Full wwPDB X-ray Structure Validation Report (i)

Nov 3, 2016 – 01:25 PM EDT

PDB ID	:	5FUO
Title	:	Extending the half-life of a Fab fragment through generation of a humanised anti-Human Serum Albumin (HSA) Fv domain: an investigation into the correlation between affinity and serum half-life
Authors	:	Adams, R.; Ceska, T.
Deposited on	:	2016-01-28
Resolution	:	3.60 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*
A user guide is available at
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

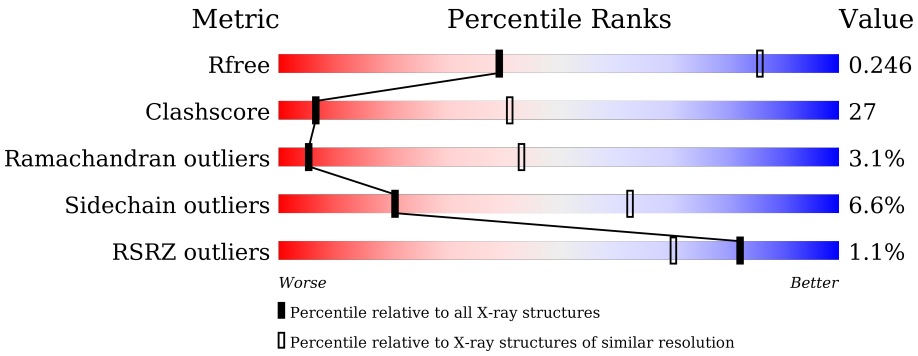
MolProbity	:	4.02b-467
Mogul	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20028320
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20028320

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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}	91344	1408 (3.80-3.40)
Clashscore	102246	1010 (3.74-3.46)
Ramachandran outliers	100387	1007 (3.76-3.44)
Sidechain outliers	100360	1007 (3.76-3.44)
RSRZ outliers	91569	1003 (3.78-3.42)

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. 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 <=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	585	2% 54% 41%
2	H	233	46% 43% 7%
3	L	217	47% 47% 7%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7869 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 SERUM ALBUMIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	579	4600	2903	777	879	41	0	0	1

- Molecule 2 is a protein called FAB HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	H	217	1623	1034	271	313	5	0	0	1

- Molecule 3 is a protein called FAB LIGHT CHAIN.

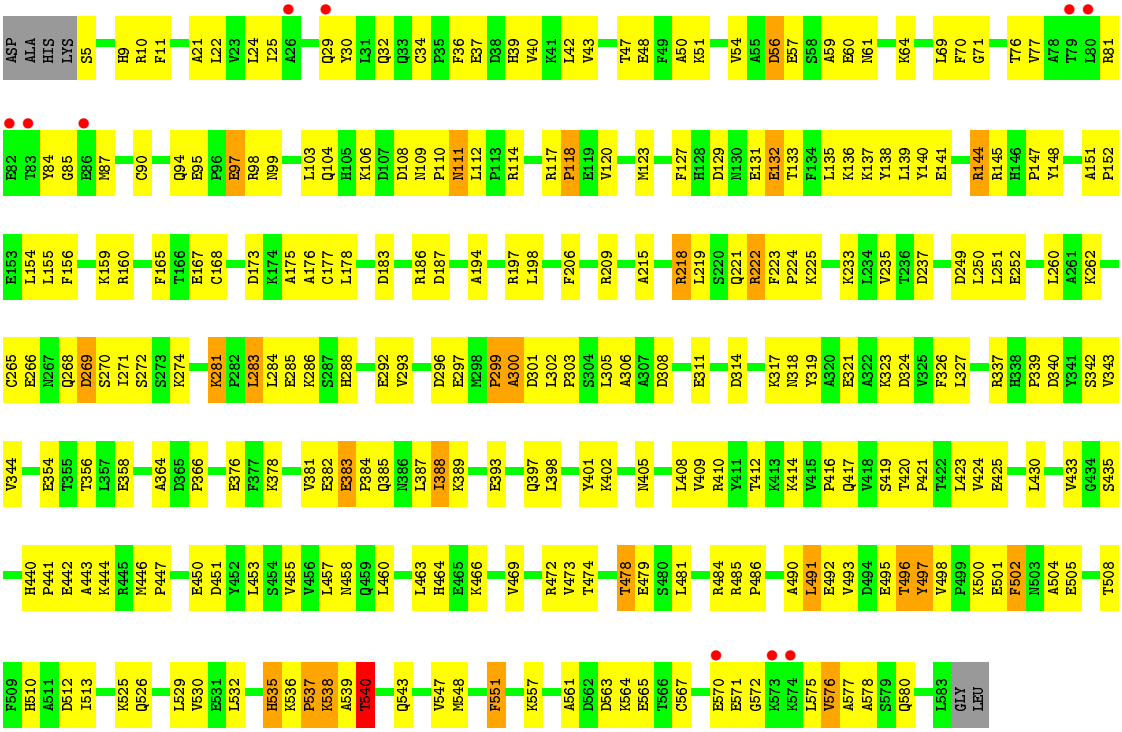
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	L	217	1646	1028	271	341	6	0	0	0

3 Residue-property plots (i)

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.

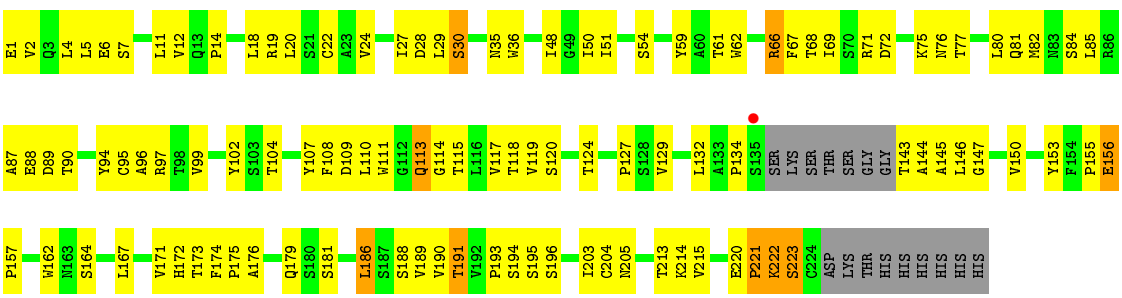
- Molecule 1: SERUM ALBUMIN

Chain A:



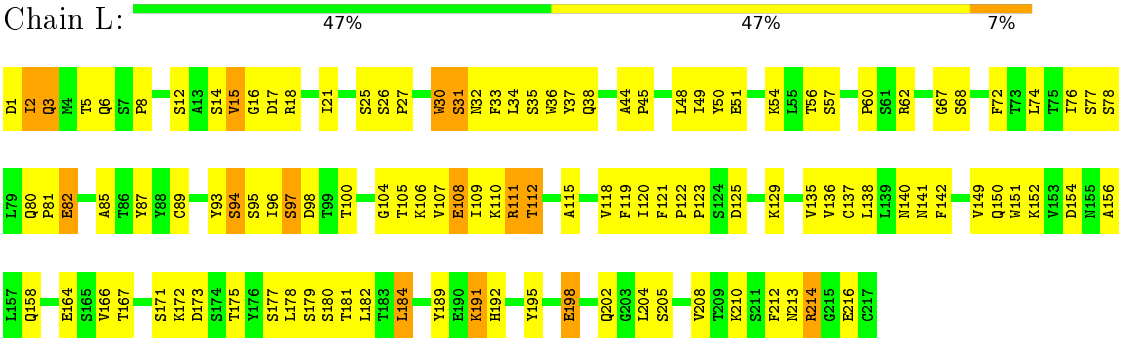
- Molecule 2: FAB HEAVY CHAIN

Chain H:



• Molecule 3: FAB LIGHT CHAIN

Chain L:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	217.64Å 217.64Å 78.65Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 3.60 28.42 – 3.60	Depositor EDS
% Data completeness (in resolution range)	97.6 (30.00-3.60) 97.7 (28.42-3.60)	Depositor EDS
R_{merge}	0.01	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.06 (at 3.65Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.215 , 0.254 0.213 , 0.246	Depositor DCC
R_{free} test set	2405 reflections (10.93%)	DCC
Wilson B-factor (Å^2)	104.4	Xtriage
Anisotropy	0.340	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 60.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.43$, $< L^2 > = 0.25$	Xtriage
Estimated twinning fraction	0.069 for -h,-k,l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	7869	wwPDB-VP
Average B, all atoms (Å^2)	115.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.50% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.52	0/4689	0.67	0/6326
2	H	0.55	0/1664	0.73	0/2275
3	L	0.53	0/1682	0.76	0/2282
All	All	0.53	0/8035	0.70	0/10883

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

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	4600	0	4518	212	0
2	H	1623	0	1601	119	0
3	L	1646	0	1590	117	0
All	All	7869	0	7709	423	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 27.

