



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:00 PM GMT

PDB ID : 3VRL
Title : Crystal structure of BMJ4 p24 capsid protein in complex with A10F9 Fab
Authors : Gu, Y.; Cao, F.; Li, S.; Yuan, Y.A.; Xia, N.
Deposited on : 2012-04-12
Resolution : 3.20 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

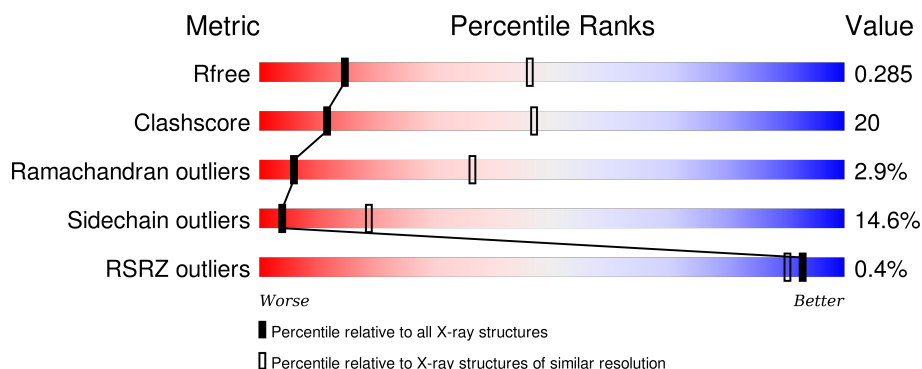
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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 $\leq 5\%$. 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	E	224	<div> <div>55%</div> <div>35%</div> <div>8%</div> <div>.</div> </div>
1	H	224	<div> <div>64%</div> <div>24%</div> <div>9%</div> <div>..</div> </div>
2	F	214	<div> <div>59%</div> <div>35%</div> <div>5%</div> </div>
2	L	214	<div> <div>59%</div> <div>35%</div> <div>6%</div> </div>
3	C	231	<div> <div>19%</div> <div>12%</div> <div>.</div> <div>68%</div> </div>

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Mol	Chain	Length	Quality of chain
3	D	231	<div> <div>18%</div> <div>12%</div> <div>2%</div> <div>68%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7806 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 A10F9 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	222	Total	C	N	O	S	0	0	0
			1669	1051	273	336	9			
1	E	222	Total	C	N	O	S	0	0	0
			1669	1051	273	336	9			

- Molecule 2 is a protein called A10F9 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	214	Total	C	N	O	S	0	0	0
			1660	1033	284	336	7			
2	F	214	Total	C	N	O	S	0	0	0
			1660	1033	284	336	7			

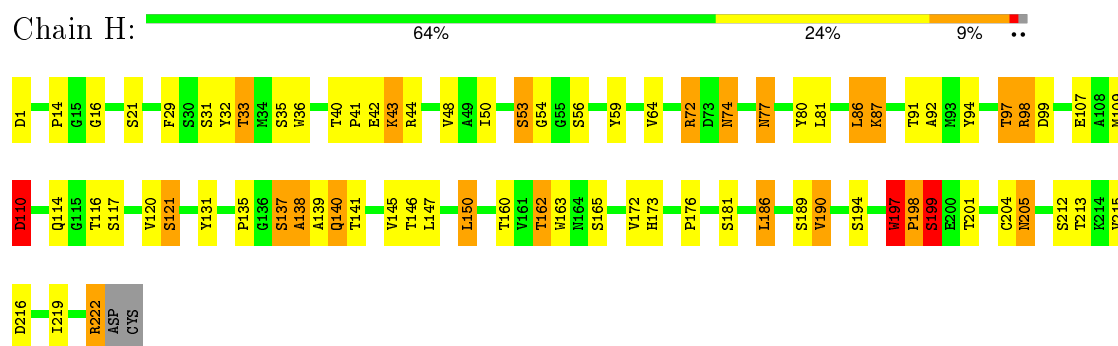
- Molecule 3 is a protein called Gag protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	73	Total	C	N	O	S	0	0	0
			574	358	99	111	6			
3	D	73	Total	C	N	O	S	0	0	0
			574	358	99	111	6			

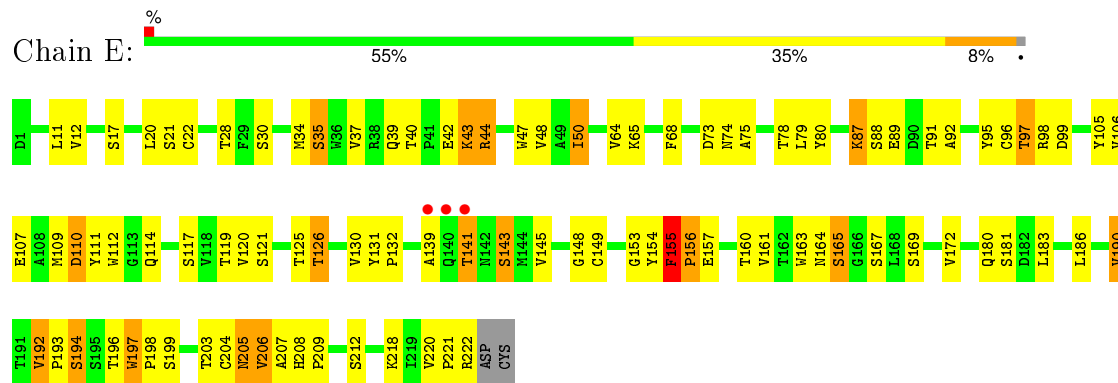
3 Residue-property plots

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.

