



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 01:12 PM GMT

PDB ID : 3MLT
Title : Crystal structure of anti-HIV-1 V3 Fab 2557 in complex with a UG1033 V3 peptide
Authors : Kong, X.-P.
Deposited on : 2010-04-18
Resolution : 2.49 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>

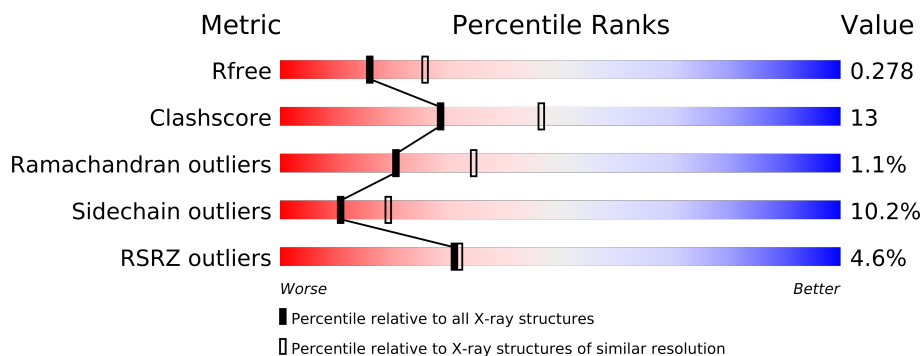
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.49 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}	66092	2784 (2.50-2.50)
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (2.50-2.50)

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	219	
1	D	219	
1	G	219	
1	L	219	
2	B	226	
2	E	226	
2	H	226	
2	I	226	
3	C	23	
3	P	23	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13628 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 Human monoclonal anti-HIV-1 gp120 V3 antibody 2557 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	216	Total	C	N	O	S	0	0	0
			1623	1019	266	333	5			
1	A	215	Total	C	N	O	S	0	0	0
			1619	1017	265	332	5			
1	D	213	Total	C	N	O	S	0	0	0
			1608	1011	263	329	5			
1	G	213	Total	C	N	O	S	0	1	0
			1614	1015	263	331	5			

- Molecule 2 is a protein called Human monoclonal anti-HIV-1 gp120 V3 antibody 2557 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	215	Total	C	N	O	S	0	0	0
			1633	1044	263	320	6			
2	B	219	Total	C	N	O	S	0	0	0
			1659	1059	267	327	6			
2	E	219	Total	C	N	O	S	0	0	0
			1659	1059	267	327	6			
2	I	219	Total	C	N	O	S	0	0	0
			1659	1059	267	327	6			

- Molecule 3 is a protein called HIV-1 gp120 third variable region (V3) crown.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	14	Total	C	N	O	0	0	0
			113	73	23	17			
3	C	14	Total	C	N	O	0	0	0
			113	73	23	17			

- Molecule 4 is water.

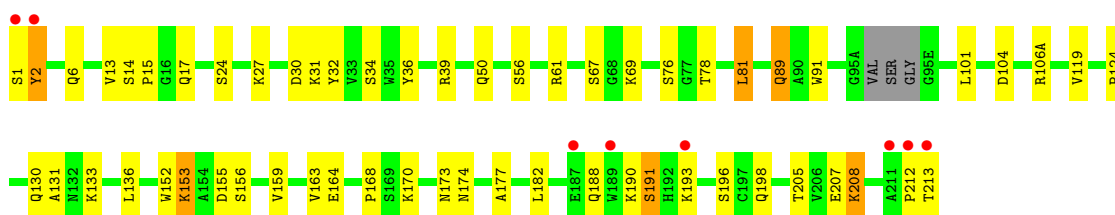
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	55	Total 55	O 55	0	0
4	H	44	Total 44	O 44	0	0
4	P	3	Total 3	O 3	0	0
4	A	12	Total 12	O 12	0	0
4	B	30	Total 30	O 30	0	0
4	C	1	Total 1	O 1	0	0
4	D	31	Total 31	O 31	0	0
4	E	49	Total 49	O 49	0	0
4	G	48	Total 48	O 48	0	0
4	I	55	Total 55	O 55	0	0

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.

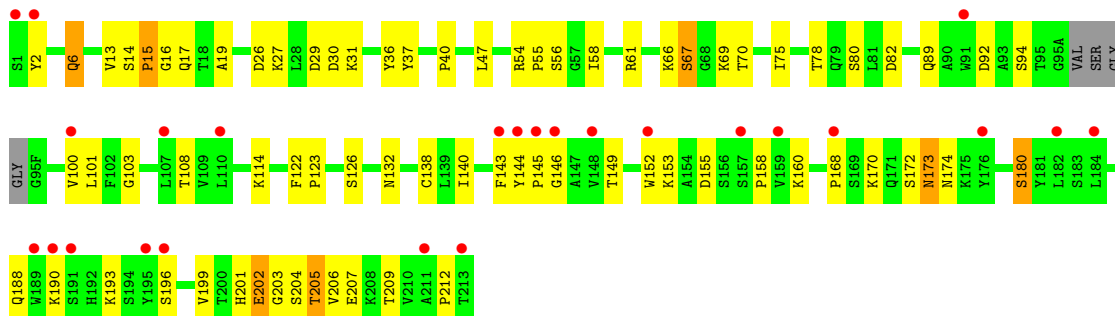
- Molecule 1: Human monoclonal anti-HIV-1 gp120 V3 antibody 2557 Fab light chain

Chain L: 



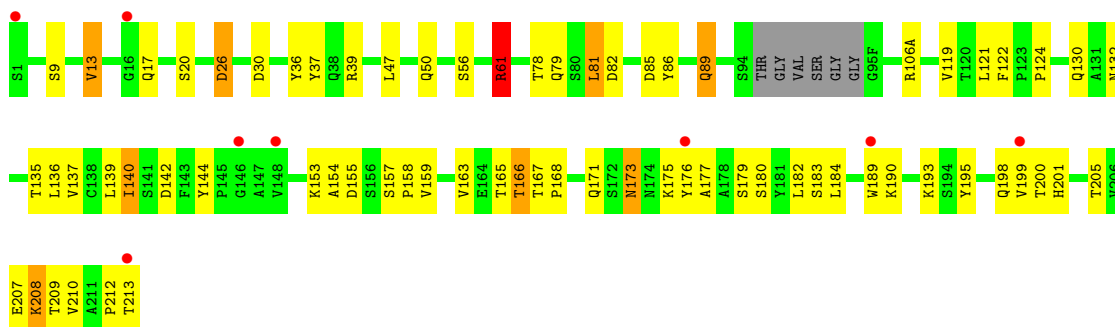
- Molecule 1: Human monoclonal anti-HIV-1 gp120 V3 antibody 2557 Fab light chain

Chain A: 