All (423) 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:A:106:LYS:HD2	1:A:148:TYR:CE1	2.00	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:485:ARG:HB3	1:A:486:PRO:HD3	1.49	0.93
3:L:12:SER:HB3	3:L:110:LYS:HE3	1.49	0.92
3:L:136:VAL:HG22	3:L:181:THR:HG23	1.51	0.91
2:H:214:LYS:HB2	2:H:214:LYS:HZ2	1.41	0.86
1:A:317:LYS:O	1:A:321:GLU:HG2	1.76	0.85
3:L:26:SER:HB3	3:L:27:PRO:HD3	1.56	0.85
2:H:2:VAL:HG13	2:H:27:ILE:HG12	1.59	0.85
1:A:302:LEU:HB3	1:A:337:ARG:NH1	1.93	0.84
3:L:178:LEU:HD23	3:L:179:SER:N	1.93	0.83
2:H:144:ALA:HB1	3:L:119:PHE:HE2	1.43	0.82
2:H:214:LYS:HB2	2:H:214:LYS:NZ	1.93	0.82
2:H:12:VAL:HG11	2:H:85:LEU:HD12	1.61	0.82
1:A:156:PHE:CE1	1:A:285:GLU:HA	2.15	0.81
1:A:120:VAL:HG21	1:A:175:ALA:HA	1.62	0.81
1:A:156:PHE:HE1	1:A:285:GLU:HA	1.46	0.79
3:L:85:ALA:O	3:L:107:VAL:HG22	1.83	0.78
2:H:127:PRO:HB3	2:H:153:TYR:HB3	1.64	0.78
1:A:151:ALA:HB3	1:A:152:PRO:HD3	1.66	0.77
2:H:143:THR:HB	2:H:193:PRO:HA	1.66	0.76
3:L:54:LYS:HZ3	3:L:54:LYS:HB3	1.49	0.76
1:A:495:GLU:OE1	1:A:538:LYS:HG3	1.85	0.76
1:A:540:THR:HA	1:A:543:GLN:NE2	2.00	0.76
1:A:490:ALA:HA	2:H:143:THR:HG21	1.67	0.76
1:A:222:ARG:HH11	1:A:222:ARG:HG3	1.50	0.76
2:H:24:VAL:HG13	2:H:76:ASN:HD21	1.51	0.75
1:A:48:GLU:HA	1:A:48:GLU:OE1	1.88	0.74
3:L:2:ILE:HD13	3:L:27:PRO:HG2	1.70	0.73
2:H:82:MET:HB3	2:H:85:LEU:HD21	1.69	0.73
2:H:90:THR:HG23	2:H:118:THR:HA	1.70	0.73
1:A:344:VAL:HG21	1:A:450:GLU:HG2	1.71	0.72
1:A:408:LEU:HD11	1:A:526:GLN:HB3	1.70	0.72
1:A:262:LYS:O	1:A:266:GLU:HG3	1.89	0.72
1:A:384:PRO:O	1:A:388:ILE:HG12	1.90	0.72
3:L:38:GLN:HB2	3:L:48:LEU:HD11	1.72	0.71
1:A:496:THR:HG22	1:A:497:TYR:N	2.03	0.71
3:L:98:ASP:O	3:L:100:THR:HG23	1.90	0.71
1:A:419:SER:OG	1:A:421:PRO:HD2	1.91	0.70
3:L:80:GLN:HB3	3:L:82:GLU:OE2	1.91	0.70
1:A:440:HIS:CD2	1:A:444:LYS:HD2	2.26	0.70
1:A:502:PHE:HA	1:A:535:HIS:HD2	1.56	0.70
3:L:12:SER:CB	3:L:110:LYS:HE3	2.21	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:430:LEU:O	1:A:433:VAL:HG12	1.91	0.69
1:A:198:LEU:HD13	1:A:458:ASN:HB2	1.73	0.69
1:A:109:ASN:ND2	1:A:463:LEU:HD21	2.08	0.68
2:H:66:ARG:HH11	2:H:66:ARG:HG3	1.59	0.68
3:L:2:ILE:O	3:L:27:PRO:HD3	1.95	0.67
1:A:209:ARG:NH1	1:A:209:ARG:HG2	2.10	0.67
3:L:95:SER:HB2	3:L:98:ASP:HB2	1.77	0.67
3:L:36:TRP:CE2	3:L:74:LEU:HB2	2.30	0.66
2:H:2:VAL:CG2	2:H:27:ILE:HG23	2.25	0.66
2:H:174:PHE:CE2	3:L:179:SER:HB3	2.32	0.65
2:H:150:VAL:HB	2:H:186:LEU:HD13	1.78	0.65
3:L:34:LEU:HG	3:L:72:PHE:CG	2.31	0.65
1:A:303:PRO:O	1:A:337:ARG:NH1	2.30	0.65
1:A:502:PHE:HA	1:A:535:HIS:CD2	2.30	0.65
1:A:398:LEU:O	1:A:402:LYS:HB2	1.96	0.65
1:A:90:CYS:O	1:A:98:ARG:HG3	1.96	0.65
1:A:420:THR:O	1:A:424:VAL:HG23	1.97	0.65
2:H:167:LEU:HD21	2:H:190:VAL:HG21	1.78	0.65
3:L:191:LYS:HG3	3:L:191:LYS:O	1.96	0.65
1:A:356:THR:HG21	1:A:376:GLU:OE2	1.97	0.64
2:H:2:VAL:HG22	2:H:27:ILE:HG23	1.79	0.64
3:L:122:PRO:HB3	3:L:212:PHE:CE1	2.33	0.64
1:A:25:ILE:HD13	1:A:154:LEU:HD23	1.78	0.64
2:H:193:PRO:HG2	2:H:196:SER:HB3	1.80	0.64
2:H:132:LEU:HB3	3:L:121:PHE:CD2	2.33	0.64
2:H:24:VAL:HG11	2:H:29:LEU:HB2	1.78	0.64
3:L:122:PRO:HB3	3:L:212:PHE:CZ	2.32	0.64
1:A:159:LYS:NZ	1:A:284:LEU:HB2	2.13	0.63
1:A:215:ALA:HB3	1:A:235:VAL:HG13	1.80	0.63
1:A:501:GLU:HG3	1:A:502:PHE:H	1.63	0.63
1:A:327:LEU:HD11	1:A:354:GLU:HG3	1.80	0.63
3:L:123:PRO:HD3	3:L:135:VAL:HG22	1.81	0.63
1:A:302:LEU:HD13	1:A:337:ARG:HG2	1.80	0.63
1:A:412:THR:OG1	1:A:423:LEU:HD13	1.98	0.63
3:L:164:GLU:HB3	3:L:180:SER:HA	1.79	0.63
1:A:218:ARG:NH1	1:A:222:ARG:HH12	1.97	0.63
2:H:222:LYS:O	2:H:223:SER:HB2	1.98	0.63
2:H:171:VAL:HG22	2:H:190:VAL:HB	1.81	0.62
3:L:8:PRO:O	3:L:105:THR:HG23	1.99	0.62
3:L:96:ILE:C	3:L:98:ASP:H	2.00	0.62
1:A:225:LYS:HG2	1:A:299:PRO:HG3	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:34:LEU:HG	3:L:72:PHE:CD2	2.35	0.62
3:L:54:LYS:HB3	3:L:54:LYS:NZ	2.14	0.62
1:A:36:PHE:O	1:A:40:VAL:HG23	1.99	0.62
1:A:151:ALA:HB2	1:A:250:LEU:HD22	1.82	0.62
1:A:504:ALA:O	1:A:508:THR:HG23	1.99	0.61
1:A:222:ARG:HG3	1:A:222:ARG:NH1	2.13	0.61
1:A:378:LYS:HE3	1:A:382:GLU:OE2	1.99	0.61
2:H:24:VAL:HG12	2:H:29:LEU:HD13	1.82	0.61
1:A:485:ARG:HB3	1:A:486:PRO:CD	2.28	0.61
2:H:150:VAL:HB	2:H:186:LEU:CD1	2.30	0.61
1:A:224:PRO:HD2	1:A:296:ASP:HB3	1.82	0.61
3:L:30:TRP:CG	3:L:31:SER:N	2.67	0.61
1:A:409:VAL:HG22	1:A:529:LEU:HD21	1.82	0.60
1:A:159:LYS:HZ2	1:A:284:LEU:HB2	1.67	0.60
1:A:218:ARG:NH1	1:A:222:ARG:NH1	2.49	0.60
3:L:192:HIS:O	3:L:214:ARG:HD3	2.00	0.60
2:H:144:ALA:CB	3:L:119:PHE:HE2	2.15	0.60
1:A:223:PHE:HB3	1:A:271:ILE:O	2.01	0.60
1:A:384:PRO:HB2	1:A:446:MET:CE	2.32	0.60
2:H:59:TYR:O	3:L:97:SER:HB2	2.02	0.60
2:H:124:THR:HG23	2:H:124:THR:O	2.02	0.60
