



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 07:47 pm BST

PDB ID : 6M87
Title : Fab 10A6 in complex with MPTS
Authors : Stanfield, R.L.; Wilson, I.A.
Deposited on : 2018-08-21
Resolution : 2.61 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
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
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)	:	2.11

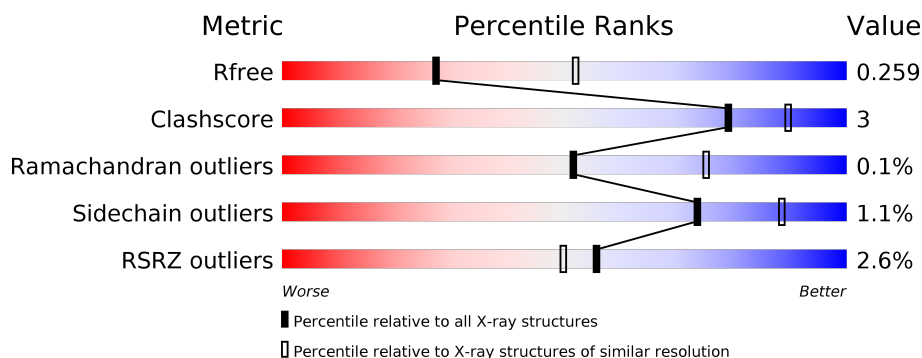
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.61 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 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	211	<div><div></div><div>92%8%</div></div>
1	C	211	<div><div>%</div><div>89%10%.</div></div>
1	E	211	<div><div>4%</div><div>92%8%</div></div>
1	G	211	<div><div>%</div><div>91%8%.</div></div>
1	I	211	<div><div>4%</div><div>91%7%.</div></div>
1	K	211	<div><div>9%</div><div>91%7%..</div></div>

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Mol	Chain	Length	Quality of chain
2	B	219	
2	D	219	
2	F	219	
2	H	219	
2	J	219	
2	L	219	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SO4	L	301	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 19716 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.

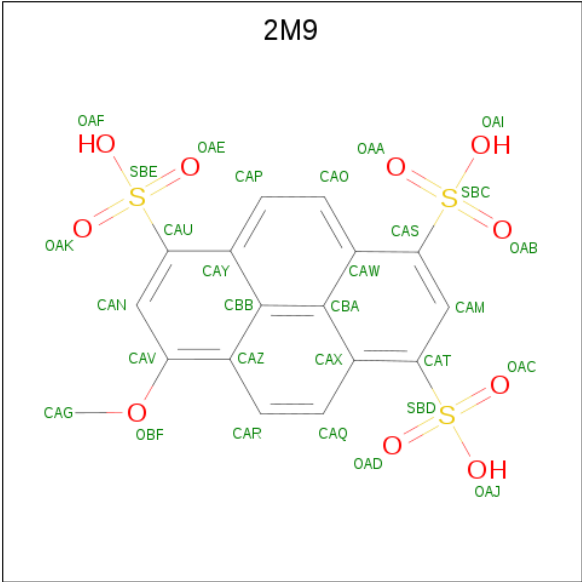
- Molecule 1 is a protein called Fab 10A6 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	211	Total	C	N	O	S	0	0	0
			1622	1009	274	331	8			
1	C	209	Total	C	N	O	S	0	0	0
			1610	1003	271	328	8			
1	E	211	Total	C	N	O	S	0	0	0
			1622	1009	274	331	8			
1	G	209	Total	C	N	O	S	0	0	0
			1610	1003	271	328	8			
1	I	208	Total	C	N	O	S	0	0	0
			1603	998	270	327	8			
1	K	207	Total	C	N	O	S	0	0	0
			1589	989	269	323	8			

- Molecule 2 is a protein called Fab 10A6 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	215	Total	C	N	O	S	0	0	0
			1613	1015	266	326	6			
2	D	216	Total	C	N	O	S	0	0	0
			1621	1020	267	328	6			
2	F	215	Total	C	N	O	S	0	0	0
			1613	1015	266	326	6			
2	H	215	Total	C	N	O	S	0	0	0
			1613	1015	266	326	6			
2	J	216	Total	C	N	O	S	0	0	0
			1621	1020	267	328	6			
2	L	216	Total	C	N	O	S	0	0	0
			1621	1020	267	328	6			

- Molecule 3 is 8-methoxypyrene-1,3,6-trisulfonic acid (three-letter code: 2M9) (formula: C₁₇H₁₂O₁₀S₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	O	S	0	0
			30	17	10	3		
3	E	1	Total	C	O	S	0	0
			30	17	10	3		
3	G	1	Total	C	O	S	0	0
			30	17	10	3		
3	K	1	Total	C	O	S	0	0
			30	17	10	3		
3	D	1	Total	C	O	S	0	0
			30	17	10	3		
3	J	1	Total	C	O	S	0	0
			30	17	10	3		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		
4	G	1	Total	O	S	0	0
			5	4	1		
4	I	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	F	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	15	Total	O	0	0
			15	15		
5	B	15	Total	O	0	0
			15	15		
5	C	8	Total	O	0	0
			8	8		
5	E	13	Total	O	0	0
			13	13		
5	G	10	Total	O	0	0
			10	10		
5	I	17	Total	O	0	0
			17	17		

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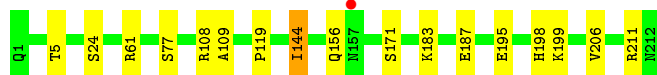
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	K	12	Total 12	O 12	0	0
5	D	12	Total 12	O 12	0	0
5	F	7	Total 7	O 7	0	0
5	H	13	Total 13	O 13	0	0
5	J	8	Total 8	O 8	0	0
5	L	18	Total 18	O 18	0	0

3 Residue-property plots [i](#)


These plots are drawn for all protein, RNA and DNA 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.