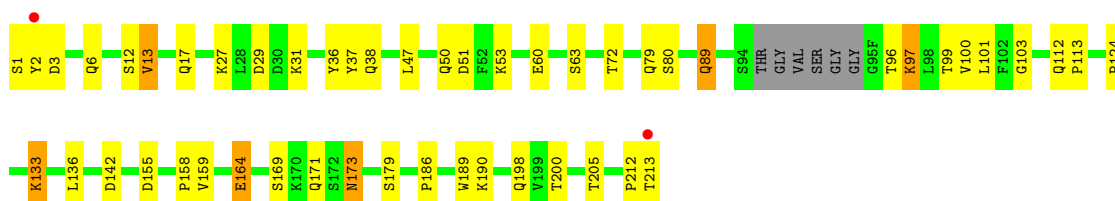
- Molecule 1: Human monoclonal anti-HIV-1 gp120 V3 antibody 2557 Fab light chain

Chain D: 



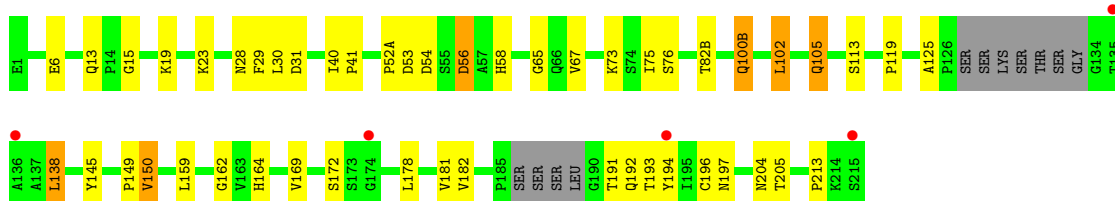
- Molecule 1: Human monoclonal anti-HIV-1 gp120 V3 antibody 2557 Fab light chain

Chain G:



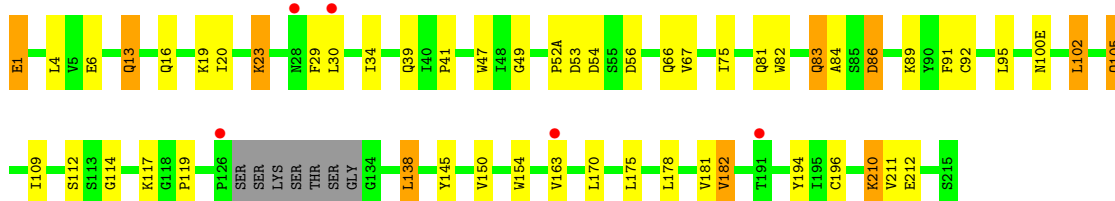
- Molecule 2: Human monoclonal anti-HIV-1 gp120 V3 antibody 2557 Fab heavy chain

Chain H:



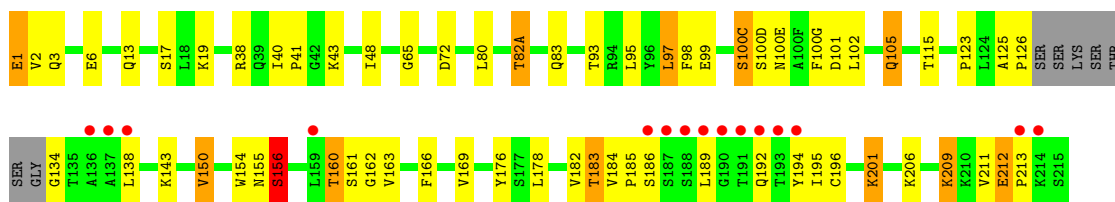
- Molecule 2: Human monoclonal anti-HIV-1 gp120 V3 antibody 2557 Fab heavy chain

Chain B:



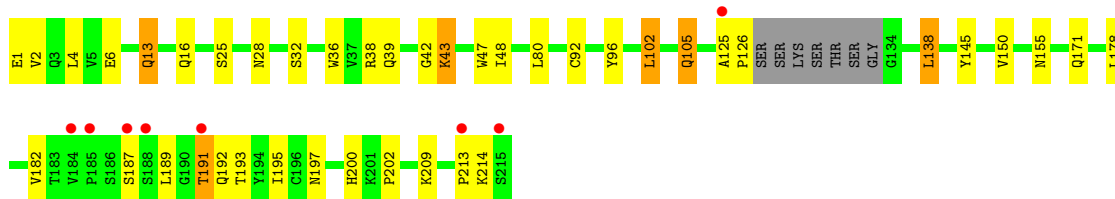
- Molecule 2: Human monoclonal anti-HIV-1 gp120 V3 antibody 2557 Fab heavy chain

Chain E:



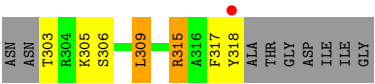
- Molecule 2: Human monoclonal anti-HIV-1 gp120 V3 antibody 2557 Fab heavy chain

Chain I:



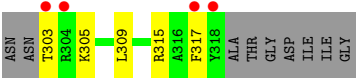
- Molecule 3: HIV-1 gp120 third variable region (V3) crown

Chain P:



- Molecule 3: HIV-1 gp120 third variable region (V3) crown

Chain C:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.68Å 142.93Å 85.01Å 90.00° 95.00° 90.00°	Depositor
Resolution (Å)	35.74 – 2.49 35.73 – 2.49	Depositor EDS
% Data completeness (in resolution range)	99.3 (35.74-2.49) 99.3 (35.73-2.49)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.202 , 0.280 0.201 , 0.278	Depositor DCC
R_{free} test set	3177 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	43.6	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 29.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 62523 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13628	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.55% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.69	0/1658	0.75	0/2261
1	D	0.76	0/1647	0.80	1/2246 (0.0%)
1	G	0.79	0/1656	0.83	0/2258
1	L	0.79	0/1662	0.82	0/2266
2	B	0.76	1/1702 (0.1%)	0.76	0/2319
2	E	0.83	0/1702	0.84	3/2319 (0.1%)
2	H	0.81	0/1675	0.83	1/2281 (0.0%)
2	I	0.83	1/1702 (0.1%)	0.81	0/2319
3	C	0.85	0/116	0.88	0/154
3	P	0.81	0/116	0.95	0/154
All	All	0.79	2/13636 (0.0%)	0.81	5/18577 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	I	92	CYS	CB-SG	-5.57	1.72	1.81
2	B	92	CYS	CB-SG	-5.17	1.73	1.81

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	101	ASP	CB-CG-OD1	6.09	123.78	118.30
2	E	101	ASP	CB-CG-OD2	-6.01	112.89	118.30
1	D	61	ARG	NE-CZ-NH2	-5.96	117.32	120.30
2	H	102	LEU	CA-CB-CG	5.09	127.00	115.30
2	E	72	ASP	CB-CG-OD2	5.06	122.86	118.30

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	1619	0	1578	56	0
1	D	1608	0	1568	48	0
1	G	1614	0	1574	37	0
1	L	1623	0	1581	38	0
2	B	1659	0	1618	36	0
2	E	1659	0	1618	52	0
2	H	1633	0	1591	38	0
2	I	1659	0	1618	30	0
3	C	113	0	115	6	0
3	P	113	0	115	8	0
4	A	12	0	0	0	0
4	B	30	0	0	0	0
4	C	1	0	0	0	0
4	D	31	0	0	1	0
4	E	49	0	0	1	0
4	G	48	0	0	2	0
4	H	44	0	0	0	0
4	I	55	0	0	1	0
4	L	55	0	0	4	0
4	P	3	0	0	0	0
All	All	13628	0	12976	329	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 13.

