Full wwPDB X-ray Structure Validation Report ⓘ

Feb 22, 2021 – 05:01 PM GMT

PDB ID	:	6SHG
Title	:	Diffraction data for RoAb13 crystal co-crystallised with PIYDIN and its RoAb13 structure
Authors	:	Saridakis, E.; Helliwell, J.R.; Govada, L.; Chain, B.; Morgan, M.; Kassen, S.C.; Chayen, N.E.
Deposited on	:	2019-08-06
Resolution	:	3.35 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:

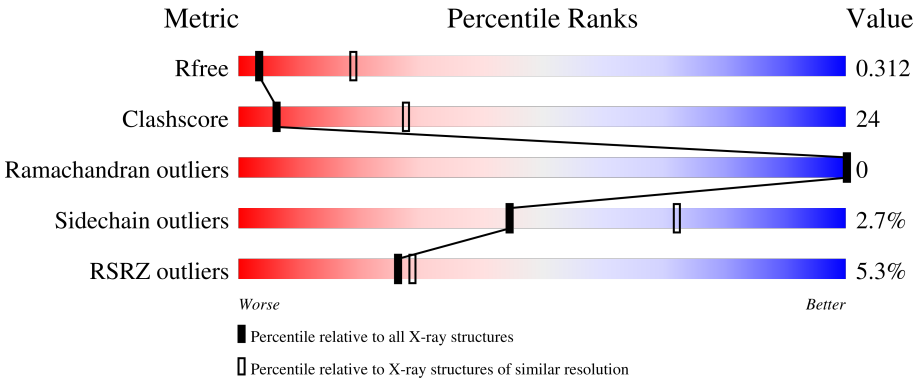
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.17.1.dev1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.17.1.dev1

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.35 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	1558 (3.42-3.30)
Clashscore	141614	1627 (3.42-3.30)
Ramachandran outliers	138981	1599 (3.42-3.30)
Sidechain outliers	138945	1598 (3.42-3.30)
RSRZ outliers	127900	1507 (3.42-3.30)

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 <=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	H	224	2% 60% 35%
2	L	220	8% 52% 42%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3280 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 Antibody RoAb13 Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	218	Total	C	N	O	S	0	1	0
			1648	1034	281	326	7			

- Molecule 2 is a protein called Antibody RoAb13 Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	214	Total	C	N	O	S	0	0	0
			1628	1018	268	333	9			

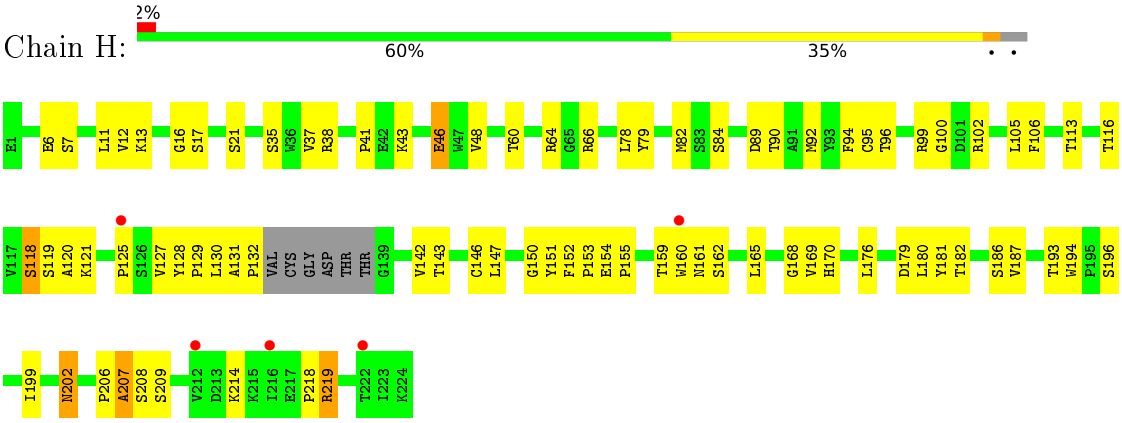
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	1	Total	O	0	0
			1	1		
3	L	3	Total	O	0	0
			3	3		