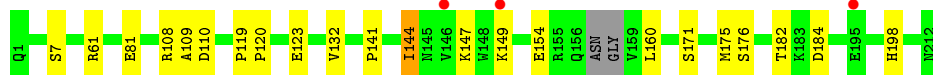
- Molecule 1: Fab 10A6 light chain

Chain A: 



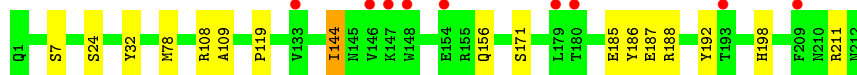
- Molecule 1: Fab 10A6 light chain

Chain C: 




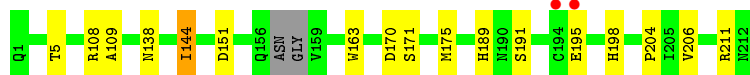
- Molecule 1: Fab 10A6 light chain

Chain E: 




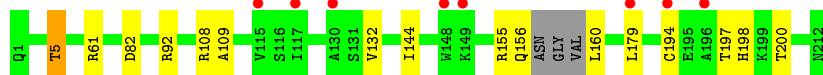
- Molecule 1: Fab 10A6 light chain

Chain G: 

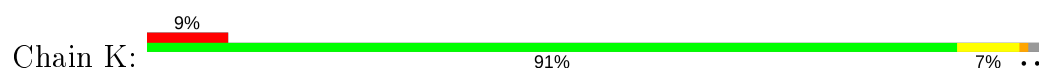


- Molecule 1: Fab 10A6 light chain

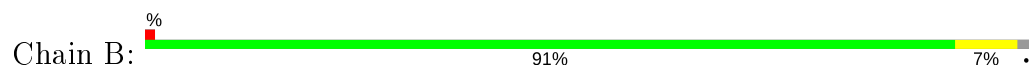
Chain I: 



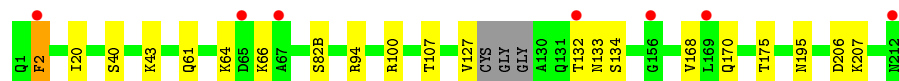
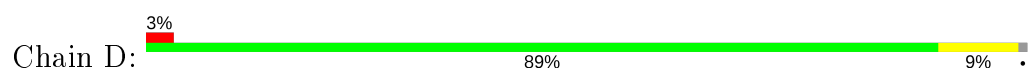
- Molecule 1: Fab 10A6 light chain



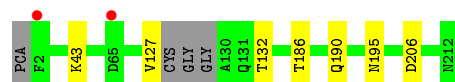
- Molecule 2: Fab 10A6 heavy chain



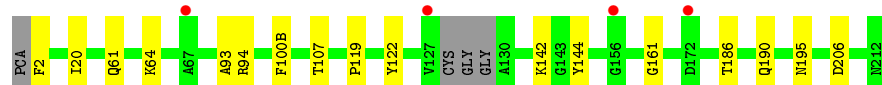
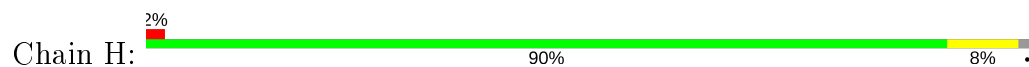
- Molecule 2: Fab 10A6 heavy chain



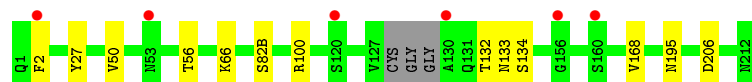
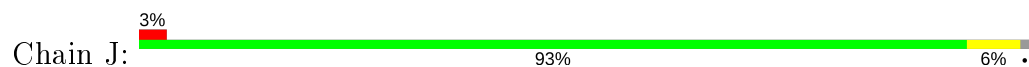
- Molecule 2: Fab 10A6 heavy chain



- Molecule 2: Fab 10A6 heavy chain



- Molecule 2: Fab 10A6 heavy chain



- Molecule 2: Fab 10A6 heavy chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	186.49 Å 186.49 Å 89.91 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.79 – 2.61 44.79 – 2.61	Depositor EDS
% Data completeness (in resolution range)	99.7 (44.79-2.61) 99.7 (44.79-2.61)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 2.61 Å)	Xtriage
Refinement program	PHENIX (1.12_2829)	Depositor
R, R_{free}	0.228 , 0.265 0.224 , 0.259	Depositor DCC
R_{free} test set	5152 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	61.3	Xtriage
Anisotropy	0.319	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.034 for -h,-k,l 0.033 for h,-h-k,-l 0.477 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	19716	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 32.06 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 9.9879e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: PCA, SO4, 2M9

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	0/1655	0.47	0/2252
1	C	0.24	0/1642	0.48	0/2233
1	E	0.25	0/1655	0.47	0/2252
1	G	0.25	0/1642	0.47	0/2233
1	I	0.24	0/1635	0.47	0/2223
1	K	0.24	0/1619	0.48	0/2200
2	B	0.25	0/1653	0.47	0/2259
2	D	0.26	0/1654	0.47	0/2262
2	F	0.25	0/1653	0.46	0/2259
2	H	0.25	0/1653	0.47	0/2259
2	J	0.26	0/1654	0.48	0/2262
2	L	0.25	0/1654	0.46	0/2262
All	All	0.25	0/19769	0.47	0/26956

There are no bond length outliers.

There are no bond angle outliers.