All (329) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:I:42:GLY:O	2:I:43:LYS:HE3	1.43	1.15
1:A:31:LYS:NZ	3:C:309:LEU:HD13	1.61	1.13
3:P:315:ARG:HH11	3:P:315:ARG:HG3	1.14	1.12
2:E:212:GLU:HB3	2:E:213:PRO:CD	1.79	1.11
2:E:212:GLU:HB3	2:E:213:PRO:HD3	1.12	1.05
2:H:192:GLN:NE2	2:H:193:THR:H	1.57	1.02
3:C:309:LEU:HD11	3:C:315:ARG:HD3	1.43	1.01

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:I:105:GLN:HE21	2:I:105:GLN:H	1.02	1.00
1:A:2:TYR:CD1	1:A:2:TYR:O	2.16	0.98
2:B:13:GLN:H	2:B:16:GLN:HE21	1.03	0.97
1:L:36:TYR:HE1	1:L:89:GLN:HG2	1.28	0.95
2:I:43:LYS:CE	2:I:43:LYS:HA	1.96	0.95
1:A:2:TYR:OH	1:A:31:LYS:HD2	1.67	0.95
2:H:192:GLN:HE21	2:H:193:THR:H	1.14	0.94
1:G:198:GLN:HG3	1:G:205:THR:CG2	1.98	0.93
2:I:13:GLN:HG2	2:I:16:GLN:NE2	1.84	0.92
1:D:166:THR:CG2	1:D:179:SER:H	1.82	0.92
2:H:149:PRO:HG3	2:E:201:LYS:HG3	1.51	0.91
2:H:105:GLN:HE21	2:H:105:GLN:H	0.99	0.91
1:A:205:THR:HG23	1:A:206:VAL:N	1.84	0.90
2:E:105:GLN:HE21	2:E:105:GLN:H	0.94	0.90
1:L:17:GLN:O	1:L:78:THR:HG23	1.72	0.89
2:I:13:GLN:H	2:I:16:GLN:HE21	1.17	0.89
2:E:105:GLN:H	2:E:105:GLN:NE2	1.70	0.88
2:H:105:GLN:NE2	2:H:105:GLN:H	1.73	0.87
1:L:2:TYR:HE2	1:L:31:LYS:HZ2	1.16	0.87
2:B:23:LYS:NZ	2:B:75:ILE:O	2.08	0.87
2:B:6:GLU:H	2:B:105:GLN:HE22	1.18	0.86
2:H:192:GLN:HE21	2:H:193:THR:N	1.75	0.85
2:B:105:GLN:H	2:B:105:GLN:HE21	1.24	0.85
1:A:31:LYS:HZ2	3:C:309:LEU:HD13	1.39	0.85
1:G:198:GLN:HG3	1:G:205:THR:HG21	1.56	0.84
2:H:105:GLN:HE21	2:H:105:GLN:N	1.76	0.83
1:D:171:GLN:HE21	1:D:177:ALA:HB2	1.43	0.83
1:G:171:GLN:HE21	1:G:173:ASN:HD21	1.26	0.83
1:G:37:TYR:HB2	1:G:47:LEU:HD11	1.61	0.83
1:G:36:TYR:HE1	1:G:89:GLN:HG2	1.43	0.82
1:L:36:TYR:CE1	1:L:89:GLN:HG2	2.13	0.82
2:E:105:GLN:HE21	2:E:105:GLN:N	1.76	0.82
2:H:40:ILE:HG23	2:H:41:PRO:HD2	1.61	0.82
1:D:36:TYR:HE1	1:D:89:GLN:HG2	1.45	0.82
2:E:1:GLU:OE2	2:E:1:GLU:N	2.12	0.81
2:I:43:LYS:HE2	2:I:43:LYS:HA	1.62	0.80
2:B:13:GLN:H	2:B:16:GLN:NE2	1.82	0.78
2:I:6:GLU:H	2:I:105:GLN:HE22	1.31	0.78
2:H:6:GLU:H	2:H:105:GLN:HE22	1.32	0.77
2:I:105:GLN:NE2	2:I:105:GLN:H	1.79	0.77
1:A:2:TYR:CG	1:A:2:TYR:O	2.38	0.76
3:P:315:ARG:HG3	3:P:315:ARG:NH1	1.95	0.76

