL:HG23	2.13	0.49
1:A:564:LYS:C	1:A:566:THR:N	2.65	0.49
1:A:36:PHE:O	1:A:40:VAL:HG23	2.13	0.49
1:B:30:TYR:HE1	1:B:103:LEU:HD23	1.78	0.49
1:B:209:ARG:HG2	2:B:4001:HLT:F2	2.03	0.49
1:B:328:GLY:HA2	2:B:4001:HLT:BR	2.69	0.48
1:B:558:CYS:HB3	1:B:568:PHE:CD2	2.48	0.48
1:B:386:ASN:O	1:B:390:GLN:HB2	2.14	0.48
1:A:168:CYS:SG	1:A:177:CYS:C	2.91	0.48
1:A:179:LEU:HB2	1:A:180:PRO:CD	2.43	0.48
1:A:49:PHE:HD1	1:A:49:PHE:C	2.17	0.48
1:A:507:PHE:HZ	1:A:576:VAL:CG2	2.26	0.48
1:B:545:LYS:HA	1:B:548:MET:HB2	1.95	0.48
1:A:50:ALA:O	1:A:54:VAL:HG23	2.13	0.48
1:B:194:ALA:HB1	1:B:455:VAL:CG1	2.43	0.48
1:B:279:CYS:HA	1:B:286:LYS:CD	2.43	0.48
1:B:32:GLN:NE2	1:B:110:PRO:HG3	2.28	0.48
1:A:81:ARG:CB	1:A:85:GLY:HA2	2.44	0.48
1:A:200:CYS:O	1:A:204:GLN:HG3	2.13	0.48
1:A:314:ASP:O	1:A:315:VAL:C	2.52	0.48
1:A:373:VAL:HG13	1:A:374:PHE:HD1	1.78	0.48
1:A:509:PHE:CZ	1:A:551:PHE:CE1	3.01	0.48
1:A:475:LYS:O	1:A:479:GLU:HB2	2.13	0.48
1:B:106:LYS:HD3	1:B:147:PRO:CB	2.37	0.48
1:B:319:TYR:CE1	1:B:323:LYS:HB2	2.49	0.47
1:A:325:VAL:HG12	1:A:329:MET:HE2	1.94	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:29:GLN:HG2	1:A:147:PRO:HA	1.95	0.47
1:B:503:ASN:OD1	1:B:504:ALA:N	2.47	0.47
1:A:51:LYS:HA	1:A:54:VAL:CG2	2.45	0.47
1:A:110:PRO:C	1:A:112:LEU:N	2.64	0.47
1:A:41:LYS:HE3	1:A:42:LEU:HD23	1.97	0.47
1:B:107:ASP:O	1:B:110:PRO:HD3	2.14	0.47
1:A:98:ARG:NH1	1:A:99:ASN:HB2	2.30	0.47
1:A:511:ALA:O	1:A:513:ILE:N	2.48	0.47
1:B:135:LEU:HD11	1:B:162:LYS:HB2	1.96	0.47
1:B:50:ALA:O	1:B:54:VAL:HG23	2.14	0.47
1:B:409:VAL:CG1	1:B:413:LYS:HE3	2.39	0.47
1:A:41:LYS:HE3	1:A:42:LEU:CD2	2.45	0.47
1:B:558:CYS:SG	1:B:567:CYS:C	2.94	0.47
1:A:563:ASP:C	1:A:564:LYS:O	2.49	0.47
1:A:42:LEU:O	1:A:46:VAL:HG23	2.15	0.47
1:B:32:GLN:NE2	1:B:110:PRO:CG	2.78	0.47
1:B:178:LEU:HG	1:B:182:LEU:HG	1.96	0.47
1:A:425:GLU:OE1	1:A:425:GLU:CA	2.60	0.46
1:B:364:ALA:O	1:B:366:PRO:HD3	2.14	0.46
1:B:206:PHE:CE2	1:B:481:LEU:HD13	2.51	0.46
1:B:572:GLY:O	1:B:576:VAL:HG23	2.15	0.46
1:A:223:PHE:CD1	1:A:272:SER:HB2	2.50	0.46
1:B:388:ILE:HG21	1:B:445:ARG:HB3	1.96	0.46
1:A:153:GLU:O	1:A:157:PHE:HD2	1.97	0.46
1:B:531:GLU:O	1:B:535:HIS:CD2	2.61	0.46