2:H:82:MET:CB	2:H:85:LEU:HD21	2.31	0.59
2:H:1:GLU:HG3	2:H:2:VAL:HG23	1.85	0.59
2:H:66:ARG:HD2	2:H:84:SER:HB2	1.85	0.59
1:A:209:ARG:HH11	1:A:209:ARG:HG2	1.68	0.58
1:A:492:GLU:OE1	2:H:143:THR:N	2.36	0.58
2:H:156:GLU:OE1	2:H:157:PRO:HA	2.02	0.58
1:A:221:GLN:HE21	1:A:339:PRO:HA	1.68	0.58
2:H:2:VAL:CG1	2:H:27:ILE:HG12	2.32	0.58
2:H:72:ASP:OD1	2:H:75:LYS:HG3	2.03	0.58
1:A:495:GLU:CB	3:L:115:ALA:HB3	2.33	0.58
3:L:135:VAL:HG12	3:L:151:TRP:CH2	2.39	0.58
1:A:5:SER:HB3	1:A:57:GLU:OE1	2.03	0.58
3:L:18:ARG:HG3	3:L:77:SER:HA	1.83	0.58
1:A:490:ALA:O	1:A:491:LEU:O	2.21	0.58
1:A:567:CYS:O	1:A:571:GLU:HB2	2.04	0.58
2:H:6:GLU:H	2:H:113:GLN:NE2	2.01	0.58
3:L:173:ASP:O	3:L:175:THR:HG23	2.05	0.57
1:A:223:PHE:CD1	1:A:272:SER:HB2	2.40	0.57
1:A:420:THR:HG23	1:A:530:VAL:HB	1.86	0.57
2:H:87:ALA:HB3	2:H:88:GLU:OE2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:525:LYS:HD2	1:A:548:MET:HE1	1.87	0.56
2:H:174:PHE:CD1	3:L:167:THR:HG23	2.40	0.56
2:H:24:VAL:CG1	2:H:29:LEU:HD13	2.36	0.56
2:H:68:THR:HB	2:H:81:GLN:HB3	1.87	0.56
2:H:94:TYR:CD1	2:H:114:GLY:HA3	2.41	0.56
3:L:87:TYR:HE1	3:L:107:VAL:HG21	1.70	0.56
2:H:134:PRO:HG2	2:H:221:PRO:HB3	1.86	0.56
2:H:6:GLU:OE2	2:H:95:CYS:HB3	2.06	0.56
1:A:168:CYS:SG	1:A:177:CYS:C	2.84	0.56
3:L:2:ILE:O	3:L:26:SER:HB3	2.05	0.56
3:L:154:ASP:OD2	3:L:192:HIS:HB3	2.06	0.56
2:H:90:THR:HA	2:H:117:VAL:O	2.06	0.56
1:A:577:ALA:HA	1:A:580:GLN:HB3	1.89	0.55
3:L:214:ARG:HH11	3:L:214:ARG:HG3	1.69	0.55
3:L:87:TYR:O	3:L:104:GLY:HA2	2.06	0.55
1:A:47:THR:O	1:A:51:LYS:HG3	2.06	0.55
1:A:458:ASN:HA	1:A:484:ARG:NH1	2.22	0.55
1:A:222:ARG:HD3	1:A:293:VAL:O	2.07	0.55
2:H:146:LEU:C	2:H:146:LEU:HD12	2.26	0.55
2:H:162:TRP:CH2	2:H:204:CYS:HB3	2.42	0.55
2:H:147:GLY:HA3	2:H:189:VAL:HA	1.89	0.55
1:A:564:LYS:O	1:A:565:GLU:HB2	2.07	0.54
1:A:29:GLN:HG2	1:A:147:PRO:HA	1.89	0.54
2:H:11:LEU:HD22	2:H:155:PRO:HD3	1.89	0.54
3:L:34:LEU:HD13	3:L:34:LEU:C	2.26	0.54
3:L:109:ILE:HG22	3:L:110:LYS:N	2.22	0.54
1:A:64:LYS:HG3	1:A:69:LEU:HG	1.90	0.54
2:H:50:ILE:HG12	2:H:51:ILE:N	2.22	0.54
1:A:112:LEU:HD21	1:A:144:ARG:NH2	2.22	0.54
1:A:265:CYS:SG	1:A:286:LYS:HD2	2.48	0.54
2:H:6:GLU:O	2:H:7:SER:HB3	2.08	0.54
2:H:28:ASP:OD1	2:H:30:SER:HB2	2.08	0.53
3:L:25:SER:HB2	3:L:93:TYR:OH	2.09	0.53
2:H:12:VAL:HB	2:H:119:VAL:HG22	1.91	0.53
1:A:50:ALA:O	1:A:54:VAL:HG23	2.09	0.53
1:A:561:ALA:C	1:A:563:ASP:N	2.62	0.53
1:A:387:LEU:HD13	2:H:195:SER:HB3	1.91	0.53
2:H:11:LEU:HD12	2:H:118:THR:O	2.08	0.53
1:A:283:LEU:HD23	1:A:284:LEU:N	2.24	0.52
2:H:127:PRO:CB	2:H:153:TYR:HB3	2.38	0.52
2:H:143:THR:CB	2:H:193:PRO:HA	2.38	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:ASP:HB3	1:A:252:GLU:CD	2.29	0.52
1:A:495:GLU:HB2	3:L:115:ALA:HB3	1.90	0.52
1:A:525:LYS:HD2	1:A:548:MET:CE	2.39	0.52
1:A:60:GLU:O	1:A:61:ASN:HB2	2.09	0.52
1:A:81:ARG:HA	1:A:85:GLY:H	1.73	0.52
3:L:60:PRO:HB3	3:L:62:ARG:NH1	2.24	0.52
2:H:144:ALA:HB1	3:L:119:PHE:CE2	2.35	0.52
1:A:383:GLU:HG3	1:A:384:PRO:N	2.24	0.52
1:A:575:LEU:HD12	1:A:578:ALA:CB	2.39	0.52
3:L:125:ASP:O	3:L:129:LYS:HG3	2.10	0.52
2:H:66:ARG:NH1	2:H:66:ARG:HG3	2.24	0.52
3:L:164:GLU:HB2	3:L:178:LEU:HD21	1.92	0.52
1:A:111:ASN:N	1:A:111:ASN:OD1	2.42	0.52
1:A:318:ASN:HB3	1:A:326:PHE:CD1	2.45	0.52
1:A:70:PHE:HZ	1:A:251:LEU:HD11	1.75	0.52
1:A:156:PHE:CE2	1:A:288:HIS:CG	2.97	0.51
3:L:135:VAL:HG12	3:L:151:TRP:HH2	1.75	0.51
1:A:510:HIS:O	1:A:513:ILE:HG22	2.11	0.51
1:A:537:PRO:O	1:A:539:ALA:N	2.44	0.51
1:A:274:LYS:NZ	1:A:297:GLU:OE2	2.39	0.51
2:H:111:TRP:CE3	3:L:45:PRO:HD2	2.45	0.51
3:L:178:LEU:HD23	3:L:178:LEU:C	2.30	0.51
3:L:93:TYR:C	3:L:95:SER:H	2.14	0.51
1:A:56:ASP:OD2	1:A:59:ALA:HB2	2.11	0.51
3:L:3:GLN:N	3:L:3:GLN:OE1	2.44	0.51
1:A:21:ALA:O	1:A:25:ILE:HG13	2.10	0.51
2:H:174:PHE:CD2	3:L:179:SER:HB3	2.46	0.51
1:A:405:ASN:HA	1:A:408:LEU:HD12	1.91	0.51
1:A:425:GLU:OE1	1:A:425:GLU:HA	2.11	0.51
1:A:505:GLU:O	1:A:505:GLU:HG2	2.11	0.51
1:A:417:GLN:CD	1:A:417:GLN:H	2.15	0.50
1:A:500:LYS:O	1:A:501:GLU:HB3	2.11	0.50
2:H:82:MET:HE1	2:H:117:VAL:HG11	1.93	0.50
1:A:39:HIS:O	1:A:43:VAL:HG23	2.11	0.50
2:H:179:GLN:C	2:H:181:SER:H	2.15	0.50
2:H:94:TYR:CE1	2:H:114:GLY:HA3	2.46	0.50
1:A:401:TYR:OH	1:A:525:LYS:HE3	2.10	0.50
1:A:194:ALA:HB1	1:A:455:VAL:CG1	2.41	0.50
1:A:496:THR:CG2	1:A:497:TYR:N	2.71	0.50
1:A:576:VAL:O	1:A:580:GLN:HB2	2.11	0.50
2:H:104:THR:HB	3:L:96:ILE:HG13	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:174:PHE:HE1	3:L:177:SER:O	1.95	0.50
3:L:152:LYS:HZ1	3:L:198:GLU:CD	2.15	0.50
3:L:2:ILE:HG22	3:L:3:GLN:N	2.26	0.50
2:H:107:TYR:CD1	3:L:35:SER:HB3	2.47	0.50
1:A:197:ARG:HG3	1:A:197:ARG:HH11	1.77	0.50
1:A:490:ALA:HA	2:H:143:THR:CG2	2.37	0.50