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	A	1622	0	1549	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1610	0	1539	15	0
1	E	1622	0	1549	11	0
1	G	1610	0	1539	12	0
1	I	1603	0	1530	12	0
1	K	1589	0	1523	11	0
2	B	1613	0	1560	11	0
2	D	1621	0	1567	14	0
2	F	1613	0	1560	3	0
2	H	1613	0	1560	8	0
2	J	1621	0	1567	9	0
2	L	1621	0	1567	4	0
3	B	30	0	10	0	0
3	D	30	0	11	0	0
3	E	30	0	11	0	0
3	G	30	0	10	0	0
3	J	30	0	10	1	0
3	K	30	0	11	0	0
4	B	5	0	0	1	0
4	D	5	0	0	1	0
4	F	5	0	0	0	0
4	G	5	0	0	0	0
4	I	5	0	0	1	0
4	L	5	0	0	2	0
5	A	15	0	0	0	0
5	B	15	0	0	0	0
5	C	8	0	0	0	0
5	D	12	0	0	0	0
5	E	13	0	0	0	0
5	F	7	0	0	0	0
5	G	10	0	0	0	0
5	H	13	0	0	0	0
5	I	17	0	0	0	0
5	J	8	0	0	0	0
5	K	12	0	0	1	0
5	L	18	0	0	0	0
All	All	19716	0	18673	107	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 3.