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:I:105:GLN:HE21	2:I:105:GLN:N	1.83	0.75
2:B:105:GLN:H	2:B:105:GLN:NE2	1.83	0.75
1:A:153:LYS:HA	1:A:158:PRO:HA	1.68	0.74
2:H:149:PRO:HG3	2:E:201:LYS:CG	2.16	0.74
1:D:37:TYR:HB2	1:D:47:LEU:HD11	1.70	0.74
1:G:38:GLN:OE1	2:I:39:GLN:NE2	2.16	0.73
2:B:82:TRP:NE1	2:B:86:ASP:OD1	2.21	0.73
1:A:36:TYR:HE1	1:A:89:GLN:HG2	1.53	0.73
1:L:153:LYS:HD2	1:L:198:GLN:HE22	1.52	0.73
1:L:188:GLN:HA	1:L:191:SER:HB3	1.72	0.72
2:E:212:GLU:CB	2:E:213:PRO:CD	2.65	0.71
1:D:142:ASP:H	1:D:171:GLN:HE22	1.36	0.71
1:A:31:LYS:HZ3	3:C:309:LEU:HD13	1.55	0.70
1:G:173:ASN:N	1:G:173:ASN:HD22	1.88	0.70
1:A:170:LYS:HE3	1:A:174:ASN:ND2	2.07	0.70
1:L:163:VAL:HG22	1:L:182:LEU:CD1	2.21	0.69
1:G:164:GLU:HG2	2:I:171:GLN:HA	1.73	0.69
2:E:163:VAL:HG22	2:E:182:VAL:HG22	1.73	0.69
2:E:126:PRO:HD3	2:E:212:GLU:O	1.92	0.69
2:B:13:GLN:HG2	2:B:16:GLN:NE2	2.07	0.69
1:G:36:TYR:CE1	1:G:89:GLN:HG2	2.27	0.69
1:D:166:THR:HG22	1:D:179:SER:O	1.93	0.68
1:D:39:ARG:NH2	1:D:81:LEU:HD22	2.09	0.68
1:A:114:LYS:HE3	1:A:202:GLU:OE2	1.94	0.68
2:E:98:PHE:H	2:E:100(E):ASN:ND2	1.91	0.68
1:A:13:VAL:HG22	1:A:14:SER:N	2.09	0.68
1:G:173:ASN:H	1:G:173:ASN:HD22	1.42	0.68
1:D:171:GLN:NE2	1:D:177:ALA:HB2	2.09	0.67
2:E:150:VAL:HG22	2:E:178:LEU:HD21	1.77	0.67
1:D:163:VAL:HG23	1:D:182:LEU:HD13	1.77	0.67
1:D:119:VAL:O	1:D:208:LYS:HE3	1.95	0.66
3:C:309:LEU:HD11	3:C:315:ARG:CD	2.23	0.66
1:A:14:SER:O	1:A:15:PRO:O	2.14	0.66
1:A:173:ASN:HD22	1:A:173:ASN:C	1.99	0.66
2:I:43:LYS:HE3	2:I:43:LYS:HA	1.75	0.66
2:I:13:GLN:H	2:I:16:GLN:NE2	1.93	0.66
2:H:192:GLN:NE2	2:H:193:THR:N	2.35	0.65
2:H:192:GLN:NE2	2:H:193:THR:OG1	2.30	0.65
2:B:119:PRO:HB3	2:B:145:TYR:HB3	1.78	0.65
1:A:201:HIS:O	1:A:202:GLU:CB	2.45	0.65
1:L:104:ASP:HB2	4:L:238:HOH:O	1.98	0.64
2:E:186:SER:HA	2:E:189:LEU:HD13	1.79	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:54:ASP:OD2	2:B:56:ASP:HB2	1.98	0.63
2:E:154:TRP:O	2:E:155:ASN:HB2	1.99	0.62
2:E:195:ILE:HA	2:E:209:LYS:O	2.00	0.62
1:A:36:TYR:CE1	1:A:89:GLN:HG2	2.33	0.62
1:A:201:HIS:C	1:A:202:GLU:HG3	2.20	0.61
1:A:67:SER:O	1:A:70:THR:HG22	1.99	0.61
1:L:173:ASN:O	1:L:174:ASN:HB2	2.00	0.61
2:B:6:GLU:H	2:B:105:GLN:NE2	1.95	0.61
1:G:124:PRO:HD3	1:G:136:LEU:HD23	1.83	0.60
2:E:201:LYS:HZ2	2:E:201:LYS:H	1.49	0.60
1:L:50:GLN:NE2	4:L:216:HOH:O	2.20	0.60
1:D:200:THR:OG1	1:D:205:THR:HG22	2.01	0.60
1:A:55:PRO:HD2	1:A:58:ILE:HG13	1.83	0.59
2:B:6:GLU:N	2:B:105:GLN:HE22	1.96	0.59
2:E:100(C):SER:HB3	2:E:100(E):ASN:ND2	2.18	0.59
1:L:106(A):ARG:NH1	4:L:247:HOH:O	2.20	0.59
1:A:201:HIS:O	1:A:202:GLU:HG3	2.02	0.58
3:P:305:LYS:HE2	3:P:318:TYR:CD2	2.38	0.58
1:G:96:THR:HG23	1:G:97:LYS:N	2.18	0.58
1:A:37:TYR:HB2	1:A:47:LEU:HD11	1.84	0.58
1:D:61:ARG:NH2	1:D:82:ASP:OD1	2.34	0.58
1:L:164:GLU:HB3	2:H:169:VAL:HG11	1.85	0.58
1:A:13:VAL:CG2	1:A:14:SER:N	2.66	0.58
1:G:72:THR:OG1	4:G:276:HOH:O	2.17	0.58
1:L:153:LYS:HB3	1:L:196:SER:HB2	1.83	0.58
2:B:1:GLU:CD	2:B:1:GLU:N	2.56	0.58
1:L:32:TYR:OH	2:H:100(B):GLN:HG3	2.04	0.58
1:G:50:GLN:HE21	1:G:53:LYS:NZ	2.02	0.58
2:E:160:THR:O	2:E:160:THR:OG1	2.22	0.57
1:G:173:ASN:H	1:G:173:ASN:ND2	2.00	0.57
1:A:201:HIS:O	1:A:202:GLU:HB2	2.02	0.57
2:E:123:PRO:HD3	2:E:209:LYS:HE2	1.87	0.57
1:G:50:GLN:HE21	1:G:53:LYS:HZ2	1.53	0.57
2:B:210:LYS:HD2	2:B:212:GLU:HG3	1.86	0.56
2:E:2:VAL:HG13	2:E:102:LEU:HD11	1.87	0.56
1:D:173:ASN:HD21	1:D:175:LYS:HB2	1.70	0.56
2:I:150:VAL:CG2	2:I:178:LEU:HD21	2.36	0.56
1:L:2:TYR:OH	1:L:31:LYS:HD2	2.06	0.56
1:D:130:GLN:C	1:D:132:ASN:H	2.09	0.56
1:L:163:VAL:HG22	1:L:182:LEU:HD11	1.88	0.55
1:L:36:TYR:HE1	1:L:89:GLN:CG	2.10	0.55
1:D:13:VAL:HG22	1:D:17:GLN:HB2	1.88	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:105:GLN:N	2:B:105:GLN:HE21	1.98	0.55
1:D:166:THR:HG23	1:D:179:SER:H	1.66	0.55
2:E:134:GLY:O	2:E:185:PRO:HA	2.07	0.55
1:A:152:TRP:HE1	1:A:180:SER:HB3	1.70	0.55
1:G:36:TYR:HE1	1:G:89:GLN:CG	2.16	0.55
2:E:100(D):SER:OG	2:E:100(D):SER:O	2.23	0.54
1:A:143:PHE:HE2	1:A:146:GLY:HA2	1.72	0.54
2:I:13:GLN:HG2	2:I:16:GLN:HE21	1.65	0.54
1:G:171:GLN:NE2	1:G:173:ASN:HD21	2.01	0.54
2:H:19:LYS:NZ	1:A:203:GLY:HA3	2.22	0.54
1:A:173:ASN:ND2	1:A:173:ASN:C	2.61	0.54
1:A:153:LYS:HB2	1:A:196:SER:HB3	1.90	0.54
2:E:138:LEU:HD23	2:E:182:VAL:O	2.08	0.54
2:E:201:LYS:NZ	2:E:201:LYS:H	2.06	0.53
1:A:144:TYR:O	1:A:201:HIS:CE1	2.61	0.53
1:G:142:ASP:OD1	1:G:171:GLN:NE2	2.40	0.53
2:E:138:LEU:H	2:E:138:LEU:HD23	1.73	0.53
1:L:2:TYR:HE2	1:L:31:LYS:NZ	1.98	0.53
1:D:140:ILE:HD12	1:D:199:VAL:HG21	1.90	0.52
2:B:1:GLU:CD	2:B:1:GLU:H3	2.12	0.52
1:L:39:ARG:NH2	1:L:81:LEU:HD22	2.24	0.52
2:H:193:THR:O	2:H:194:TYR:HB2	2.08	0.52
1:A:114:LYS:CE	1:A:202:GLU:OE2	2.56	0.52
1:D:173:ASN:ND2	1:D:175:LYS:H	2.08	0.52
2:H:28:ASN:HD21	2:H:31:ASP:CG	2.12	0.52
2:H:149:PRO:CG	2:E:201:LYS:HG3	2.32	0.52
2:H:162:GLY:O	2:H:182:VAL:HA	2.10	0.52
1:A:2:TYR:OH	1:A:31:LYS:CD	2.48	0.51
1:A:201:HIS:C	1:A:202:GLU:CG	2.77	0.51
1:G:171:GLN:HE21	1:G:173:ASN:ND2	2.02	0.51
1:A:61:ARG:NH2	1:A:82:ASP:OD1	2.37	0.51
1:D:163:VAL:HG23	1:D:182:LEU:CD1	2.40	0.51
2:I:126:PRO:HG3	2:I:138:LEU:HD23	1.93	0.51
1:G:171:GLN:HB2	1:G:173:ASN:HD21	1.76	0.51
2:E:138:LEU:HD12	2:E:211:VAL:HB	1.92	0.51
1:D:184:LEU:HD11	1:D:195:TYR:CZ	2.46	0.51
1:G:171:GLN:HB2	1:G:173:ASN:ND2	2.26	0.51
1:A:17:GLN:O	1:A:78:THR:HG23	2.11	0.51
1:G:2:TYR:CZ	1:G:27:LYS:HB2	2.46	0.51
2:B:89:LYS:HD3	2:B:91:PHE:CZ	2.45	0.50
1:A:13:VAL:CG2	1:A:14:SER:H	2.23	0.50
1:D:140:ILE:HD12	1:D:199:VAL:CG2	2.42	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:119:VAL:O	1:L:208:LYS:NZ	2.45	0.50
2:I:4:LEU:HG	2:I:102:LEU:HD13	1.92	0.50
1:D:153:LYS:HE2	1:D:198:GLN:NE2	2.27	0.50
1:L:2:TYR:HB2	4:L:226:HOH:O	2.12	0.50
1:D:130:GLN:C	1:D:132:ASN:N	2.65	0.49
1:L:124:PRO:HD3	1:L:136:LEU:HD23	1.94	0.49
2:I:214:LYS:NZ	4:I:230:HOH:O	2.44	0.49
2:H:28:ASN:ND2	2:H:31:ASP:CG	2.65	0.49
1:D:144:TYR:O	1:D:201:HIS:HE1	1.95	0.49
1:D:36:TYR:CE1	1:D:89:GLN:HG2	2.36	0.49
2:H:138:LEU:O	2:H:181:VAL:HG23	2.12	0.49
2:B:4:LEU:HG	2:B:102:LEU:HD13	1.94	0.49
1:D:198:GLN:HG2	1:D:207:GLU:HB2	1.94	0.49
2:I:1:GLU:HG2	2:I:2:VAL:N	2.28	0.49
