



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 19, 2026 – 03:50 AM UTC

PDB ID : 6PDS / pdb_00006pds
Title : Vaccine-elicited NHP FP-targeting antibody 0PV-a.04 in complex with HIV fusion peptide (residue 512-519)
Authors : Xu, K.; Liu, K.; Wang, Y.; Kwong, P.D.
Deposited on : 2019-06-19
Resolution : 1.89 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

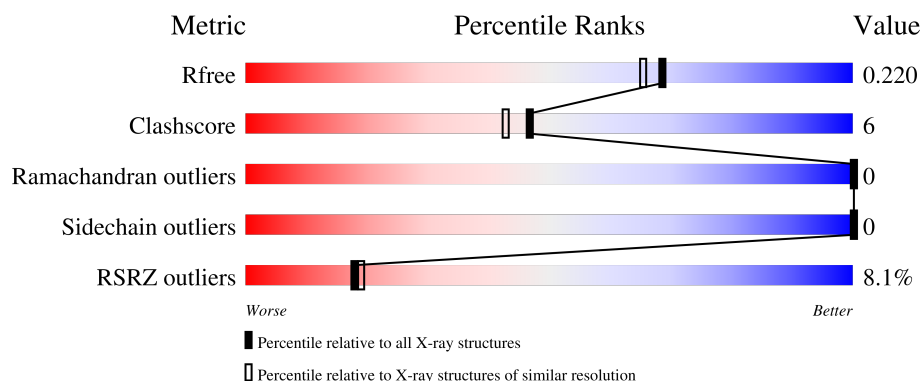
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	212	14% 87% 13%
1	L	212	4% 88% 12%
2	B	220	7% 88% 12%
2	H	220	7% 87% 12%
3	C	8	12% 88% 12%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	G	8	88% 12%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7418 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 0PV-a.04 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	212	Total	C	N	O	S	0	0	0
			1618	1004	276	331	7			
1	L	212	Total	C	N	O	S	0	0	0
			1618	1004	276	331	7			

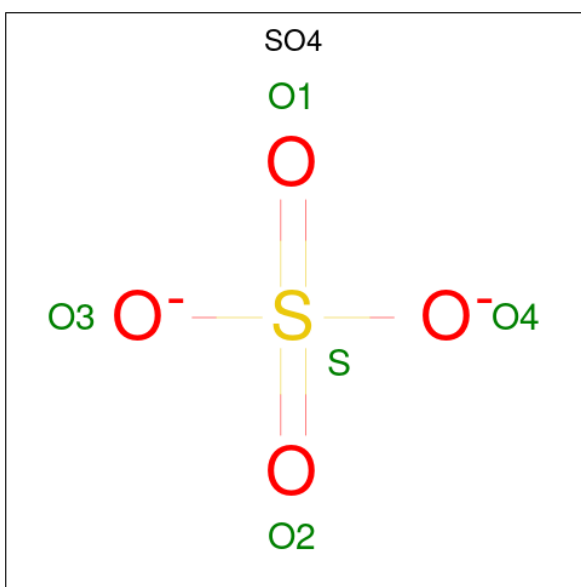
- Molecule 2 is a protein called antibody 0PV-a.04 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	220	Total	C	N	O	S	0	0	0
			1678	1059	284	331	4			
2	H	218	Total	C	N	O	S	0	0	0
			1665	1051	282	328	4			

- Molecule 3 is a protein called HIV-1 fusion peptide residue 512-519.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	G	7	Total	C	N	O	0	0	0
			40	26	7	7			
3	C	8	Total	C	N	O	0	0	0
			51	35	8	8			

