



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:00 PM UTC

PDB ID : 2IH3 / pdb_00002ih3
Title : Ion selectivity in a semi-synthetic K⁺ channel locked in the conductive conformation
Authors : Valiyaveetil, F.I.; Leonetti, M.; Muir, T.W.; MacKinnon, R.
Deposited on : 2006-09-25
Resolution : 1.72 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

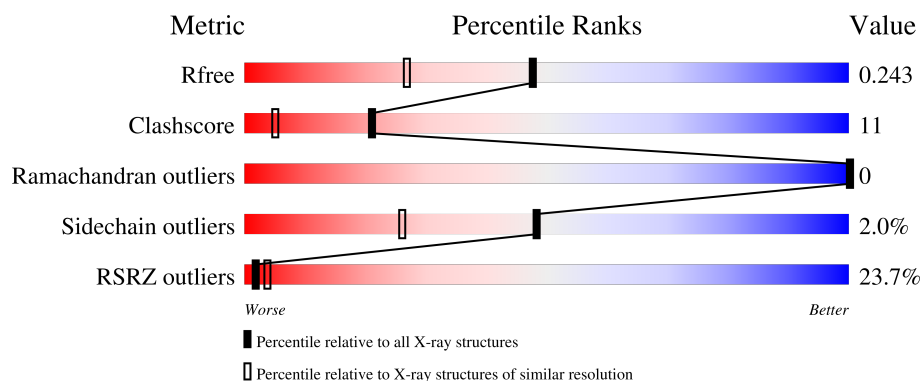
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.72 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	1039 (1.72-1.72)
Clashscore	190562	1049 (1.72-1.72)
Ramachandran outliers	187476	1041 (1.72-1.72)
Sidechain outliers	187428	1041 (1.72-1.72)
RSRZ outliers	180081	1039 (1.72-1.72)

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	219	26% 82% 15% .
2	B	212	24% 75% 24% .
3	C	122	13% 65% 13% . 20%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4338 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 Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	0	0	0
			1648	1042	275	325	6			

- Molecule 2 is a protein called FAB Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	212	Total	C	N	O	S	0	0	0
			1649	1023	283	338	5			

- Molecule 3 is a protein called Voltage-gated potassium channel.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	98	Total	C	N	O	S	0	0	0
			737	487	125	123	2			

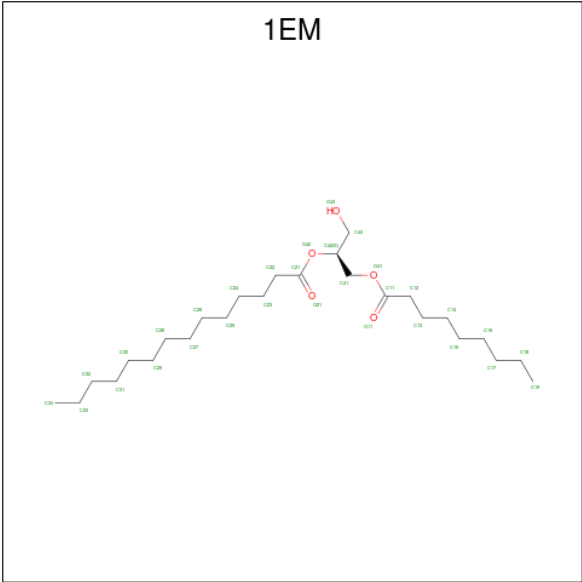
There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1	MET	-	cloning artifact	UNP P0A334
C	2	ALA	-	cloning artifact	UNP P0A334
C	69	ALA	SER	engineered mutation	UNP P0A334
C	70	CYS	VAL	engineered mutation	UNP P0A334
C	77	DAL	GLY	modified residue	UNP P0A334

- Molecule 4 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	7	Total	K	0	0
			7	7		

- Molecule 5 is (1S)-2-HYDROXY-1-[(NONANOYLOXY)METHYL]ETHYL MYRISTATE (CCD ID: 1EM) (formula: C₂₆H₅₀O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			31	26	5		