1:B:472:ARG:HH12	1:B:494:ASP:CA	2.27	0.46
1:A:66:LEU:HB3	1:A:70:PHE:HE2	1.80	0.46
1:A:403:PHE:O	1:A:406:ALA:HB3	2.16	0.46
1:B:504:ALA:C	1:B:506:THR:N	2.69	0.46
1:B:376:GLU:O	1:B:379:PRO:HD2	2.15	0.46
1:B:63:ASP:OD1	1:B:63:ASP:O	2.34	0.46
1:B:472:ARG:NH1	1:B:494:ASP:CB	2.79	0.46
1:B:31:LEU:HG	1:B:74:LEU:HD22	1.98	0.46
1:B:73:LYS:O	1:B:76:THR:HG23	2.16	0.46
1:A:507:PHE:CZ	1:A:576:VAL:HG22	2.48	0.46
1:B:137:LYS:O	1:B:141:GLU:HG2	2.16	0.46
1:B:44:ASN:HA	1:B:47:THR:HG22	1.98	0.46
1:A:459:GLN:O	1:A:462:VAL:HG22	2.16	0.46
1:B:441:PRO:O	1:B:442:GLU:C	2.52	0.46
1:A:373:VAL:HG13	1:A:374:PHE:N	2.31	0.46
1:A:532:LEU:HD11	1:A:547:VAL:HG11	1.98	0.46
1:A:310:VAL:O	1:A:370:TYR:HE1	1.98	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:61:ASN:C	1:A:63:ASP:N	2.69	0.45
1:A:360:CYS:SG	1:A:370:TYR:N	2.90	0.45
1:A:409:VAL:HG13	1:A:410:ARG:N	2.31	0.45
1:B:57:GLU:O	1:B:59:ALA:N	2.50	0.45
1:B:51:LYS:C	1:B:53:CYS:H	2.20	0.45
1:B:372:LYS:O	1:B:375:ASP:HB2	2.17	0.45
1:B:381:VAL:O	1:B:385:GLN:HG3	2.15	0.45
1:A:90:CYS:HA	1:A:93:LYS:HG3	1.98	0.45
1:B:393:GLU:HA	1:B:396:GLU:HG3	1.98	0.45
1:B:342:SER:HA	1:B:447:PRO:HA	1.99	0.45
1:B:472:ARG:NH1	1:B:494:ASP:HA	2.32	0.45
1:A:554:PHE:CZ	1:A:568:PHE:HD1	2.35	0.45
1:A:81:ARG:CA	1:A:85:GLY:HA2	2.47	0.45
1:A:98:ARG:HG2	1:A:99:ASN:N	2.30	0.45
1:A:509:PHE:CE1	1:A:551:PHE:CZ	3.04	0.45
1:A:554:PHE:CE1	1:A:572:GLY:HA2	2.52	0.45
1:A:120:VAL:HG21	1:A:175:ALA:CA	2.44	0.45
1:B:141:GLU:OE1	1:B:144:ARG:HD3	2.16	0.45
1:A:351:LYS:CD	3:A:2016:HOH:O	2.62	0.45
1:B:472:ARG:NH1	1:B:494:ASP:CA	2.80	0.45
1:A:413:LYS:NZ	1:A:537:PRO:O	2.50	0.45
1:A:117:ARG:HG3	1:A:123:MET:HE3	1.97	0.44
1:A:492:GLU:HG3	1:A:493:VAL:H	1.81	0.44
1:A:67:HIS:CE1	1:A:99:ASN:HD21	2.35	0.44
1:A:538:LYS:O	1:A:539:ALA:C	2.55	0.44
1:B:456:VAL:O	1:B:459:GLN:HB3	2.17	0.44
1:A:434:GLY:O	1:A:438:CYS:HB2	2.17	0.44
1:B:139:LEU:HD22	1:B:154:LEU:HG	2.00	0.44
1:B:87:MET:HE3	1:B:105:HIS:HB3	1.99	0.44
1:A:507:PHE:CD1	1:A:509:PHE:HE2	2.34	0.44
1:B:139:LEU:HD21	1:B:158:ALA:HB2	1.99	0.44
1:A:566:THR:O	1:A:569:ALA:N	2.48	0.44
1:A:61:ASN:O	1:A:63:ASP:N	2.50	0.44