All (107) 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:E:144:ILE:HG13	1:E:198:HIS:HD2	1.37	0.86
1:E:144:ILE:HG13	1:E:198:HIS:CD2	2.25	0.70
1:A:144:ILE:HG13	1:A:198:HIS:HD2	1.59	0.68
1:K:156:GLN:NE2	5:K:401:HOH:O	2.26	0.67
1:C:144:ILE:HG13	1:C:198:HIS:HD2	1.59	0.67
1:I:144:ILE:HD11	1:I:198:HIS:CD2	2.31	0.65
2:L:100:ARG:NH2	4:L:301:SO4:O4	2.29	0.65
1:K:182:THR:HG22	1:K:183:LYS:H	1.61	0.64
1:C:144:ILE:HG13	1:C:198:HIS:CD2	2.33	0.62
1:K:108:ARG:HD3	1:K:109:ALA:O	2.00	0.62
1:K:187:GLU:O	1:K:211:ARG:NH1	2.32	0.61
2:B:158:LEU:N	3:J:301:2M9:OAE	2.33	0.61
1:K:53:LYS:NZ	4:L:301:SO4:O2	2.35	0.59
2:H:122:TYR:HE2	2:H:142:LYS:HD3	1.68	0.59
1:I:61:ARG:NH1	1:I:82:ASP:OD2	2.37	0.58
1:A:108:ARG:NH1	1:A:109:ALA:O	2.37	0.57
1:I:108:ARG:NH1	1:I:109:ALA:O	2.37	0.57
1:G:189:HIS:O	1:G:211:ARG:HD2	2.03	0.57
1:A:119:PRO:HG3	2:B:127:VAL:HG21	1.86	0.57
2:H:61:GLN:HG3	2:H:64:LYS:HE3	1.86	0.56
2:L:40:SER:O	2:L:43:LYS:HG2	2.06	0.56
1:C:182:THR:OG1	1:C:184:ASP:OD1	2.24	0.55
1:E:185:GLU:HA	1:E:188:ARG:NH1	2.20	0.55
1:A:144:ILE:HG13	1:A:198:HIS:CD2	2.39	0.54
1:G:108:ARG:NH1	1:G:109:ALA:O	2.40	0.54
1:G:144:ILE:HG13	1:G:198:HIS:HD2	1.72	0.54
1:E:185:GLU:HA	1:E:188:ARG:HH11	1.72	0.54
2:H:195:ASN:ND2	2:H:206:ASP:OD1	2.40	0.53
1:C:175:MET:HG2	1:C:176:SER:N	2.23	0.53
1:A:195:GLU:HG3	1:A:206:VAL:HG22	1.91	0.53
1:C:108:ARG:NH1	1:C:109:ALA:O	2.42	0.53
2:J:2:PHE:CD1	2:J:27:TYR:HB3	2.44	0.52
1:K:182:THR:HG22	1:K:183:LYS:N	2.25	0.52
2:B:195:ASN:ND2	2:B:206:ASP:OD1	2.41	0.52
1:E:187:GLU:OE1	1:E:211:ARG:NH2	2.43	0.52
2:F:195:ASN:ND2	2:F:206:ASP:OD1	2.43	0.52
2:L:195:ASN:ND2	2:L:206:ASP:OD1	2.43	0.52
1:K:195:GLU:HG2	1:K:206:VAL:HG22	1.93	0.51
1:G:144:ILE:HG13	1:G:198:HIS:CD2	2.46	0.50
2:J:195:ASN:ND2	2:J:206:ASP:OD1	2.43	0.50
4:I:301:SO4:O2	2:J:100:ARG:NE	2.43	0.49
1:I:61:ARG:NH2	1:I:82:ASP:OD1	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:GLU:OE1	1:A:211:ARG:NH2	2.45	0.49
1:C:120:PRO:HD3	1:C:132:VAL:HG22	1.94	0.48
1:E:119:PRO:HG3	2:F:127:VAL:HG21	1.96	0.48
2:D:133:ASN:OD1	2:D:134:SER:N	2.43	0.48
1:G:195:GLU:HG2	1:G:204:PRO:HB2	1.95	0.48
1:G:151:ASP:OD1	1:G:191:SER:N	2.45	0.48
1:E:108:ARG:NH1	1:E:109:ALA:O	2.47	0.47
2:D:61:GLN:HG3	2:D:64:LYS:HE3	1.96	0.47
1:A:24:SER:HB2	1:K:5:THR:HB	1.95	0.47
2:B:40:SER:O	2:B:43:LYS:HG2	2.14	0.47
1:I:198:HIS:CE1	1:I:200:THR:HG23	2.50	0.47
2:D:40:SER:O	2:D:43:LYS:HG2	2.15	0.47
1:C:123:GLU:OE1	2:D:207:LYS:NZ	2.43	0.47
2:J:66:LYS:NZ	2:J:82(B):SER:O	2.48	0.47
2:L:61:GLN:O	2:L:64:LYS:HG2	2.15	0.46
2:B:158:LEU:HG	2:J:56:THR:HG23	1.97	0.46
2:B:66:LYS:NZ	2:B:82(B):SER:O	2.49	0.46
2:B:158:LEU:HD22	2:J:50:VAL:HG13	1.98	0.46
2:H:20:ILE:HD12	2:H:107:THR:HG21	1.98	0.46
1:K:78:MET:HE1	1:K:106:ILE:HG13	1.98	0.45
2:B:20:ILE:HD12	2:B:107:THR:HG21	1.98	0.45
1:C:119:PRO:HG3	2:D:127:VAL:HG21	1.97	0.45
1:I:144:ILE:HD13	1:I:144:ILE:HA	1.83	0.45
1:K:147:LYS:HG3	1:K:195:GLU:HB2	1.99	0.45
1:C:160:LEU:HD11	2:D:168:VAL:HB	1.99	0.45
2:H:93:ALA:HB1	2:H:100(B):PHE:HB3	1.97	0.45
1:G:195:GLU:HG3	1:G:206:VAL:HG22	1.99	0.44
1:I:132:VAL:HG13	1:I:179:LEU:HB3	2.00	0.44
1:E:186:TYR:O	1:E:192:TYR:OH	2.33	0.44
2:D:2:PHE:HD1	2:D:94:ARG:NH1	2.16	0.44
2:D:20:ILE:HD12	2:D:107:THR:HG21	2.00	0.43
1:C:61:ARG:NH2	1:C:81:GLU:OE1	2.51	0.43
2:F:186:THR:O	2:F:190:GLN:N	2.51	0.43
1:G:138:ASN:ND2	1:G:170:ASP:OD2	2.50	0.43
2:H:186:THR:O	2:H:190:GLN:N	2.52	0.43
2:D:195:ASN:ND2	2:D:206:ASP:OD1	2.52	0.43
4:B:302:SO4:O1	1:G:108:ARG:NH2	2.52	0.43
1:C:7:SER:HB3	1:G:5:THR:HG22	2.00	0.43
2:D:170:GLN:NE2	2:D:175:THR:OG1	2.45	0.43
1:K:128:GLY:O	1:K:182:THR:HG23	2.18	0.43
2:B:186:THR:O	2:B:190:GLN:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:163:TRP:CD1	1:G:175:MET:HG3	2.54	0.42
2:D:100:ARG:NE	4:D:302:SO4:O4	2.53	0.42
1:E:24:SER:HB2	1:I:5:THR:HB	1.99	0.42
1:E:108:ARG:HD2	1:E:171:SER:HB2	2.02	0.42
2:D:2:PHE:HD2	2:D:2:PHE:H	1.68	0.42
2:J:133:ASN:OD1	2:J:134:SER:N	2.48	0.42
1:I:160:LEU:HD21	2:J:168:VAL:CG2	2.50	0.41
2:D:66:LYS:NZ	2:D:82(B):SER:O	2.53	0.41
1:A:61:ARG:HD2	1:A:77:SER:O	2.20	0.41
2:B:61:GLN:HG3	2:B:64:LYS:HE3	2.01	0.41
1:C:147:LYS:HD3	1:C:154:GLU:OE2	2.20	0.41
2:H:2:PHE:HD1	2:H:94:ARG:NH1	2.18	0.41
1:C:108:ARG:HD2	1:C:171:SER:HB2	2.02	0.41
1:I:160:LEU:HD11	2:J:168:VAL:HB	2.02	0.41
1:G:108:ARG:HD2	1:G:171:SER:HB2	2.02	0.41
2:H:119:PRO:HB3	2:H:144:TYR:HB3	2.02	0.41
1:A:183:LYS:O	1:A:187:GLU:HG2	2.20	0.41
2:B:2:PHE:N	2:B:2:PHE:CD1	2.89	0.41
2:D:2:PHE:N	2:D:2:PHE:CD2	2.90	0.40
1:I:155:ARG:NE	1:I:156:GLN:O	2.54	0.40
1:E:7:SER:HB2	1:I:5:THR:HG23	2.03	0.40
1:C:110:ASP:OD1	1:C:141:PRO:HD3	2.22	0.40
1:A:108:ARG:HD2	1:A:171:SER:HB2	2.04	0.40
1:C:149:LYS:HG2	1:C:154:GLU:HA	2.03	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	A	209/211 (99%)	197 (94%)	12 (6%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	205/211 (97%)	195 (95%)	10 (5%)	0	100	100
1	E	209/211 (99%)	199 (95%)	10 (5%)	0	100	100
1	G	205/211 (97%)	195 (95%)	10 (5%)	0	100	100
1	I	204/211 (97%)	194 (95%)	10 (5%)	0	100	100
1	K	201/211 (95%)	189 (94%)	12 (6%)	0	100	100
2	B	211/219 (96%)	201 (95%)	10 (5%)	0	100	100
2	D	212/219 (97%)	202 (95%)	10 (5%)	0	100	100
2	F	211/219 (96%)	202 (96%)	8 (4%)	1 (0%)	29	52
2	H	211/219 (96%)	201 (95%)	9 (4%)	1 (0%)	29	52
2	J	212/219 (97%)	201 (95%)	11 (5%)	0	100	100
2	L	212/219 (97%)	201 (95%)	11 (5%)	0	100	100
All	All	2502/2580 (97%)	2377 (95%)	123 (5%)	2 (0%)	51	75

