



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 23, 2024 – 04:01 AM EDT

PDB ID : 4PLJ
Title : Hepatitis E Virus E2s domain (Genotype IV) in complex with a neutralizing antibody 8G12
Authors : Tang, X.H.; Li, S.W.; Sivaraman, J.
Deposited on : 2014-05-18
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
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)	:	2.37.1

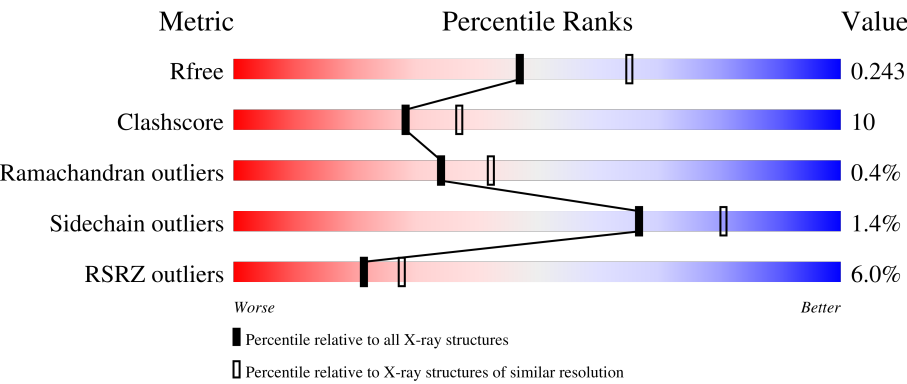
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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

Mol	Chain	Length	Quality of chain
1	A	148	<div><div>0%</div><div>91%</div><div>9%</div></div>
1	B	148	<div><div>5%</div><div>89%</div><div>10%</div><div>.</div></div>
2	C	212	<div><div>15%</div><div>68%</div><div>26%</div><div>5%</div></div>
2	L	212	<div><div>3%</div><div>79%</div><div>19%</div><div>.</div></div>
3	D	229	<div><div>7%</div><div>76%</div><div>17%</div><div>6%</div></div>

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Mol	Chain	Length	Quality of chain
3	H	229	<div><div></div><div>3%</div><div>83%</div><div>12%</div><div>5%</div></div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9128 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 Capsid protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	148	Total	C	N	O	S	0	0	0
			1129	716	189	222	2			
1	B	147	Total	C	N	O	S	0	0	0
			1123	713	188	220	2			

- Molecule 2 is a protein called 8G12 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	212	Total	C	N	O	S	0	0	0
			1632	1018	274	334	6			
2	C	201	Total	C	N	O	S	0	0	0
			1553	975	255	317	6			

- Molecule 3 is a protein called 8G12 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	218	Total	C	N	O	S	0	0	0
			1669	1064	274	324	7			
3	D	216	Total	C	N	O	S	0	0	0
			1658	1057	272	322	7			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	79	Total	O	0	0
			79	79		
4	B	76	Total	O	0	0
			76	76		
4	L	57	Total	O	0	0
			57	57		
4	H	74	Total	O	0	0
			74	74		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	37	Total	O	0	0
			37	37		
4	D	41	Total	O	0	0
			41	41		

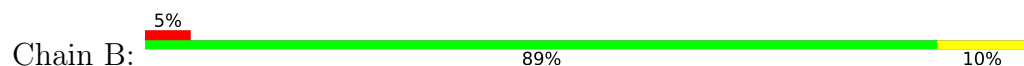
3 Residue-property plots [i](#)

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.

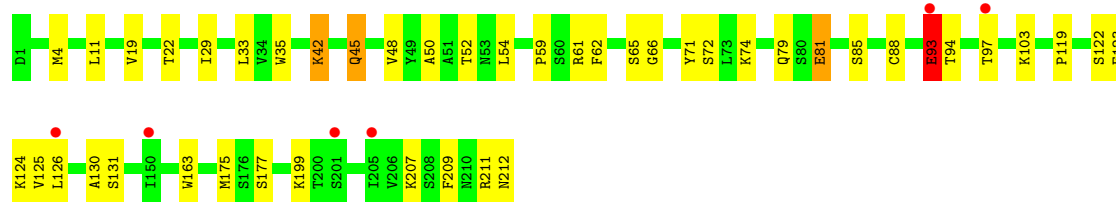
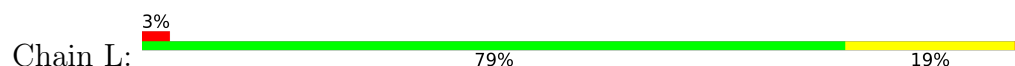
- Molecule 1: Capsid protein



