



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:18 PM UTC

PDB ID : 3O45 / pdb_00003o45
Title : Crystal Structure of 101F Fab Bound to 17-mer Peptide Epitope
Authors : McLellan, J.S.; Chen, M.; Chang, J.S.; Yang, Y.; Kim, A.; Graham, B.S.; Kwong, P.D.
Deposited on : 2010-07-26
Resolution : 2.87 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

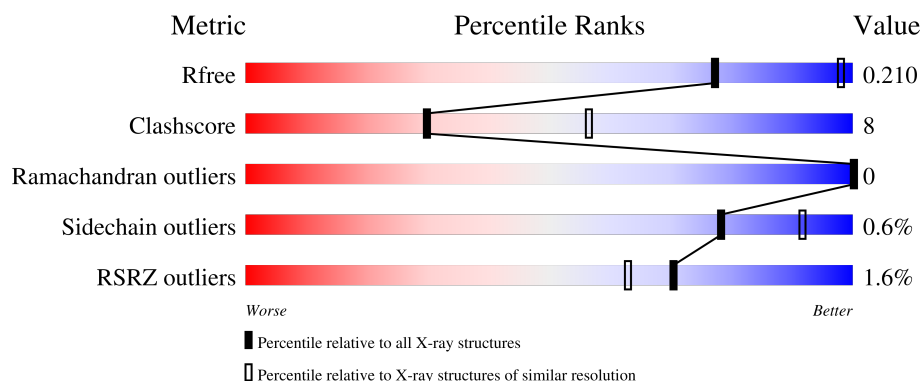
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.87 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}	180053	3557 (2.90-2.86)
Clashscore	190562	3801 (2.90-2.86)
Ramachandran outliers	187476	3699 (2.90-2.86)
Sidechain outliers	187428	3702 (2.90-2.86)
RSRZ outliers	180081	3558 (2.90-2.86)

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	B	218	% 79% 21%
1	L	218	2% 76% 24%
2	A	220	2% 86% 14%
2	H	220	% 82% 18%
3	C	19	5% 53% 47%

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Mol	Chain	Length	Quality of chain
3	P	19	 A horizontal bar chart showing the quality of chain 3. The bar is divided into three segments: a green segment representing 63%, a yellow segment representing 5%, and a grey segment representing 32%. The percentages are labeled below the bar.

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6953 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 Mouse monoclonal antibody 101F 101F Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	218	Total	C	N	O	S	0	0	0
			1692	1059	284	342	7			
1	B	217	Total	C	N	O	S	0	0	0
			1686	1056	283	341	6			

- Molecule 2 is a protein called Mouse monoclonal antibody 101F 101F Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	220	Total	C	N	O	S	0	0	0
			1669	1064	274	325	6			
2	A	220	Total	C	N	O	S	0	0	0
			1669	1064	274	325	6			

- Molecule 3 is a protein called Fusion glycoprotein F1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	13	Total	C	N	O	0	0	1
			94	58	20	16			
3	C	10	Total	C	N	O	0	0	0
			80	50	16	14			

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	L	1	Total	O	S	0	0
			5	4	1		
4	L	1	Total	O	S	0	0
			5	4	1		
4	L	1	Total	O	S	0	0
			5	4	1		
4	H	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

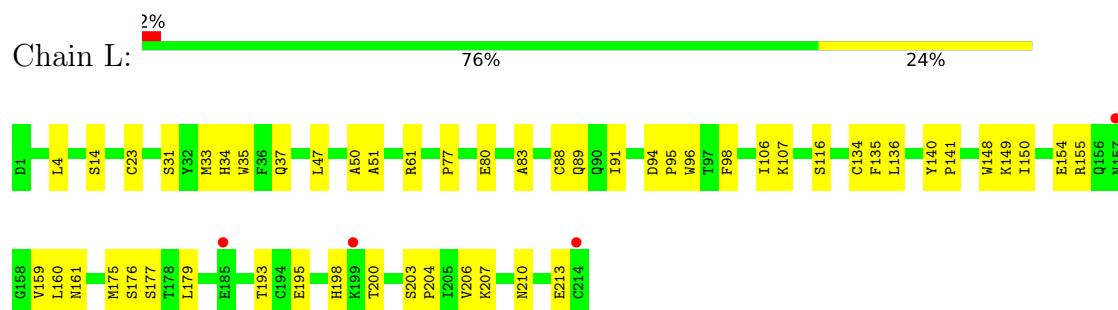
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	L	8	Total	O	0	0
			8	8		
5	H	6	Total	O	0	0
			6	6		
5	A	8	Total	O	0	0
			8	8		
5	B	5	Total	O	0	0
			5	5		
5	P	1	Total	O	0	0
			1	1		