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	43	LYS
2	H	161	GLY

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	184/184 (100%)	180 (98%)	4 (2%)	52	76
1	C	183/184 (100%)	182 (100%)	1 (0%)	88	96
1	E	184/184 (100%)	180 (98%)	4 (2%)	52	76
1	G	183/184 (100%)	182 (100%)	1 (0%)	88	96
1	I	182/184 (99%)	178 (98%)	4 (2%)	52	76
1	K	181/184 (98%)	178 (98%)	3 (2%)	60	81
2	B	186/187 (100%)	185 (100%)	1 (0%)	88	96

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	186/187 (100%)	184 (99%)	2 (1%)	73	88
2	F	186/187 (100%)	185 (100%)	1 (0%)	88	96
2	H	186/187 (100%)	186 (100%)	0	100	100
2	J	186/187 (100%)	185 (100%)	1 (0%)	88	96
2	L	186/187 (100%)	183 (98%)	3 (2%)	62	82
All	All	2213/2226 (99%)	2188 (99%)	25 (1%)	73	88

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	THR
1	A	144	ILE
1	A	156	GLN
1	A	199	LYS
2	B	113	SER
1	C	144	ILE
1	E	32	TYR
1	E	78	MET
1	E	144	ILE
1	E	156	GLN
1	G	144	ILE
1	I	5	THR
1	I	92	ARG
1	I	194	CYS
1	I	197	THR
1	K	5	THR
1	K	108	ARG
1	K	160	LEU
2	D	2	PHE
2	D	132	THR
2	F	132	THR
2	J	132	THR
2	L	2	PHE
2	L	56	THR
2	L	64	LYS

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

Mol	Chain	Res	Type
2	B	212	ASN
2	H	33	ASN
2	L	212	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

9 non-standard protein/DNA/RNA residues 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
1	PCA	A	1	1	7,8,9	1.79	1 (14%)	9,10,12	2.14	5 (55%)
1	PCA	G	1	1	7,8,9	1.79	1 (14%)	9,10,12	2.14	5 (55%)
2	PCA	L	1	2	7,8,9	1.80	1 (14%)	9,10,12	2.15	5 (55%)
2	PCA	J	1	2	7,8,9	1.79	1 (14%)	9,10,12	2.13	5 (55%)
2	PCA	D	1	2	7,8,9	1.81	1 (14%)	9,10,12	2.09	5 (55%)
1	PCA	K	1	1	7,8,9	1.79	1 (14%)	9,10,12	2.15	5 (55%)
1	PCA	C	1	1	7,8,9	1.80	1 (14%)	9,10,12	2.15	5 (55%)
1	PCA	I	1	1	7,8,9	1.79	1 (14%)	9,10,12	2.15	5 (55%)
1	PCA	E	1	1	7,8,9	1.79	1 (14%)	9,10,12	2.15	5 (55%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	PCA	A	1	1	-	0/0/11/13	0/1/1/1
1	PCA	G	1	1	-	0/0/11/13	0/1/1/1
2	PCA	L	1	2	-	0/0/11/13	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PCA	J	1	2	-	0/0/11/13	0/1/1/1
2	PCA	D	1	2	-	0/0/11/13	0/1/1/1
1	PCA	K	1	1	-	0/0/11/13	0/1/1/1
1	PCA	C	1	1	-	0/0/11/13	0/1/1/1
1	PCA	I	1	1	-	0/0/11/13	0/1/1/1
1	PCA	E	1	1	-	0/0/11/13	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	PCA	CD-N	4.67	1.46	1.34
1	C	1	PCA	CD-N	4.65	1.46	1.34
2	L	1	PCA	CD-N	4.64	1.46	1.34
1	I	1	PCA	CD-N	4.63	1.46	1.34
1	K	1	PCA	CD-N	4.61	1.46	1.34
1	G	1	PCA	CD-N	4.61	1.46	1.34
1	E	1	PCA	CD-N	4.61	1.46	1.34
1	A	1	PCA	CD-N	4.61	1.46	1.34
2	J	1	PCA	CD-N	4.61	1.46	1.34