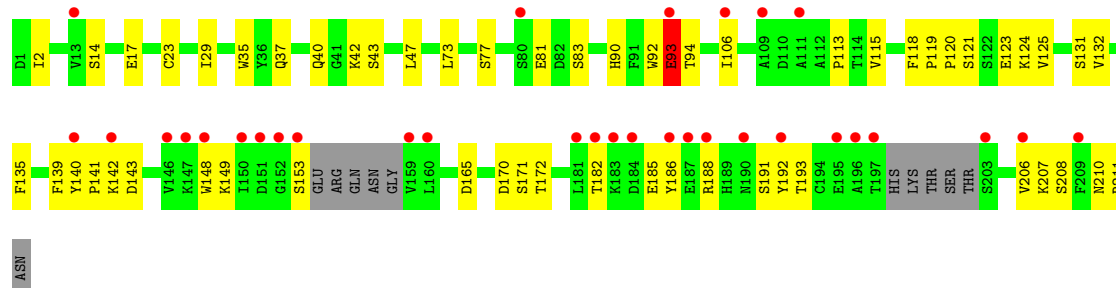
- Molecule 1: Capsid protein




- Molecule 2: 8G12 light chain

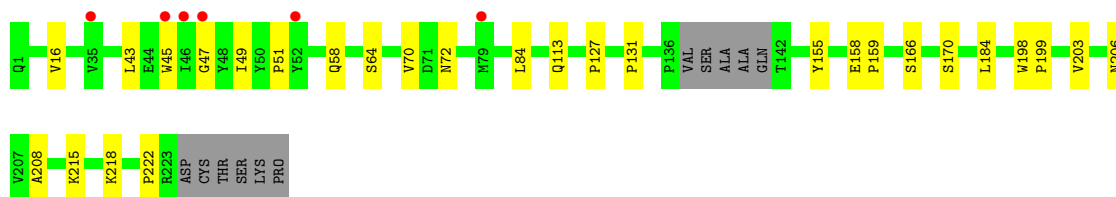


- Molecule 2: 8G12 light chain




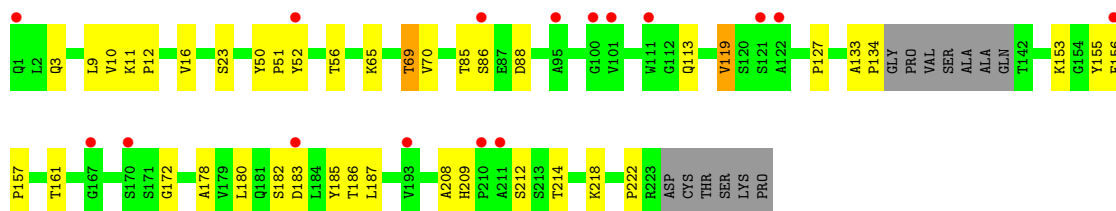
- Molecule 3: 8G12 heavy chain

Chain H:  3% 83% 12% 5%



- Molecule 3: 8G12 heavy chain

Chain D:  7% 76% 17% 6%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.77Å 89.46Å 137.64Å 90.00° 95.90° 90.00°	Depositor
Resolution (Å)	29.14 – 2.30 29.14 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.14-2.30) 97.3 (29.14-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.187 , 0.242 0.194 , 0.243	Depositor DCC
R_{free} test set	2000 reflections (3.27%)	wwPDB-VP
Wilson B-factor (Å ²)	32.8	Xtriage
Anisotropy	0.684	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 48.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9128	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.22% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