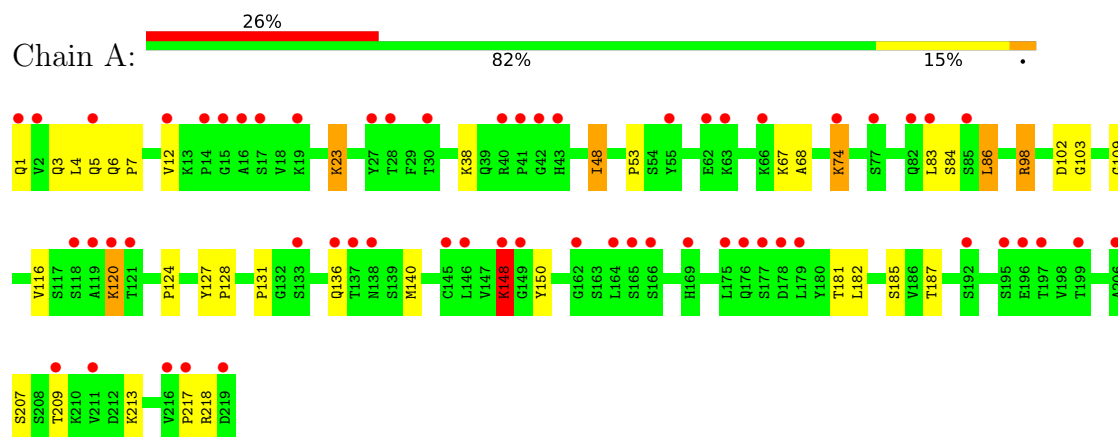
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	110	Total	O	0	0
			110	110		
6	B	120	Total	O	0	0
			120	120		
6	C	36	Total	O	0	0
			36	36		

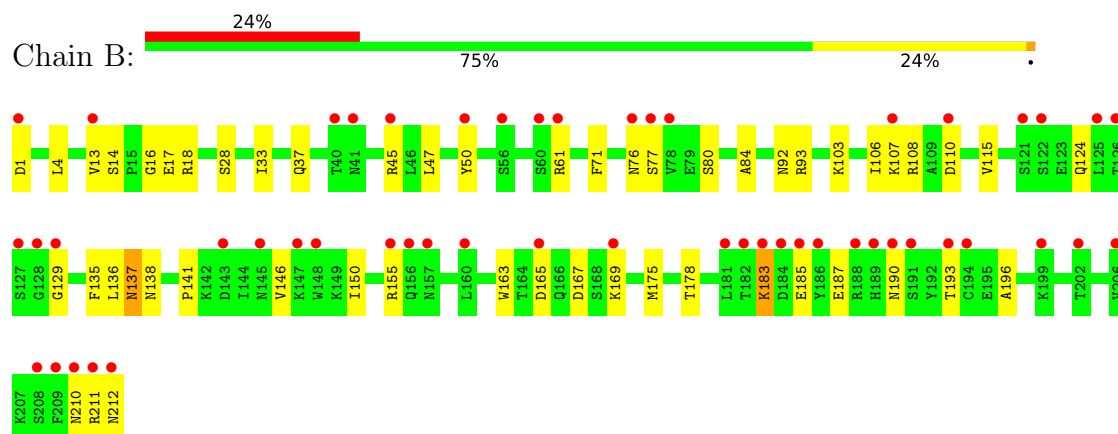
3 Residue-property plots [i](#)

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.

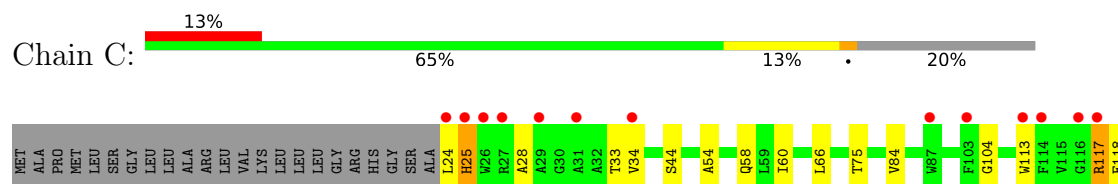
• Molecule 1: FAB Heavy Chain



• Molecule 2: FAB Light Chain



• Molecule 3: Voltage-gated potassium channel





4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	155.72Å 155.72Å 75.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.20 – 1.72 31.20 – 1.72	Depositor EDS
% Data completeness (in resolution range)	93.4 (31.20-1.72) 93.5 (31.20-1.72)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 1.72Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.233 , 0.242 0.229 , 0.243	Depositor DCC
R_{free} test set	9442 reflections (9.96%)	wwPDB-VP
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.302	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 47.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.022 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4338	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% 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: DAL, K, 1EM

