



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2022 – 02:07 PM EDT

PDB ID : 7REW
Title : Crystal Structure of IL-13 in complex with MMAb3 Fab
Authors : Sudom, A.; Min, X.
Deposited on : 2021-07-13
Resolution : 2.10 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 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 : 2.28.1
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) : 2.28.1

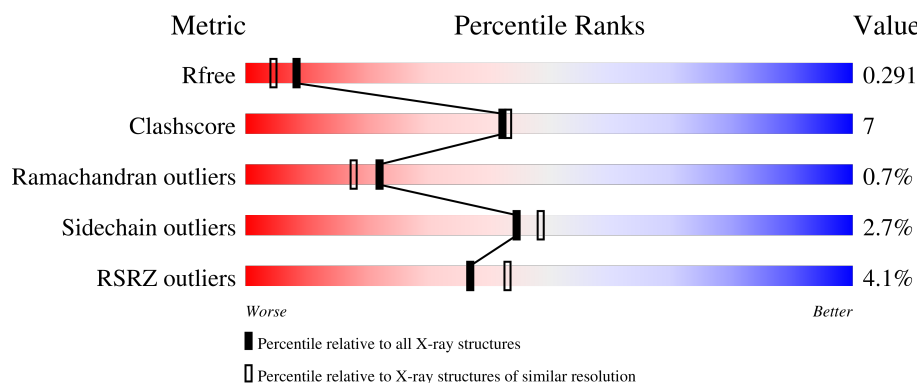
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	227	<div> <div>6%</div> <div>74%</div> <div>15%</div> <div>11%</div> </div>
1	H	227	<div> <div>3%</div> <div>74%</div> <div>12%</div> <div>13%</div> </div>
2	B	212	<div> <div>%</div> <div>78%</div> <div>19%</div> <div>..</div> </div>
2	L	212	<div> <div>4%</div> <div>80%</div> <div>17%</div> <div>.</div> </div>
3	G	112	<div> <div>4%</div> <div>73%</div> <div>8%</div> <div>18%</div> </div>

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Mol	Chain	Length	Quality of chain
3	I	112	<div><div></div><div>4%</div><div>66%</div><div>16%</div><div>•</div><div>17%</div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7712 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 anti-cyno interleukin 13 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	197	Total	C	N	O	S	0	0	0
			1492	951	245	289	7			
1	A	203	Total	C	N	O	S	0	0	0
			1529	975	250	297	7			

- Molecule 2 is a protein called anti-cyno interleukin 13 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	206	Total	C	N	O	S	0	0	0
			1557	979	256	317	5			
2	B	208	Total	C	N	O	S	0	0	0
			1570	986	258	321	5			

- Molecule 3 is a protein called IL13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	92	Total	C	N	O	S	0	0	0
			716	455	123	132	6			
3	I	93	Total	C	N	O	S	0	0	0
			717	456	124	131	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	2	GLY	-	expression tag	UNP Q0PW92
I	2	GLY	-	expression tag	UNP Q0PW92

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	29	Total	O	0	0
			29	29		

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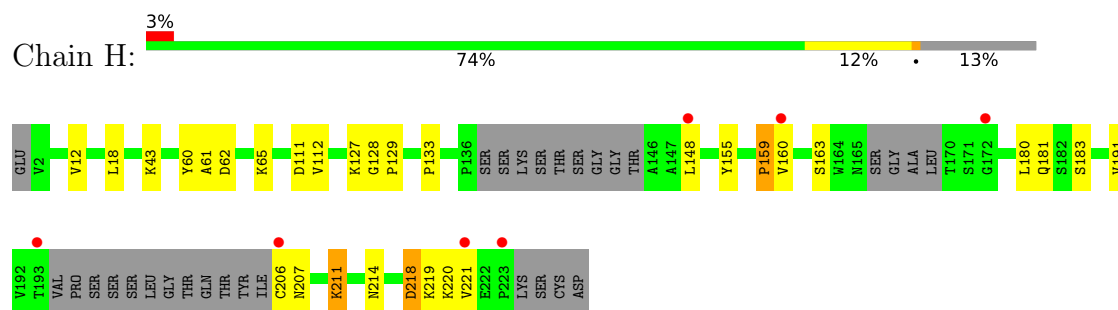
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	38	Total 38	O 38	0	0
4	A	29	Total 29	O 29	0	0
4	B	20	Total 20	O 20	0	0
4	G	10	Total 10	O 10	0	0
4	I	5	Total 5	O 5	0	0