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	0.54	1/1157 (0.1%)	0.60	0/1584
1	B	0.54	1/1151 (0.1%)	0.62	1/1577 (0.1%)
2	C	0.49	1/1588 (0.1%)	0.65	3/2156 (0.1%)
2	L	0.44	0/1670	0.59	1/2268 (0.0%)
3	D	0.40	0/1705	0.55	0/2329
3	H	0.46	0/1717	0.56	0/2346
All	All	0.47	3/8988 (0.0%)	0.59	5/12260 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
2	C	0	2
2	L	0	1
All	All	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	93	GLU	C-N	8.99	1.54	1.34
1	B	537	PHE	C-N	-6.95	1.18	1.34
1	A	537	PHE	C-N	-5.75	1.20	1.34

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	92	TRP	C-N-CA	-7.04	104.11	121.70
2	C	93	GLU	CA-C-N	-6.22	103.52	117.20
2	L	93	GLU	C-N-CA	5.84	136.29	121.70
1	B	537	PHE	O-C-N	-5.81	113.41	122.70
2	C	29	ILE	N-CA-C	-5.25	96.82	111.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	93	GLU	Peptide,Mainchain
2	L	93	GLU	Peptide

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	1129	0	1103	22	0
1	B	1123	0	1097	24	0
2	C	1553	0	1485	49	0
2	L	1632	0	1543	56	0
3	D	1658	0	1614	30	0
3	H	1669	0	1624	17	0
4	A	79	0	0	3	0
4	B	76	0	0	8	0
4	C	37	0	0	4	0
4	D	41	0	0	5	0
4	H	74	0	0	4	0
4	L	57	0	0	5	0
All	All	9128	0	8466	168	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 10.