1:B:70:PHE:CD1	1:B:70:PHE:N	2.86	0.44
1:A:525:LYS:O	1:A:528:ALA:HB3	2.17	0.44
1:A:265:CYS:SG	1:A:286:LYS:HD2	2.58	0.44
1:A:420:THR:HG23	1:A:530:VAL:CG1	2.47	0.44
1:A:325:VAL:HG12	1:A:329:MET:CE	2.47	0.44
1:A:179:LEU:O	1:A:180:PRO:C	2.56	0.44
1:A:14:LEU:HD13	1:A:22:LEU:CD1	2.42	0.44
1:B:141:GLU:HA	1:B:141:GLU:OE1	2.18	0.44
1:A:204:GLN:HE21	1:A:246:CYS:HB3	1.82	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:19:PHE:C	1:A:19:PHE:CD1	2.91	0.44
1:A:500:LYS:O	1:A:535:HIS:CD2	2.71	0.43
1:B:135:LEU:HD11	1:B:162:LYS:HD3	2.00	0.43
1:B:540:THR:CB	1:B:544:LEU:HG	2.47	0.43
1:B:30:TYR:CE1	1:B:103:LEU:HD23	2.53	0.43
1:B:373:VAL:C	1:B:375:ASP:H	2.21	0.43
1:B:9:HIS:O	1:B:13:ASP:HB2	2.18	0.43
1:A:512:ASP:O	1:A:515:THR:HG22	2.19	0.43
1:A:394:LEU:CD1	1:A:398:LEU:HD11	2.47	0.43
1:B:6:GLU:HG2	1:B:66:LEU:HD11	2.01	0.43
1:A:360:CYS:SG	1:A:369:CYS:C	2.97	0.43
1:A:274:LYS:HE3	1:A:296:ASP:HA	2.01	0.43
1:A:95:GLU:O	1:A:96:PRO:C	2.57	0.43
1:A:451:ASP:O	1:A:454:SER:HB2	2.18	0.43
1:B:422:THR:HG23	1:B:463:LEU:CD1	2.49	0.43
1:B:209:ARG:CG	2:B:4001:HLT:F2	2.56	0.43
1:A:509:PHE:CE1	1:A:551:PHE:HZ	2.36	0.43
1:B:272:SER:HB3	1:B:275:LEU:HG	2.01	0.43
1:B:310:VAL:CG1	1:B:374:PHE:CE1	3.01	0.43
1:B:367:HIS:HA	1:B:370:TYR:CZ	2.54	0.43
1:A:78:ALA:C	1:A:80:LEU:H	2.21	0.43
1:B:66:LEU:HD22	1:B:251:LEU:HD12	2.00	0.42
1:A:95:GLU:O	1:A:97:GLU:N	2.52	0.42
1:B:442:GLU:HA	1:B:445:ARG:HD2	2.00	0.42
1:B:56:ASP:N	1:B:56:ASP:OD1	2.52	0.42
1:B:98:ARG:HG2	1:B:102:PHE:CE2	2.54	0.42
1:B:117:ARG:HA	1:B:118:PRO:HD3	1.85	0.42
1:A:536:LYS:N	1:A:537:PRO:HD3	2.33	0.42
1:B:75:CYS:HA	1:B:78:ALA:HB3	2.01	0.42
1:A:240:LYS:HE2	1:A:244:GLU:OE2	2.19	0.42
1:A:194:ALA:HB1	1:A:455:VAL:CG1	2.50	0.42
1:B:374:PHE:CD1	1:B:374:PHE:N	2.87	0.42
1:B:532:LEU:HA	1:B:532:LEU:HD23	1.88	0.42
1:B:531:GLU:HA	1:B:531:GLU:OE1	2.19	0.42
1:A:408:LEU:HD11	1:A:530:VAL:CG2	2.49	0.42
1:A:61:ASN:C	1:A:63:ASP:H	2.21	0.42
1:B:179:LEU:HB2	1:B:180:PRO:HD3	2.01	0.42
1:B:168:CYS:SG	1:B:177:CYS:C	2.98	0.42
1:A:100:GLU:HA	1:A:100:GLU:OE1	2.20	0.42
1:B:30:TYR:OH	1:B:103:LEU:HD21	2.20	0.42
1:A:420:THR:HG23	1:A:530:VAL:HG11	2.02	0.42