1:D:85:ASP:OD1	1:D:106(A):ARG:HD3	2.13	0.49
1:G:112:GLN:HB2	1:G:113:PRO:HD2	1.95	0.49
1:L:168:PRO:HA	1:L:177:ALA:O	2.13	0.49
2:E:38:ARG:HB2	2:E:48:ILE:HD11	1.93	0.49
1:L:153:LYS:HG2	1:L:156:SER:HA	1.95	0.49
2:E:184:VAL:HG11	2:E:194:TYR:CE1	2.49	0.48
2:H:119:PRO:HD2	2:H:205:THR:HG21	1.94	0.48
2:H:40:ILE:CG2	2:H:41:PRO:HD2	2.40	0.48
2:B:210:LYS:CD	2:B:212:GLU:HG3	2.42	0.48
2:H:54:ASP:OD2	2:H:56:ASP:HB2	2.13	0.48
1:L:152:TRP:CD2	1:L:182:LEU:HD13	2.49	0.48
1:G:50:GLN:O	1:G:51:ASP:HB2	2.13	0.48
2:B:154:TRP:CH2	2:B:196:CYS:HB3	2.49	0.48
1:G:99:THR:HG22	1:G:101:LEU:HD23	1.95	0.48
1:A:14:SER:O	1:A:15:PRO:C	2.52	0.48
1:D:163:VAL:CG2	1:D:182:LEU:HD13	2.43	0.48
2:H:28:ASN:ND2	2:H:31:ASP:HB2	2.28	0.48
2:I:145:TYR:CE1	2:I:150:VAL:HG13	2.48	0.47
3:C:305:LYS:O	3:C:305:LYS:HG2	2.12	0.47
2:E:17:SER:HB3	2:E:82(A):THR:O	2.14	0.47
1:D:166:THR:HG21	1:D:179:SER:H	1.74	0.47
2:B:30:LEU:HD13	2:B:53:ASP:OD1	2.14	0.47
1:A:26:ASP:OD2	1:A:69:LYS:NZ	2.46	0.47
1:D:124:PRO:HD3	1:D:136:LEU:HD23	1.96	0.47
3:P:315:ARG:CG	3:P:315:ARG:HH11	2.01	0.47
1:G:13:VAL:HG22	1:G:17:GLN:HB2	1.97	0.47
2:B:29:PHE:CE2	2:B:52(A):PRO:HB3	2.49	0.47
2:E:95:LEU:HD21	2:E:97:LEU:HG	1.97	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:E:2:VAL:O	2:E:3:GLN:HB3	2.15	0.47
2:B:83:GLN:HG2	2:B:84:ALA:N	2.30	0.47
1:D:122:PHE:HE1	2:E:125:ALA:O	1.97	0.46
1:G:186:PRO:O	1:G:190:LYS:HG3	2.15	0.46
1:A:201:HIS:O	1:A:202:GLU:CG	2.62	0.46
2:B:210:LYS:NZ	2:B:211:VAL:O	2.43	0.46
2:B:163:VAL:HG22	2:B:182:VAL:HG13	1.97	0.46
1:D:26:ASP:N	1:D:26:ASP:OD1	2.41	0.46
1:A:54:ARG:NH1	1:A:58:ILE:O	2.44	0.46
2:I:191:THR:HG22	2:I:192:GLN:N	2.31	0.46
1:L:27:LYS:HB2	1:L:31:LYS:HE3	1.98	0.46
2:E:2:VAL:CG1	2:E:102:LEU:HD11	2.46	0.46
2:B:150:VAL:CG2	2:B:178:LEU:HD21	2.46	0.46
2:H:75:ILE:O	2:H:76:SER:HB2	2.15	0.46
2:B:82:TRP:CD1	2:B:86:ASP:OD1	2.69	0.46
1:A:2:TYR:CZ	1:A:31:LYS:HD2	2.48	0.46
1:A:205:THR:O	1:A:206:VAL:HG23	2.16	0.46
1:D:171:GLN:NE2	1:D:176:TYR:O	2.49	0.46
2:B:20:ILE:HG21	2:B:109:ILE:HD11	1.98	0.46
4:D:222:HOH:O	2:E:143:LYS:HG2	2.16	0.46
1:G:189:TRP:CZ2	1:G:212:PRO:HA	2.51	0.46
1:A:100:VAL:HB	2:B:47:TRP:CG	2.51	0.45
1:L:17:GLN:O	1:L:78:THR:CG2	2.54	0.45
1:D:189:TRP:CZ2	1:D:212:PRO:HA	2.51	0.45
1:A:140:ILE:HG12	1:A:199:VAL:HG21	1.97	0.45
1:A:19:ALA:N	1:A:75:ILE:O	2.40	0.45
2:I:200:HIS:CD2	2:I:202:PRO:HD2	2.50	0.45
1:G:89:GLN:HA	1:G:101:LEU:O	2.17	0.45
2:E:82(A):THR:N	4:E:227:HOH:O	2.49	0.45
1:A:144:TYR:CG	1:A:145:PRO:HA	2.51	0.45
2:E:19:LYS:HA	2:E:80:LEU:O	2.17	0.45
2:E:154:TRP:O	2:E:156:SER:N	2.50	0.45
1:D:166:THR:CG2	1:D:179:SER:N	2.66	0.45
1:G:96:THR:CG2	1:G:97:LYS:N	2.79	0.45
2:I:38:ARG:HB2	2:I:48:ILE:HD11	1.99	0.45
1:G:133:LYS:HA	1:G:133:LYS:HE3	1.99	0.45
1:D:165:THR:HG1	1:D:180:SER:HG	1.65	0.45
2:B:95:LEU:HD11	2:B:100(E):ASN:HB3	1.99	0.45
1:L:152:TRP:CG	1:L:182:LEU:HD13	2.52	0.44
2:H:15:GLY:HA2	2:H:82(B):THR:HG23	1.99	0.44
2:I:43:LYS:CA	2:I:43:LYS:CE	2.82	0.44
2:H:178:LEU:C	2:H:178:LEU:HD12	2.38	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:E:40:ILE:HG23	2:E:41:PRO:HD2	1.99	0.44
2:E:138:LEU:CD2	2:E:184:VAL:HG22	2.47	0.44
2:E:155:ASN:N	2:E:195:ILE:O	2.34	0.44
1:D:79:GLN:NE2	1:D:82:ASP:OD2	2.51	0.44
1:G:169:SER:HB3	4:G:245:HOH:O	2.18	0.44
1:A:144:TYR:CD1	1:A:145:PRO:HA	2.53	0.43
1:L:14:SER:O	1:L:15:PRO:C	2.55	0.43
1:L:198:GLN:HG3	1:L:207:GLU:HG3	2.00	0.43
1:D:13:VAL:HG11	1:D:78:THR:HG21	2.00	0.43
2:B:178:LEU:C	2:B:178:LEU:HD12	2.38	0.43
1:D:39:ARG:HH21	1:D:81:LEU:HD22	1.77	0.43
1:L:177:ALA:HB3	2:H:164:HIS:CE1	2.54	0.43
2:E:212:GLU:CB	2:E:213:PRO:HD3	2.07	0.43
2:H:54:ASP:OD1	3:P:305:LYS:NZ	2.50	0.43
2:H:19:LYS:HZ2	1:A:203:GLY:HA3	1.83	0.43
2:I:36:TRP:CE3	2:I:80:LEU:HD22	2.54	0.43
1:D:139:LEU:HD22	2:E:166:PHE:CE1	2.53	0.43
1:D:154:ALA:HB2	1:D:195:TYR:CE2	2.53	0.43
1:A:190:LYS:HA	1:A:212:PRO:HG3	2.00	0.43
2:E:93:THR:HG23	2:E:100(G):PHE:CD1	2.54	0.43
1:A:92:ASP:HB2	1:A:101:LEU:HD11	2.01	0.43
2:H:125:ALA:HB1	2:H:213:PRO:HA	1.99	0.42
1:D:155:ASP:OD1	1:D:193:LYS:HB2	2.19	0.42
1:A:31:LYS:O	1:A:66:LYS:NZ	2.51	0.42
2:E:154:TRP:CZ3	2:E:196:CYS:HB3	2.54	0.42
2:B:47:TRP:CZ2	2:B:49:GLY:HA2	2.55	0.42
2:I:209:LYS:HD2	2:I:209:LYS:HA	1.85	0.42
1:L:2:TYR:OH	1:L:31:LYS:CD	2.67	0.42
2:E:100(C):SER:HB3	2:E:100(E):ASN:HD21	1.84	0.42
1:A:14:SER:O	1:A:17:GLN:HB3	2.19	0.42
1:G:31:LYS:HA	1:G:31:LYS:HD2	1.84	0.42
2:H:30:LEU:HA	2:H:52(A):PRO:HB2	2.01	0.42
1:L:61:ARG:HD2	1:L:76:SER:O	2.19	0.42
2:E:162:GLY:O	2:E:183:THR:N	2.41	0.42
2:H:145:TYR:CZ	2:H:150:VAL:HG13	2.55	0.42
3:P:309:LEU:N	3:P:309:LEU:HD23	2.34	0.42
2:I:150:VAL:HG22	2:I:178:LEU:HD21	2.01	0.42
1:D:157:SER:HA	1:D:158:PRO:HD3	1.94	0.42
1:D:36:TYR:O	1:D:86:TYR:HA	2.19	0.42
3:P:309:LEU:HG	3:P:317:PHE:HE1	1.84	0.42
1:G:100:VAL:HB	2:I:47:TRP:CD2	2.54	0.42
1:L:31:LYS:HD3	1:L:91:TRP:O	2.20	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:193:LYS:O	1:L:212:PRO:HD2	2.20	0.41
2:H:58:HIS:NE2	3:P:318:TYR:OH	2.45	0.41
2:H:29:PHE:CE2	2:H:52(A):PRO:HB3	2.54	0.41
1:A:138:CYS:HB2	1:A:152:TRP:CZ2	2.56	0.41
1:A:61:ARG:HH22	1:A:82:ASP:CG	2.20	0.41
2:I:125:ALA:HA	2:I:126:PRO:HD3	1.87	0.41
1:A:122:PHE:HA	1:A:123:PRO:HD3	1.90	0.41
2:B:138:LEU:O	2:B:181:VAL:HA	2.20	0.41
1:L:89:GLN:HA	1:L:101:LEU:O	2.21	0.41
2:B:194:TYR:H	2:B:210:LYS:HZ1	1.66	0.41
1:G:89:GLN:HB2	1:G:101:LEU:O	2.20	0.41
1:A:6:GLN:OE1	1:A:103:GLY:HA3	2.21	0.41
2:E:169:VAL:O	2:E:176:TYR:HA	2.21	0.41
1:D:144:TYR:O	1:D:201:HIS:CE1	2.72	0.41
2:E:6:GLU:H	2:E:105:GLN:HE22	1.69	0.41
1:L:212:PRO:HA	1:L:213:THR:HA	1.61	0.41
1:D:167:THR:O	1:D:168:PRO:C	2.59	0.41
1:D:135:THR:OG1	2:E:143:LYS:HE2	2.20	0.40
1:G:6:GLN:OE1	1:G:103:GLY:HA3	2.21	0.40
2:I:32:SER:HB3	2:I:96:TYR:CE2	2.56	0.40
2:E:209:LYS:HD3	2:E:209:LYS:HA	1.85	0.40
1:D:121:LEU:HD12	1:D:137:VAL:O	2.20	0.40
1:L:34:SER:HB2	1:L:89:GLN:HG3	2.03	0.40
1:A:89:GLN:HA	1:A:101:LEU:O	2.20	0.40
2:B:83:GLN:HG2	2:B:84:ALA:H	1.86	0.40
2:H:53:ASP:O	2:H:73:LYS:NZ	2.47	0.40
2:H:40:ILE:HG23	2:H:41:PRO:CD	2.40	0.40
2:B:154:TRP:CZ3	2:B:196:CYS:HB3	2.56	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	211/219 (96%)	187 (89%)	19 (9%)	5 (2%)	9 13