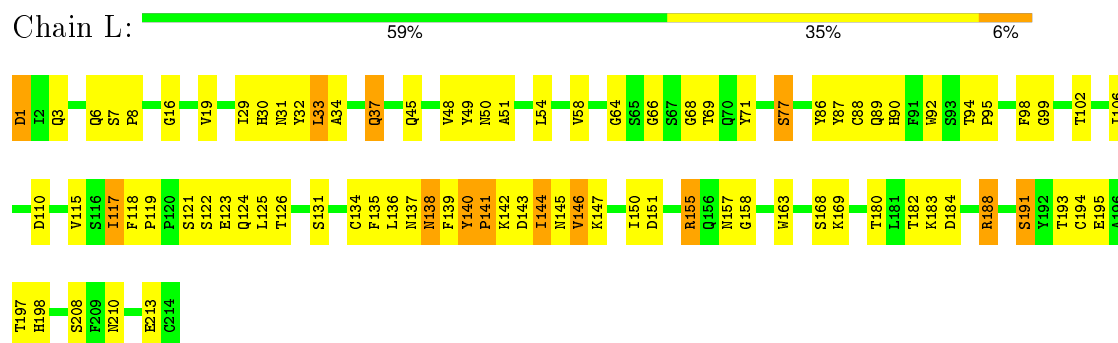
• Molecule 1: A10F9 Fab heavy chain



• Molecule 1: A10F9 Fab heavy chain

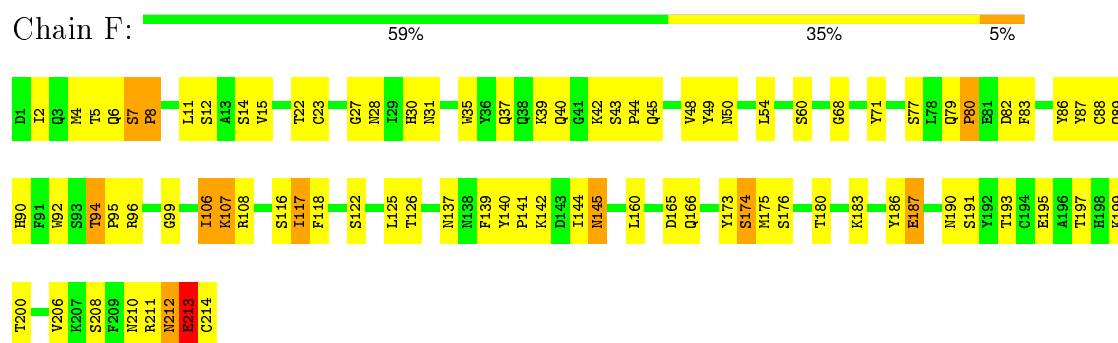


• Molecule 2: A10F9 Fab light chain



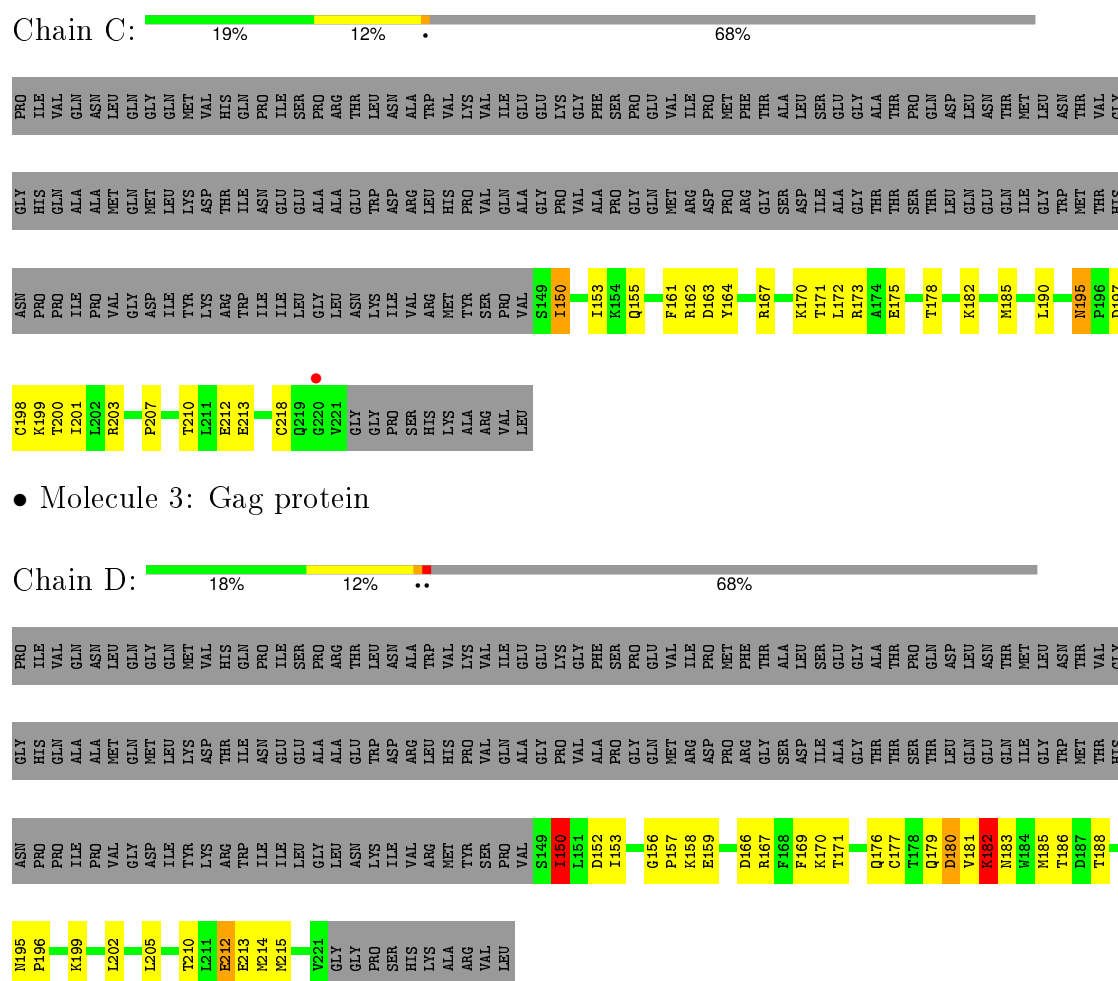
- Molecule 2: A10F9 Fab light chain

Chain F:



- Molecule 3: Gag protein

Chain C:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.11Å 124.06Å 150.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.02 – 3.20 37.02 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (37.02-3.20) 99.6 (37.02-3.20)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.82 (at 3.18Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.211 , 0.292 0.207 , 0.285	Depositor DCC
R_{free} test set	1339 reflections (5.36%)	DCC
Wilson B-factor (Å ²)	67.7	Xtriage
Anisotropy	0.024	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 18.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 26331 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7806	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.64% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	E	0.65	0/1710	0.83	0/2333
1	H	0.70	1/1710 (0.1%)	0.85	2/2333 (0.1%)
2	F	0.75	1/1699 (0.1%)	0.80	2/2306 (0.1%)
2	L	0.71	0/1699	0.79	0/2306
3	C	0.56	0/582	0.80	0/785
3	D	0.77	2/582 (0.3%)	0.79	1/785 (0.1%)
All	All	0.70	4/7982 (0.1%)	0.82	5/10848 (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
1	E	0	1
1	H	0	1
2	F	0	1
2	L	0	1
All	All	0	4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	214	CYS	C-O	13.93	1.49	1.23
3	D	182	LYS	CE-NZ	7.73	1.68	1.49
1	H	87	LYS	CD-CE	7.04	1.68	1.51
3	D	158	LYS	CE-NZ	6.56	1.65	1.49

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	186	LEU	CA-CB-CG	7.18	131.82	115.30
1	H	87	LYS	CD-CE-NZ	6.22	126.01	111.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	214	CYS	CA-C-O	-6.12	107.24	120.10
2	F	94	THR	C-N-CD	5.76	140.50	128.40
3	D	150	ILE	CB-CA-C	-5.05	101.50	111.60

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	155	PHE	Peptide
2	F	7	SER	Peptide
1	H	197	TRP	Peptide
2	L	140	TYR	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	E	1669	0	1633	72	0
1	H	1669	0	1631	78	0
2	F	1660	0	1585	55	0
2	L	1660	0	1587	71	0
3	C	574	0	570	30	0
3	D	574	0	570	19	0
All	All	7806	0	7576	308	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 20.