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	1.06	0/1692	1.11	7/2312 (0.3%)
2	B	1.01	0/1686	1.09	2/2287 (0.1%)
3	C	1.39	3/750 (0.4%)	1.29	10/1031 (1.0%)
All	All	1.11	3/4128 (0.1%)	1.14	19/5630 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	54	ALA	CA-CB	7.17	1.63	1.53
3	C	75	THR	C-O	5.39	1.32	1.24
3	C	75	THR	CA-CB	5.06	1.61	1.53

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	137	ASN	N-CA-C	6.80	120.77	109.95
3	C	119	GLN	CB-CA-C	-6.51	100.40	110.81
2	B	80	SER	N-CA-C	6.48	119.17	111.33
1	A	68	ALA	N-CA-C	6.01	119.52	109.95
1	A	103	GLY	N-CA-C	5.92	122.90	115.21
3	C	119	GLN	N-CA-C	-5.72	104.24	111.11
3	C	120	GLU	N-CA-C	5.65	117.11	111.07
3	C	60	ILE	N-CA-C	5.56	118.53	113.20
3	C	25	HIS	N-CA-C	5.52	117.73	111.11
1	A	148	LYS	N-CA-C	5.40	117.89	109.52
1	A	84	SER	N-CA-C	5.36	117.83	109.52
3	C	84	VAL	N-CA-C	-5.35	108.07	113.47
1	A	48	ILE	CB-CA-C	-5.25	105.61	111.80
3	C	104	GLY	N-CA-C	-5.22	106.47	112.73
3	C	119	GLN	CA-CB-CG	5.15	124.40	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	86	LEU	N-CA-C	5.11	117.28	110.53
3	C	28	ALA	N-CA-C	-5.10	105.81	111.36
3	C	117	ARG	CA-CB-CG	-5.09	103.91	114.10
1	A	98	ARG	NE-CZ-NH2	-5.01	114.69	119.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1648	0	1616	32	0
2	B	1649	0	1576	50	0
3	C	737	0	756	11	0
4	C	7	0	0	0	0
5	C	31	0	50	0	0
6	A	110	0	0	4	0
6	B	120	0	0	14	0
6	C	36	0	0	6	0
All	All	4338	0	3998	89	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 11.