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	1	PCA	OE-CD-CG	-3.14	121.28	126.76
2	J	1	PCA	OE-CD-CG	-3.13	121.30	126.76
1	A	1	PCA	OE-CD-CG	-3.13	121.31	126.76
1	K	1	PCA	OE-CD-CG	-3.13	121.31	126.76
1	C	1	PCA	OE-CD-CG	-3.12	121.31	126.76
1	I	1	PCA	OE-CD-CG	-3.12	121.33	126.76
2	L	1	PCA	OE-CD-CG	-3.10	121.35	126.76
1	G	1	PCA	OE-CD-CG	-3.10	121.35	126.76
2	D	1	PCA	OE-CD-CG	-3.08	121.40	126.76
2	D	1	PCA	CA-N-CD	-2.90	103.65	113.58
2	L	1	PCA	CA-N-CD	-2.86	103.78	113.58
2	J	1	PCA	CA-N-CD	-2.83	103.89	113.58
1	K	1	PCA	CA-N-CD	-2.81	103.96	113.58
1	I	1	PCA	CA-N-CD	-2.81	103.97	113.58
1	E	1	PCA	CA-N-CD	-2.80	103.98	113.58
1	C	1	PCA	CA-N-CD	-2.80	103.99	113.58
1	G	1	PCA	CA-N-CD	-2.80	104.00	113.58
1	A	1	PCA	CA-N-CD	-2.78	104.05	113.58
1	K	1	PCA	CB-CA-N	2.70	111.06	103.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	I	1	PCA	CB-CA-N	2.69	111.03	103.30
1	G	1	PCA	CB-CA-N	2.68	110.99	103.30
1	C	1	PCA	CB-CA-N	2.68	110.99	103.30
1	E	1	PCA	CB-CA-N	2.68	110.99	103.30
2	L	1	PCA	CB-CA-N	2.67	110.97	103.30
1	A	1	PCA	CB-CA-N	2.66	110.93	103.30
2	D	1	PCA	CB-CA-N	2.65	110.92	103.30
2	J	1	PCA	CB-CA-N	2.63	110.86	103.30
2	J	1	PCA	CB-CA-C	-2.60	109.13	112.70
1	I	1	PCA	CB-CA-C	-2.59	109.14	112.70
1	K	1	PCA	CB-CA-C	-2.59	109.14	112.70
1	E	1	PCA	CB-CA-C	-2.59	109.14	112.70
1	G	1	PCA	CB-CA-C	-2.57	109.17	112.70
1	C	1	PCA	CB-CA-C	-2.57	109.17	112.70
1	A	1	PCA	CB-CA-C	-2.56	109.18	112.70
1	E	1	PCA	CG-CD-N	2.51	114.89	108.39
1	I	1	PCA	CG-CD-N	2.49	114.85	108.39
1	K	1	PCA	CG-CD-N	2.49	114.84	108.39
1	G	1	PCA	CG-CD-N	2.49	114.83	108.39
1	C	1	PCA	CG-CD-N	2.48	114.82	108.39
1	A	1	PCA	CG-CD-N	2.48	114.82	108.39
2	L	1	PCA	CG-CD-N	2.48	114.81	108.39
2	J	1	PCA	CG-CD-N	2.48	114.80	108.39
2	L	1	PCA	CB-CA-C	-2.45	109.33	112.70
2	D	1	PCA	CG-CD-N	2.43	114.67	108.39
2	D	1	PCA	CB-CA-C	-2.16	109.73	112.70

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

12 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	$\# Z > 2$
3	2M9	J	301	-	30,33,33	0.36	0	46,55,55	0.51	0
4	SO4	G	302	-	4,4,4	0.14	0	6,6,6	0.05	0
4	SO4	D	302	-	4,4,4	0.14	0	6,6,6	0.04	0
3	2M9	G	301	-	30,33,33	0.35	0	46,55,55	0.52	0
3	2M9	E	301	-	30,33,33	0.35	0	46,55,55	0.48	0
4	SO4	L	301	-	4,4,4	0.15	0	6,6,6	0.06	0
4	SO4	F	301	-	4,4,4	0.14	0	6,6,6	0.06	0
3	2M9	B	301	-	30,33,33	0.33	0	46,55,55	0.50	0
3	2M9	K	301	-	30,33,33	0.33	0	46,55,55	0.46	0
4	SO4	I	301	-	4,4,4	0.14	0	6,6,6	0.05	0
4	SO4	B	302	-	4,4,4	0.14	0	6,6,6	0.07	0
3	2M9	D	301	-	30,33,33	0.34	0	46,55,55	0.47	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	2M9	J	301	-	-	0/20/20/20	0/4/4/4
3	2M9	G	301	-	-	1/20/20/20	0/4/4/4
3	2M9	E	301	-	-	0/20/20/20	0/4/4/4
3	2M9	B	301	-	-	0/20/20/20	0/4/4/4
3	2M9	K	301	-	-	0/20/20/20	0/4/4/4
3	2M9	D	301	-	-	0/20/20/20	0/4/4/4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