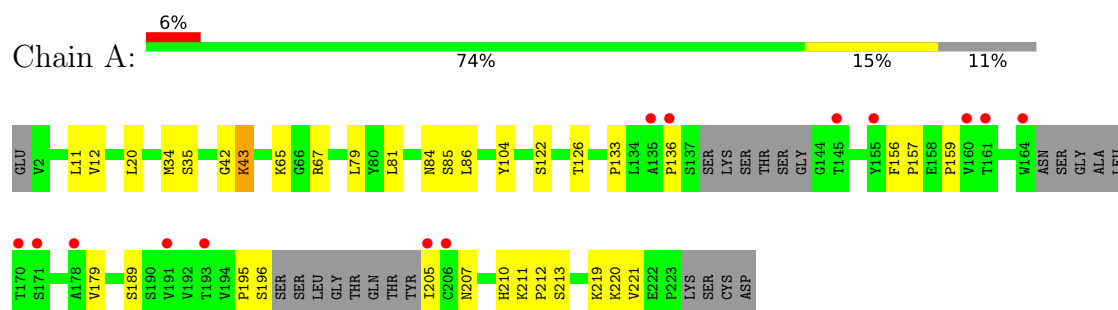
3 Residue-property plots [i](#)

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.

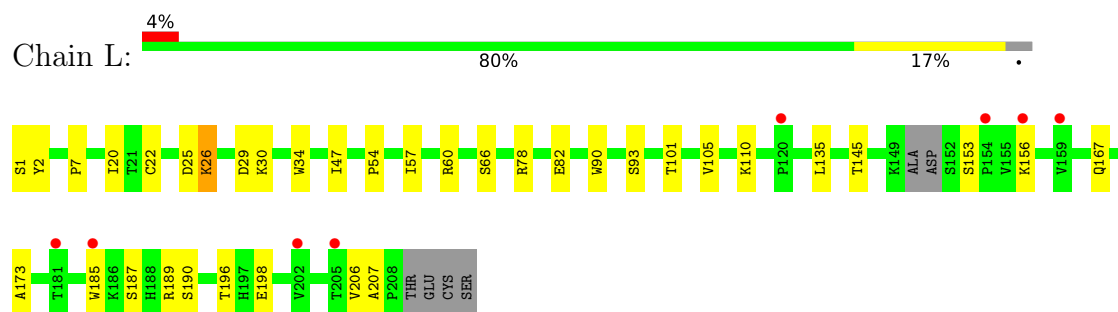
- Molecule 1: anti-cyno interleukin 13 Fab heavy chain



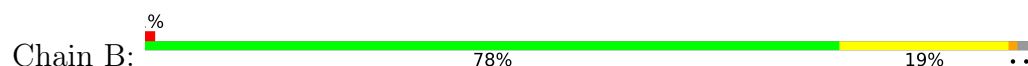
- Molecule 1: anti-cyno interleukin 13 Fab heavy chain