All (89) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:34:VAL:HA	6:C:335:HOH:O	1.41	1.18
2:B:107:LYS:C	6:B:274:HOH:O	2.18	0.85
1:A:53:PRO:HB3	1:A:74:LYS:HE2	1.66	0.79
3:C:34:VAL:CA	6:C:335:HOH:O	2.10	0.78
1:A:23:LYS:HE3	1:A:23:LYS:H	1.51	0.75
2:B:108:ARG:N	6:B:274:HOH:O	2.20	0.75
3:C:33:THR:C	6:C:335:HOH:O	2.32	0.72
2:B:167:ASP:OD2	2:B:169:LYS:HB2	1.89	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:34:VAL:N	6:C:335:HOH:O	2.23	0.71
2:B:108:ARG:CA	6:B:274:HOH:O	2.39	0.70
2:B:76:ASN:OD1	6:B:262:HOH:O	2.08	0.70
2:B:178:THR:O	6:B:297:HOH:O	2.10	0.69
1:A:140:MET:HE3	1:A:187:THR:HG22	1.75	0.68
2:B:61:ARG:HG3	2:B:76:ASN:O	1.93	0.68
1:A:23:LYS:HE3	1:A:23:LYS:N	2.08	0.68
3:C:113:TRP:HE1	3:C:117:ARG:NH1	1.92	0.67
2:B:45:ARG:HD2	6:B:330:HOH:O	1.97	0.63
2:B:106:ILE:HG22	6:B:274:HOH:O	1.98	0.62
2:B:137:ASN:HB3	2:B:138:ASN:HD22	1.64	0.61
2:B:183:LYS:O	2:B:187:GLU:HG2	1.99	0.61
2:B:13:VAL:CG2	2:B:17:GLU:HB3	2.30	0.61
3:C:24:LEU:N	6:C:317:HOH:O	2.34	0.60
2:B:92:ASN:HD21	3:C:58:GLN:HE21	1.50	0.60
1:A:67:LYS:NZ	6:A:299:HOH:O	2.10	0.59
2:B:193:THR:HG23	6:B:248:HOH:O	2.02	0.59
1:A:120:LYS:O	1:A:120:LYS:HD2	2.04	0.57
2:B:108:ARG:HA	6:B:274:HOH:O	2.03	0.56
3:C:25:HIS:ND1	6:C:330:HOH:O	2.33	0.55
2:B:92:ASN:HD22	2:B:93:ARG:HG3	1.71	0.55
1:A:53:PRO:CB	1:A:74:LYS:HE2	2.35	0.55
2:B:190:ASN:HD21	2:B:210:ASN:HD22	1.55	0.55
2:B:28:SER:OG	6:B:269:HOH:O	2.18	0.54
3:C:118:GLU:OE2	3:C:121:ARG:HD3	2.08	0.54
2:B:190:ASN:ND2	2:B:210:ASN:HD22	2.06	0.53
2:B:150:ILE:HD12	2:B:155:ARG:HH11	1.74	0.52
2:B:146:VAL:HB	2:B:175:MET:HE1	1.93	0.51
1:A:128:PRO:HG3	1:A:213:LYS:HD2	1.92	0.51
2:B:183:LYS:HB3	2:B:183:LYS:NZ	2.25	0.50
2:B:115:VAL:HG22	2:B:136:LEU:CD1	2.41	0.50
2:B:211:ARG:HH11	2:B:211:ARG:HG2	1.76	0.50
2:B:37:GLN:HB2	2:B:47:LEU:HD11	1.93	0.50
3:C:118:GLU:O	3:C:119:GLN:C	2.55	0.50
1:A:148:LYS:HB3	1:A:181:THR:HG23	1.94	0.50
1:A:83:LEU:HB3	1:A:86:LEU:HD21	1.94	0.50
3:C:44:SER:OG	3:C:66:LEU:HA	2.12	0.50
1:A:1:GLN:OE1	1:A:1:GLN:N	2.39	0.49
2:B:1:ASP:HA	6:B:237:HOH:O	2.12	0.49
1:A:5:GLN:HG3	6:A:287:HOH:O	2.13	0.49
1:A:6:GLN:HE21	1:A:109:GLY:HA3	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:110:ASP:OD1	2:B:141:PRO:HD3	2.12	0.48
1:A:128:PRO:HG3	1:A:213:LYS:CG	2.42	0.48
2:B:13:VAL:HG22	2:B:14:SER:N	2.28	0.48
2:B:137:ASN:HB3	2:B:138:ASN:ND2	2.29	0.48
2:B:33:ILE:HG21	2:B:71:PHE:CD2	2.49	0.47
1:A:182:LEU:C	1:A:182:LEU:HD12	2.38	0.47
2:B:136:LEU:CD1	2:B:196:ALA:HB2	2.45	0.47
2:B:185:GLU:HG2	6:B:292:HOH:O	2.15	0.47
1:A:207:SER:O	1:A:209:THR:HG23	2.15	0.46
1:A:185:SER:HB2	2:B:135:PHE:CE2	2.51	0.46
2:B:212:ASN:ND2	2:B:212:ASN:OXT	2.49	0.46
2:B:18:ARG:NH1	2:B:76:ASN:OD1	2.49	0.45
2:B:167:ASP:HB2	6:B:320:HOH:O	2.17	0.45
1:A:120:LYS:HD2	1:A:120:LYS:C	2.41	0.44
1:A:148:LYS:HB2	1:A:148:LYS:NZ	2.33	0.44
2:B:13:VAL:HG22	2:B:17:GLU:HB3	1.98	0.44
1:A:3:GLN:HE21	1:A:98:ARG:NH1	2.16	0.44
1:A:102:ASP:O	2:B:50:TYR:OH	2.36	0.44
2:B:136:LEU:HD11	2:B:196:ALA:HB2	1.98	0.44
1:A:3:GLN:HG3	6:A:317:HOH:O	2.18	0.44
1:A:102:ASP:HB2	2:B:50:TYR:OH	2.18	0.43
1:A:12:VAL:O	1:A:116:VAL:HA	2.19	0.43
1:A:127:TYR:HA	1:A:128:PRO:HD3	1.87	0.43
2:B:124:GLN:HG2	2:B:129:GLY:O	2.19	0.43
2:B:18:ARG:NH1	2:B:76:ASN:ND2	2.68	0.42
2:B:16:GLY:C	2:B:77:SER:HA	2.44	0.42
2:B:16:GLY:O	2:B:77:SER:HA	2.19	0.42
2:B:136:LEU:HD23	2:B:175:MET:HE2	2.00	0.42
2:B:103:LYS:HE2	6:B:268:HOH:O	2.18	0.42
1:A:217:PRO:HA	6:A:263:HOH:O	2.19	0.42
1:A:131:PRO:O	1:A:218:ARG:HD3	2.19	0.42
1:A:7:PRO:HD3	1:A:23:LYS:HZ1	1.85	0.41
2:B:136:LEU:HD22	2:B:136:LEU:N	2.35	0.41
2:B:163:TRP:CD1	2:B:175:MET:HG3	2.56	0.41
1:A:38:LYS:HB2	1:A:48:ILE:HD11	2.01	0.41
1:A:124:PRO:HB3	1:A:150:TYR:HB3	2.03	0.41
1:A:3:GLN:HE21	1:A:98:ARG:HH11	1.68	0.41
2:B:84:ALA:O	2:B:103:LYS:HD2	2.21	0.41
1:A:128:PRO:HG3	1:A:213:LYS:CD	2.51	0.40
2:B:167:ASP:OD2	2:B:169:LYS:CB	2.66	0.40