All (308) 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)
3:D:182:LYS:NZ	3:D:182:LYS:CE	1.68	1.55
1:H:53:SER:H	3:C:200:THR:HG21	1.14	1.07
2:F:137:ASN:HA	2:F:174:SER:HB3	1.38	1.05
2:L:94:THR:HG23	2:L:95:PRO:HD3	1.35	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:155:ARG:HG2	2:L:155:ARG:HH11	1.22	1.04
1:H:40:THR:HG22	1:H:42:GLU:H	1.26	1.01
1:H:162:THR:HG22	1:H:205:ASN:ND2	1.74	1.00
2:L:29:ILE:HG23	2:L:92:TRP:HB2	1.44	0.99
1:H:72:ARG:HD3	1:H:74:ASN:ND2	1.86	0.90
2:F:193:THR:HA	2:F:208:SER:HB3	1.55	0.89
1:H:97:THR:HG21	1:H:109:MET:CE	2.04	0.87
2:L:94:THR:CG2	2:L:95:PRO:HD3	2.04	0.86
1:E:64:VAL:HG13	1:E:68:PHE:HB2	1.58	0.85
2:L:137:ASN:HB3	2:L:138:ASN:ND2	1.91	0.84
1:H:42:GLU:O	1:H:43:LYS:HB2	1.77	0.84
2:L:94:THR:HG23	2:L:95:PRO:CD	2.09	0.83
1:H:197:TRP:HZ2	1:H:219:ILE:O	1.62	0.82
1:H:72:ARG:HD3	1:H:74:ASN:HD21	1.45	0.82
1:E:197:TRP:O	1:E:198:PRO:C	2.17	0.80
1:H:197:TRP:CZ2	1:H:219:ILE:O	2.34	0.80
2:L:155:ARG:HH11	2:L:155:ARG:CG	1.93	0.80
1:H:97:THR:HG21	1:H:109:MET:HE1	1.62	0.79
2:L:140:TYR:CD2	2:L:141:PRO:HD3	2.18	0.79
2:L:90:HIS:HD2	2:L:92:TRP:H	1.30	0.79
1:E:221:PRO:O	1:E:222:ARG:HB2	1.81	0.78
2:L:117:ILE:HD11	2:L:194:CYS:HB2	1.63	0.78
1:H:72:ARG:HH11	1:H:74:ASN:HD21	1.30	0.78
1:H:53:SER:N	3:C:200:THR:HG21	1.97	0.77
1:H:97:THR:CG2	1:H:109:MET:CE	2.61	0.77
2:L:29:ILE:CG2	2:L:92:TRP:HB2	2.15	0.77
1:H:162:THR:CG2	1:H:205:ASN:ND2	2.48	0.77
1:E:97:THR:HG21	1:E:109:MET:CE	2.14	0.77
1:H:40:THR:HG22	1:H:42:GLU:N	2.00	0.76
1:H:98:ARG:NH1	1:H:110:ASP:OD2	2.18	0.76
2:F:6:GLN:HE22	2:F:87:TYR:HA	1.51	0.75
1:E:165:SER:HA	1:E:205:ASN:ND2	2.02	0.75
1:H:162:THR:CG2	1:H:205:ASN:HD22	2.00	0.75
3:C:153:ILE:H	3:C:153:ILE:HD12	1.52	0.74
2:L:150:ILE:HD12	2:L:155:ARG:HD3	1.69	0.74
1:E:97:THR:CG2	1:E:109:MET:CE	2.66	0.74
2:L:155:ARG:HH12	2:L:158:GLY:H	1.37	0.73
1:H:162:THR:HG22	1:H:205:ASN:HD22	1.54	0.73
2:L:155:ARG:CZ	2:L:157:ASN:HB2	2.18	0.72
1:H:97:THR:CG2	1:H:109:MET:HE1	2.18	0.72
1:E:35:SER:HB3	1:E:50:ILE:HG23	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:139:PHE:CE2	2:F:174:SER:HA	2.26	0.71
1:E:148:GLY:HA2	1:E:163:TRP:CH2	2.26	0.71
2:F:144:ILE:HG22	2:F:145:ASN:N	2.06	0.70
2:L:151:ASP:HA	2:L:191:SER:HB3	1.74	0.69
1:E:163:TRP:CE3	1:E:204:CYS:HB3	2.28	0.68
1:E:97:THR:HG21	1:E:109:MET:SD	2.33	0.68
1:H:35:SER:HB2	1:H:109:MET:HE1	1.76	0.68
2:L:136:LEU:HD11	2:L:146:VAL:HG13	1.76	0.67
2:F:6:GLN:NE2	2:F:87:TYR:HA	2.10	0.66
1:E:194:SER:O	1:E:198:PRO:HD2	1.95	0.66
1:H:162:THR:HG22	1:H:205:ASN:HD21	1.57	0.66
1:E:197:TRP:O	1:E:199:SER:N	2.28	0.66
3:C:150:ILE:HD12	3:C:185:MET:CE	2.25	0.66
3:D:153:ILE:HD13	3:D:171:THR:HG21	1.76	0.66
1:H:21:SER:HB3	1:H:80:TYR:CD1	2.30	0.66
3:C:150:ILE:HG12	3:C:175:GLU:HG3	1.78	0.66
2:L:144:ILE:HG23	2:L:198:HIS:HB2	1.77	0.65
1:H:173:HIS:HE1	2:L:138:ASN:OD1	1.79	0.65
1:E:91:THR:HG23	1:E:119:THR:HA	1.77	0.65
2:L:188:ARG:NH1	2:L:188:ARG:HB3	2.11	0.65
1:H:145:VAL:HB	1:H:194:SER:HB3	1.80	0.64
1:H:107:GLU:OE1	2:L:50:ASN:ND2	2.29	0.64
2:L:137:ASN:HB3	2:L:138:ASN:HD22	1.61	0.64
1:H:33:THR:HG22	1:H:99:ASP:CG	2.18	0.64
1:H:222:ARG:HB2	1:H:222:ARG:HH11	1.62	0.64
1:H:42:GLU:O	1:H:43:LYS:CB	2.47	0.63
2:F:94:THR:HG23	2:F:95:PRO:HD3	1.79	0.63
1:E:207:ALA:O	1:E:209:PRO:HD3	1.99	0.63
2:L:136:LEU:HD11	2:L:146:VAL:CG1	2.29	0.62
3:D:153:ILE:CD1	3:D:171:THR:HG21	2.29	0.62
1:E:180:GLN:NE2	2:F:160:LEU:HD11	2.14	0.62
2:L:139:PHE:CZ	2:L:144:ILE:HD11	2.35	0.62
2:F:195:GLU:HG2	2:F:206:VAL:HG22	1.80	0.62
1:H:97:THR:HG21	1:H:109:MET:SD	2.40	0.62
2:F:31:ASN:HA	2:F:71:TYR:HE1	1.65	0.61
2:L:90:HIS:CD2	2:L:92:TRP:H	2.17	0.61
3:C:153:ILE:N	3:C:153:ILE:HD12	2.16	0.61
3:D:205:LEU:HD11	3:D:214:MET:HA	1.83	0.61
1:H:72:ARG:HH11	1:H:74:ASN:ND2	1.97	0.60
1:E:64:VAL:CG1	1:E:68:PHE:HB2	2.30	0.60
2:L:110:ASP:HA	2:L:140:TYR:HB3	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:35:SER:OG	1:E:109:MET:HE1	2.01	0.60
1:E:97:THR:CG2	1:E:109:MET:HE2	2.30	0.60
1:E:155:PHE:O	1:E:208:HIS:CE1	2.55	0.60
1:H:72:ARG:NH1	1:H:74:ASN:HD21	2.00	0.60
2:L:188:ARG:HH11	2:L:188:ARG:CG	2.14	0.60
3:C:150:ILE:CG1	3:C:175:GLU:HG3	2.33	0.59
2:L:142:LYS:HD3	2:L:142:LYS:O	2.02	0.59
2:L:6:GLN:HE21	2:L:99:GLY:HA3	1.67	0.59
2:F:2:ILE:HG23	2:F:27:GLY:H	1.66	0.59
1:H:40:THR:CG2	1:H:42:GLU:OE1	2.50	0.59
1:H:139:ALA:O	1:H:140:GLN:HB2	2.03	0.59
2:F:40:GLN:HE22	2:F:165:ASP:HB3	1.68	0.58
3:C:212:GLU:OE1	3:C:212:GLU:HA	2.01	0.58
2:L:122:SER:HA	2:L:125:LEU:HD12	1.85	0.58
2:F:2:ILE:HD12	2:F:27:GLY:HA3	1.84	0.58
3:D:166:ASP:O	3:D:170:LYS:HG3	2.04	0.58
3:D:159:GLU:OE2	3:D:167:ARG:HD3	2.04	0.57