2:H:12:VAL:HG11	2:H:85:LEU:CD1	2.36	0.50
3:L:14:SER:HA	3:L:110:LYS:HB2	1.92	0.50
1:A:441:PRO:C	1:A:443:ALA:H	2.15	0.50
1:A:408:LEU:HD11	1:A:526:GLN:CB	2.40	0.49
1:A:420:THR:HB	1:A:421:PRO:HD3	1.93	0.49
2:H:5:LEU:HG	2:H:113:GLN:OE1	2.13	0.49
1:A:383:GLU:CG	1:A:384:PRO:HD3	2.43	0.49
1:A:571:GLU:HG3	1:A:575:LEU:HD22	1.94	0.49
1:A:268:GLN:C	1:A:270:SER:H	2.16	0.49
2:H:127:PRO:HB2	2:H:150:VAL:HG12	1.95	0.49
1:A:135:LEU:O	1:A:138:TYR:HB3	2.13	0.49
2:H:48:ILE:HG12	2:H:62:TRP:HH2	1.77	0.49
2:H:203:ILE:HG22	2:H:205:ASN:ND2	2.28	0.49
3:L:120:ILE:CD1	3:L:137:CYS:HB2	2.43	0.48
3:L:204:LEU:HD13	3:L:208:VAL:HG22	1.94	0.48
1:A:32:GLN:NE2	1:A:110:PRO:HG3	2.28	0.48
2:H:18:LEU:HB3	2:H:82:MET:HE2	1.95	0.48
3:L:189:TYR:HA	3:L:195:TYR:OH	2.14	0.48
3:L:142:PHE:N	3:L:175:THR:HB	2.28	0.48
1:A:209:ARG:CG	1:A:209:ARG:HH11	2.25	0.48
1:A:302:LEU:HD13	1:A:337:ARG:CG	2.42	0.48
3:L:95:SER:O	3:L:97:SER:N	2.46	0.48
1:A:25:ILE:HG22	1:A:29:GLN:NE2	2.28	0.48
1:A:127:PHE:CE1	1:A:131:GLU:HG3	2.49	0.48
2:H:28:ASP:O	2:H:29:LEU:HB3	2.14	0.48
3:L:6:GLN:HG3	3:L:89:CYS:SG	2.54	0.48
3:L:96:ILE:C	3:L:98:ASP:N	2.67	0.48
1:A:22:LEU:HD21	1:A:155:LEU:HD11	1.95	0.47
1:A:283:LEU:HD23	1:A:284:LEU:H	1.79	0.47
1:A:495:GLU:HB3	3:L:115:ALA:HB3	1.95	0.47
1:A:464:HIS:CG	1:A:473:VAL:HG11	2.49	0.47
1:A:319:TYR:OH	1:A:358:GLU:HG2	2.14	0.47
1:A:538:LYS:HE3	3:L:204:LEU:HD22	1.96	0.47
2:H:173:THR:HA	2:H:188:SER:HA	1.95	0.47
1:A:223:PHE:HD1	1:A:272:SER:HB2	1.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:GLU:OE1	1:A:99:ASN:HB2	2.14	0.47
3:L:32:ASN:O	3:L:34:LEU:N	2.48	0.47
1:A:314:ASP:HB3	1:A:318:ASN:ND2	2.30	0.47
1:A:364:ALA:O	1:A:366:PRO:HD3	2.14	0.47
1:A:167:GLU:HG2	1:A:167:GLU:O	2.14	0.47
1:A:249:ASP:HB3	1:A:252:GLU:CG	2.45	0.47
2:H:113:GLN:HB2	2:H:113:GLN:HE21	1.49	0.47
2:H:67:PHE:HA	2:H:81:GLN:O	2.14	0.47
1:A:319:TYR:CE1	1:A:323:LYS:HD3	2.50	0.47
1:A:420:THR:HG23	1:A:530:VAL:CG1	2.45	0.47
1:A:563:ASP:OD1	1:A:564:LYS:N	2.48	0.47
1:A:441:PRO:O	1:A:443:ALA:N	2.48	0.47
1:A:71:GLY:HA3	1:A:98:ARG:NH2	2.30	0.47
1:A:570:GLU:C	1:A:572:GLY:H	2.18	0.47
2:H:143:THR:CG2	2:H:193:PRO:HA	2.44	0.47
3:L:44:ALA:HB1	3:L:45:PRO:CD	2.45	0.47
1:A:474:THR:O	1:A:478:THR:OG1	2.25	0.46
2:H:48:ILE:HG12	2:H:62:TRP:CH2	2.50	0.46
1:A:384:PRO:HB2	1:A:446:MET:HE1	1.96	0.46
3:L:182:LEU:HD21	3:L:184:LEU:HD21	1.97	0.46
1:A:120:VAL:HG21	1:A:175:ALA:CA	2.39	0.46
1:A:183:ASP:O	1:A:186:ARG:HB3	2.16	0.46
2:H:127:PRO:HD2	2:H:213:THR:HG21	1.98	0.46
1:A:206:PHE:CE2	1:A:481:LEU:HB2	2.51	0.46
1:A:571:GLU:HB3	1:A:575:LEU:CD2	2.46	0.46
1:A:76:THR:OG1	1:A:77:VAL:HG23	2.15	0.46
1:A:159:LYS:HD3	1:A:285:GLU:OE2	2.15	0.46
3:L:111:ARG:O	3:L:112:THR:C	2.53	0.46
3:L:95:SER:HB2	3:L:98:ASP:CB	2.44	0.46
1:A:24:LEU:HA	1:A:43:VAL:HG21	1.96	0.46
1:A:490:ALA:O	1:A:491:LEU:C	2.54	0.46
1:A:34:CYS:SG	1:A:84:TYR:CE1	3.08	0.46
2:H:36:TRP:HD1	2:H:69:ILE:CD1	2.29	0.46
2:H:6:GLU:OE2	2:H:95:CYS:N	2.48	0.46
3:L:85:ALA:H	3:L:107:VAL:HG23	1.79	0.46
1:A:420:THR:HG23	1:A:530:VAL:CB	2.46	0.46
2:H:186:LEU:C	2:H:186:LEU:HD22	2.37	0.46
3:L:118:VAL:HB	3:L:210:LYS:HG3	1.97	0.46
1:A:11:PHE:CD1	1:A:54:VAL:HG21	2.51	0.45
2:H:179:GLN:C	2:H:181:SER:N	2.68	0.45
1:A:156:PHE:CE2	1:A:160:ARG:NE	2.84	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:492:GLU:HG3	1:A:493:VAL:H	1.81	0.45
3:L:179:SER:O	3:L:179:SER:OG	2.33	0.45
1:A:106:LYS:HD2	1:A:148:TYR:HE1	1.66	0.45
1:A:306:ALA:O	1:A:311:GLU:HG3	2.17	0.45
3:L:50:TYR:OH	3:L:54:LYS:NZ	2.38	0.45
2:H:127:PRO:HB3	2:H:153:TYR:CB	2.39	0.45
1:A:457:LEU:CD1	1:A:485:ARG:HA	2.46	0.45
2:H:189:VAL:HG21	3:L:138:LEU:HD22	1.98	0.45
2:H:51:ILE:HG23	2:H:51:ILE:O	2.17	0.45
3:L:6:GLN:OE1	3:L:104:GLY:HA2	2.15	0.45
3:L:50:TYR:HD2	3:L:51:GLU:HG2	1.80	0.45
1:A:30:TYR:O	1:A:87:MET:HE3	2.16	0.45
3:L:120:ILE:HD12	3:L:137:CYS:HA	1.97	0.45
3:L:48:LEU:C	3:L:49:ILE:HG13	2.37	0.45
1:A:106:LYS:HD3	1:A:147:PRO:HB2	1.98	0.45
1:A:537:PRO:O	1:A:538:LYS:C	2.55	0.45
1:A:496:THR:O	1:A:497:TYR:O	2.35	0.45
1:A:260:LEU:O	1:A:260:LEU:HD12	2.16	0.45
1:A:496:THR:HG22	1:A:497:TYR:H	1.81	0.45
2:H:164:SER:HA	2:H:205:ASN:OD1	2.17	0.45
1:A:221:GLN:NE2	1:A:339:PRO:HA	2.31	0.44
1:A:433:VAL:HG11	1:A:453:LEU:HD21	1.99	0.44
2:H:143:THR:HG22	2:H:194:SER:H	1.82	0.44
2:H:24:VAL:HG11	2:H:29:LEU:CB	2.45	0.44
3:L:16:GLY:O	3:L:78:SER:HA	2.17	0.44
1:A:381:VAL:O	1:A:385:GLN:HG3	2.17	0.44
3:L:111:ARG:O	3:L:112:THR:O	2.35	0.44
3:L:18:ARG:HG3	3:L:76:ILE:O	2.16	0.44
2:H:102:TYR:N	2:H:102:TYR:CD1	2.84	0.44
3:L:120:ILE:HD11	3:L:151:TRP:CH2	2.52	0.44
3:L:120:ILE:HD13	3:L:137:CYS:SG	2.57	0.44
2:H:97:ARG:NH2	2:H:109:ASP:OD2	2.48	0.44
2:H:20:LEU:HD12	2:H:80:LEU:HD23	1.98	0.44