There are no symmetry-related clashes.

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	217/219 (99%)	210 (97%)	7 (3%)	0	100	100
2	B	210/212 (99%)	203 (97%)	7 (3%)	0	100	100
3	C	95/122 (78%)	95 (100%)	0	0	100	100
All	All	522/553 (94%)	508 (97%)	14 (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	185/185 (100%)	179 (97%)	6 (3%)	34	11
2	B	190/190 (100%)	187 (98%)	3 (2%)	55	34
3	C	70/89 (79%)	70 (100%)	0	100	100
All	All	445/464 (96%)	436 (98%)	9 (2%)	48	26

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	23	LYS
1	A	74	LYS
1	A	120	LYS
1	A	136	GLN

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Mol	Chain	Res	Type
1	A	148	LYS
2	B	4	LEU
2	B	165	ASP
2	B	183	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	5	GLN
1	A	6	GLN
2	B	37	GLN
2	B	41	ASN
2	B	85	ASN
2	B	92	ASN
2	B	137	ASN
2	B	138	ASN
2	B	190	ASN
2	B	210	ASN
2	B	212	ASN
3	C	58	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 7 are monoatomic - leaving 1 for Mogul analysis.

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	$\# Z > 2$
5	1EM	C	301	-	30,30,30	1.20	3 (10%)	32,32,32	1.13	2 (6%)

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
5	1EM	C	301	-	-	10/32/32/32	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	301	1EM	O41-C11	3.83	1.44	1.33
5	C	301	1EM	O42-C21	3.56	1.44	1.34
5	C	301	1EM	O41-C41	-2.41	1.39	1.45

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	301	1EM	O42-C21-C22	3.70	119.49	111.48
5	C	301	1EM	O41-C11-C12	2.01	117.96	111.83

There are no chirality outliers.