3:C:150:ILE:HD12	3:C:185:MET:HE2	1.85	0.57
3:D:205:LEU:CD1	3:D:214:MET:HA	2.34	0.57
2:L:33:LEU:HD11	2:L:88:CYS:HB2	1.87	0.57
2:L:155:ARG:NH1	2:L:155:ARG:CG	2.62	0.56
1:E:165:SER:HA	1:E:205:ASN:HD21	1.69	0.56
2:F:37:GLN:HG3	2:F:86:TYR:CE2	2.39	0.56
1:H:86:LEU:HB3	1:H:120:VAL:HG21	1.87	0.56
2:L:115:VAL:HA	2:L:135:PHE:O	2.06	0.56
1:E:148:GLY:HA2	1:E:163:TRP:HH2	1.70	0.56
2:L:140:TYR:CD2	2:L:141:PRO:CD	2.88	0.55
2:L:29:ILE:HG22	2:L:32:TYR:HB2	1.86	0.55
2:L:188:ARG:HH11	2:L:188:ARG:CB	2.20	0.55
2:L:7:SER:HB2	2:L:8:PRO:HA	1.90	0.54
2:L:155:ARG:NH1	2:L:157:ASN:HB2	2.23	0.54
3:C:170:LYS:HA	3:C:173:ARG:HH21	1.72	0.54
2:F:31:ASN:HA	2:F:71:TYR:CE1	2.42	0.54
1:E:131:TYR:O	1:E:149:CYS:HB2	2.07	0.53
2:F:186:TYR:CE2	2:F:211:ARG:HD2	2.43	0.53
3:D:212:GLU:HA	3:D:212:GLU:OE2	2.07	0.53
2:L:6:GLN:HE22	2:L:87:TYR:HA	1.73	0.53
1:H:150:LEU:O	1:H:150:LEU:HD23	2.08	0.53
3:C:150:ILE:HA	3:C:153:ILE:HD13	1.91	0.53
2:L:8:PRO:O	2:L:102:THR:HG23	2.09	0.53
2:L:155:ARG:NH1	2:L:155:ARG:HG2	2.04	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:97:THR:HG21	1:E:109:MET:HE1	1.91	0.52
3:C:172:LEU:HD21	3:C:182:LYS:HB3	1.91	0.52
2:F:193:THR:HA	2:F:208:SER:CB	2.35	0.52
1:E:192:VAL:HG12	1:E:196:THR:OG1	2.09	0.52
2:L:140:TYR:O	2:L:198:HIS:CE1	2.62	0.52
1:H:135:PRO:HD3	1:H:147:LEU:CD2	2.39	0.52
1:H:40:THR:HG21	1:H:42:GLU:OE1	2.09	0.52
1:E:180:GLN:HE21	2:F:160:LEU:HD11	1.73	0.52
2:L:51:ALA:O	2:L:64:GLY:HA3	2.10	0.52
2:L:188:ARG:NH1	2:L:188:ARG:CB	2.73	0.52
2:F:37:GLN:NE2	2:F:39:LYS:HE2	2.25	0.52
2:F:90:HIS:O	2:F:96:ARG:HB2	2.11	0.51
2:L:188:ARG:HH11	2:L:188:ARG:HG3	1.76	0.51
2:L:123:GLU:O	2:L:126:THR:HB	2.10	0.51
2:F:106:ILE:O	2:F:166:GLN:NE2	2.41	0.51
2:L:54:LEU:HD13	2:L:58:VAL:HG12	1.93	0.51
1:E:155:PHE:HB3	1:E:156:PRO:CD	2.41	0.51
2:F:7:SER:HB3	2:F:8:PRO:HD2	1.93	0.51
1:E:39:GLN:HA	1:E:44:ARG:O	2.12	0.50
1:H:41:PRO:HD3	1:H:91:THR:O	2.11	0.50
1:H:197:TRP:CG	1:H:198:PRO:HD3	2.46	0.50
1:H:16:GLY:O	1:H:86:LEU:HD23	2.12	0.50
2:L:48:VAL:HG12	2:L:49:TYR:N	2.26	0.50
1:E:22:CYS:SG	1:E:96:CYS:CB	2.99	0.50
3:D:182:LYS:NZ	3:D:182:LYS:CD	2.64	0.50
1:H:189:SER:HB3	2:L:135:PHE:CE1	2.47	0.50
2:L:37:GLN:HG3	2:L:86:TYR:CE2	2.46	0.50
3:C:210:THR:OG1	3:C:213:GLU:HG3	2.11	0.50
2:F:144:ILE:CG2	2:F:145:ASN:N	2.75	0.50
3:D:210:THR:H	3:D:213:GLU:HG3	1.77	0.50
1:H:33:THR:CG2	1:H:99:ASP:OD2	2.60	0.49
2:L:29:ILE:HG23	2:L:92:TRP:CB	2.28	0.49
1:E:20:LEU:O	1:E:80:TYR:HA	2.12	0.49
1:H:72:ARG:CD	1:H:74:ASN:HD21	2.22	0.49
1:H:213:THR:HG22	1:H:215:VAL:HG23	1.93	0.49
2:F:6:GLN:HG3	2:F:99:GLY:HA3	1.95	0.49
1:E:125:THR:HA	1:E:155:PHE:HB3	1.95	0.49
1:H:36:TRP:NE1	1:H:81:LEU:HB2	2.28	0.49
1:H:53:SER:OG	3:C:200:THR:CG2	2.61	0.49
2:F:79:GLN:O	2:F:82:ASP:N	2.45	0.49
1:H:35:SER:CB	1:H:109:MET:HE1	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:42:GLU:O	1:E:43:LYS:HB2	2.12	0.48
1:E:88:SER:HA	1:E:120:VAL:O	2.13	0.48
3:C:150:ILE:HD11	3:C:175:GLU:HG3	1.95	0.48
2:F:79:GLN:O	2:F:80:PRO:C	2.51	0.48
2:L:89:GLN:HB2	2:L:98:PHE:CD2	2.47	0.48
3:C:155:GLN:HB3	3:C:195:ASN:H	1.77	0.48
3:C:161:PHE:HB2	3:C:218:CYS:HB3	1.94	0.48
3:C:150:ILE:CD1	3:C:175:GLU:HG3	2.44	0.48
1:E:87:LYS:HG3	1:E:89:GLU:HB2	1.94	0.48
1:H:163:TRP:CZ3	1:H:204:CYS:HB3	2.48	0.48
1:E:37:VAL:HG13	1:E:47:TRP:HA	1.95	0.48
2:L:33:LEU:HG	2:L:34:ALA:N	2.27	0.48
2:L:29:ILE:O	2:L:29:ILE:HG22	2.14	0.47
2:F:183:LYS:O	2:F:186:TYR:HB3	2.13	0.47
1:E:11:LEU:HD12	1:E:12:VAL:N	2.29	0.47
3:D:179:GLN:O	3:D:180:ASP:C	2.53	0.47
1:E:153:GLY:HA2	1:E:183:LEU:HB3	1.96	0.47
1:H:135:PRO:HD3	1:H:147:LEU:HD22	1.95	0.47
3:D:181:VAL:O	3:D:185:MET:HG3	2.15	0.47
1:E:98:ARG:HG2	1:E:99:ASP:N	2.30	0.47
1:H:48:VAL:HG13	1:H:64:VAL:HG21	1.97	0.47
1:H:165:SER:H	1:H:205:ASN:HD21	1.63	0.47
1:H:36:TRP:CD1	1:H:81:LEU:HB2	2.49	0.47
2:F:7:SER:HB3	2:F:8:PRO:CD	2.45	0.46
2:F:117:ILE:HG12	2:F:118:PHE:N	2.30	0.46
2:L:16:GLY:O	2:L:77:SER:HA	2.15	0.46
2:L:193:THR:HA	2:L:208:SER:HB3	1.97	0.46
2:F:144:ILE:HG22	2:F:145:ASN:H	1.80	0.46
1:E:130:VAL:HG21	1:E:206:VAL:HG11	1.96	0.46
1:E:73:ASP:OD1	1:E:75:ALA:HB3	2.15	0.46
2:L:1:ASP:HB3	2:L:3:GLN:NE2	2.31	0.46
2:F:140:TYR:CG	2:F:141:PRO:HA	2.50	0.46
2:F:83:PHE:CE2	2:F:106:ILE:HB	2.51	0.46
2:F:183:LYS:O	2:F:187:GLU:HG3	2.16	0.46
1:E:197:TRP:C	1:E:199:SER:N	2.67	0.46
2:L:139:PHE:CE2	2:L:144:ILE:HD11	2.51	0.46
1:H:131:TYR:CE1	2:L:124:GLN:HG3	2.51	0.46
3:D:150:ILE:HG13	3:D:150:ILE:H	1.47	0.45
2:L:29:ILE:O	2:L:30:HIS:C	2.54	0.45
1:E:145:VAL:HG23	1:E:194:SER:HB3	1.97	0.45
3:D:202:LEU:CD2	3:D:214:MET:HB3	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:97:THR:CG2	1:H:109:MET:HE2	2.42	0.45