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



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

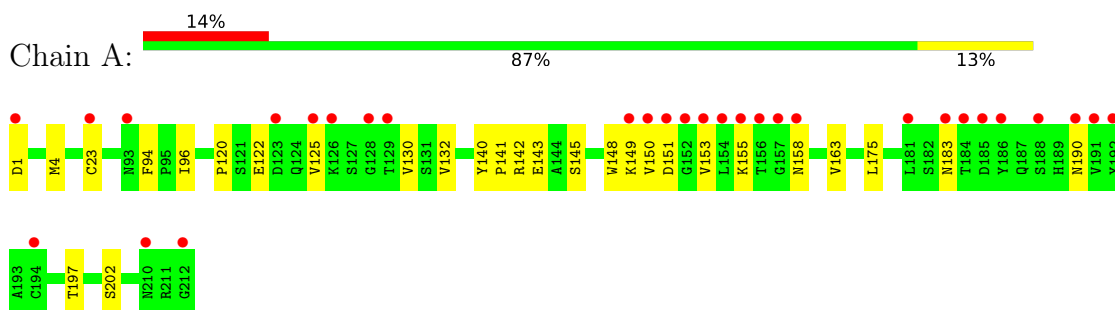
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	139	Total	O	0	0
			139	139		
5	B	196	Total	O	0	0
			196	196		
5	H	215	Total	O	0	0
			215	215		
5	L	164	Total	O	0	0
			164	164		
5	G	8	Total	O	0	0
			8	8		
5	C	6	Total	O	0	0
			6	6		

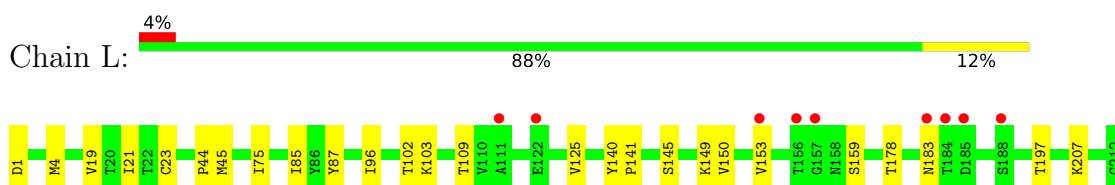
3 Residue-property plots

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.

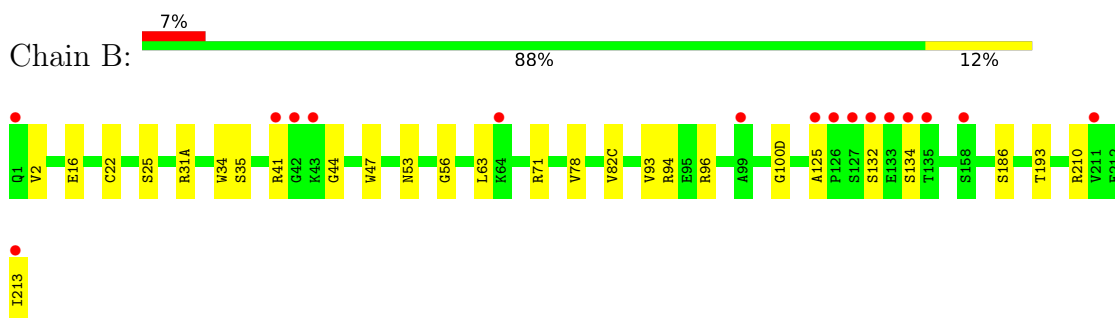
- Molecule 1: antibody 0PV-a.04 light chain



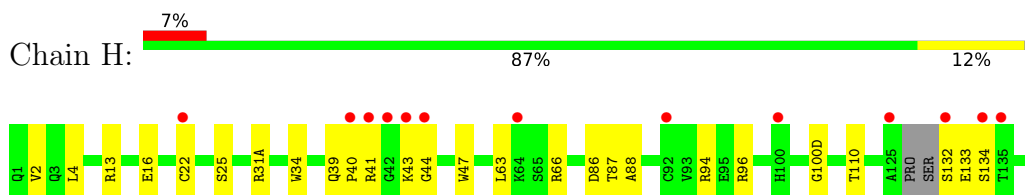
- Molecule 1: antibody 0PV-a.04 light chain



- Molecule 2: antibody 0PV-a.04 heavy chain



- Molecule 2: antibody 0PV-a.04 heavy chain



- Molecule 3: HIV-1 fusion peptide residue 512-519

Chain G:  88% 12%



- Molecule 3: HIV-1 fusion peptide residue 512-519

Chain C:  12% 88% 12%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.86Å 74.38Å 187.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.45 – 1.89 39.45 – 1.89	Depositor EDS
% Data completeness (in resolution range)	96.0 (39.45-1.89) 96.0 (39.45-1.89)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 1.89Å)	Xtriage
Refinement program	PHENIX 1.14_3228	Depositor
R, R_{free}	0.193 , 0.220 0.193 , 0.220	Depositor DCC
R_{free} test set	3794 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	22.9	Xtriage
Anisotropy	0.367	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7418	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.67% 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