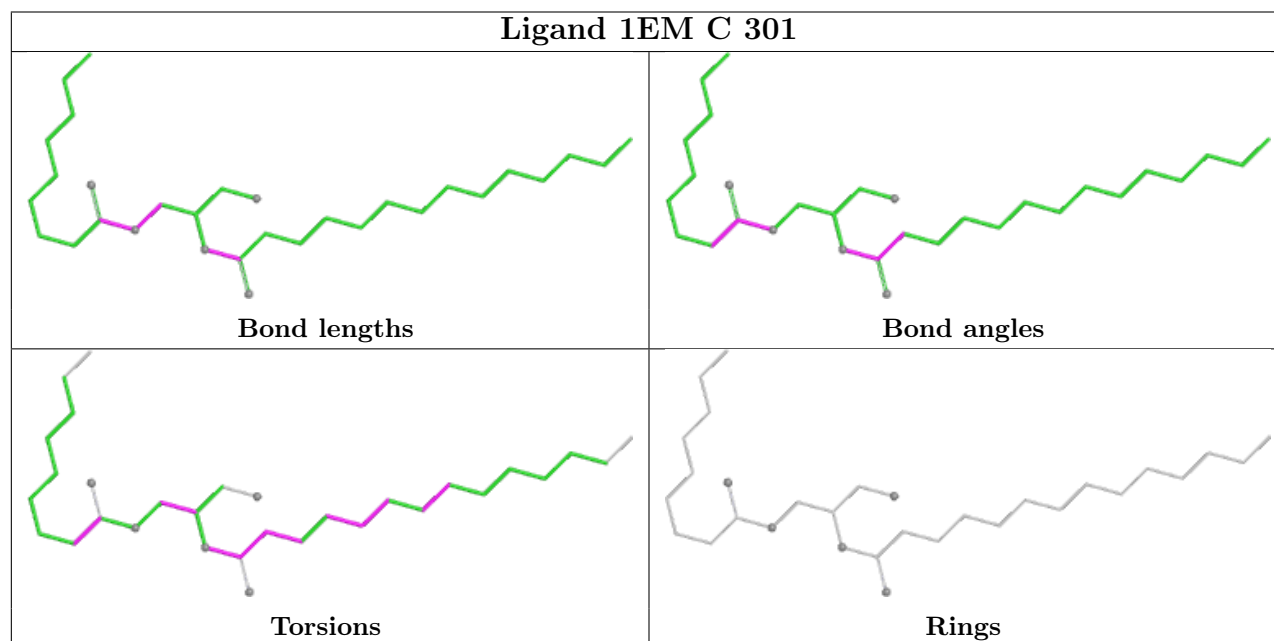
All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	301	1EM	O21-C21-O42-C42
5	C	301	1EM	C22-C21-O42-C42
5	C	301	1EM	C23-C24-C25-C26
5	C	301	1EM	C21-C22-C23-C24
5	C	301	1EM	C24-C25-C26-C27
5	C	301	1EM	C26-C27-C28-C29
5	C	301	1EM	O41-C11-C12-C13
5	C	301	1EM	O41-C41-C42-O42
5	C	301	1EM	O11-C11-C12-C13
5	C	301	1EM	O42-C21-C22-C23

There are no ring outliers.

No monomer is involved in short contacts.

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

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	219/219 (100%)	1.38	58 (26%) 1 3	18, 36, 51, 64	0
2	B	212/212 (100%)	1.16	51 (24%) 2 3	16, 34, 55, 67	0
3	C	97/122 (79%)	0.51	16 (16%) 4 6	14, 21, 57, 64	0
All	All	528/553 (95%)	1.13	125 (23%) 2 4	14, 34, 55, 67	0

All (125) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	125	LEU	5.2
1	A	74	LYS	5.1
3	C	24	LEU	5.0
2	B	212	ASN	4.9
1	A	178	ASP	4.7
1	A	219	ASP	4.6
3	C	120	GLU	4.6
3	C	119	GLN	4.5
1	A	138	ASN	4.3
3	C	27	ARG	4.3
1	A	133	SER	4.2
1	A	137	THR	4.1
1	A	120	LYS	4.0
3	C	26	TRP	4.0
3	C	117	ARG	3.8
1	A	145	CYS	3.8
2	B	127	SER	3.8
2	B	190	ASN	3.8
3	C	121	ARG	3.7
2	B	50	TYR	3.6
1	A	136	GLN	3.6
1	A	177	SER	3.6
2	B	76	ASN	3.5