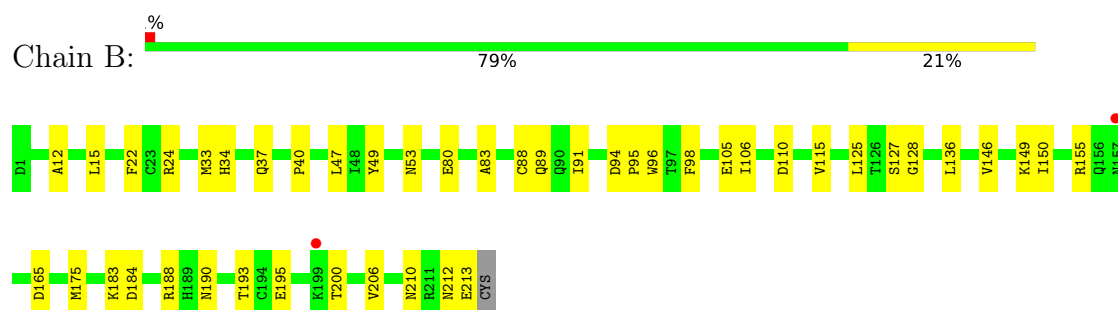
3 Residue-property plots [i](#)

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.

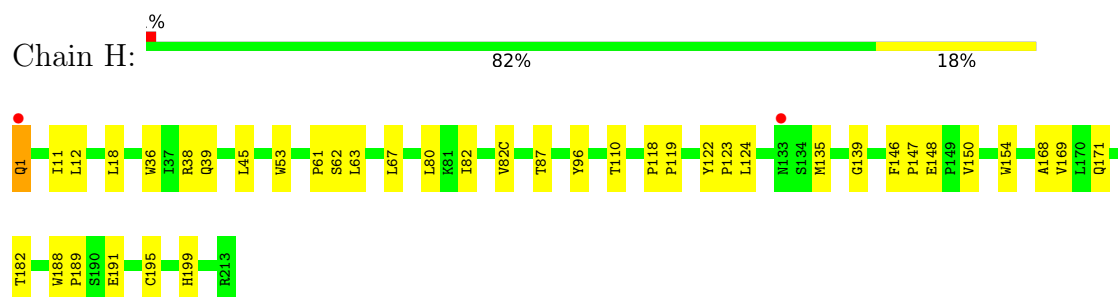
- Molecule 1: Mouse monoclonal antibody 101F 101F Fab light chain



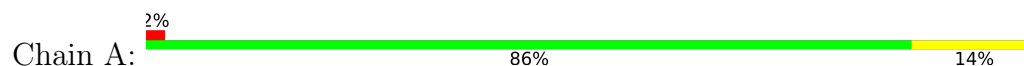
- Molecule 1: Mouse monoclonal antibody 101F 101F Fab light chain