All (168) 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:B:554:LYS:NZ	2:C:93:GLU:HB3	1.45	1.31
4:B:702:HOH:O	2:C:93:GLU:HB2	1.31	1.28
2:L:93:GLU:HG3	2:L:94:THR:O	1.19	1.25
2:L:93:GLU:HA	2:L:93:GLU:OE1	1.59	1.02
2:L:93:GLU:CG	2:L:94:THR:O	2.10	0.99
1:B:554:LYS:HZ2	2:C:93:GLU:HB3	1.19	0.98
1:B:554:LYS:HZ3	2:C:93:GLU:HB3	1.27	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:554:LYS:NZ	2:C:93:GLU:CB	2.30	0.95
1:A:554:LYS:HZ2	2:L:93:GLU:CD	1.71	0.93
1:B:554:LYS:HZ2	2:C:93:GLU:CB	1.84	0.89
3:D:88:ASP:OD1	4:D:301:HOH:O	1.91	0.87
3:H:166:SER:H	3:H:206:ASN:HD21	1.21	0.86
2:L:45:GLN:NE2	4:L:302:HOH:O	2.08	0.84
1:A:554:LYS:HZ2	2:L:93:GLU:CG	1.91	0.83
1:A:588:LEU:HA	2:L:93:GLU:OE2	1.80	0.82
2:L:93:GLU:HG3	2:L:94:THR:C	2.00	0.81
3:H:184:LEU:O	4:H:301:HOH:O	2.01	0.78
2:L:93:GLU:OE1	2:L:93:GLU:CA	2.32	0.78
2:L:66:GLY:O	4:L:301:HOH:O	2.01	0.78
2:C:123:GLU:OE2	4:C:301:HOH:O	2.00	0.78
1:A:554:LYS:NZ	2:L:93:GLU:CB	2.49	0.76
2:C:42:LYS:HD2	2:C:43:SER:H	1.51	0.76
1:A:554:LYS:NZ	2:L:93:GLU:CG	2.51	0.74
1:A:468:ASN:HD22	1:B:472:TRP:HH2	1.33	0.73
1:A:588:LEU:CA	2:L:93:GLU:OE2	2.39	0.70
3:D:85:THR:O	3:D:119:VAL:HG11	1.93	0.69
1:A:554:LYS:HZ1	2:L:93:GLU:HB2	1.57	0.68
3:D:182:SER:OG	3:D:183:ASP:N	2.27	0.68
1:A:554:LYS:NZ	2:L:93:GLU:HB2	2.08	0.67
2:C:77:SER:OG	4:C:302:HOH:O	2.10	0.67
1:A:554:LYS:HZ2	2:L:93:GLU:CB	2.08	0.66
2:L:124:LYS:HE3	2:L:130:ALA:HA	1.76	0.66
2:L:163:TRP:CD1	2:L:175:MET:HG3	2.30	0.66
1:A:472:TRP:HH2	1:B:468:ASN:HD22	1.44	0.66
1:B:554:LYS:HZ3	2:C:93:GLU:CB	2.02	0.64
2:L:123:GLU:OE1	3:H:218:LYS:NZ	2.24	0.64
2:C:210:ASN:O	2:C:211:ARG:HG2	1.98	0.64
1:A:588:LEU:C	2:L:93:GLU:OE2	2.37	0.64
2:L:123:GLU:HA	2:L:126:LEU:HD22	1.80	0.63
3:D:3:GLN:OE1	3:D:113:GLN:NE2	2.32	0.62
1:B:459:SER:OG	4:B:701:HOH:O	2.14	0.61
3:H:58:GLN:NE2	4:H:305:HOH:O	2.33	0.61
3:H:208:ALA:HB2	3:H:215:LYS:HD3	1.81	0.61
1:A:469:ASP:HB2	1:A:601:LEU:HD12	1.83	0.60
3:H:166:SER:H	3:H:206:ASN:ND2	1.97	0.60
2:C:142:LYS:HG3	2:C:143:ASP:H	1.66	0.60
1:B:518:LYS:HE2	4:B:709:HOH:O	2.02	0.59
1:B:549:GLU:OE1	4:B:702:HOH:O	2.17	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:211:ARG:O	4:L:303:HOH:O	2.17	0.59
2:C:120:PRO:HD3	2:C:132:VAL:HG22	1.85	0.59
1:B:554:LYS:CE	2:C:93:GLU:HB3	2.32	0.58
2:C:186:TYR:CE1	2:C:192:TYR:HE2	2.21	0.58
1:B:470:VAL:HG22	1:B:600:VAL:HG22	1.86	0.58
2:C:182:THR:HG23	2:C:185:GLU:H	1.68	0.57
1:A:554:LYS:NZ	2:L:93:GLU:CD	2.53	0.57
3:D:161:THR:HG22	3:D:208:ALA:HB3	1.86	0.56
3:D:178:ALA:HB2	3:D:187:LEU:HD23	1.87	0.56