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	209/219 (95%)	196 (94%)	13 (6%)	0	100	100
1	G	210/219 (96%)	199 (95%)	9 (4%)	2 (1%)	22	38
1	L	212/219 (97%)	198 (93%)	10 (5%)	4 (2%)	12	19
2	B	215/226 (95%)	202 (94%)	11 (5%)	2 (1%)	25	42
2	E	215/226 (95%)	196 (91%)	16 (7%)	3 (1%)	16	27
2	H	209/226 (92%)	194 (93%)	13 (6%)	2 (1%)	22	38
2	I	215/226 (95%)	203 (94%)	11 (5%)	1 (0%)	38	60
3	C	12/23 (52%)	12 (100%)	0	0	100	100
3	P	12/23 (52%)	12 (100%)	0	0	100	100
All	All	1720/1826 (94%)	1599 (93%)	102 (6%)	19 (1%)	21	34

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	2	TYR
2	E	156	SER
2	E	212	GLU
1	L	131	ALA
1	A	15	PRO
1	A	16	GLY
1	A	168	PRO
2	B	114	GLY
2	H	65	GLY
1	A	67	SER
2	B	41	PRO
1	G	155	ASP
2	I	213	PRO
1	L	155	ASP
1	A	202	GLU
1	L	191	SER
2	H	172	SER
2	E	65	GLY
1	G	158	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	184/186 (99%)	161 (88%)	23 (12%)	7	12
1	D	183/186 (98%)	163 (89%)	20 (11%)	9	16
1	G	184/186 (99%)	166 (90%)	18 (10%)	12	21
1	L	184/186 (99%)	166 (90%)	18 (10%)	12	21
2	B	188/194 (97%)	168 (89%)	20 (11%)	10	18
2	E	188/194 (97%)	169 (90%)	19 (10%)	11	20
2	H	184/194 (95%)	169 (92%)	15 (8%)	17	29
2	I	188/194 (97%)	173 (92%)	15 (8%)	17	31
3	C	11/17 (65%)	9 (82%)	2 (18%)	2	4
3	P	11/17 (65%)	7 (64%)	4 (36%)	0	0
All	All	1505/1554 (97%)	1351 (90%)	154 (10%)	11	19