1:B:57:GLU:HB3	1:B:58:SER:H	1.52	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:485:ARG:O	1:B:486:PRO:C	2.58	0.42
1:B:30:TYR:HA	1:B:30:TYR:HD1	1.74	0.42
1:B:70:PHE:N	1:B:70:PHE:HD1	2.17	0.42
1:A:222:ARG:C	1:A:224:PRO:HD3	2.40	0.42
1:A:99:ASN:O	1:A:103:LEU:HG	2.19	0.42
1:B:178:LEU:O	1:B:179:LEU:C	2.58	0.42
1:A:492:GLU:O	1:A:493:VAL:C	2.57	0.42
1:A:539:ALA:O	1:A:540:THR:OG1	2.33	0.42
1:B:151:ALA:HB3	1:B:152:PRO:CD	2.39	0.42
1:B:373:VAL:C	1:B:375:ASP:N	2.73	0.42
1:A:535:HIS:O	1:A:535:HIS:ND1	2.52	0.42
1:A:417:GLN:H	1:A:417:GLN:CD	2.23	0.42
1:B:34:CYS:HA	1:B:35:PRO:HD3	1.90	0.41
1:B:374:PHE:N	1:B:374:PHE:HD1	2.18	0.41
1:B:408:LEU:HD23	1:B:408:LEU:HA	1.94	0.41
1:B:278:CYS:HB3	1:B:289:CYS:HB3	1.91	0.41
1:A:370:TYR:HD1	1:A:370:TYR:C	2.22	0.41
1:A:110:PRO:HB3	1:A:112:LEU:HG	2.01	0.41
1:A:308:ASP:N	1:A:308:ASP:OD1	2.53	0.41
1:A:97:GLU:CB	1:A:100:GLU:CG	2.98	0.41
1:B:18:ASN:O	1:B:22:LEU:HG	2.21	0.41
1:B:109:ASN:O	1:B:110:PRO:C	2.58	0.41
1:B:29:GLN:HG3	1:B:143:ALA:HB1	2.01	0.41
1:A:302:LEU:HA	1:A:302:LEU:HD23	1.95	0.41
1:B:231:VAL:O	1:B:235:VAL:HG23	2.21	0.41
1:B:333:GLU:O	1:B:337:ARG:HG2	2.20	0.41
1:A:164:ALA:O	1:A:178:LEU:HD12	2.20	0.41
1:A:433:VAL:HG22	1:A:452:TYR:HD2	1.83	0.41
1:A:139:LEU:HD22	1:A:154:LEU:HG	2.03	0.41
1:A:394:LEU:O	1:A:398:LEU:HG	2.20	0.41
1:B:472:ARG:NH1	1:B:494:ASP:HB2	2.36	0.41
1:A:563:ASP:O	1:A:564:LYS:C	2.58	0.41
1:A:344:VAL:HG23	1:A:450:GLU:HG2	2.02	0.41
1:A:411:TYR:HA	1:A:414:LYS:HD3	2.01	0.41
1:B:408:LEU:HD22	1:B:530:VAL:CG2	2.51	0.41
1:B:32:GLN:HE21	1:B:110:PRO:HG2	1.86	0.41
2:B:4001:HLT:CL	2:B:4002:HLT:F1	2.66	0.41
1:A:315:VAL:HG12	1:A:316:CYS:N	2.35	0.41
1:B:38:ASP:O	1:B:41:LYS:HB3	2.21	0.41
1:A:378:LYS:HB3	1:A:379:PRO:CD	2.46	0.40
1:B:29:GLN:HG2	1:B:147:PRO:HA	2.02	0.40
1:A:507:PHE:CD1	1:A:509:PHE:CD2	3.07	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:179:LEU:HD22	1:B:179:LEU:HA	1.91	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	574/585 (98%)	498 (87%)	58 (10%)	18 (3%)	7 5
1	B	574/585 (98%)	493 (86%)	67 (12%)	14 (2%)	9 8
All	All	1148/1170 (98%)	991 (86%)	125 (11%)	32 (3%)	8 6