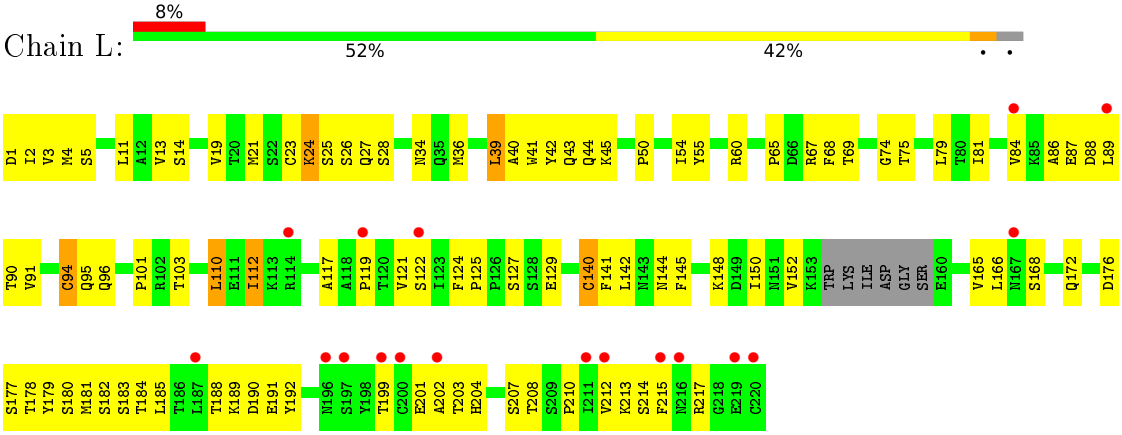
3 Residue-property plots (i)

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: Antibody RoAb13 Heavy Chain



- Molecule 2: Antibody RoAb13 Light Chain



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	76.61Å 76.61Å 270.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	67.52 – 3.35 73.70 – 3.09	Depositor EDS
% Data completeness (in resolution range)	86.5 (67.52-3.35) 66.9 (73.70-3.09)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.76 (at 3.07Å)	Xtriage
Refinement program	PHENIX 1.12_2829, REFMAC 5	Depositor
R, R_{free}	0.261 , 0.312 0.261 , 0.312	Depositor DCC
R_{free} test set	554 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å^2)	79.4	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 64.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.86	EDS
Total number of atoms	3280	wwPDB-VP
Average B, all atoms (Å^2)	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.59% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|>5$	RMSZ	#$\|Z\|>5$
1	H	0.74	2/1689 (0.1%)	0.87	3/2297 (0.1%)
2	L	0.73	3/1664 (0.2%)	0.83	1/2264 (0.0%)
All	All	0.74	5/3353 (0.1%)	0.85	4/4561 (0.1%)

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	H	0	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	95	CYS	CB-SG	-7.52	1.69	1.82
2	L	94	CYS	CB-SG	5.90	1.92	1.82
2	L	23	CYS	CB-SG	5.17	1.91	1.82
1	H	46	GLU	CG-CD	5.05	1.59	1.51
2	L	42	TYR	CD1-CE1	5.01	1.46	1.39

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	H	219	ARG	NE-CZ-NH1	-6.35	117.12	120.30
1	H	95	CYS	CA-CB-SG	-5.67	103.80	114.00
2	L	112	ILE	CG1-CB-CG2	-5.53	99.23	111.40
1	H	147	LEU	CA-CB-CG	5.01	126.83	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	H	207	ALA	Peptide

5.2 Too-close contacts (i)

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	H	1648	0	1609	74	0
2	L	1628	0	1530	95	0
3	H	1	0	0	0	0
3	L	3	0	0	0	0
All	All	3280	0	3139	156	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 24.