1:A:459:SER:OG	4:A:701:HOH:O	2.12	0.56
3:D:11:LYS:HG3	3:D:12:PRO:HD2	1.87	0.56
2:L:61:ARG:NE	2:L:79:GLN:HE21	2.04	0.55
2:C:42:LYS:CD	2:C:43:SER:H	2.20	0.55
2:L:124:LYS:NZ	2:L:131:SER:H	2.04	0.55
2:C:37:GLN:HB2	2:C:47:LEU:HD11	1.88	0.55
3:H:127:PRO:HB3	3:H:155:TYR:HB3	1.89	0.54
3:H:72:ASN:ND2	4:H:303:HOH:O	2.23	0.54
1:B:554:LYS:NZ	2:C:93:GLU:CG	2.70	0.53
1:B:554:LYS:HZ2	2:C:93:GLU:CG	2.20	0.53
2:L:212:ASN:HA	4:L:303:HOH:O	2.08	0.53
2:C:148:TRP:O	2:C:149:LYS:HD2	2.09	0.53
3:H:16:VAL:HG12	3:H:84:LEU:HD11	1.91	0.53
2:L:42:LYS:HE2	3:H:113:GLN:O	2.09	0.53
3:D:9:LEU:HB2	3:D:157:PRO:HG3	1.89	0.53
3:D:127:PRO:HB3	3:D:155:TYR:HB3	1.91	0.52
1:A:554:LYS:HD3	2:L:93:GLU:OE1	2.08	0.52
2:L:124:LYS:HE3	2:L:131:SER:H	1.74	0.52
1:B:511:SER:HB3	4:B:735:HOH:O	2.09	0.52
2:L:61:ARG:CZ	2:L:79:GLN:HE21	2.22	0.52
1:B:524:ARG:HH21	1:B:524:ARG:HB3	1.75	0.51
2:L:124:LYS:CE	2:L:131:SER:H	2.24	0.51
2:L:122:SER:O	2:L:126:LEU:HD13	2.11	0.50
2:L:175:MET:HE1	2:L:177:SER:HB2	1.94	0.50
2:L:74:LYS:HE3	4:L:305:HOH:O	2.10	0.50
2:C:186:TYR:O	2:C:192:TYR:OH	2.29	0.50
2:C:193:THR:HG23	2:C:208:SER:HB3	1.94	0.50
1:A:554:LYS:CD	2:L:93:GLU:OE1	2.61	0.49
3:D:153:LYS:HA	3:D:186:THR:HG22	1.92	0.49
3:D:178:ALA:HA	3:D:187:LEU:HB3	1.92	0.49
2:C:113:PRO:HB3	2:C:139:PHE:HB3	1.94	0.49
2:L:124:LYS:HZ1	2:L:131:SER:H	1.59	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:35:TRP:CD2	2:C:73:LEU:HB2	2.48	0.49
2:C:142:LYS:HG3	2:C:143:ASP:N	2.26	0.49
3:D:212:SER:HB3	3:D:214:THR:HG23	1.95	0.49
3:D:10:VAL:HG21	3:D:16:VAL:HB	1.95	0.48
3:H:64:SER:N	4:H:302:HOH:O	2.16	0.48
2:C:149:LYS:HG3	2:C:153:SER:C	2.33	0.48
2:C:186:TYR:HE1	2:C:192:TYR:HE2	1.61	0.48
2:L:4:MET:HG2	2:L:97:THR:HG22	1.96	0.48
2:L:85:SER:OG	2:L:103:LYS:HD3	2.14	0.48
2:L:22:THR:HG22	2:L:72:SER:OG	2.13	0.48
3:D:172:GLY:O	4:D:302:HOH:O	2.20	0.47
4:B:702:HOH:O	2:C:93:GLU:CB	2.14	0.47
2:C:140:TYR:CG	2:C:141:PRO:HA	2.49	0.47
3:D:161:THR:CG2	3:D:208:ALA:HB3	2.44	0.47
2:C:93:GLU:CG	2:C:94:THR:O	2.63	0.47
2:L:59:PRO:HG2	2:L:62:PHE:CE2	2.50	0.47
3:D:86:SER:HA	3:D:119:VAL:HG13	1.96	0.46
3:H:45:TRP:CH2	3:H:47:GLY:HA2	2.51	0.46
2:L:124:LYS:HB3	2:L:124:LYS:HE2	1.54	0.46
2:C:2:ILE:HD12	2:C:90:HIS:CE1	2.50	0.46
2:C:93:GLU:HG3	2:C:94:THR:C	2.36	0.46
2:L:52:THR:HG22	2:L:65:SER:HA	1.97	0.46
1:B:589:GLY:N	2:C:93:GLU:OE1	2.50	0.45
2:C:124:LYS:HE3	2:C:131:SER:OG	2.16	0.45
3:D:23:SER:N	4:D:303:HOH:O	2.48	0.45
2:L:35:TRP:CZ3	2:L:88:CYS:HB3	2.51	0.45
2:C:170:ASP:OD1	2:C:172:THR:HG22	2.17	0.45
2:C:206:VAL:O	2:C:207:LYS:HG2	2.17	0.45