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

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	0.25	0/1652	0.55	2/2245 (0.1%)
1	L	0.19	0/1652	0.50	0/2245
2	B	0.22	0/1718	0.54	1/2346 (0.0%)
2	H	0.21	0/1704	0.51	0/2326
3	C	0.10	0/51	0.23	0/68
3	G	0.13	0/39	0.29	0/52
All	All	0.22	0/6816	0.52	3/9282 (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
2	B	0	1
2	H	0	1
All	All	0	2

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	B	44	GLY	N-CA-C	-5.36	100.48	113.18
1	A	149	LYS	CD-CE-NZ	-5.31	94.92	111.90
1	A	149	LYS	CG-CD-CE	5.27	123.43	111.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	63	LEU	Peptide
2	H	63	LEU	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	A	1618	0	1560	20	0
1	L	1618	0	1560	20	0
2	B	1678	0	1645	18	0
2	H	1665	0	1633	21	0
3	C	51	0	53	1	0
3	G	40	0	44	0	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
4	H	5	0	0	0	0
4	L	5	0	0	0	0
5	A	139	0	0	3	0
5	B	196	0	0	5	1
5	C	6	0	0	0	0
5	G	8	0	0	0	0
5	H	215	0	0	9	0
5	L	164	0	0	4	0
All	All	7418	0	6495	75	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 6.

All (75) 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:L:207:LYS:NZ	5:L:401:HOH:O	1.91	1.04
1:A:125:VAL:HA	1:A:183:ASN:HD21	1.29	0.95
1:A:1:ASP:OD2	5:A:401:HOH:O	2.02	0.77
1:A:125:VAL:HA	1:A:183:ASN:ND2	2.01	0.76
1:L:1:ASP:OD2	5:L:402:HOH:O	2.05	0.74
2:H:13:ARG:HB2	2:H:16:GLU:HG3	1.75	0.69