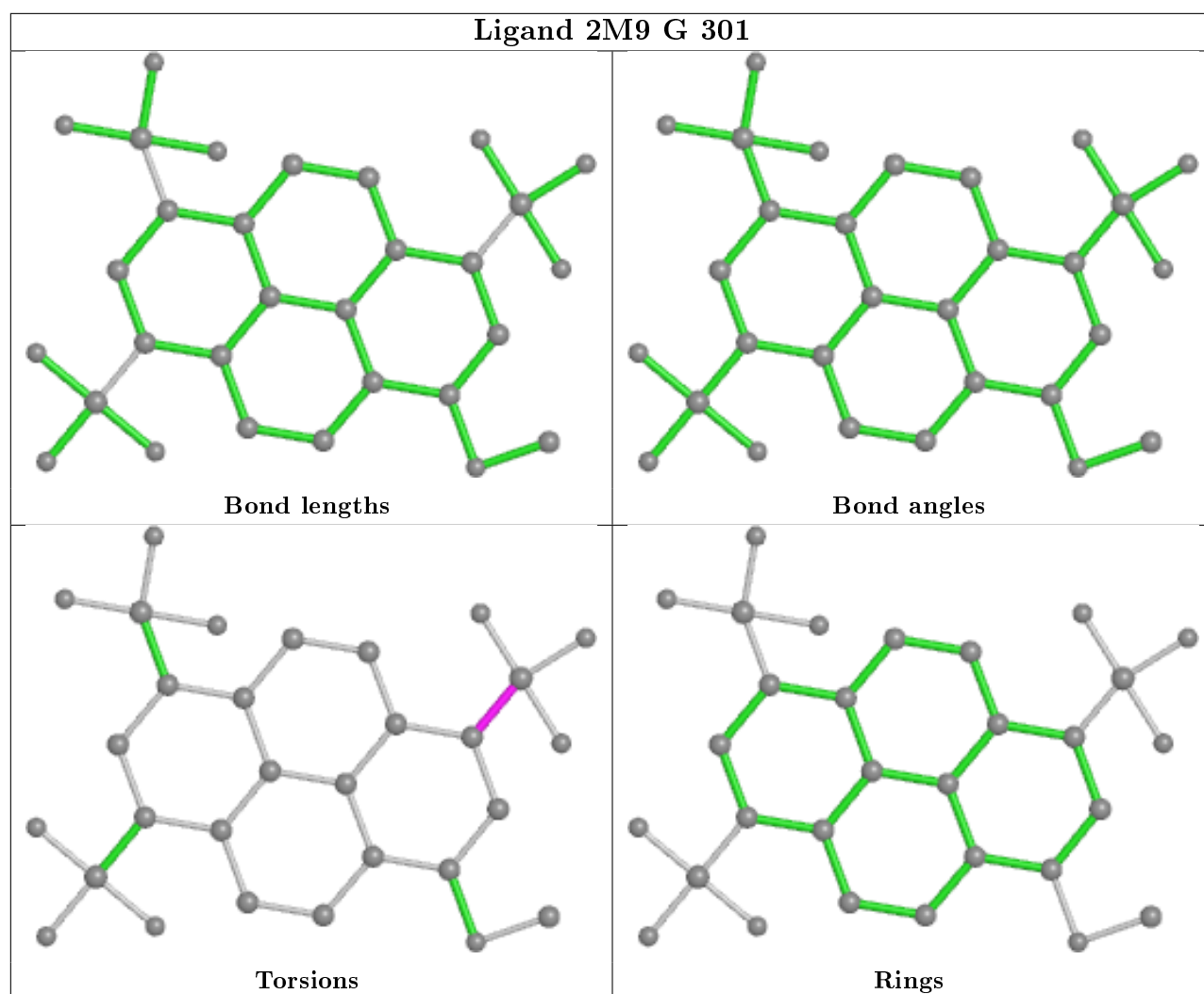
Mol	Chain	Res	Type	Atoms
3	G	301	2M9	CAY-CAU-SBE-OAE

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	J	301	2M9	1	0
4	D	302	SO4	1	0
4	L	301	SO4	2	0
4	I	301	SO4	1	0
4	B	302	SO4	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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 ⓘ

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	210/211 (99%)	0.28	1 (0%) 91 89	39, 66, 109, 138	0
1	C	208/211 (98%)	0.28	3 (1%) 75 71	41, 68, 108, 126	0
1	E	210/211 (99%)	0.42	9 (4%) 35 28	38, 67, 118, 176	0
1	G	208/211 (98%)	0.29	2 (0%) 82 80	39, 69, 96, 123	0
1	I	207/211 (98%)	0.28	8 (3%) 39 32	38, 70, 114, 133	0
1	K	206/211 (97%)	0.62	18 (8%) 10 7	40, 68, 135, 275	0
2	B	215/219 (98%)	0.29	3 (1%) 75 71	39, 63, 85, 257	0
2	D	215/219 (98%)	0.24	7 (3%) 46 39	46, 70, 105, 211	0
2	F	215/219 (98%)	0.25	2 (0%) 84 82	39, 67, 92, 166	0
2	H	215/219 (98%)	0.18	4 (1%) 66 62	38, 68, 90, 152	0
2	J	215/219 (98%)	0.33	6 (2%) 53 46	45, 73, 106, 140	0
2	L	215/219 (98%)	0.23	4 (1%) 66 62	43, 63, 100, 151	0
All	All	2539/2580 (98%)	0.31	67 (2%) 56 50	38, 67, 111, 275	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	K	148	TRP	20.5
2	B	172	ASP	8.0
1	K	130	ALA	7.9
1	K	182	THR	6.5
1	C	146	VAL	6.2
1	E	146	VAL	6.1
1	K	181	LEU	6.0
2	L	172	ASP	5.3
2	L	156	GLY	4.6
2	F	2	PHE	4.2
2	D	2	PHE	4.0

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Mol	Chain	Res	Type	RSRZ
2	H	67	ALA	3.7
1	K	155	ARG	3.5
1	E	179	LEU	3.4
2	B	156	GLY	3.4
1	G	194	CYS	3.3
1	I	149	LYS	3.2
1	I	194	CYS	3.2
1	K	205	ILE	3.2
2	J	160	SER	3.1
2	F	65	ASP	3.1
2	D	156	GLY	3.0
1	E	147	LYS	3.0
1	E	209	PHE	2.9
1	I	130	ALA	2.9
1	K	195	GLU	2.8
2	H	156	GLY	2.8
1	K	147	LYS	2.8
1	I	196	ALA	2.7
1	A	157	ASN	2.7
1	I	148	TRP	2.7
1	I	117	ILE	2.7
1	C	149	LYS	2.7
1	C	195	GLU	2.7
1	K	199	LYS	2.7
1	K	149	LYS	2.6
1	K	143	ASP	2.6
1	I	115	VAL	2.6
1	E	154	GLU	2.6
1	K	159	VAL	2.5
1	E	148	TRP	2.5
2	J	130	ALA	2.5
2	H	172	ASP	2.5
2	D	212	ASN	2.5
2	L	82	LEU	2.4
1	E	180	THR	2.4
1	E	193	THR	2.4
2	D	169	LEU	2.4
2	J	156	GLY	2.3
1	K	146	VAL	2.3
1	G	195	GLU	2.3
2	J	2	PHE	2.3
2	L	98	PHE	2.3