All (156) 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:H:99:ARG:NH1	2:L:55:TYR:CE2	2.08	1.20
1:H:99:ARG:NH2	2:L:55:TYR:OH	1.82	1.11
2:L:112:ILE:HG22	2:L:172:GLN:OE1	1.50	1.10
1:H:99:ARG:NH2	2:L:55:TYR:CZ	2.20	1.08
1:H:150:GLY:HA2	1:H:180:LEU:HB3	1.47	0.95
1:H:90:THR:HG23	1:H:116:THR:HA	1.50	0.92
1:H:99:ARG:CZ	2:L:55:TYR:CZ	2.55	0.89
1:H:99:ARG:NH1	2:L:55:TYR:CZ	2.45	0.84
2:L:119:PRO:HB2	2:L:142:LEU:HD12	1.61	0.82
2:L:201:GLU:HG3	2:L:212:VAL:HG12	1.62	0.82
1:H:160:TRP:HB3	1:H:165:LEU:HD12	1.62	0.79
1:H:152:PHE:CD1	1:H:153:PRO:HA	2.18	0.78
2:L:204:HIS:H	2:L:207:SER:HB3	1.47	0.78
1:H:99:ARG:HH12	2:L:55:TYR:HE2	1.32	0.76
2:L:152:VAL:HG22	2:L:202:ALA:HA	1.68	0.76
2:L:202:ALA:O	2:L:210:PRO:HB3	1.87	0.74
1:H:13:LYS:HE2	1:H:119:SER:O	1.88	0.73
2:L:112:ILE:CG2	2:L:172:GLN:OE1	2.35	0.72
1:H:99:ARG:HH22	2:L:55:TYR:HH	1.38	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:41:TRP:HB2	2:L:54:ILE:HB	1.74	0.69
1:H:131:ALA:O	1:H:219:ARG:NH2	2.26	0.68
1:H:130:LEU:HD22	2:L:124:PHE:HB3	1.75	0.67
1:H:176:LEU:HB2	1:H:181:TYR:CE1	2.30	0.67
2:L:86:ALA:O	2:L:89:LEU:HG	1.96	0.64
1:H:102:ARG:NH2	1:H:105:LEU:HB2	2.13	0.63
2:L:65:PRO:HD2	2:L:68:PHE:HD2	1.64	0.63
1:H:102:ARG:HH21	1:H:105:LEU:HB2	1.63	0.63
2:L:140:CYS:N	2:L:183:SER:OG	2.25	0.62
1:H:154:GLU:HG2	1:H:155:PRO:HA	1.81	0.61
2:L:5:SER:HB2	2:L:24:LYS:HE3	1.83	0.60
2:L:3:VAL:H	2:L:26:SER:HG	1.49	0.60
2:L:150:ILE:HB	2:L:204:HIS:HD2	1.66	0.60
2:L:165:VAL:HG13	2:L:185:LEU:HA	1.83	0.60
2:L:28:SER:HA	2:L:74:GLY:O	2.02	0.60
1:H:6:GLU:OE2	1:H:94:PHE:HA	2.02	0.60
2:L:179:TYR:O	2:L:180:SER:OG	2.20	0.58
1:H:168:GLY:O	1:H:187:VAL:HA	2.04	0.58
1:H:161:ASN:OD1	1:H:199:ILE:HA	2.02	0.58
2:L:190:ASP:N	2:L:190:ASP:OD1	2.34	0.58
2:L:4:MET:CE	2:L:96:GLN:HG2	2.34	0.57
2:L:43:GLN:HG3	2:L:90:THR:HG21	1.86	0.57
2:L:172:GLN:HG3	2:L:179:TYR:CE1	2.40	0.57
2:L:122:SER:O	2:L:140:CYS:HA	2.04	0.57
2:L:3:VAL:N	2:L:26:SER:OG	2.24	0.56
1:H:99:ARG:HH11	1:H:99:ARG:HG2	1.70	0.56
2:L:44:GLN:HB2	2:L:50:PRO:HG3	1.86	0.56
1:H:13:LYS:HA	1:H:118:SER:O	2.05	0.56
1:H:48:VAL:O	1:H:60:THR:HG22	2.06	0.56
1:H:127:VAL:O	1:H:214:LYS:HE3	2.06	0.55
2:L:150:ILE:HG23	2:L:181:MET:HE3	1.88	0.55
1:H:35:SER:OG	1:H:96:THR:OG1	2.25	0.54
2:L:178:THR:OG1	2:L:179:TYR:N	2.41	0.54
2:L:41:TRP:CZ3	2:L:94:CYS:HB3	2.43	0.54
2:L:2:ILE:O	2:L:103:THR:HG21	2.07	0.54