All (32) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	268	GLN
1	A	300	ALA
1	A	315	VAL
1	A	538	LYS
1	B	54	VAL
1	B	57	GLU
1	B	58	SER
1	B	442	GLU
1	A	58	SER
1	A	59	ALA
1	A	512	ASP
1	A	540	THR
1	B	60	GLU
1	B	85	GLY
1	B	300	ALA
1	B	555	VAL
1	B	565	GLU
1	A	479	GLU
1	A	62	CYS

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Mol	Chain	Res	Type
1	A	97	GLU
1	A	150	TYR
1	A	179	LEU
1	A	565	GLU
1	A	147	PRO
1	A	321	GLU
1	B	77	VAL
1	B	86	GLU
1	B	443	ALA
1	B	505	GLU
1	B	179	LEU
1	A	113	PRO
1	A	537	PRO

5.3.2 Protein sidechains ⓘ

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	425/511 (83%)	381 (90%)	44 (10%)	10	13
1	B	415/511 (81%)	377 (91%)	38 (9%)	13	18
All	All	840/1022 (82%)	758 (90%)	82 (10%)	12	15

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	GLU
1	A	38	ASP
1	A	49	PHE
1	A	68	THR
1	A	79	THR
1	A	86	GLU
1	A	87	MET
1	A	98	ARG
1	A	109	ASN
1	A	117	ARG
1	A	137	LYS
1	A	174	LYS

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Mol	Chain	Res	Type
1	A	208	GLU
1	A	232	SER
1	A	236	THR
1	A	245	CYS
1	A	268	GLN
1	A	273	SER
1	A	281	LYS
1	A	283	LEU
1	A	284	LEU
1	A	292	GLU
1	A	308	ASP
1	A	323	LYS
1	A	336	ARG
1	A	337	ARG
1	A	344	VAL
1	A	370	TYR
1	A	375	ASP
1	A	396	GLU
1	A	414	LYS
1	A	425	GLU
1	A	435	SER
1	A	442	GLU
1	A	444	LYS
1	A	445	ARG
1	A	467	THR
1	A	471	ASP
1	A	480	SER
1	A	489	SER
1	A	508	THR
1	A	516	LEU
1	A	532	LEU
1	A	568	PHE
1	B	10	ARG
1	B	30	TYR
1	B	38	ASP
1	B	56	ASP
1	B	57	GLU
1	B	58	SER
1	B	63	ASP
1	B	167	GLU
1	B	179	LEU
1	B	184	GLU