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:41:ARG:HG2	5:B:425:HOH:O	1.94	0.68
2:H:41:ARG:HG3	5:H:582:HOH:O	1.94	0.66
2:B:193:THR:HG23	2:B:210:ARG:HE	1.63	0.63
1:A:158:ASN:HB2	5:A:411:HOH:O	1.97	0.63
1:L:109:THR:HG22	3:C:17:VAL:HG21	1.82	0.59
2:B:193:THR:CG2	2:B:210:ARG:HE	2.16	0.59
2:H:133:GLU:N	5:H:401:HOH:O	2.29	0.58
2:H:34:TRP:CZ2	2:H:94:ARG:HD3	2.42	0.55
1:A:120:PRO:HG3	1:A:130:VAL:HG13	1.89	0.54
2:B:71:ARG:HD2	5:B:569:HOH:O	2.08	0.54
1:L:145:SER:HB3	1:L:197:THR:OG1	2.08	0.54
2:B:34:TRP:CZ2	2:B:94:ARG:HD3	2.43	0.53
2:H:39:GLN:NE2	5:H:403:HOH:O	2.29	0.53
2:B:96:ARG:HD2	2:B:100(D):GLY:C	2.33	0.52
1:L:85:ILE:HG12	1:L:103:LYS:HD3	1.92	0.52
2:B:134:SER:O	2:B:186:SER:HB2	2.11	0.50
2:H:44:GLY:HA2	1:L:87:TYR:CE1	2.47	0.49
1:A:120:PRO:HG3	1:A:130:VAL:CG1	2.42	0.49
2:B:31(A):ARG:HG2	2:B:53:ASN:HB3	1.94	0.49
5:H:529:HOH:O	1:L:45:MET:SD	2.60	0.49
1:L:1:ASP:N	5:L:407:HOH:O	2.44	0.48
1:L:150:VAL:O	1:L:153:VAL:HG22	2.13	0.48
1:A:151:ASP:OD2	1:A:190:ASN:HB2	2.13	0.48
1:A:145:SER:HB3	1:A:197:THR:OG1	2.14	0.48
2:B:213:ILE:H	2:B:213:ILE:HD12	1.77	0.48
1:L:4:MET:HE3	1:L:23:CYS:SG	2.53	0.48
1:L:44:PRO:O	1:L:45:MET:HE2	2.13	0.47
1:A:122:GLU:HG3	1:A:125:VAL:HB	1.97	0.47
2:H:40:PRO:HG3	2:H:88:ALA:HB2	1.97	0.47
2:B:22:CYS:HB3	2:B:78:VAL:HG22	1.97	0.46
2:B:35:SER:HB2	2:B:93:VAL:HG13	1.98	0.46
2:H:2:VAL:HA	2:H:25:SER:O	2.15	0.46
2:B:56:GLY:O	2:B:71:ARG:NH1	2.49	0.46
1:A:163:VAL:HG22	1:A:175:LEU:HD12	1.98	0.45
1:A:120:PRO:HD3	1:A:132:VAL:HG22	1.99	0.45
2:B:2:VAL:HA	2:B:25:SER:O	2.17	0.45
2:H:134:SER:N	5:H:401:HOH:O	2.09	0.45
1:L:21:ILE:HG12	1:L:102:THR:HG21	1.98	0.45
2:H:44:GLY:HA2	1:L:87:TYR:HE1	1.82	0.44
1:A:96:ILE:HB	2:B:47:TRP:CG	2.52	0.44
2:H:31(A):ARG:NH2	5:H:402:HOH:O	2.12	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:4:LEU:HB3	2:H:22:CYS:SG	2.58	0.43
1:A:143:GLU:CD	1:A:143:GLU:H	2.26	0.43
2:H:40:PRO:HG2	5:H:409:HOH:O	2.19	0.43
2:H:66:ARG:NH1	2:H:86:ASP:OD2	2.50	0.43
2:H:47:TRP:CG	1:L:96:ILE:HB	2.54	0.43
1:A:142:ARG:HE	1:A:142:ARG:HB3	1.71	0.43
1:L:140:TYR:CG	1:L:141:PRO:HA	2.54	0.43
1:A:202:SER:HB3	5:A:518:HOH:O	2.19	0.42
2:H:132:SER:N	5:H:401:HOH:O	2.52	0.42
1:A:4:MET:HE3	1:A:23:CYS:SG	2.59	0.42
1:A:140:TYR:CG	1:A:141:PRO:HA	2.54	0.42
1:A:150:VAL:O	1:A:153:VAL:HG22	2.19	0.42
2:B:125:ALA:HB2	5:B:465:HOH:O	2.19	0.42
2:B:132:SER:N	5:B:407:HOH:O	2.51	0.42
2:H:43:LYS:NZ	2:H:43:LYS:HB3	2.34	0.42
1:L:183:ASN:HD22	1:L:183:ASN:HA	1.67	0.42
2:H:40:PRO:HG3	5:H:412:HOH:O	2.20	0.42
2:H:44:GLY:HA3	5:L:452:HOH:O	2.19	0.41
1:L:159:SER:HA	1:L:178:THR:O	2.21	0.41
2:H:96:ARG:HD3	2:H:100(D):GLY:C	2.46	0.41
1:A:94:PHE:HB3	5:B:463:HOH:O	2.19	0.41
1:L:19:VAL:HG22	1:L:75:ILE:HB	2.03	0.41
2:H:87:THR:HG23	2:H:110:THR:HA	2.01	0.41
1:L:149:LYS:HA	1:L:153:VAL:O	2.20	0.41
1:L:125:VAL:O	1:L:183:ASN:OD1	2.39	0.40
2:B:35:SER:HB2	2:B:93:VAL:CG1	2.51	0.40
2:B:16:GLU:O	2:B:82(C):VAL:HG22	2.21	0.40
1:A:148:TRP:HB2	1:A:155:LYS:O	2.21	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)
5:B:542:HOH:O	5:B:575:HOH:O[4_545]	2.18	0.02

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	210/212 (99%)	205 (98%)	5 (2%)	0	100	100
1	L	210/212 (99%)	205 (98%)	5 (2%)	0	100	100
2	B	216/220 (98%)	211 (98%)	5 (2%)	0	100	100
2	H	214/220 (97%)	207 (97%)	7 (3%)	0	100	100
3	C	6/8 (75%)	6 (100%)	0	0	100	100
3	G	5/8 (62%)	5 (100%)	0	0	100	100
All	All	861/880 (98%)	839 (97%)	22 (3%)	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	A	187/187 (100%)	187 (100%)	0	100	100
1	L	187/187 (100%)	187 (100%)	0	100	100
2	B	194/194 (100%)	194 (100%)	0	100	100
2	H	192/194 (99%)	192 (100%)	0	100	100
3	C	4/4 (100%)	4 (100%)	0	100	100
3	G	3/4 (75%)	3 (100%)	0	100	100
All	All	767/770 (100%)	767 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