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Mol	Chain	Res	Type	RSRZ
1	K	207	LYS	2.2
2	B	5	LEU	2.2
2	D	132	THR	2.2
2	J	120	SER	2.2
2	D	67	ALA	2.2
1	K	198	HIS	2.1
2	D	65	ASP	2.1
2	J	53	ASN	2.1
1	K	194	CYS	2.1
1	K	131	SER	2.0
1	K	204	PRO	2.0
1	E	133	VAL	2.0
2	H	127	VAL	2.0
1	I	179	LEU	2.0

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

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(Å ²)	Q<0.9
2	PCA	J	1	8/9	0.81	0.38	101,106,125,126	0
2	PCA	L	1	8/9	0.82	0.34	83,92,121,130	0
1	PCA	A	1	8/9	0.82	0.51	155,177,213,227	0
1	PCA	G	1	8/9	0.83	0.36	72,102,132,152	0
2	PCA	D	1	8/9	0.85	0.25	71,82,84,85	0
1	PCA	K	1	8/9	0.88	0.20	55,60,77,84	0
1	PCA	I	1	8/9	0.91	0.38	77,107,137,155	0
1	PCA	C	1	8/9	0.96	0.18	48,54,70,79	0
1	PCA	E	1	8/9	0.96	0.26	39,46,62,92	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

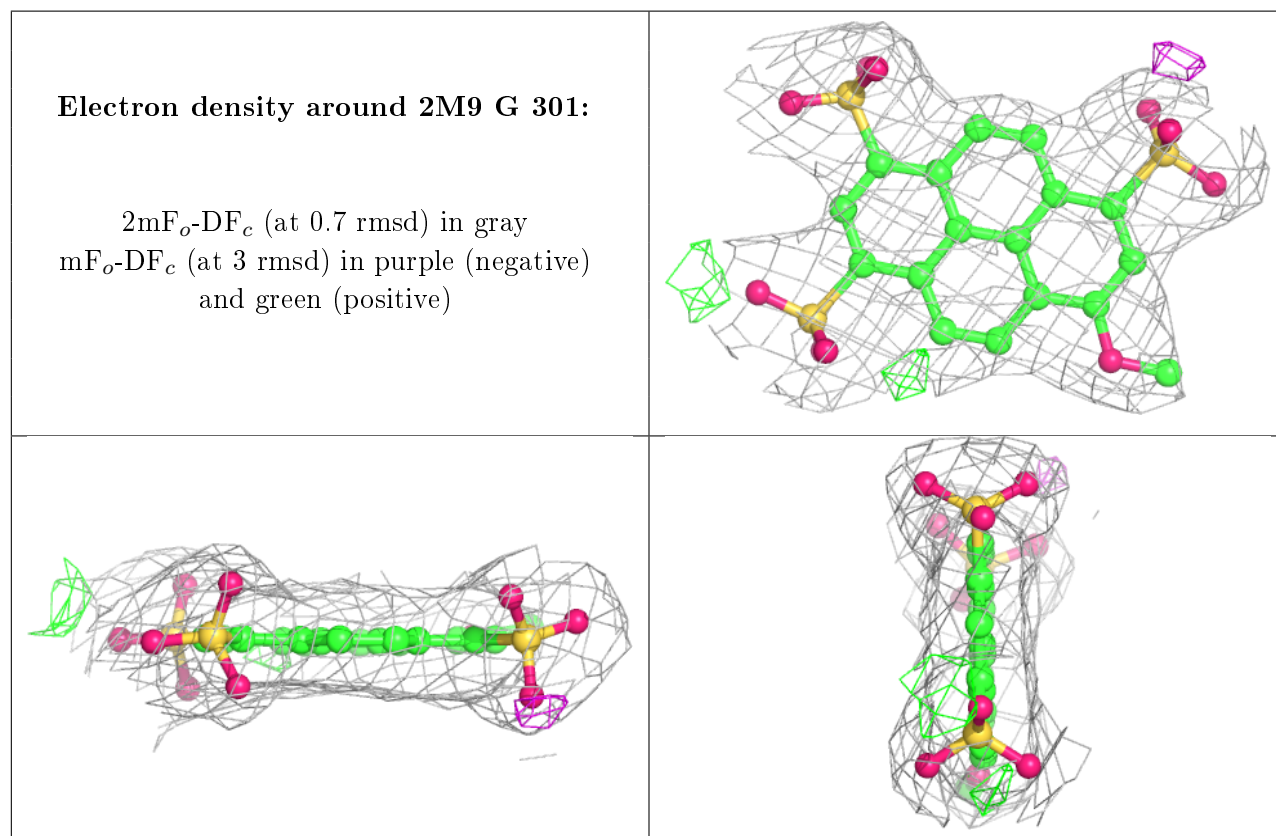
6.4 Ligands ⓘ

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(Å ²)	Q<0.9
4	SO4	D	302	5/5	0.84	0.13	133,136,136,137	0
4	SO4	I	301	5/5	0.92	0.14	76,80,97,99	0
4	SO4	F	301	5/5	0.94	0.15	75,75,80,87	0
4	SO4	L	301	5/5	0.94	0.13	72,78,93,93	0
4	SO4	B	302	5/5	0.95	0.17	63,71,75,77	0
3	2M9	J	301	30/30	0.97	0.17	46,62,73,84	0
3	2M9	K	301	30/30	0.97	0.17	42,54,66,70	0
3	2M9	D	301	30/30	0.97	0.18	54,66,81,96	0
3	2M9	E	301	30/30	0.98	0.17	41,51,66,67	0
4	SO4	G	302	5/5	0.98	0.12	66,77,78,80	0
3	2M9	G	301	30/30	0.98	0.16	45,61,73,80	0
3	2M9	B	301	30/30	0.98	0.19	40,48,61,74	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers

There are no such residues in this entry.