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Mol	Chain	Res	Type
1	B	204	GLN
1	B	212	LYS
1	B	232	SER
1	B	245	CYS
1	B	274	LYS
1	B	284	LEU
1	B	287	SER
1	B	295	ASN
1	B	305	LEU
1	B	323	LYS
1	B	324	ASP
1	B	336	ARG
1	B	337	ARG
1	B	344	VAL
1	B	360	CYS
1	B	370	TYR
1	B	375	ASP
1	B	398	LEU
1	B	414	LYS
1	B	435	SER
1	B	445	ARG
1	B	467	THR
1	B	471	ASP
1	B	480	SER
1	B	489	SER
1	B	532	LEU
1	B	566	THR
1	B	568	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	HIS
1	A	109	ASN
1	A	146	HIS
1	A	204	GLN
1	A	318	ASN
1	A	385	GLN
1	A	458	ASN
1	A	483	ASN
1	B	32	GLN
1	B	196	GLN

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Mol	Chain	Res	Type
1	B	386	ASN
1	B	483	ASN
1	B	535	HIS

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	HLT	A	4001	-	6,6,6	1.00	1 (16%)	9,9,9	1.62	1 (11%)
2	HLT	A	4002	-	6,6,6	0.70	0	9,9,9	1.30	0
2	HLT	A	4003	-	6,6,6	0.95	0	9,9,9	1.55	1 (11%)
2	HLT	B	4001	-	6,6,6	0.81	0	9,9,9	1.40	1 (11%)
2	HLT	B	4002	-	6,6,6	0.59	0	9,9,9	1.30	0
2	HLT	B	4003	-	6,6,6	0.80	0	9,9,9	1.61	1 (11%)

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	HLT	A	4001	-	-	0/6/6/6	0/0/0/0
2	HLT	A	4002	-	-	0/6/6/6	0/0/0/0
2	HLT	A	4003	-	-	0/6/6/6	0/0/0/0
2	HLT	B	4001	-	-	0/6/6/6	0/0/0/0
2	HLT	B	4002	-	-	0/6/6/6	0/0/0/0
2	HLT	B	4003	-	-	0/6/6/6	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	4001	HLT	BR-C1	-2.08	1.87	1.95

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	4003	HLT	BR-C1-CL	-3.33	108.72	111.13
2	A	4001	HLT	BR-C1-CL	-3.13	108.86	111.13
2	B	4003	HLT	BR-C1-CL	-3.01	108.95	111.13
2	B	4001	HLT	BR-C1-CL	-2.28	109.48	111.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	578/585 (98%)	0.57	58 (10%) 8 8	25, 69, 132, 144	0
1	B	576/585 (98%)	0.68	71 (12%) 5 5	34, 78, 137, 150	0
All	All	1154/1170 (98%)	0.62	129 (11%) 6 7	25, 73, 135, 150	0

All (129) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	539	ALA	10.2
1	B	539	ALA	9.6
1	B	563	ASP	8.7
1	B	562	ASP	7.6
1	A	96	PRO	6.1
1	A	85	GLY	5.8
1	A	506	THR	5.5
1	B	83	THR	5.4
1	A	568	PHE	5.1
1	B	558	CYS	5.0
1	B	508	THR	4.9
1	B	577	ALA	4.6
1	B	507	PHE	4.5
1	A	516	LEU	4.4
1	A	565	GLU	4.4
1	B	82	GLU	4.3
1	A	300	ALA	4.3
1	A	84	TYR	4.3
1	B	364	ALA	4.3
1	B	165	PHE	4.2
1	A	505	GLU	4.1
1	B	572	GLY	4.0
1	B	172	ALA	4.0
1	A	364	ALA	4.0