2:L:65:PRO:HD2	2:L:68:PHE:CD2	2.43	0.54
1:H:99:ARG:NH1	1:H:99:ARG:HG2	2.22	0.53
1:H:99:ARG:NH1	2:L:55:TYR:CD2	2.69	0.53
1:H:125:PRO:HB3	1:H:151:TYR:HB3	1.90	0.53
2:L:13:VAL:HG22	2:L:14:SER:H	1.73	0.53
2:L:21:MET:HE3	2:L:79:LEU:HD23	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:99:ARG:NH2	2:L:55:TYR:CE1	2.75	0.53
1:H:120:ALA:HB3	1:H:152:PHE:CE2	2.44	0.53
1:H:37:VAL:HG13	1:H:46:GLU:O	2.09	0.52
2:L:39:LEU:HG	2:L:40:ALA:N	2.24	0.52
2:L:88:ASP:O	2:L:110:LEU:HD11	2.09	0.52
1:H:152:PHE:CG	1:H:153:PRO:HA	2.44	0.52
1:H:118:SER:HB3	1:H:120:ALA:H	1.73	0.52
1:H:96:THR:HG1	1:H:106:PHE:HD1	1.58	0.51
1:H:169:VAL:HA	1:H:186:SER:O	2.10	0.51
2:L:112:ILE:HG22	2:L:172:GLN:CD	2.29	0.51
1:H:11:LEU:HB2	1:H:153:PRO:HG3	1.92	0.51
1:H:38:ARG:HA	1:H:92:MET:O	2.11	0.51
2:L:165:VAL:HG12	2:L:166:LEU:H	1.76	0.51
2:L:117:ALA:HB3	2:L:145:PHE:HA	1.91	0.50
2:L:176:ASP:OD2	2:L:178:THR:HB	2.12	0.50
1:H:99:ARG:NH2	2:L:55:TYR:HH	2.03	0.50
1:H:146:CYS:HB2	1:H:160:TRP:CH2	2.47	0.50
1:H:160:TRP:O	1:H:165:LEU:HG	2.12	0.50
1:H:194:TRP:CH2	1:H:218:PRO:HG3	2.48	0.49
2:L:19:VAL:HB	2:L:81:ILE:HB	1.93	0.49
2:L:145:PHE:N	2:L:178:THR:OG1	2.45	0.49
2:L:208:THR:O	2:L:210:PRO:HD3	2.12	0.49
1:H:100:GLY:O	1:H:102:ARG:N	2.45	0.49
2:L:189:LYS:O	2:L:192:TYR:HB3	2.11	0.49
1:H:160:TRP:HB2	1:H:165:LEU:HB2	1.95	0.48
2:L:110:LEU:HD23	2:L:110:LEU:H	1.76	0.48
1:H:89:ASP:OD1	1:H:89:ASP:N	2.46	0.48
1:H:169:VAL:O	1:H:170:HIS:HD2	1.96	0.48
1:H:193:THR:O	1:H:196:SER:HB2	2.13	0.48
2:L:13:VAL:HG11	2:L:84:VAL:HG21	1.96	0.48
1:H:12:VAL:CG1	1:H:16:GLY:HA3	2.44	0.47
2:L:24:LYS:HD2	2:L:25:SER:N	2.30	0.47
2:L:152:VAL:HG13	2:L:201:GLU:O	2.14	0.47
1:H:64:ARG:HH11	1:H:64:ARG:HG2	1.80	0.47
2:L:43:GLN:OE1	2:L:45:LYS:HE3	2.15	0.47
2:L:121:VAL:HG22	2:L:142:LEU:CD1	2.44	0.47
2:L:145:PHE:HE2	2:L:180:SER:HA	1.79	0.47
2:L:5:SER:HB2	2:L:24:LYS:HB3	1.97	0.46
1:H:17:SER:HA	1:H:82:MET:O	2.15	0.46
1:H:128:TYR:HA	1:H:129:PRO:HD3	1.77	0.46
2:L:60:ARG:NH1	2:L:69:THR:HG23	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:121:VAL:HG22	2:L:142:LEU:HD13	1.98	0.46
1:H:12:VAL:HG11	1:H:16:GLY:HA3	1.97	0.46
1:H:161:ASN:HB2	1:H:165:LEU:HG	1.97	0.45
2:L:34:ASN:O	2:L:36:MET:HG2	2.16	0.45
1:H:125:PRO:HB3	1:H:151:TYR:CB	2.46	0.45
2:L:214:SER:OG	2:L:215:PHE:N	2.49	0.45
2:L:27:GLN:O	2:L:75:THR:HG22	2.17	0.45
2:L:203:THR:HG22	2:L:210:PRO:HG3	1.98	0.45