1:A:48:GLU:O	1:A:51:LYS:N	2.50	0.44
3:L:136:VAL:HG22	3:L:181:THR:CG2	2.35	0.44
1:A:225:LYS:CG	1:A:299:PRO:HG3	2.46	0.44
1:A:94:GLN:O	1:A:97:GLU:HB2	2.17	0.44
3:L:120:ILE:O	3:L:120:ILE:CG2	2.66	0.44
1:A:378:LYS:HE2	1:A:378:LYS:HB3	1.70	0.44
1:A:538:LYS:HE3	3:L:204:LEU:CD2	2.47	0.44
1:A:71:GLY:HA3	1:A:98:ARG:CZ	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:416:PRO:HB2	1:A:497:TYR:CD1	2.52	0.44
2:H:20:LEU:HD22	2:H:115:THR:HG21	2.00	0.44
3:L:156:ALA:O	3:L:158:GLN:NE2	2.50	0.44
1:A:118:PRO:HG2	1:A:123:MET:CG	2.47	0.44
1:A:414:LYS:O	1:A:472:ARG:NH1	2.51	0.44
1:A:393:GLU:O	1:A:397:GLN:HG3	2.17	0.43
2:H:97:ARG:O	2:H:108:PHE:HA	2.18	0.43
1:A:173:ASP:HB3	1:A:176:ALA:HB3	2.00	0.43
1:A:300:ALA:O	1:A:302:LEU:N	2.52	0.43
2:H:22:CYS:O	2:H:77:THR:HG23	2.18	0.43
3:L:18:ARG:HA	3:L:76:ILE:O	2.18	0.43
1:A:356:THR:HG21	1:A:376:GLU:CD	2.39	0.43
1:A:197:ARG:HG3	1:A:197:ARG:NH1	2.33	0.43
1:A:233:LYS:NZ	1:A:237:ASP:OD2	2.46	0.43
2:H:36:TRP:CE2	2:H:80:LEU:HB2	2.54	0.43
1:A:500:LYS:O	1:A:501:GLU:CB	2.66	0.43
2:H:36:TRP:CD1	2:H:69:ILE:HG12	2.54	0.43
1:A:151:ALA:CB	1:A:250:LEU:HD22	2.46	0.43
1:A:571:GLU:HB3	1:A:575:LEU:HD23	2.00	0.43
3:L:26:SER:HB3	3:L:27:PRO:CD	2.39	0.43
3:L:67:GLY:HA3	3:L:72:PHE:HA	2.01	0.43
2:H:24:VAL:O	2:H:76:ASN:ND2	2.52	0.43
3:L:150:GLN:HB2	3:L:198:GLU:HB3	1.99	0.43
3:L:26:SER:CB	3:L:27:PRO:HD3	2.38	0.43
3:L:6:GLN:HE21	3:L:21:ILE:CG2	2.32	0.43
1:A:108:ASP:O	1:A:110:PRO:HD3	2.19	0.43
1:A:485:ARG:NH2	1:A:486:PRO:HG3	2.34	0.43
1:A:30:TYR:HE1	1:A:103:LEU:HD23	1.84	0.43
2:H:24:VAL:CG1	2:H:76:ASN:HD21	2.25	0.42
3:L:87:TYR:HE1	3:L:107:VAL:CG2	2.31	0.42
1:A:496:THR:O	1:A:497:TYR:C	2.58	0.42
2:H:4:LEU:HD23	2:H:24:VAL:HB	2.01	0.42
3:L:172:LYS:HG2	3:L:172:LYS:HZ2	1.78	0.42
1:A:165:PHE:CE1	1:A:178:LEU:HD21	2.54	0.42
1:A:495:GLU:OE1	1:A:538:LYS:CG	2.64	0.42
2:H:5:LEU:HA	2:H:113:GLN:OE1	2.19	0.42
3:L:107:VAL:O	3:L:107:VAL:HG23	2.19	0.42
3:L:213:ASN:O	3:L:214:ARG:C	2.58	0.42
3:L:2:ILE:O	3:L:27:PRO:CD	2.67	0.42
1:A:118:PRO:HG2	1:A:123:MET:HG3	2.02	0.42
1:A:141:GLU:OE1	1:A:145:ARG:NH1	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:LEU:CD2	1:A:155:LEU:HD11	2.50	0.42
1:A:342:SER:HA	1:A:447:PRO:HA	2.02	0.42
1:A:457:LEU:O	1:A:460:LEU:HB3	2.19	0.42
1:A:537:PRO:HB2	1:A:538:LYS:H	1.68	0.42
3:L:109:ILE:CG2	3:L:110:LYS:N	2.83	0.42
3:L:93:TYR:C	3:L:95:SER:N	2.73	0.42
1:A:340:ASP:O	1:A:447:PRO:HD3	2.20	0.42
2:H:147:GLY:HA3	2:H:189:VAL:HG12	2.02	0.42
2:H:19:ARG:HG3	2:H:81:GLN:OE1	2.20	0.42
2:H:12:VAL:HG21	2:H:18:LEU:HD22	2.02	0.41
2:H:145:ALA:HB2	2:H:191:THR:HA	2.02	0.41
1:A:133:THR:O	1:A:137:LYS:HB2	2.20	0.41
2:H:172:HIS:CD2	3:L:140:ASN:OD1	2.73	0.41
2:H:175:PRO:HG2	3:L:166:VAL:O	2.20	0.41
2:H:85:LEU:HA	2:H:89:ASP:OD2	2.21	0.41
3:L:56:THR:HG22	3:L:57:SER:N	2.35	0.41
1:A:9:HIS:HE1	1:A:10:ARG:NH1	2.18	0.41
2:H:174:PHE:CE1	3:L:167:THR:HG23	2.54	0.41
1:A:384:PRO:HB2	1:A:446:MET:HE2	2.01	0.41
2:H:129:VAL:HG21	2:H:215:VAL:HB	2.02	0.41
2:H:153:TYR:OH	2:H:176:ALA:HB2	2.19	0.41
1:A:219:LEU:HD22	1:A:223:PHE:CE2	2.56	0.41
1:A:551:PHE:HA	1:A:551:PHE:HD1	1.77	0.41
2:H:85:LEU:HB3	2:H:119:VAL:HG21	2.01	0.41
2:H:35:ASN:OD1	2:H:50:ILE:HB	2.20	0.41
3:L:35:SER:OG	3:L:37:TYR:HE1	2.04	0.41
3:L:15:VAL:CG1	3:L:81:PRO:HD3	2.50	0.41
1:A:42:LEU:HG	1:A:42:LEU:H	1.64	0.41
3:L:35:SER:O	3:L:89:CYS:HA	2.21	0.41
1:A:25:ILE:HD11	1:A:139:LEU:HD22	2.02	0.41
2:H:222:LYS:O	2:H:223:SER:CB	2.68	0.41
2:H:97:ARG:CZ	2:H:110:LEU:HD12	2.51	0.41
1:A:498:VAL:HG13	1:A:498:VAL:O	2.21	0.41
1:A:561:ALA:C	1:A:563:ASP:H	2.21	0.41
2:H:14:PRO:HD3	2:H:120:SER:C	2.41	0.41
3:L:166:VAL:CG2	3:L:178:LEU:HG	2.50	0.41
1:A:132:GLU:OE1	1:A:136:LYS:HE3	2.21	0.41
1:A:140:TYR:O	1:A:144:ARG:HG2	2.21	0.41
1:A:433:VAL:HG21	1:A:453:LEU:HG	2.02	0.41
1:A:543:GLN:HG2	1:A:543:GLN:H	1.63	0.41
2:H:96:ALA:HB1	2:H:108:PHE:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:LYS:HB3	1:A:285:GLU:CB	2.51	0.40
3:L:12:SER:HA	3:L:108:GLU:O	2.21	0.40
3:L:120:ILE:CD1	3:L:137:CYS:CB	2.99	0.40
1:A:70:PHE:HZ	1:A:251:LEU:CD1	2.33	0.40
1:A:458:ASN:CB	1:A:484:ARG:HH12	2.35	0.40
1:A:543:GLN:O	1:A:547:VAL:HG23	2.20	0.40
2:H:146:LEU:C	2:H:146:LEU:CD1	2.88	0.40
1:A:410:ARG:O	1:A:414:LYS:HG3	2.22	0.40
2:H:11:LEU:HB2	2:H:155:PRO:HG3	2.03	0.40
1:A:186:ARG:NH1	1:A:187:ASP:OD1	2.55	0.40
1:A:343:VAL:HG23	1:A:451:ASP:OD2	2.21	0.40
3:L:149:VAL:HG11	3:L:180:SER:CB	2.52	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	577/585 (99%)	503 (87%)	57 (10%)	17 (3%)	6 44
2	H	213/233 (91%)	191 (90%)	20 (9%)	2 (1%)	21 67
3	L	215/217 (99%)	182 (85%)	21 (10%)	12 (6%)	2 26
All	All	1005/1035 (97%)	876 (87%)	98 (10%)	31 (3%)	5 43