All (154) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	1	SER
1	L	6	GLN
1	L	13	VAL
1	L	24	SER
1	L	30	ASP
1	L	56	SER
1	L	67	SER
1	L	69	LYS
1	L	81	LEU
1	L	89	GLN
1	L	130	GLN
1	L	133	LYS
1	L	153	LYS
1	L	159	VAL
1	L	170	LYS
1	L	190	LYS
1	L	205	THR
1	L	208	LYS
2	H	13	GLN
2	H	23	LYS
2	H	56	ASP
2	H	67	VAL

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Mol	Chain	Res	Type
2	H	100(B)	GLN
2	H	102	LEU
2	H	105	GLN
2	H	113	SER
2	H	138	LEU
2	H	150	VAL
2	H	159	LEU
2	H	191	THR
2	H	196	CYS
2	H	197	ASN
2	H	204	ASN
3	P	303	THR
3	P	306	SER
3	P	309	LEU
3	P	315	ARG
1	A	6	GLN
1	A	27	LYS
1	A	29	ASP
1	A	30	ASP
1	A	40	PRO
1	A	56	SER
1	A	80	SER
1	A	94	SER
1	A	108	THR
1	A	126	SER
1	A	132	ASN
1	A	149	THR
1	A	155	ASP
1	A	160	LYS
1	A	172	SER
1	A	173	ASN
1	A	180	SER
1	A	188	GLN
1	A	193	LYS
1	A	204	SER
1	A	205	THR
1	A	207	GLU
1	A	209	THR
2	B	1	GLU
2	B	13	GLN
2	B	19	LYS
2	B	23	LYS

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Mol	Chain	Res	Type
2	B	34	ILE
2	B	39	GLN
2	B	66	GLN
2	B	67	VAL
2	B	81	GLN
2	B	83	GLN
2	B	86	ASP
2	B	102	LEU
2	B	105	GLN
2	B	112	SER
2	B	117	LYS
2	B	138	LEU
2	B	170	LEU
2	B	175	LEU
2	B	182	VAL
2	B	210	LYS
3	C	303	THR
3	C	317	PHE
1	D	9	SER
1	D	13	VAL
1	D	20	SER
1	D	26	ASP
1	D	30	ASP
1	D	50	GLN
1	D	56	SER
1	D	61	ARG
1	D	81	LEU
1	D	89	GLN
1	D	140	ILE
1	D	159	VAL
1	D	166	THR
1	D	173	ASN
1	D	183	SER
1	D	190	LYS
1	D	208	LYS
1	D	209	THR
1	D	210	VAL
1	D	213	THR
2	E	1	GLU
2	E	13	GLN
2	E	43	LYS
2	E	82(A)	THR