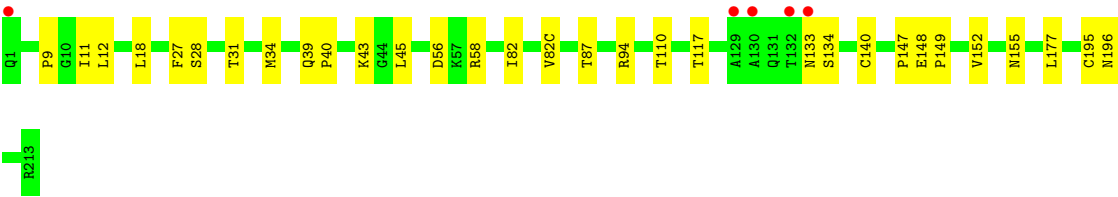


- Molecule 2: Mouse monoclonal antibody 101F 101F Fab heavy chain

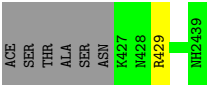


- Molecule 2: Mouse monoclonal antibody 101F 101F Fab heavy chain

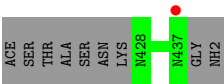




• Molecule 3: Fusion glycoprotein F1



• Molecule 3: Fusion glycoprotein F1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.54Å 92.97Å 140.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.54 – 2.87 25.54 – 2.87	Depositor EDS
% Data completeness (in resolution range)	98.0 (25.54-2.87) 97.7 (25.54-2.87)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 2.89Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.200 , 0.254 (Not available) , 0.210	Depositor DCC
R_{free} test set	1228 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	32.3	Xtrriage
Anisotropy	0.530	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 28.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6953	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.12% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, NH2

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	B	0.25	0/1728	0.66	0/2352
1	L	0.25	0/1734	0.68	0/2360
2	A	0.25	0/1715	0.70	2/2348 (0.1%)
2	H	0.26	0/1715	0.72	2/2348 (0.1%)
3	C	0.22	0/80	0.62	0/105
3	P	0.26	0/93	0.63	0/121
All	All	0.25	0/7065	0.69	4/9634 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	199	HIS	CA-C-N	5.31	125.37	119.32
2	H	199	HIS	C-N-CA	5.31	125.37	119.32
2	A	117	THR	CA-C-N	5.25	123.44	119.66
2	A	117	THR	C-N-CA	5.25	123.44	119.66

There are no chirality outliers.

There are no planarity outliers.

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	B	1686	0	1605	27	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	L	1692	0	1610	36	0
2	A	1669	0	1647	24	0
2	H	1669	0	1647	34	0
3	C	80	0	83	0	0
3	P	94	0	99	2	0
4	A	5	0	0	0	0
4	B	10	0	0	0	0
4	H	5	0	0	0	0
4	L	15	0	0	0	0
5	A	8	0	0	0	0
5	B	5	0	0	0	0
5	H	6	0	0	0	0
5	L	8	0	0	0	0
5	P	1	0	0	0	0
All	All	6953	0	6691	115	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 8.