1:H:78:LEU:HD12	1:H:79:TYR:H	1.81	0.45
2:L:24:LYS:HE3	2:L:24:LYS:HB3	1.46	0.45
2:L:60:ARG:HH12	2:L:69:THR:HG23	1.82	0.44
1:H:176:LEU:HB2	1:H:181:TYR:CZ	2.52	0.44
2:L:121:VAL:HG12	2:L:122:SER:N	2.33	0.44
1:H:66:ARG:NH2	1:H:89:ASP:OD2	2.49	0.44
2:L:144:ASN:HA	2:L:178:THR:OG1	2.18	0.44
2:L:11:LEU:HD12	2:L:11:LEU:HA	1.72	0.44
1:H:142:VAL:HG22	1:H:143:THR:H	1.83	0.44
2:L:39:LEU:HG	2:L:40:ALA:H	1.83	0.44
2:L:21:MET:CB	2:L:79:LEU:HB3	2.47	0.43
2:L:67:ARG:NH2	2:L:87:GLU:OE2	2.49	0.43
1:H:152:PHE:HB2	1:H:179:ASP:O	2.19	0.43
2:L:127:SER:C	2:L:129:GLU:H	2.22	0.43
1:H:162:SER:N	1:H:202:ASN:OD1	2.51	0.43
1:H:206:PRO:O	1:H:208:SER:N	2.46	0.43
2:L:90:THR:OG1	2:L:91:VAL:N	2.52	0.43
1:H:11:LEU:HG	1:H:12:VAL:N	2.34	0.42
2:L:121:VAL:HA	2:L:141:PHE:O	2.19	0.42
1:H:41:PRO:C	1:H:43:LYS:H	2.21	0.42
1:H:6:GLU:HB3	1:H:113:THR:CG2	2.49	0.42
2:L:150:ILE:HD11	2:L:202:ALA:HB1	2.00	0.42
1:H:16:GLY:O	1:H:84:SER:N	2.47	0.42
2:L:217:ARG:HH11	2:L:217:ARG:HG3	1.84	0.42
2:L:188:THR:HB	2:L:191:GLU:HB2	2.00	0.42
1:H:159:THR:OG1	1:H:202:ASN:HB2	2.20	0.42
2:L:11:LEU:HB3	2:L:110:LEU:HA	2.02	0.42
1:H:132:PRO:O	1:H:219:ARG:NE	2.44	0.42
2:L:1:ASP:HA	2:L:101:PRO:HD2	2.02	0.42
2:L:21:MET:HB2	2:L:79:LEU:HB3	2.00	0.42
1:H:207:ALA:N	1:H:209:SER:H	2.18	0.42
1:H:102:ARG:HH21	1:H:105:LEU:CD2	2.32	0.41
2:L:166:LEU:HB2	2:L:184:THR:HB	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:7:SER:HB2	1:H:21:SER:HB2	2.02	0.41
2:L:4:MET:HE3	2:L:96:GLN:HG2	2.02	0.41
1:H:121:LYS:HB2	1:H:121:LYS:HE3	1.86	0.41
2:L:212:VAL:O	2:L:213:LYS:HD2	2.20	0.41
2:L:2:ILE:H	2:L:2:ILE:HG13	1.68	0.41
2:L:86:ALA:C	2:L:88:ASP:H	2.23	0.41
1:H:182:THR:HG21	2:L:184:THR:HG21	2.03	0.41
1:H:186:SER:HB2	2:L:141:PHE:CD2	2.56	0.41
2:L:124:PHE:HA	2:L:125:PRO:HD3	1.89	0.41
2:L:54:ILE:HD13	2:L:60:ARG:HA	2.03	0.40
2:L:150:ILE:HD12	2:L:204:HIS:CD2	2.57	0.40
2:L:199:THR:HG23	2:L:214:SER:HB2	2.02	0.40
1:H:142:VAL:HG22	1:H:143:THR:N	2.36	0.40
2:L:112:ILE:CG2	2:L:177:SER:HB3	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	215/224 (96%)	182 (85%)	33 (15%)	0	100	100
2	L	210/220 (96%)	179 (85%)	31 (15%)	0	100	100
All	All	425/444 (96%)	361 (85%)	64 (15%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	H	186/191 (97%)	184 (99%)	2 (1%)	73 86
2	L	182/196 (93%)	174 (96%)	8 (4%)	28 59
All	All	368/387 (95%)	358 (97%)	10 (3%)	44 72