All (31) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	300	ALA
1	A	491	LEU
1	A	497	TYR
1	A	538	LYS
2	H	223	SER

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Mol	Chain	Res	Type
3	L	17	ASP
1	A	301	ASP
1	A	442	GLU
1	A	479	GLU
1	A	537	PRO
1	A	540	THR
3	L	3	GLN
3	L	30	TRP
3	L	33	PHE
3	L	112	THR
3	L	141	ASN
3	L	214	ARG
1	A	56	ASP
1	A	129	ASP
1	A	269	ASP
1	A	281	LYS
1	A	502	PHE
1	A	512	ASP
3	L	31	SER
3	L	97	SER
3	L	216	GLU
1	A	118	PRO
3	L	94	SER
2	H	221	PRO
3	L	2	ILE
1	A	388	ILE

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	506/511 (99%)	475 (94%)	31 (6%)	23	65
2	H	180/195 (92%)	168 (93%)	12 (7%)	20	62
3	L	190/190 (100%)	175 (92%)	15 (8%)	15	54
All	All	876/896 (98%)	818 (93%)	58 (7%)	21	63

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	GLU
1	A	97	GLU
1	A	104	GLN
1	A	111	ASN
1	A	114	ARG
1	A	117	ARG
1	A	132	GLU
1	A	144	ARG
1	A	218	ARG
1	A	222	ARG
1	A	269	ASP
1	A	283	LEU
1	A	292	GLU
1	A	299	PRO
1	A	305	LEU
1	A	308	ASP
1	A	324	ASP
1	A	383	GLU
1	A	389	LYS
1	A	435	SER
1	A	466	LYS
1	A	469	VAL
1	A	478	THR
1	A	496	THR
1	A	532	LEU
1	A	535	HIS
1	A	536	LYS
1	A	540	THR
1	A	551	PHE
1	A	557	LYS
1	A	576	VAL
2	H	30	SER
2	H	54	SER
2	H	61	THR
2	H	66	ARG
2	H	71	ARG
2	H	99	VAL
2	H	113	GLN
2	H	156	GLU
2	H	186	LEU
2	H	191	THR
2	H	220	GLU