All (115) 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)
2:H:12:LEU:HD11	2:H:82(C):VAL:HG21	1.67	0.76
1:B:127:SER:N	1:B:128:GLY:HA2	2.00	0.76
1:L:195:GLU:HG2	1:L:206:VAL:HG22	1.68	0.72
1:L:14:SER:HA	1:L:107:LYS:HB2	1.72	0.72
2:A:11:ILE:HD12	2:A:147:PRO:HG3	1.72	0.71
2:H:11:ILE:HD12	2:H:147:PRO:HG3	1.76	0.66
1:L:155:ARG:HD3	1:L:179:LEU:HD11	1.77	0.66
1:B:146:VAL:HG21	1:B:175:MET:HE1	1.76	0.66
1:B:127:SER:H	1:B:128:GLY:HA2	1.61	0.65
2:A:27:PHE:CD1	2:A:34:MET:HE2	2.32	0.65
2:H:96:TYR:CZ	3:P:429:ARG:HD2	2.33	0.64
1:B:115:VAL:HG22	1:B:136:LEU:HD22	1.81	0.62
1:L:95:PRO:HG3	2:H:61:PRO:HD2	1.82	0.62
1:L:161:ASN:HB3	1:L:175:MET:HE3	1.82	0.61
1:B:37:GLN:HB2	1:B:47:LEU:HD11	1.84	0.60
1:L:150:ILE:HD11	1:L:179:LEU:HD21	1.85	0.58
2:A:11:ILE:HB	2:A:147:PRO:HG2	1.86	0.58
2:A:12:LEU:CD1	2:A:82(C):VAL:HG21	2.34	0.57
2:A:40:PRO:HG2	2:A:43:LYS:HD3	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:12:LEU:HD11	2:A:82(C):VAL:HG21	1.87	0.56
1:L:94:ASP:HA	1:L:95:PRO:C	2.30	0.56
2:H:39:GLN:HB2	2:H:45:LEU:HD23	1.88	0.55
1:L:37:GLN:HB2	1:L:47:LEU:HD11	1.87	0.54
1:B:195:GLU:HG2	1:B:206:VAL:HG22	1.90	0.54
1:B:125:LEU:O	1:B:183:LYS:HD3	2.08	0.53
2:H:96:TYR:CE2	3:P:429:ARG:HD2	2.43	0.53
2:A:31:THR:HB	2:A:34:MET:HG3	1.90	0.52
1:L:210:ASN:HB2	1:L:213:GLU:HG2	1.91	0.52
1:B:22:PHE:CE2	1:B:24:ARG:HB2	2.45	0.52
2:A:148:GLU:HB3	2:A:149:PRO:HA	1.92	0.51
2:A:39:GLN:HB2	2:A:45:LEU:HD23	1.92	0.51
1:B:149:LYS:HB2	1:B:193:THR:HB	1.93	0.51
1:L:33:MET:SD	1:L:88:CYS:HB2	2.50	0.51
2:A:87:THR:HG23	2:A:110:THR:HA	1.92	0.51
1:B:127:SER:N	1:B:128:GLY:CA	2.73	0.50
2:H:12:LEU:CD1	2:H:82(C):VAL:HG21	2.39	0.50
1:B:22:PHE:HE2	1:B:24:ARG:HB2	1.77	0.50
1:L:175:MET:HG2	1:L:176:SER:N	2.27	0.50
2:H:36:TRP:CD1	2:H:80:LEU:HB2	2.47	0.50
2:A:18:LEU:HD23	2:A:82:ILE:HD12	1.94	0.50
2:A:34:MET:HE1	2:A:94:ARG:CD	2.42	0.50
2:H:135:MET:HE3	2:H:182:THR:HG22	1.94	0.49
2:H:148:GLU:OE2	2:H:168:ALA:HB3	2.11	0.49
1:B:210:ASN:HB2	1:B:213:GLU:HG2	1.94	0.49
2:H:154:TRP:CZ3	2:H:195:CYS:HB3	2.48	0.49
2:H:12:LEU:C	2:H:12:LEU:HD12	2.38	0.49
2:H:87:THR:HG23	2:H:110:THR:HA	1.95	0.49
1:L:31:SER:O	1:L:50:ALA:HA	2.12	0.48
2:A:34:MET:HE1	2:A:94:ARG:CG	2.42	0.48
1:L:4:LEU:HD13	1:L:33:MET:HE1	1.95	0.48
1:B:150:ILE:HD12	1:B:155:ARG:HD2	1.96	0.48
1:L:160:LEU:HD11	2:H:171:GLN:HB2	1.95	0.48
1:B:184:ASP:O	1:B:188:ARG:HG3	2.14	0.47
1:L:23:CYS:HB2	1:L:35:TRP:CH2	2.50	0.47
1:B:136:LEU:HD11	1:B:146:VAL:HG22	1.96	0.47
2:H:38:ARG:NH2	2:H:63:LEU:HD21	2.30	0.47
1:L:160:LEU:HD22	2:H:169:VAL:HG11	1.97	0.46
1:L:175:MET:HE2	1:L:177:SER:HB2	1.97	0.46
2:H:18:LEU:HD23	2:H:82:ILE:HD12	1.96	0.46
2:A:11:ILE:HG12	2:A:12:LEU:N	2.29	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:33:MET:SD	1:B:88:CYS:HB2	2.56	0.46
1:L:148:TRP:O	1:L:154:GLU:HA	2.16	0.45
1:B:190:ASN:HD21	1:B:212:ASN:ND2	2.14	0.45
1:L:61:ARG:HD2	1:L:77:PRO:O	2.16	0.45
