



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:54 AM UTC

PDB ID : 5CCB / pdb_00005ccb
Title : Crystal structure of human m1A58 methyltransferase in a complex with tRNA³Lys and SAH
Authors : Finer-Moore, J.; Czudnochowski, N.; O'Connell III, J.D.; Wang, A.L.; Stroud, R.M.
Deposited on : 2015-07-01
Resolution : 2.00 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)
Xtrriage (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

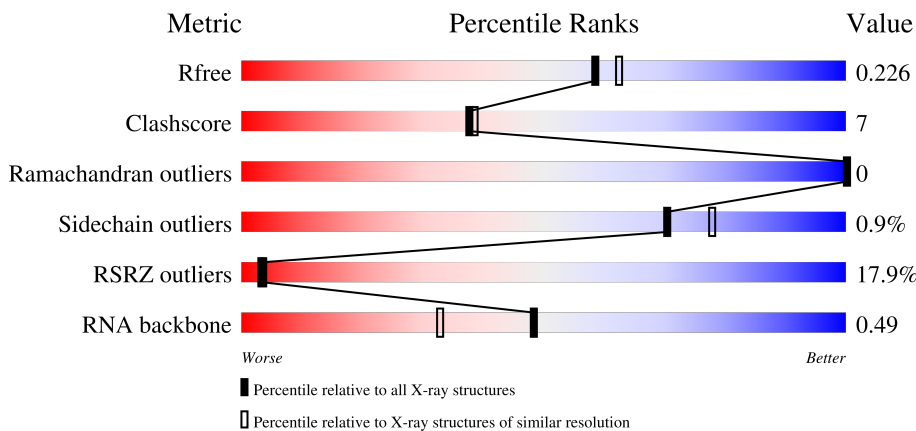
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)
RNA backbone	3983	1000 (2.36-1.64)

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	289	 9% 90% 8%
2	B	497	 19% 64% 10% 25%
3	N	77	 16% 57% 34% 9%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 13154 atoms, of which 5959 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 tRNA (adenine(58)-N(1))-methyltransferase catalytic subunit TRMT61A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	282	4305	1362	2144	390	398	11	0	0	0

- Molecule 2 is a protein called tRNA (adenine(58)-N(1))-methyltransferase non-catalytic subunit TRM6.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	B	371	5882	1864	2955	517	532	14	0	0	0

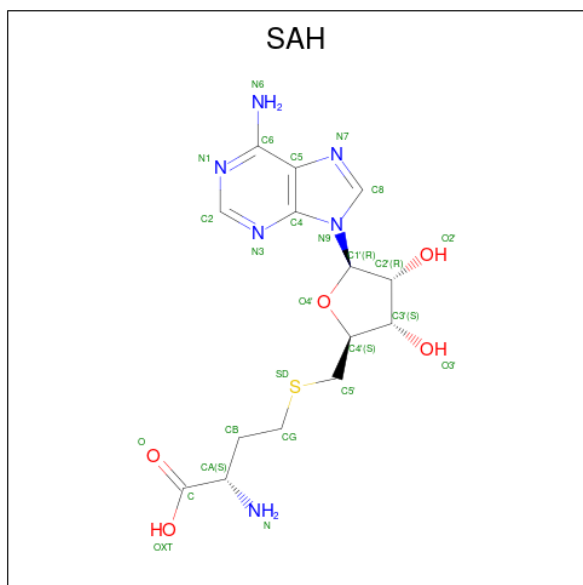
- Molecule 3 is a RNA chain called tRNA3Lys.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	P			
3	N	77	2505	742	841	296	548	78	0	1	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
N	74	C	-	insertion	GB 339572
N	75	C	-	insertion	GB 339572
N	76	A	-	insertion	GB 339572

- Molecule 4 is S-ADENOSYL-L-HOMOCYSTEINE (CCD ID: SAH) (formula: C₁₄H₂₀N₆O₅S).