3:C:210:THR:H	3:C:213:GLU:HG3	1.82	0.45
2:F:212:ASN:OD1	2:F:212:ASN:N	2.48	0.45
1:E:105:TYR:O	1:E:107:GLU:HG3	2.16	0.45
2:L:147:LYS:HB3	2:L:195:GLU:HB2	1.99	0.45
2:F:139:PHE:CE1	2:F:144:ILE:HG13	2.52	0.45
1:H:172:VAL:HG22	1:H:190:VAL:HG13	1.98	0.45
3:C:153:ILE:H	3:C:153:ILE:CD1	2.27	0.45
1:H:32:TYR:O	1:H:72:ARG:NH2	2.50	0.45
1:H:197:TRP:O	1:H:199:SER:N	2.50	0.45
3:C:175:GLU:O	3:C:182:LYS:HE2	2.16	0.45
1:E:130:VAL:CG2	1:E:206:VAL:HG11	2.46	0.45
1:H:197:TRP:CD1	1:H:198:PRO:HD3	2.52	0.45
1:E:180:GLN:N	1:E:183:LEU:O	2.49	0.45
1:H:54:GLY:HA3	3:C:197:ASP:CG	2.37	0.45
3:D:150:ILE:C	3:D:152:ASP:H	2.20	0.44
1:H:53:SER:OG	3:C:200:THR:HG21	2.18	0.44
1:E:34:MET:HB3	1:E:79:LEU:HD22	2.00	0.44
2:F:31:ASN:O	2:F:50:ASN:HA	2.18	0.44
2:F:30:HIS:CD2	2:F:92:TRP:CZ3	3.06	0.44
1:H:44:ARG:HB2	1:H:44:ARG:HE	1.56	0.44
1:E:47:TRP:HZ2	1:E:50:ILE:HG12	1.81	0.44
1:H:40:THR:HG23	1:H:41:PRO:HD2	1.99	0.44
2:F:165:ASP:O	2:F:166:GLN:C	2.54	0.44
3:D:195:ASN:HB2	3:D:196:PRO:HD2	2.00	0.44
1:E:22:CYS:O	1:E:78:THR:HA	2.18	0.44
1:E:143:SER:O	1:E:193:PRO:HA	2.18	0.44
1:E:155:PHE:CB	1:E:156:PRO:CD	2.96	0.43
1:E:95:TYR:HE2	2:F:43:SER:HB3	1.83	0.43
2:L:66:GLY:HA3	2:L:71:TYR:HA	2.00	0.43
1:H:29:PHE:CE2	1:H:72:ARG:NH1	2.86	0.43
2:L:118:PHE:HA	2:L:119:PRO:HD3	1.81	0.43
1:E:79:LEU:HA	1:E:79:LEU:HD12	1.81	0.43
3:C:150:ILE:HA	3:C:171:THR:HG21	2.00	0.43
1:H:176:PRO:HG2	2:L:163:TRP:O	2.19	0.43
1:H:54:GLY:HA3	3:C:197:ASP:OD1	2.19	0.43
2:F:142:LYS:HA	2:F:173:TYR:CD2	2.54	0.43
1:E:163:TRP:CZ3	1:E:204:CYS:HB3	2.53	0.43
1:H:98:ARG:HD3	1:H:110:ASP:OD2	2.19	0.43
1:E:180:GLN:OE1	1:E:180:GLN:HA	2.19	0.43
2:F:90:HIS:HD2	2:F:92:TRP:H	1.67	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:40:THR:OG1	1:E:44:ARG:HD2	2.19	0.43
1:E:154:TYR:CG	1:E:154:TYR:O	2.72	0.43
3:C:198:CYS:O	3:C:201:ILE:HG22	2.19	0.42
1:E:197:TRP:CH2	1:E:221:PRO:HB3	2.54	0.42
1:E:97:THR:CG2	1:E:109:MET:HE1	2.45	0.42
1:E:39:GLN:O	1:E:92:ALA:HB1	2.19	0.42
1:E:203:THR:HA	1:E:218:LYS:HA	2.00	0.42
1:E:125:THR:HG22	1:E:126:THR:N	2.34	0.42
1:H:197:TRP:CH2	1:H:219:ILE:O	2.73	0.42
2:L:188:ARG:CZ	2:L:188:ARG:HB3	2.48	0.42
1:H:131:TYR:HB3	2:L:121:SER:OG	2.19	0.42
1:E:110:ASP:HB3	1:E:111:TYR:HD1	1.85	0.42
2:F:4:MET:HE2	2:F:23:CYS:SG	2.58	0.42
2:F:40:GLN:HE22	2:F:165:ASP:CB	2.32	0.42
2:L:48:VAL:CG1	2:L:49:TYR:N	2.82	0.42
2:L:31:ASN:O	2:L:50:ASN:HA	2.20	0.42
1:E:155:PHE:HB3	1:E:156:PRO:HD2	2.01	0.42
2:F:37:GLN:HE22	2:F:39:LYS:HE2	1.85	0.42
1:H:50:ILE:HG22	1:H:59:TYR:HB3	2.00	0.42
2:F:28:ASN:OD1	2:F:28:ASN:C	2.58	0.42
3:D:183:ASN:O	3:D:186:THR:HB	2.20	0.42
1:H:92:ALA:HB3	1:H:94:TYR:CE1	2.55	0.42
2:F:23:CYS:HB3	2:F:35:TRP:CH2	2.55	0.42
2:F:4:MET:CE	2:F:23:CYS:SG	3.08	0.42
1:E:172:VAL:HG22	1:E:190:VAL:HG12	2.02	0.42
1:H:172:VAL:CG1	1:H:173:HIS:N	2.83	0.42
2:F:40:GLN:C	2:F:42:LYS:H	2.23	0.42
1:H:14:PRO:HD3	1:H:121:SER:C	2.40	0.42
2:F:35:TRP:CZ3	2:F:88:CYS:HB3	2.55	0.42
1:E:35:SER:OG	1:E:109:MET:CE	2.66	0.41
1:E:110:ASP:HB3	1:E:111:TYR:CD1	2.55	0.41
1:E:209:PRO:O	1:E:212:SER:N	2.36	0.41
2:F:40:GLN:OE1	2:F:165:ASP:OD2	2.37	0.41
2:L:7:SER:HA	2:L:8:PRO:C	2.41	0.41
2:F:37:GLN:O	2:F:44:PRO:HA	2.21	0.41
2:L:182:THR:HG22	2:L:184:ASP:H	1.84	0.41
2:L:210:ASN:O	2:L:213:GLU:HG2	2.19	0.41
3:D:202:LEU:HD22	3:D:214:MET:HB3	2.01	0.41
1:E:164:ASN:HB2	1:E:167:SER:OG	2.21	0.41
3:C:172:LEU:HG	3:C:182:LYS:HD3	2.03	0.41
1:H:33:THR:HG22	1:H:99:ASP:OD2	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:33:THR:HG21	1:H:99:ASP:OD2	2.20	0.41
1:E:125:THR:HA	1:E:156:PRO:HD2	2.03	0.41
1:H:77:ASN:HD22	1:H:77:ASN:HA	1.63	0.41
1:E:42:GLU:O	1:E:43:LYS:CB	2.69	0.41
1:E:30:SER:HA	1:E:74:ASN:OD1	2.21	0.41
1:H:33:THR:OG1	3:C:200:THR:HG22	2.21	0.41
2:L:32:TYR:CE1	3:C:207:PRO:HG3	2.55	0.41
2:F:48:VAL:HG12	2:F:49:TYR:N	2.36	0.41
1:E:112:TRP:CD1	1:E:112:TRP:N	2.89	0.40
2:F:210:ASN:O	2:F:213:GLU:HG3	2.21	0.40
2:F:83:PHE:CZ	2:F:106:ILE:HB	2.56	0.40
2:F:107:LYS:HA	2:F:140:TYR:OH	2.21	0.40
3:C:164:TYR:CE2	3:C:190:LEU:HD12	2.57	0.40
3:D:150:ILE:C	3:D:152:ASP:N	2.75	0.40
1:H:204:CYS:O	1:H:216:ASP:HA	2.21	0.40
1:E:132:PRO:HA	1:E:149:CYS:CB	2.52	0.40
1:H:137:SER:O	1:H:138:ALA:HB2	2.22	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	E	220/224 (98%)	188 (86%)	22 (10%)	10 (4%)	3	24
1	H	220/224 (98%)	194 (88%)	17 (8%)	9 (4%)	3	27
2	F	212/214 (99%)	186 (88%)	21 (10%)	5 (2%)	7	43
2	L	212/214 (99%)	190 (90%)	19 (9%)	3 (1%)	14	57
3	C	71/231 (31%)	69 (97%)	2 (3%)	0	100	100
3	D	71/231 (31%)	66 (93%)	3 (4%)	2 (3%)	6	37
All	All	1006/1338 (75%)	893 (89%)	84 (8%)	29 (3%)	6	36