1:B:110:ASP:HB3	1:B:200:THR:HG22	1.98	0.45
2:H:188:TRP:CG	2:H:189:PRO:HA	2.52	0.45
2:A:34:MET:HE1	2:A:94:ARG:HD2	1.99	0.45
1:L:198:HIS:CE1	1:L:200:THR:HG23	2.52	0.45
1:B:89:GLN:HB2	1:B:98:PHE:CD2	2.51	0.45
2:H:191:GLU:OE2	2:A:28:SER:HB2	2.16	0.45
1:L:135:PHE:C	1:L:136:LEU:HD12	2.42	0.45
1:B:15:LEU:HD21	1:B:80:GLU:HG3	1.99	0.45
1:B:91:ILE:HG22	1:B:96:TRP:HB3	1.99	0.44
1:L:149:LYS:HB2	1:L:193:THR:HB	1.99	0.44
2:A:31:THR:HB	2:A:34:MET:CG	2.47	0.44
2:H:122:TYR:HA	2:H:123:PRO:HD3	1.85	0.44
1:B:40:PRO:HB3	1:B:165:ASP:OD2	2.18	0.44
2:A:9:PRO:HG2	2:A:12:LEU:HD23	2.00	0.44
2:H:11:ILE:HB	2:H:147:PRO:HG2	2.00	0.44
2:H:146:PHE:HA	2:H:147:PRO:HA	1.79	0.44
1:B:34:HIS:HE2	1:B:91:ILE:HD11	1.84	0.43
1:L:203:SER:HA	1:L:204:PRO:HD3	1.76	0.43
2:H:1:GLN:HG3	2:A:155:ASN:HB3	1.99	0.43
2:H:62:SER:HB3	2:H:63:LEU:HD12	2.00	0.43
2:A:133:ASN:CG	2:A:134:SER:H	2.26	0.43
1:L:206:VAL:O	1:L:207:LYS:HD3	2.18	0.43
2:H:53:TRP:H	2:H:53:TRP:HE3	1.66	0.43
2:A:12:LEU:HD12	2:A:12:LEU:C	2.44	0.43
1:B:49:TYR:O	1:B:53:ASN:HB2	2.19	0.43
2:H:11:ILE:HB	2:H:147:PRO:CG	2.49	0.43
2:H:1:GLN:CD	2:H:1:GLN:N	2.76	0.42
1:L:89:GLN:HB2	1:L:98:PHE:CD2	2.54	0.42
1:B:83:ALA:HB2	1:B:106:ILE:HG12	2.02	0.42
1:L:95:PRO:HG3	2:H:61:PRO:CD	2.47	0.42
1:L:150:ILE:HD12	1:L:155:ARG:HD2	2.02	0.42
2:H:124:LEU:HB2	2:H:139:GLY:C	2.45	0.42
2:A:56:ASP:CG	2:A:58:ARG:HH12	2.28	0.42
1:L:136:LEU:HD12	1:L:136:LEU:N	2.35	0.42
1:L:83:ALA:HB2	1:L:106:ILE:HG12	2.01	0.41
1:L:159:VAL:HG22	1:L:179:LEU:HD13	2.02	0.41
2:A:152:VAL:HA	2:A:196:ASN:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:116:SER:O	1:L:134:CYS:HA	2.21	0.41
2:H:118:PRO:HA	2:H:119:PRO:HD3	1.85	0.41
1:L:50:ALA:O	1:L:51:ALA:HB3	2.21	0.41
2:H:12:LEU:HD12	2:H:12:LEU:O	2.21	0.41
2:H:67:LEU:HD23	2:H:82:ILE:HG12	2.02	0.41
2:H:188:TRP:CD1	2:H:189:PRO:HA	2.55	0.41
2:A:177:LEU:HD23	2:A:177:LEU:C	2.45	0.41
1:L:140:TYR:O	1:L:198:HIS:HE1	2.03	0.41
1:L:91:ILE:HA	1:L:96:TRP:CD1	2.56	0.41
1:B:94:ASP:HA	1:B:95:PRO:C	2.45	0.41
2:H:63:LEU:HD12	2:H:63:LEU:N	2.37	0.40
1:L:140:TYR:CG	1:L:141:PRO:HA	2.57	0.40
1:B:12:ALA:HA	1:B:105:GLU:O	2.21	0.40
1:L:34:HIS:HE2	1:L:91:ILE:HD11	1.85	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	B	215/218 (99%)	206 (96%)	9 (4%)	0	100	100
1	L	216/218 (99%)	206 (95%)	10 (5%)	0	100	100
2	A	218/220 (99%)	212 (97%)	6 (3%)	0	100	100
2	H	218/220 (99%)	208 (95%)	10 (5%)	0	100	100
3	C	8/19 (42%)	8 (100%)	0	0	100	100
3	P	11/19 (58%)	10 (91%)	1 (9%)	0	100	100
All	All	886/914 (97%)	850 (96%)	36 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	B	189/190 (100%)	189 (100%)	0	100	100
1	L	190/190 (100%)	189 (100%)	1 (0%)	81	93
2	A	191/191 (100%)	189 (99%)	2 (1%)	68	87
2	H	191/191 (100%)	189 (99%)	2 (1%)	68	87
3	C	9/14 (64%)	9 (100%)	0	100	100
3	P	10/14 (71%)	10 (100%)	0	100	100
All	All	780/790 (99%)	775 (99%)	5 (1%)	78	92