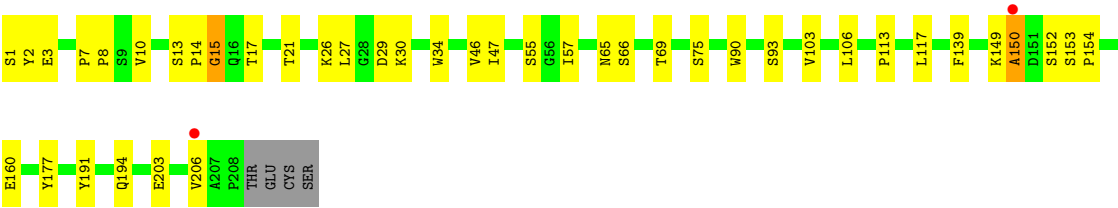


- Molecule 2: anti-cyno interleukin 13 Fab light chain



- Molecule 2: anti-cyno interleukin 13 Fab light chain

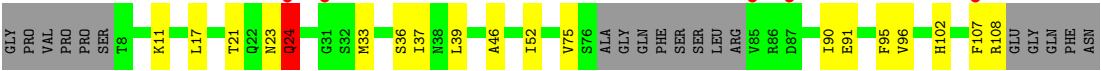




● Molecule 3: IL13



● Molecule 3: IL13



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	57.79Å 74.35Å 76.53Å 96.85° 109.89° 112.49°	Depositor
Resolution (Å)	35.96 – 2.10 35.96 – 2.10	Depositor EDS
% Data completeness (in resolution range)	87.9 (35.96-2.10) 87.9 (35.96-2.10)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.97 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.14 _3260	Depositor
R, R_{free}	0.245 , 0.291 0.245 , 0.291	Depositor DCC
R_{free} test set	2617 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	39.0	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 39.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.022 for h,-h-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7712	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.64% 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, 0.333 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	0.26	0/1568	0.47	0/2136
1	H	0.27	0/1530	0.50	0/2083
2	B	0.25	0/1612	0.47	1/2202 (0.0%)
2	L	0.28	0/1598	0.51	1/2181 (0.0%)
3	G	0.25	0/723	0.44	0/972
3	I	0.28	0/725	0.46	0/978
All	All	0.27	0/7756	0.48	2/10552 (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
3	I	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	189	ARG	NE-CZ-NH1	-8.92	115.84	120.30
2	B	152	SER	C-N-CA	6.34	137.54	121.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	I	23	ASN	Peptide
3	I	24	GLN	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	1529	0	1475	23	0
1	H	1492	0	1434	21	1
2	B	1570	0	1511	24	0
2	L	1557	0	1501	22	0
3	G	716	0	740	7	0
3	I	717	0	741	12	1
4	A	29	0	0	0	0
4	B	20	0	0	0	0
4	G	10	0	0	0	0
4	H	29	0	0	0	0
4	I	5	0	0	0	0
4	L	38	0	0	0	0
All	All	7712	0	7402	104	1

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 7.