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	118	SER
1	H	202	ASN
2	L	24	LYS
2	L	39	LEU
2	L	95	GLN
2	L	110	LEU
2	L	140	CYS
2	L	148	LYS
2	L	168	SER
2	L	182	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	39	GLN
1	H	170	HIS
2	L	44	GLN
2	L	95	GLN
2	L	204	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

There are no monosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	H	218/224 (97%)	0.43	5 (2%)	60	63	54, 82, 112, 132	0
2	L	214/220 (97%)	0.63	18 (8%)	11	13	60, 104, 149, 170	0
All	All	432/444 (97%)	0.53	23 (5%)	26	28	54, 89, 136, 170	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	187	LEU	5.4
2	L	202	ALA	4.9
2	L	211	ILE	4.2
2	L	220	CYS	4.2
1	H	222	THR	4.0
2	L	212	VAL	3.6
2	L	196	ASN	3.6
2	L	114	ARG	3.3
2	L	219	GLU	3.3
2	L	200	CYS	3.2
2	L	122	SER	2.9
1	H	160	TRP	2.9
2	L	119	PRO	2.6
1	H	216	ILE	2.5
2	L	84	VAL	2.4
2	L	167	ASN	2.4
2	L	89	LEU	2.4
2	L	215	PHE	2.3
1	H	212	VAL	2.2
2	L	199	THR	2.1
2	L	216	ASN	2.1
2	L	197	SER	2.1
1	H	125	PRO	2.1

6.2 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no 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.