3:D:157:PRO:O	3:D:209:HIS:HE1	1.99	0.45
3:H:199:PRO:CB	3:H:222:PRO:HG3	2.46	0.45
3:D:209:HIS:HD2	3:D:212:SER:CB	2.30	0.45
2:L:122:SER:HA	2:L:125:VAL:HG22	1.99	0.44
2:C:81:GLU:OE1	2:C:81:GLU:N	2.47	0.44
1:A:460:ARG:O	4:A:702:HOH:O	2.21	0.44
3:D:65:LYS:N	4:D:311:HOH:O	2.46	0.44
3:D:69:THR:HB	4:D:318:HOH:O	2.18	0.44
1:B:524:ARG:HB3	1:B:524:ARG:NH2	2.32	0.44
2:L:11:LEU:HD13	2:L:19:VAL:HG23	1.99	0.44
2:C:118:PHE:HA	2:C:119:PRO:HD3	1.88	0.44
2:C:171:SER:HA	4:C:310:HOH:O	2.17	0.44
3:H:49:ILE:O	3:H:51:PRO:HD3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:10:VAL:O	3:D:119:VAL:HA	2.18	0.43
3:D:209:HIS:HB3	3:D:214:THR:OG1	2.18	0.43
1:B:468:ASN:O	4:B:703:HOH:O	2.20	0.43
1:B:468:ASN:HA	1:B:600:VAL:HG13	2.01	0.43
2:L:33:LEU:O	2:L:50:ALA:O	2.37	0.43
2:C:188:ARG:HD2	2:C:188:ARG:HA	1.91	0.43
1:B:534:LYS:HE2	4:B:714:HOH:O	2.19	0.43
2:L:119:PRO:HB3	2:L:209:PHE:CE2	2.54	0.43
2:C:115:VAL:HA	2:C:135:PHE:O	2.18	0.43
3:D:218:LYS:HD3	3:D:218:LYS:HA	1.87	0.43
1:A:461:PRO:HG2	1:A:464:VAL:HG12	2.01	0.42
2:C:93:GLU:HG3	2:C:94:THR:O	2.18	0.42
2:L:11:LEU:HD13	2:L:19:VAL:CG2	2.49	0.42
2:L:74:LYS:HE2	2:L:74:LYS:HB3	1.85	0.42
2:L:81:GLU:H	2:L:81:GLU:CD	2.22	0.42
2:L:42:LYS:HD3	2:L:42:LYS:HA	1.67	0.42
2:L:48:VAL:HG22	2:L:54:LEU:HD12	2.02	0.42
3:H:131:PRO:HD3	3:H:218:LYS:HE2	2.02	0.42
3:D:50:TYR:CZ	3:D:52:TYR:HB3	2.55	0.42
1:A:549:GLU:OE1	2:L:93:GLU:OE1	2.38	0.41
2:C:172:THR:HG23	4:C:312:HOH:O	2.19	0.41
3:D:209:HIS:HD2	3:D:212:SER:OG	2.03	0.41
2:C:35:TRP:CE2	2:C:73:LEU:HB2	2.55	0.41
1:A:534:LYS:HE2	4:A:710:HOH:O	2.20	0.41
2:C:83:SER:HB2	2:C:106:ILE:HG12	2.03	0.41
1:A:554:LYS:HZ1	2:L:93:GLU:CB	2.23	0.41
2:C:14:SER:N	2:C:17:GLU:OE1	2.48	0.41
1:B:554:LYS:NZ	2:C:93:GLU:CD	2.74	0.41
3:D:51:PRO:HB3	3:D:70:VAL:HG21	2.02	0.41
3:D:133:ALA:HA	3:D:134:PRO:HD3	1.93	0.41
2:L:29:ILE:HD11	2:L:71:TYR:CE2	2.56	0.41
2:L:207:LYS:HA	2:L:207:LYS:HD3	1.64	0.40
2:C:121:SER:O	2:C:125:VAL:HG22	2.21	0.40
1:B:548:TRP:CE2	1:B:596:SER:HB2	2.56	0.40
3:H:198:TRP:HD1	3:H:203:VAL:HG23	1.86	0.40
3:D:156:PHE:HA	3:D:157:PRO:HA	1.89	0.40
3:D:180:LEU:HD13	3:D:185:TYR:CE1	2.57	0.40
3:H:158:GLU:HA	3:H:159:PRO:HA	1.89	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	146/148 (99%)	140 (96%)	6 (4%)	0	100	100
1	B	145/148 (98%)	140 (97%)	5 (3%)	0	100	100
2	C	195/212 (92%)	183 (94%)	11 (6%)	1 (0%)	29	35
2	L	210/212 (99%)	200 (95%)	8 (4%)	2 (1%)	15	17
3	D	212/229 (93%)	201 (95%)	10 (5%)	1 (0%)	29	35
3	H	214/229 (93%)	209 (98%)	5 (2%)	0	100	100
All	All	1122/1178 (95%)	1073 (96%)	45 (4%)	4 (0%)	34	42