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	80	GLU
2	H	1	GLN
2	H	150	VAL
2	A	140	CYS
2	A	195	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	37	GLN
1	L	53	ASN
1	L	74	ASN
1	L	137	ASN
1	L	138	ASN
1	L	210	ASN
2	H	16	GLN
2	H	77	GLN
2	H	131	GLN
2	H	164	HIS
2	A	77	GLN
2	A	164	HIS

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Mol	Chain	Res	Type
1	B	37	GLN
1	B	42	GLN
1	B	137	ASN
1	B	161	ASN
1	B	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	B	216	-	4,4,4	0.24	0	6,6,6	0.12	0
4	SO4	L	217	-	4,4,4	0.24	0	6,6,6	0.08	0
4	SO4	H	214	-	4,4,4	0.25	0	6,6,6	0.07	0
4	SO4	B	215	-	4,4,4	0.24	0	6,6,6	0.07	0
4	SO4	L	216	-	4,4,4	0.23	0	6,6,6	0.07	0
4	SO4	A	214	-	4,4,4	0.23	0	6,6,6	0.08	0
4	SO4	L	215	-	4,4,4	0.23	0	6,6,6	0.10	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	B	217/218 (99%)	-0.10	2 (0%) 81 75	13, 33, 69, 82	4 (1%)
1	L	218/218 (100%)	-0.16	4 (1%) 67 60	14, 32, 50, 79	4 (1%)
2	A	220/220 (100%)	-0.28	5 (2%) 61 52	12, 26, 58, 105	0
2	H	220/220 (100%)	-0.28	2 (0%) 81 75	12, 25, 54, 84	0
3	C	10/19 (52%)	-0.09	1 (10%) 12 11	16, 27, 44, 71	0
3	P	12/19 (63%)	-0.05	0 100 100	17, 29, 66, 66	0
All	All	897/914 (98%)	-0.20	14 (1%) 70 63	12, 29, 62, 105	8 (0%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	199	LYS	3.7
1	B	157	ASN	3.4
2	A	1	GLN	3.3
2	A	129	ALA	2.9
1	L	157	ASN	2.7
1	L	214	CYS	2.6
3	C	437	ASN	2.5
2	A	130	ALA	2.4
2	A	132	THR	2.4
1	L	199	LYS	2.3
2	H	133	ASN	2.2
2	H	1	GLN	2.1
2	A	133	ASN	2.1
1	L	185	GLU	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 [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	SO4	B	216	5/5	0.72	0.19	80,84,85,86	0
4	SO4	L	216	5/5	0.75	0.12	85,86,87,88	0
4	SO4	L	217	5/5	0.84	0.10	82,82,83,83	0
4	SO4	H	214	5/5	0.84	0.12	73,73,75,75	0
4	SO4	A	214	5/5	0.84	0.12	71,71,73,74	0
4	SO4	L	215	5/5	0.84	0.11	63,65,68,70	0
4	SO4	B	215	5/5	0.90	0.08	68,68,69,71	0

6.5 Other polymers [i](#)

There are no such residues in this entry.