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Mol	Chain	Res	Type
2	H	222	LYS
3	L	1	ASP
3	L	5	THR
3	L	15	VAL
3	L	68	SER
3	L	82	GLU
3	L	94	SER
3	L	106	LYS
3	L	108	GLU
3	L	111	ARG
3	L	171	SER
3	L	184	LEU
3	L	191	LYS
3	L	198	GLU
3	L	202	GLN
3	L	205	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	HIS
1	A	29	GLN
1	A	61	ASN
1	A	67	HIS
1	A	109	ASN
1	A	221	GLN
1	A	440	HIS
1	A	483	ASN
1	A	543	GLN
2	H	3	GLN
2	H	113	GLN
3	L	155	ASN
3	L	163	GLN
3	L	202	GLN

5.3.3 RNA

There are no RNA molecules 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

There are no ligands in this entry.

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, 95^{th} 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	579/585 (98%)	-0.25	10 (1%) 73 59	75, 111, 185, 211	0
2	H	217/233 (93%)	-0.30	1 (0%) 91 86	81, 103, 131, 151	0
3	L	217/217 (100%)	-0.33	0 100 100	70, 103, 131, 163	0
All	All	1013/1035 (97%)	-0.27	11 (1%) 82 70	70, 107, 177, 211	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	79	THR	5.3
2	H	135	SER	4.1
1	A	83	THR	3.8
1	A	573	LYS	3.8
1	A	26	ALA	2.8
1	A	570	GLU	2.7
1	A	82	GLU	2.6
1	A	86	GLU	2.6
1	A	574	LYS	2.1
1	A	80	LEU	2.1
1	A	29	GLN	2.1

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

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.