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	43	LYS
1	H	110	ASP
1	H	138	ALA
1	H	181	SER
1	H	197	TRP
2	L	141	PRO
1	E	156	PRO
1	E	165	SER
2	F	8	PRO
1	H	199	SER
2	L	77	SER
1	E	181	SER
2	F	68	GLY
3	D	156	GLY
1	H	137	SER
1	E	141	THR
2	F	77	SER
1	H	140	GLN
1	E	139	ALA
1	E	43	LYS
1	E	65	LYS
1	E	106	VAL
2	L	68	GLY
2	F	213	GLU
3	D	157	PRO
1	E	155	PHE
1	E	197	TRP
1	H	198	PRO
2	F	80	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	E	191/193 (99%)	163 (85%)	28 (15%)	4 18

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	191/193 (99%)	162 (85%)	29 (15%)	3	17
2	F	188/188 (100%)	157 (84%)	31 (16%)	3	13
2	L	188/188 (100%)	165 (88%)	23 (12%)	6	27
3	C	64/197 (32%)	56 (88%)	8 (12%)	6	26
3	D	64/197 (32%)	54 (84%)	10 (16%)	3	15
All	All	886/1156 (77%)	757 (85%)	129 (15%)	4	19

All (129) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	1	ASP
1	H	31	SER
1	H	33	THR
1	H	53	SER
1	H	56	SER
1	H	72	ARG
1	H	74	ASN
1	H	77	ASN
1	H	86	LEU
1	H	87	LYS
1	H	97	THR
1	H	98	ARG
1	H	110	ASP
1	H	114	GLN
1	H	116	THR
1	H	117	SER
1	H	121	SER
1	H	141	THR
1	H	146	THR
1	H	150	LEU
1	H	160	THR
1	H	162	THR
1	H	186	LEU
1	H	190	VAL
1	H	199	SER
1	H	201	THR
1	H	205	ASN
1	H	212	SER
1	H	222	ARG
2	L	1	ASP