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Mol	Chain	Res	Type
2	E	83	GLN
2	E	97	LEU
2	E	99	GLU
2	E	100(C)	SER
2	E	105	GLN
2	E	115	THR
2	E	150	VAL
2	E	156	SER
2	E	160	THR
2	E	161	SER
2	E	183	THR
2	E	192	GLN
2	E	201	LYS
2	E	206	LYS
2	E	209	LYS
1	G	1	SER
1	G	3	ASP
1	G	12	SER
1	G	13	VAL
1	G	29	ASP
1	G	60	GLU
1	G	63	SER
1	G	79	GLN
1	G	80	SER
1	G	89	GLN
1	G	97	LYS
1	G	133	LYS
1	G	159	VAL
1	G	164	GLU
1	G	173	ASN
1	G	179	SER
1	G	200	THR
1	G	213	THR
2	I	13	GLN
2	I	25	SER
2	I	28	ASN
2	I	43	LYS
2	I	102	LEU
2	I	105	GLN
2	I	138	LEU
2	I	155	ASN
2	I	182	VAL

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Mol	Chain	Res	Type
2	I	187	SER
2	I	189	LEU
2	I	191	THR
2	I	193	THR
2	I	195	ILE
2	I	197	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (46) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	50	GLN
1	L	89	GLN
1	L	112	GLN
1	L	198	GLN
2	H	3	GLN
2	H	28	ASN
2	H	81	GLN
2	H	105	GLN
2	H	164	HIS
2	H	192	GLN
1	A	79	GLN
1	A	132	ASN
1	A	173	ASN
1	A	174	ASN
1	A	201	HIS
2	B	3	GLN
2	B	16	GLN
2	B	81	GLN
2	B	83	GLN
2	B	105	GLN
2	B	192	GLN
2	B	204	ASN
1	D	171	GLN
1	D	173	ASN
1	D	174	ASN
1	D	192	HIS
1	D	201	HIS
2	E	3	GLN
2	E	13	GLN
2	E	16	GLN
2	E	81	GLN
2	E	100(E)	ASN

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Mol	Chain	Res	Type
2	E	105	GLN
2	E	155	ASN
2	E	192	GLN
1	G	38	GLN
1	G	50	GLN
1	G	79	GLN
1	G	173	ASN
1	G	198	GLN
2	I	16	GLN
2	I	28	ASN
2	I	39	GLN
2	I	100(E)	ASN
2	I	105	GLN
2	I	155	ASN

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 ⓘ

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, 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	215/219 (98%)	0.60	25 (11%) 5 5	26, 58, 87, 92	0
1	D	213/219 (97%)	0.17	8 (3%) 38 40	25, 46, 67, 74	0
1	G	213/219 (97%)	-0.09	2 (0%) 81 82	22, 41, 60, 67	0
1	L	216/219 (98%)	-0.03	8 (3%) 39 41	21, 37, 73, 76	0
2	B	219/226 (96%)	0.17	5 (2%) 57 60	25, 43, 59, 69	0
2	E	219/226 (96%)	0.15	15 (6%) 17 17	23, 37, 78, 88	0
2	H	215/226 (95%)	0.11	5 (2%) 57 60	22, 39, 71, 80	0
2	I	219/226 (96%)	-0.03	8 (3%) 39 41	18, 34, 66, 80	0
3	C	14/23 (60%)	1.73	4 (28%) 1 1	64, 74, 87, 88	0
3	P	14/23 (60%)	0.43	1 (7%) 16 15	43, 51, 62, 65	0
All	All	1757/1826 (96%)	0.15	81 (4%) 31 32	18, 42, 74, 92	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	191	THR	6.9
1	A	189	TRP	5.9
3	C	318	TYR	5.7
2	E	188	SER	5.5
2	E	190	GLY	5.0
3	C	317	PHE	5.0
1	A	184	LEU	4.8
1	A	2	TYR	4.7
1	A	146	GLY	4.7
1	L	2	TYR	4.7
2	E	189	LEU	4.5
2	E	193	THR	4.4
1	L	1	SER	4.4

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Mol	Chain	Res	Type	RSRZ
3	C	303	THR	4.3
1	A	213	THR	4.2
2	E	137	ALA	4.1
2	E	194	TYR	3.8
2	E	136	ALA	3.6
2	B	30	LEU	3.4
1	A	145	PRO	3.4
2	B	191	THR	3.2
1	D	146	GLY	3.2
1	A	176	TYR	3.2
2	E	159	LEU	3.2
1	A	195	TYR	3.2
2	H	136	ALA	3.1
2	E	187	SER	3.1
1	D	1	SER	3.1
2	I	188	SER	3.1
2	I	191	THR	3.0
1	A	144	TYR	3.0
1	G	213	THR	3.0
2	I	184	VAL	2.9
3	C	304	ARG	2.8
1	G	2	TYR	2.8
1	D	148	VAL	2.8
1	A	107	LEU	2.8
1	A	143	PHE	2.8
1	A	159	VAL	2.8
2	E	192	GLN	2.7
2	E	213	PRO	2.7
2	E	138	LEU	2.7
1	L	189	TRP	2.7
2	I	187	SER	2.6
1	A	211	ALA	2.6
1	A	196	SER	2.6
1	D	16	GLY	2.6
2	E	214	LYS	2.6
1	D	199	VAL	2.6
1	A	1	SER	2.6
2	B	163	VAL	2.5
1	A	190	LYS	2.5
1	L	193	LYS	2.4
1	A	152	TRP	2.4
2	E	186	SER	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	28	ASN	2.3
1	A	191	SER	2.3
1	L	211	ALA	2.2
1	A	91	TRP	2.2
2	B	126	PRO	2.2
2	I	215	SER	2.2
1	D	189	TRP	2.2
1	D	176	TYR	2.2
1	L	187	GLU	2.1
1	A	148	VAL	2.1
2	I	125	ALA	2.1
2	I	213	PRO	2.1
1	A	110	LEU	2.1
2	H	174	GLY	2.1
2	I	185	PRO	2.1
3	P	318	TYR	2.1
1	L	213	THR	2.1
1	A	100	VAL	2.1
2	H	135	THR	2.1
2	H	215	SER	2.1
1	L	212	PRO	2.1
1	A	157	SER	2.1
2	H	194	TYR	2.0
1	A	182	LEU	2.0
1	D	213	THR	2.0
1	A	168	PRO	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 ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.