All (104) 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:G:22:GLN:NE2	3:G:24:GLN:O	1.90	1.04
1:A:211:LYS:HG3	1:A:212:PRO:HD3	1.64	0.79
1:H:62:ASP:HA	1:H:65:LYS:HD2	1.63	0.78
1:A:179:VAL:HG11	2:B:160:GLU:HB3	1.69	0.74
3:I:37:ILE:HG12	3:I:46:ALA:HB1	1.72	0.69
2:L:1:SER:HA	2:L:30:LYS:HE3	1.75	0.69
1:A:220:LYS:HD2	1:A:221:VAL:H	1.59	0.68
2:L:1:SER:OG	2:L:2:TYR:N	2.32	0.61
1:A:20:LEU:HD12	1:A:81:LEU:HD23	1.83	0.61
3:I:24:GLN:NE2	3:I:24:GLN:HA	2.16	0.60
2:B:117:LEU:HD13	2:B:206:VAL:HG23	1.84	0.60
1:H:60:TYR:HB2	1:H:65:LYS:HG2	1.85	0.59
3:I:24:GLN:HA	3:I:24:GLN:HE21	1.68	0.58
1:H:133:PRO:HB3	1:H:219:LYS:HZ1	1.68	0.57
2:L:185:TRP:HH2	2:L:206:VAL:HG12	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:211:LYS:O	2:B:26:LYS:NZ	2.37	0.57
2:L:26:LYS:HD2	2:L:29:ASP:HB2	1.86	0.57
2:B:1:SER:OG	2:B:2:TYR:N	2.38	0.57
1:H:211:LYS:HD3	1:H:214:ASN:HA	1.87	0.56
1:H:211:LYS:HA	1:H:214:ASN:H	1.70	0.56
3:G:22:GLN:O	3:G:22:GLN:HG3	2.05	0.56
2:B:46:VAL:HA	2:B:57:ILE:HD13	1.89	0.55
1:A:210:HIS:CD2	1:A:212:PRO:HD2	2.42	0.55
2:L:185:TRP:CH2	2:L:206:VAL:HG12	2.41	0.54
2:L:7:PRO:O	2:L:101:THR:HG22	2.07	0.54
2:B:34:TRP:HB2	2:B:47:ILE:HB	1.91	0.52
2:L:60:ARG:CZ	2:L:78:ARG:HD3	2.39	0.51
3:I:17:LEU:O	3:I:21:THR:HG23	2.11	0.51
1:H:220:LYS:HG3	1:H:221:VAL:H	1.77	0.49
2:L:20:ILE:HG23	2:L:101:THR:HG21	1.94	0.49
3:G:40:THR:OG1	3:G:41:ALA:N	2.43	0.49
1:A:12:VAL:HG11	1:A:86:LEU:HD13	1.94	0.49
2:B:13:SER:O	2:B:15:GLY:N	2.41	0.49
2:L:26:LYS:HE3	2:L:30:LYS:HE2	1.95	0.49
2:B:113:PRO:HB3	2:B:139:PHE:HB3	1.95	0.49
1:H:127:LYS:NZ	1:H:128:GLY:O	2.45	0.49
2:L:110:LYS:HD2	2:L:198:GLU:HG3	1.96	0.48
3:I:33:MET:SD	3:I:91:GLU:HG2	2.54	0.48
3:I:52:ILE:HD13	3:I:75:VAL:HG23	1.95	0.48
3:I:24:GLN:NE2	3:I:24:GLN:CA	2.77	0.47
3:I:90:ILE:HD11	3:I:95:PHE:HD2	1.77	0.47
2:B:1:SER:HA	2:B:30:LYS:HE3	1.95	0.47
1:H:191:VAL:HG21	2:L:135:LEU:HD13	1.96	0.47
2:B:194:GLN:NE2	2:B:203:GLU:OE2	2.36	0.47
1:A:42:GLY:H	1:A:43:LYS:HE2	1.80	0.47
2:L:82:GLU:HG3	2:L:105:VAL:HG23	1.96	0.47
2:B:150:ALA:HA	2:B:191:TYR:CE1	2.50	0.47
1:A:159:PRO:O	1:A:210:HIS:HD2	1.98	0.46
2:L:90:TRP:CZ2	2:L:93:SER:HA	2.50	0.46
2:L:190:SER:OG	2:L:207:ALA:HB2	2.15	0.46
1:A:67:ARG:CG	1:A:85:SER:HB2	2.46	0.46
2:B:17:THR:HG22	2:B:75:SER:HA	1.97	0.46
2:L:54:PRO:HG2	2:L:57:ILE:HG13	1.98	0.46