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Mol	Chain	Res	Type	RSRZ
2	B	194	CYS	3.5
3	C	113	TRP	3.4
2	B	183	LYS	3.4
1	A	63	LYS	3.4
2	B	157	ASN	3.4
2	B	210	ASN	3.3
2	B	209	PHE	3.3
2	B	1	ASP	3.2
2	B	193	THR	3.1
1	A	17	SER	3.1
2	B	78	VAL	3.0
2	B	13	VAL	3.0
1	A	166	SER	3.0
2	B	165	ASP	3.0
1	A	43	HIS	3.0
2	B	147	LYS	3.0
1	A	195	SER	2.9
1	A	211	VAL	2.9
2	B	186	TYR	2.9
1	A	40	ARG	2.9
2	B	206	VAL	2.9
1	A	196	GLU	2.9
2	B	107	LYS	2.9
1	A	121	THR	2.8
3	C	31	ALA	2.8
2	B	56	SER	2.8
2	B	181	LEU	2.8
1	A	118	SER	2.8
2	B	143	ASP	2.8
1	A	15	GLY	2.8
3	C	103	PHE	2.7
1	A	148	LYS	2.7
1	A	165	SER	2.7
1	A	42	GLY	2.7
1	A	28	THR	2.6
2	B	202	THR	2.6
3	C	87	TRP	2.6
1	A	197	THR	2.6
1	A	169	HIS	2.6
1	A	41	PRO	2.6
1	A	55	TYR	2.6
2	B	60	SER	2.6

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Mol	Chain	Res	Type	RSRZ
2	B	208	SER	2.6
1	A	179	LEU	2.5
1	A	5	GLN	2.5
2	B	121	SER	2.5
2	B	160	LEU	2.5
2	B	128	GLY	2.5
2	B	40	THR	2.5
1	A	19	LYS	2.5
2	B	41	ASN	2.5
1	A	176	GLN	2.4
2	B	184	ASP	2.4
2	B	148	TRP	2.4
2	B	77	SER	2.4
1	A	30	THR	2.4
1	A	199	THR	2.4
2	B	126	THR	2.4
2	B	122	SER	2.4
2	B	45	ARG	2.4
2	B	169	LYS	2.4
1	A	1	GLN	2.3
1	A	162	GLY	2.3
1	A	164	LEU	2.3
1	A	146	LEU	2.3
1	A	206	ALA	2.3
2	B	156	GLN	2.3
3	C	116	GLY	2.3
3	C	34	VAL	2.3
3	C	25	HIS	2.2
1	A	16	ALA	2.2
1	A	119	ALA	2.2
1	A	175	LEU	2.2
2	B	211	ARG	2.2
2	B	191	SER	2.2
1	A	149	GLY	2.2
1	A	62	GLU	2.2
2	B	185	GLU	2.2
1	A	12	VAL	2.2
1	A	77	SER	2.2
2	B	155	ARG	2.2
2	B	188	ARG	2.2
1	A	14	PRO	2.2
1	A	217	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	82	GLN	2.1
3	C	29	ALA	2.1
2	B	189	HIS	2.1
1	A	66	LYS	2.1
2	B	145	ASN	2.1
1	A	2	VAL	2.1
1	A	83	LEU	2.1
3	C	114	PHE	2.1
2	B	129	GLY	2.1
2	B	199	LYS	2.1
2	B	110	ASP	2.1
2	B	61	ARG	2.1
1	A	27	TYR	2.1
1	A	85	SER	2.0
1	A	192	SER	2.0
1	A	209	THR	2.0
2	B	182	THR	2.0
1	A	216	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
3	DAL	C	77	5/6	0.98	0.05	12,14,14,19	0