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Mol	Chain	Res	Type	RSRZ
1	B	569	ALA	3.9
1	A	83	THR	3.9
1	A	513	ILE	3.8
1	B	116	VAL	3.8
1	B	566	THR	3.8
1	B	516	LEU	3.8
1	B	515	THR	3.8
1	B	512	ASP	3.7
1	B	567	CYS	3.7
1	B	546	ALA	3.7
1	B	540	THR	3.6
1	A	56	ASP	3.6
1	B	578	ALA	3.5
1	B	96	PRO	3.5
1	A	116	VAL	3.4
1	A	515	THR	3.3
1	A	563	ASP	3.3
1	B	60	GLU	3.3
1	B	502	PHE	3.3
1	A	271	ILE	3.2
1	A	581	ALA	3.2
1	B	509	PHE	3.2
1	A	90	CYS	3.2
1	A	357	LEU	3.2
1	A	558	CYS	3.2
1	B	511	ALA	3.2
1	A	223	PHE	3.2
1	A	510	HIS	3.1
1	B	5	SER	3.1
1	B	568	PHE	3.1
1	A	178	LEU	3.1
1	A	95	GLU	3.1
1	B	81	ARG	3.1
1	B	95	GLU	3.0
1	A	87	MET	3.0
1	B	58	SER	3.0
1	A	91	CYS	3.0
1	B	169	CYS	3.0
1	B	555	VAL	3.0
1	A	59	ALA	3.0
1	B	570	GLU	2.9
1	A	507	PHE	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	169	CYS	2.8
1	A	7	VAL	2.8
1	B	124	CYS	2.8
1	B	77	VAL	2.8
1	A	559	CYS	2.8
1	B	223	PHE	2.7
1	B	565	GLU	2.7
1	A	275	LEU	2.7
1	A	165	PHE	2.7
1	A	561	ALA	2.7
1	B	168	CYS	2.6
1	A	122	VAL	2.6
1	A	172	ALA	2.6
1	B	365	ASP	2.6
1	B	86	GLU	2.6
1	B	97	GLU	2.6
1	A	517	SER	2.6
1	A	511	ALA	2.5
1	A	111	ASN	2.5
1	A	562	ASP	2.5
1	B	513	ILE	2.5
1	B	573	LYS	2.5
1	A	473	VAL	2.5
1	B	579	SER	2.5
1	B	392	CYS	2.5
1	A	508	THR	2.4
1	A	224	PRO	2.4
1	B	55	ALA	2.4
1	A	113	PRO	2.4
1	B	554	PHE	2.4
1	B	560	LYS	2.4
1	A	5	SER	2.4
1	B	357	LEU	2.4
1	A	86	GLU	2.4
1	B	501	GLU	2.4
1	A	502	PHE	2.4
1	B	50	ALA	2.3
1	B	62	CYS	2.3
1	B	179	LEU	2.3
1	A	279	CYS	2.3
1	B	59	ALA	2.3
1	B	335	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	319	TYR	2.2
1	A	124	CYS	2.2
1	B	559	CYS	2.2
1	A	555	VAL	2.2
1	B	271	ILE	2.2
1	B	170	GLN	2.2
1	B	166	THR	2.2
1	A	8	ALA	2.2
1	B	178	LEU	2.2
1	B	114	ARG	2.1
1	A	540	THR	2.1
1	B	80	LEU	2.1
1	B	57	GLU	2.1
1	B	63	ASP	2.1
1	A	247	HIS	2.1
1	A	92	ALA	2.1
1	B	561	ALA	2.1
1	A	278	CYS	2.0
1	A	110	PRO	2.0
1	B	54	VAL	2.0
1	B	53	CYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	HLT	A	4003	7/7	0.55	17.29	84,86,87,95	7

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	HLT	B	4003	7/7	0.60	14.10	85,86,90,94	7
2	HLT	B	4002	7/7	0.28	4.34	75,76,77,81	7
2	HLT	B	4001	7/7	0.28	4.20	91,91,92,95	7
2	HLT	A	4001	7/7	0.31	3.78	73,73,74,81	7
2	HLT	A	4002	7/7	0.22	1.43	70,70,72,78	7

6.5 Other polymers ⓘ

There are no such residues in this entry.