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	L	93	GLU
2	L	199	LYS
2	C	40	GLN
3	D	222	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	125/125 (100%)	124 (99%)	1 (1%)	81	91
1	B	125/125 (100%)	125 (100%)	0	100	100
2	C	179/190 (94%)	176 (98%)	3 (2%)	60	76

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	L	185/190 (97%)	181 (98%)	4 (2%)	52	69
3	D	189/199 (95%)	186 (98%)	3 (2%)	62	78
3	H	190/199 (96%)	187 (98%)	3 (2%)	62	78
All	All	993/1028 (97%)	979 (99%)	14 (1%)	67	81

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	528	THR
2	L	42	LYS
2	L	45	GLN
2	L	81	GLU
2	L	93	GLU
3	H	43	LEU
3	H	70	VAL
3	H	170	SER
2	C	23	CYS
2	C	165	ASP
2	C	191	SER
3	D	56	THR
3	D	69	THR
3	D	119	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	468	ASN
1	B	468	ASN
2	L	32	ASN
2	L	70	GLN
2	L	79	GLN
3	H	58	GLN
3	H	60	GLN
3	H	143	ASN
3	H	181	GLN
3	H	206	ASN
3	D	209	HIS

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	537:PHE	C	538:VAL	N	1.18

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	148/148 (100%)	0.02	1 (0%) 87 91	29, 42, 60, 78	0
1	B	147/148 (99%)	0.14	7 (4%) 30 37	27, 40, 55, 72	0
2	C	201/212 (94%)	0.78	32 (15%) 1 2	34, 72, 107, 125	0
2	L	212/212 (100%)	0.20	6 (2%) 53 60	27, 55, 75, 88	0
3	D	216/229 (94%)	0.49	16 (7%) 14 19	38, 62, 84, 105	0
3	H	218/229 (95%)	0.01	6 (2%) 53 60	28, 44, 64, 91	0
All	All	1142/1178 (96%)	0.29	68 (5%) 21 28	27, 52, 91, 125	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	197	THR	6.1
2	C	188	ARG	6.1
2	C	196	ALA	5.7
2	L	93	GLU	5.6
2	C	150	ILE	4.7
2	C	153	SER	4.2
2	C	206	VAL	4.2
2	C	93	GLU	4.0
2	C	146	VAL	3.8
2	C	109	ALA	3.6
2	C	186	TYR	3.5
3	D	167	GLY	3.3
2	C	152	GLY	3.3
2	C	80	SER	3.3
1	B	598	VAL	3.3
2	C	13	VAL	3.2
2	C	142	LYS	3.2
1	B	546	SER	3.2
2	C	195	GLU	3.2

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Mol	Chain	Res	Type	RSRZ
3	D	52	TYR	3.1
3	H	45	TRP	3.1
2	C	190	ASN	3.0
2	C	192	TYR	3.0
2	C	184	ASP	3.0
1	B	597	ALA	2.9
2	C	111	ALA	2.9
2	C	160	LEU	2.9
2	C	182	THR	2.8
3	D	193	VAL	2.8
3	D	95	ALA	2.8
2	C	181	LEU	2.7
1	B	548	TRP	2.7
3	D	183	ASP	2.7
2	C	209	PHE	2.6
3	D	210	PRO	2.6
1	B	547	PHE	2.6
1	B	545	LEU	2.6
2	L	126	LEU	2.5
3	H	52	TYR	2.5
3	H	46	ILE	2.5
3	D	211	ALA	2.4
2	C	147	LYS	2.4
2	C	187	GLU	2.4
1	A	597	ALA	2.3
3	H	47	GLY	2.3
2	C	183	LYS	2.3
2	C	148	TRP	2.3
2	C	159	VAL	2.3
3	D	1	GLN	2.2
3	H	79	MET	2.2
3	D	122	ALA	2.2
2	L	205	ILE	2.2
3	D	156	PHE	2.2
1	B	595	ILE	2.2
3	D	111	TRP	2.2
3	H	35	VAL	2.2
2	C	151	ASP	2.2
2	C	203	SER	2.1
3	D	121	SER	2.1
3	D	101	VAL	2.1
2	C	140	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
3	D	100	GLY	2.1
3	D	170	SER	2.1
2	L	150	ILE	2.0
2	L	201	SER	2.0
2	C	106	ILE	2.0
2	L	97	THR	2.0
3	D	86	SER	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.