1:H:12:VAL:HG21	1:H:18:LEU:HG	1.98	0.45
2:L:190:SER:HG	2:L:207:ALA:HB2	1.81	0.45
1:A:210:HIS:CE1	1:A:213:SER:HG	2.35	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:21:THR:HG23	2:B:69:THR:HG23	1.98	0.45
2:L:82:GLU:CG	2:L:105:VAL:H	2.29	0.45
1:A:34:MET:HB3	1:A:79:LEU:HD22	1.98	0.45
1:A:67:ARG:HG2	1:A:84:ASN:O	2.17	0.45
1:H:43:LYS:HE3	1:H:43:LYS:HB3	1.87	0.45
1:H:133:PRO:CB	1:H:219:LYS:HZ1	2.30	0.44
3:I:37:ILE:HD12	3:I:37:ILE:HA	1.88	0.44
2:B:10:VAL:O	2:B:103:VAL:HA	2.18	0.44
2:B:26:LYS:HB3	2:B:29:ASP:OD1	2.18	0.43
1:H:180:LEU:HD23	1:H:181:GLN:O	2.19	0.43
1:A:189:SER:HB3	2:B:177:TYR:OH	2.18	0.43
2:B:27:LEU:O	2:B:27:LEU:HD13	2.18	0.43
2:L:82:GLU:HG3	2:L:105:VAL:H	1.83	0.43
3:G:45:CYS:SG	3:G:72:PRO:HD2	2.59	0.43
1:A:205:ILE:HA	1:A:220:LYS:HD3	2.00	0.43
1:H:206:CYS:O	1:H:218:ASP:HA	2.19	0.43
2:L:167:GLN:OE1	2:L:173:ALA:HB2	2.18	0.43
2:B:149:LYS:HE2	2:B:154:PRO:HG3	2.01	0.43
3:I:17:LEU:HD22	3:I:96:VAL:HG13	1.99	0.43
1:H:129:PRO:HB3	1:H:155:TYR:HB3	2.01	0.43
2:B:150:ALA:HB3	2:B:153:SER:HB2	1.99	0.43
3:I:39:LEU:HD13	3:I:102:HIS:HB2	2.01	0.43
1:A:205:ILE:N	1:A:220:LYS:HZ3	2.17	0.43
1:A:122:SER:OG	1:A:156:PHE:HZ	2.02	0.43
1:A:157:PRO:HD2	1:A:212:PRO:HB2	2.01	0.42
1:H:163:SER:HB2	1:H:207:ASN:HB2	2.00	0.42
3:I:37:ILE:HG12	3:I:46:ALA:CB	2.45	0.42
2:B:90:TRP:CZ2	2:B:93:SER:HA	2.55	0.42
1:H:61:ALA:O	1:H:65:LYS:HG3	2.20	0.42
2:L:25:ASP:O	2:L:26:LYS:HG3	2.19	0.42
2:L:34:TRP:HB2	2:L:47:ILE:HB	2.02	0.42
1:H:111:ASP:HB3	1:H:112:VAL:H	1.70	0.42
2:B:14:PRO:HD3	2:B:106:LEU:O	2.19	0.42
1:A:156:PHE:CD1	1:A:157:PRO:HA	2.55	0.41
1:H:148:LEU:HD12	1:H:148:LEU:HA	1.77	0.41
1:H:159:PRO:HB2	1:H:160:VAL:H	1.69	0.41
2:B:7:PRO:HA	2:B:8:PRO:HD3	1.92	0.41
1:H:133:PRO:HD3	1:H:219:LYS:HE2	2.03	0.41
1:A:11:LEU:HD11	1:A:122:SER:OG	2.21	0.41
1:A:126:THR:HG22	1:A:213:SER:OG	2.20	0.41
1:A:133:PRO:HD3	1:A:219:LYS:HZ3	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:19:ASN:O	3:G:22:GLN:HG2	2.21	0.41
2:B:27:LEU:HD13	2:B:65:ASN:OD1	2.20	0.41
3:G:60:ILE:H	3:G:60:ILE:HG13	1.75	0.41
2:B:27:LEU:CD1	2:B:65:ASN:OD1	2.69	0.40
2:L:145:THR:OG1	2:L:196:THR:HB	2.22	0.40
1:A:104:TYR:CZ	3:G:72:PRO:HB3	2.56	0.40
1:A:133:PRO:HD3	1:A:219:LYS:NZ	2.36	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:183:SER:N	3:I:24:GLN:OE1[1_565]	1.94	0.26