6.3 Carbohydrates [i](#)

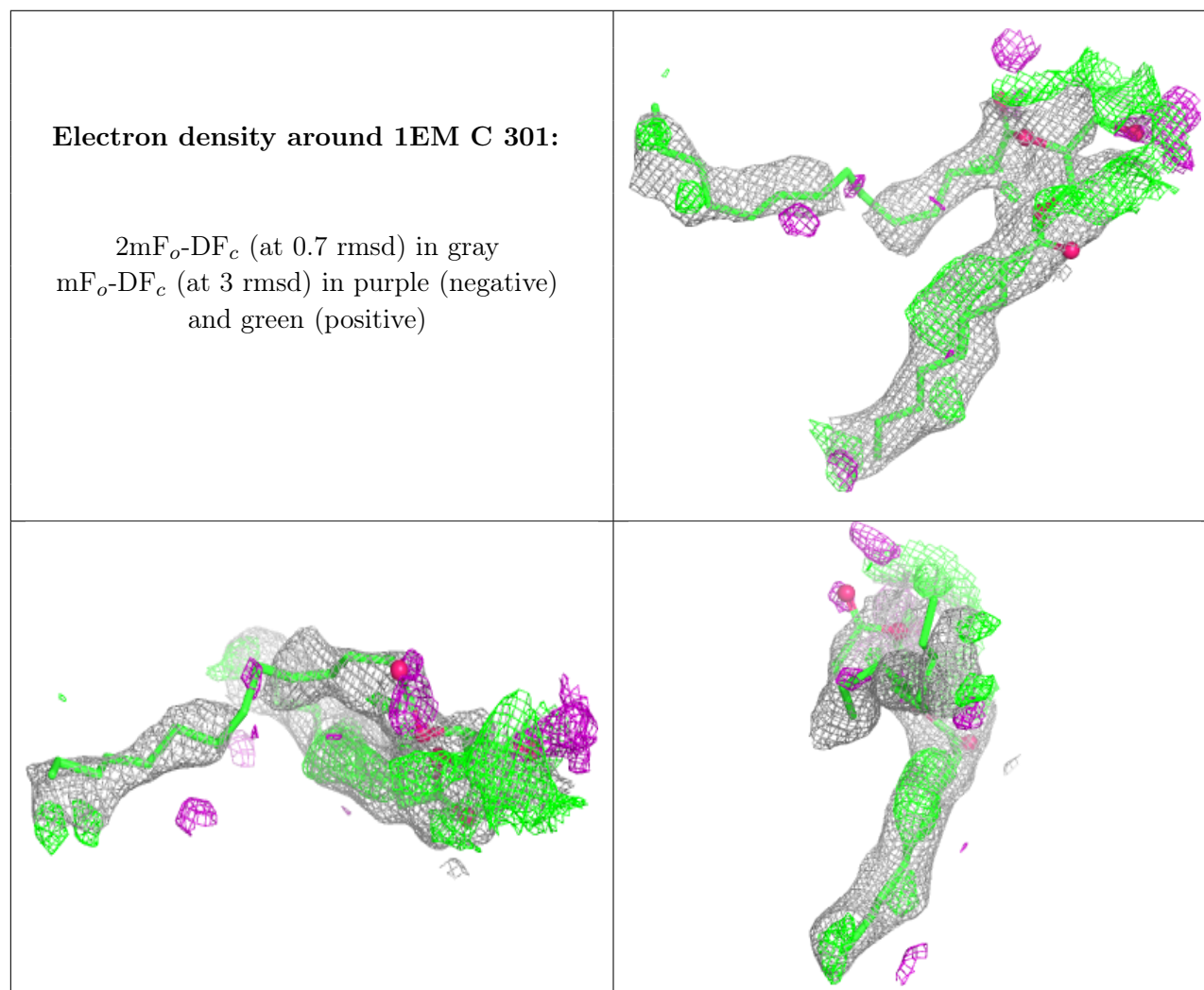
There are no oligosaccharides in this entry.

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	K	C	207	1/1	0.64	0.31	61,61,61,61	1
5	1EM	C	301	31/31	0.66	0.33	64,71,76,79	0
4	K	C	206	1/1	0.82	0.28	74,74,74,74	1
4	K	C	205	1/1	0.95	0.53	61,61,61,61	1
4	K	C	201	1/1	1.00	0.02	15,15,15,15	1
4	K	C	202	1/1	1.00	0.02	16,16,16,16	1
4	K	C	203	1/1	1.00	0.03	17,17,17,17	1
4	K	C	204	1/1	1.00	0.01	17,17,17,17	1

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 [i](#)

There are no such residues in this entry.