Mol	Chain	Res	Type
1	A	31	ASN
1	A	183	ASN
2	H	171	GLN
1	L	187	GLN
1	L	210	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](#)

4 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	L	301	-	4,4,4	0.21	0	6,6,6	0.21	0
4	SO4	B	301	-	4,4,4	0.23	0	6,6,6	0.10	0
4	SO4	H	301	-	4,4,4	0.25	0	6,6,6	0.10	0
4	SO4	A	301	-	4,4,4	0.24	0	6,6,6	0.09	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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	127:SER	C	132:SER	N	13.13

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	212/212 (100%)	0.76	30 (14%) 6 6	17, 30, 61, 75	0
1	L	212/212 (100%)	0.41	9 (4%) 40 43	15, 28, 48, 62	0
2	B	220/220 (100%)	0.33	16 (7%) 21 22	14, 25, 55, 76	0
2	H	218/220 (99%)	0.13	15 (6%) 23 24	11, 22, 52, 78	0
3	C	8/8 (100%)	0.86	1 (12%) 8 8	21, 27, 41, 50	0
3	G	7/8 (87%)	-0.22	0 100 100	16, 17, 23, 30	0
All	All	877/880 (99%)	0.40	71 (8%) 18 19	11, 26, 55, 78	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	156	THR	5.4
2	H	92	CYS	5.0
1	A	153	VAL	4.9
2	H	42	GLY	4.7
1	A	190	ASN	4.7
1	A	191	VAL	4.6
1	A	154	LEU	4.6
2	H	213	ILE	4.4
1	L	157	GLY	4.2
1	A	157	GLY	4.0
2	H	43	LYS	3.9
1	A	181	LEU	3.8
2	B	126	PRO	3.7
1	A	212	GLY	3.5
2	B	41	ARG	3.5
2	B	133	GLU	3.4
2	B	127	SER	3.4
3	C	18	PHE	3.3
2	H	134	SER	3.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	B	135	THR	3.3
2	B	132	SER	3.3
2	B	1	GLN	3.2
2	B	134	SER	3.2
2	H	22	CYS	3.2
2	H	40	PRO	3.2
2	H	132	SER	3.0
1	L	156	THR	2.9
2	B	213	ILE	2.8
1	L	184	THR	2.8
2	B	125	ALA	2.8
2	H	44	GLY	2.8
1	A	149	LYS	2.7
1	A	123	ASP	2.7
1	A	151	ASP	2.7
1	A	192	TYR	2.7
2	H	41	ARG	2.7
2	H	187	SER	2.7
2	B	99	ALA	2.6
1	A	183	ASN	2.6
1	A	150	VAL	2.6
1	A	128	GLY	2.6
1	A	93	ASN	2.6
2	B	42	GLY	2.6
2	H	125	ALA	2.6
1	L	188	SER	2.6
1	A	194	CYS	2.5
2	B	43	LYS	2.5
1	L	183	ASN	2.5
1	L	153	VAL	2.5
1	A	125	VAL	2.4
1	A	155	LYS	2.4
1	A	184	THR	2.4
2	B	211	VAL	2.3
1	A	23	CYS	2.3
1	L	122	GLU	2.2
1	L	185	ASP	2.2
1	A	210	ASN	2.2
2	H	64	LYS	2.2
1	A	1	ASP	2.1
2	H	135	THR	2.1
2	B	158	SER	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	126	LYS	2.1
2	B	64	LYS	2.1
1	A	152	GLY	2.1
1	A	158	ASN	2.1
1	L	111	ALA	2.1
1	A	185	ASP	2.1
2	H	100	HIS	2.0
1	A	188	SER	2.0
1	A	129	THR	2.0
1	A	186	TYR	2.0

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 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	A	301	5/5	0.85	0.09	45,53,56,58	0
4	SO4	H	301	5/5	0.90	0.08	41,53,64,65	0
4	SO4	B	301	5/5	0.95	0.11	38,41,46,46	0
4	SO4	L	301	5/5	0.98	0.07	26,26,28,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.