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Mol	Chain	Res	Type
2	L	19	VAL
2	L	33	LEU
2	L	37	GLN
2	L	45	GLN
2	L	69	THR
2	L	106	ILE
2	L	117	ILE
2	L	131	SER
2	L	134	CYS
2	L	138	ASN
2	L	143	ASP
2	L	144	ILE
2	L	145	ASN
2	L	146	VAL
2	L	155	ARG
2	L	168	SER
2	L	169	LYS
2	L	180	THR
2	L	183	LYS
2	L	188	ARG
2	L	191	SER
2	L	197	THR
3	C	150	ILE
3	C	162	ARG
3	C	163	ASP
3	C	167	ARG
3	C	178	THR
3	C	195	ASN
3	C	199	LYS
3	C	203	ARG
1	E	17	SER
1	E	21	SER
1	E	28	THR
1	E	35	SER
1	E	44	ARG
1	E	48	VAL
1	E	50	ILE
1	E	87	LYS
1	E	97	THR
1	E	110	ASP
1	E	114	GLN
1	E	117	SER

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Mol	Chain	Res	Type
1	E	121	SER
1	E	126	THR
1	E	141	THR
1	E	143	SER
1	E	155	PHE
1	E	157	GLU
1	E	160	THR
1	E	161	VAL
1	E	169	SER
1	E	186	LEU
1	E	190	VAL
1	E	192	VAL
1	E	194	SER
1	E	205	ASN
1	E	206	VAL
1	E	220	VAL
2	F	5	THR
2	F	11	LEU
2	F	12	SER
2	F	14	SER
2	F	15	VAL
2	F	22	THR
2	F	45	GLN
2	F	54	LEU
2	F	60	SER
2	F	89	GLN
2	F	106	ILE
2	F	107	LYS
2	F	108	ARG
2	F	116	SER
2	F	117	ILE
2	F	122	SER
2	F	125	LEU
2	F	126	THR
2	F	145	ASN
2	F	174	SER
2	F	175	MET
2	F	176	SER
2	F	180	THR
2	F	187	GLU
2	F	190	ASN
2	F	191	SER

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Mol	Chain	Res	Type
2	F	197	THR
2	F	199	LYS
2	F	200	THR
2	F	212	ASN
2	F	213	GLU
3	D	150	ILE
3	D	169	PHE
3	D	176	GLN
3	D	177	CYS
3	D	180	ASP
3	D	182	LYS
3	D	188	THR
3	D	199	LYS
3	D	212	GLU
3	D	215	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	74	ASN
1	H	77	ASN
1	H	173	HIS
1	H	205	ASN
2	L	3	GLN
2	L	6	GLN
2	L	30	HIS
2	L	89	GLN
2	L	90	HIS
2	L	138	ASN
3	C	195	ASN
1	E	205	ASN
1	E	208	HIS
2	F	6	GLN
2	F	30	HIS
2	F	37	GLN
2	F	40	GLN
2	F	76	ASN
2	F	90	HIS
2	F	138	ASN
2	F	145	ASN
2	F	161	ASN
2	F	190	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	E	222/224 (99%)	-0.20	3 (1%) 78 65	51, 70, 110, 127	0
1	H	222/224 (99%)	-0.38	0 100 100	44, 59, 92, 110	0
2	F	214/214 (100%)	-0.29	0 100 100	52, 69, 85, 121	0
2	L	214/214 (100%)	-0.33	0 100 100	47, 66, 85, 118	0
3	C	73/231 (31%)	-0.13	1 (1%) 78 65	50, 73, 97, 108	0
3	D	73/231 (31%)	-0.24	0 100 100	56, 79, 106, 114	0
All	All	1018/1338 (76%)	-0.29	4 (0%) 93 90	44, 68, 97, 127	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	141	THR	4.1
3	C	220	GLY	2.2
1	E	139	ALA	2.1
1	E	140	GLN	2.1

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 carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.