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	195/227 (86%)	176 (90%)	17 (9%)	2 (1%)	15	11
1	H	189/227 (83%)	177 (94%)	10 (5%)	2 (1%)	14	9
2	B	206/212 (97%)	196 (95%)	8 (4%)	2 (1%)	15	11
2	L	202/212 (95%)	195 (96%)	7 (4%)	0	100	100
3	G	86/112 (77%)	80 (93%)	6 (7%)	0	100	100
3	I	89/112 (80%)	84 (94%)	4 (4%)	1 (1%)	14	9
All	All	967/1102 (88%)	908 (94%)	52 (5%)	7 (1%)	22	18

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	150	ALA

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Mol	Chain	Res	Type
3	I	24	GLN
2	B	15	GLY
1	A	195	PRO
1	H	211	LYS
1	H	159	PRO
1	A	136	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	170/190 (90%)	165 (97%)	5 (3%)	42	46
1	H	165/190 (87%)	164 (99%)	1 (1%)	86	90
2	B	178/182 (98%)	175 (98%)	3 (2%)	60	67
2	L	177/182 (97%)	171 (97%)	6 (3%)	37	39
3	G	82/97 (84%)	79 (96%)	3 (4%)	34	35
3	I	81/97 (84%)	76 (94%)	5 (6%)	18	15
All	All	853/938 (91%)	830 (97%)	23 (3%)	44	48

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	218	ASP
2	L	22	CYS
2	L	26	LYS
2	L	66	SER
2	L	153	SER
2	L	156	LYS
2	L	187	SER
1	A	35	SER
1	A	43	LYS
1	A	65	LYS
1	A	196	SER
1	A	207	ASN

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Mol	Chain	Res	Type
2	B	3	GLU
2	B	55	SER
2	B	66	SER
3	G	22	GLN
3	G	58	SER
3	G	68	ASN
3	I	11	LYS
3	I	24	GLN
3	I	36	SER
3	I	107	PHE
3	I	108	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	98	GLN
3	G	68	ASN
3	I	24	GLN

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 ⓘ

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	203/227 (89%)	0.40	14 (6%) 16 21	30, 48, 83, 91	0
1	H	197/227 (86%)	0.24	7 (3%) 42 49	29, 47, 78, 92	0
2	B	208/212 (98%)	0.17	2 (0%) 82 85	33, 52, 70, 75	0
2	L	206/212 (97%)	0.20	8 (3%) 39 45	30, 51, 83, 95	0
3	G	92/112 (82%)	0.15	5 (5%) 25 31	35, 53, 78, 93	0
3	I	93/112 (83%)	0.32	5 (5%) 25 31	34, 56, 87, 92	0
All	All	999/1102 (90%)	0.25	41 (4%) 37 43	29, 51, 80, 95	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	136	PRO	5.3
1	A	170	THR	4.6
2	L	159	VAL	3.9
3	I	85	VAL	3.7
1	A	164	TRP	3.4
1	H	172	GLY	3.2
1	A	160	VAL	2.7
2	L	181	THR	2.7
1	A	161	THR	2.7
2	L	154	PRO	2.6
3	G	30	ASN	2.6
1	H	221	VAL	2.5
1	H	223	PRO	2.5
1	H	148	LEU	2.5
2	L	120	PRO	2.5
3	G	40	THR	2.5
3	I	87	ASP	2.4
2	B	150	ALA	2.4
1	H	206	CYS	2.4

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Mol	Chain	Res	Type	RSRZ
2	L	185	TRP	2.4
3	I	107	PHE	2.3
3	G	22	GLN	2.3
1	A	155	TYR	2.3
2	L	202	VAL	2.3
3	G	24	GLN	2.3
1	A	178	ALA	2.3
2	L	156	LYS	2.2
1	H	160	VAL	2.2
2	B	206	VAL	2.2
1	A	171	SER	2.2
1	A	135	ALA	2.2
2	L	205	THR	2.2
1	H	193	THR	2.1
3	I	24	GLN	2.1
3	G	85	VAL	2.1
1	A	206	CYS	2.1
1	A	193	THR	2.1
3	I	31	GLY	2.1
1	A	191	VAL	2.0
1	A	205	ILE	2.0
1	A	145	THR	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 monosaccharides 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.