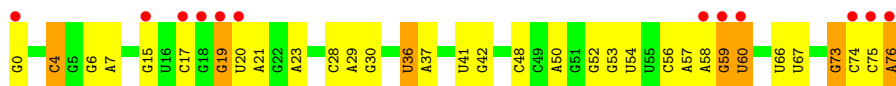
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	H	N	O			S
4	A	1	45	14	19	6	5	1	0	0

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Na		
5	N	1	1	1	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	164	164	164	0	0
6	B	144	144	144	0	0
6	N	108	108	108	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	137.18Å 137.18Å 177.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	74.37 – 2.00 74.37 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (74.37-2.00) 90.8 (74.37-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.12 (at 1.84Å)	Xtrriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, R_{free}	0.199 , 0.222 0.203 , 0.226	Depositor DCC
R_{free} test set	7193 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	33.0	Xtrriage
Anisotropy	0.201	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13154	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.92% 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: SAH, NA

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	0.39	0/2214	0.71	0/3012
2	B	0.35	0/2981	0.71	0/4009
3	N	0.22	0/1859	0.41	0/2896
All	All	0.33	0/7054	0.64	0/9917

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	2161	2144	2131	17	0
2	B	2927	2955	2949	41	0
3	N	1664	841	841	32	0
4	A	26	19	19	0	0
5	N	1	0	0	0	0
6	A	164	0	0	0	1
6	B	144	0	0	2	0
6	N	108	0	0	5	0
All	All	7195	5959	5940	89	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 7.

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)
1:A:1:MET:HB3	1:A:2:SER:HA	1.53	0.88
3:N:7:A:O2'	6:N:201:HOH:O	2.09	0.70
2:B:384:LEU:HD12	2:B:410:CYS:SG	2.37	0.65
3:N:17:C:OP2	3:N:19:G:N2	2.30	0.64
2:B:99:ASP:OD1	3:N:53:G:O2'	2.15	0.62
3:N:21:A:N3	3:N:21:A:O4'	2.33	0.61
1:A:43:LEU:HD11	1:A:60:CYS:SG	2.40	0.60
2:B:57:VAL:HG23	2:B:66:PHE:CZ	2.36	0.59
3:N:19:G:H2'	3:N:19:G:N3	2.16	0.59
3:N:76:A:H3'	3:N:76:A:N3	2.18	0.59
2:B:354:ARG:O	2:B:358:LEU:HD13	2.04	0.58
2:B:35:LYS:NZ	2:B:46:THR:O	2.32	0.58
2:B:176:GLU:HB3	2:B:179:LYS:HG3	1.86	0.56
3:N:17:C:C5	3:N:19:G:C6	2.94	0.56
2:B:157:ILE:C	2:B:158:ILE:HD12	2.31	0.55
1:A:1:MET:CB	1:A:2:SER:HA	2.25	0.54
2:B:46:THR:HG23	2:B:51:TRP:CZ3	2.43	0.54
2:B:132:SER:O	6:B:501:HOH:O	2.18	0.54
1:A:108:CYS:HB3	1:A:169:PHE:CD1	2.44	0.53
3:N:56:C:C5	3:N:58[A]:A:N1	2.78	0.52
1:A:243:VAL:HG23	1:A:265:PHE:CZ	2.44	0.52
2:B:65:ALA:O	2:B:66:PHE:CD2	2.63	0.52
3:N:48:C:N3	6:N:203:HOH:O	2.34	0.52
3:N:36:U:H2'	3:N:36:U:O2	2.09	0.52
3:N:36:U:O2	3:N:36:U:C2'	2.59	0.51
2:B:46:THR:HG23	2:B:51:TRP:CH2	2.46	0.51
3:N:4:C:H5''	3:N:4:C:H6	1.75	0.51
1:A:37:GLN:OE1	1:A:42:VAL:HG22	2.11	0.51
3:N:20:U:O2	3:N:20:U:O4'	2.30	0.49
3:N:59:G:C1'	3:N:60:U:OP1	2.61	0.49
3:N:0:G:C5	3:N:74:C:C6	3.00	0.49
3:N:17:C:H3'	3:N:17:C:OP1	2.12	0.49
3:N:28:C:O2'	3:N:29:A:H5'	2.12	0.49
2:B:385:LEU:O	6:B:502:HOH:O	2.20	0.48
3:N:0:G:O6	3:N:75:C:N4	2.46	0.48
3:N:75:C:H3'	3:N:75:C:O2	2.13	0.48
2:B:66:PHE:HA	2:B:76:PRO:HD3	1.95	0.48
3:N:59:G:H1'	3:N:60:U:OP1	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:345:GLU:OE2	2:B:348:ARG:NH2	2.46	0.48
2:B:64:THR:HG22	2:B:66:PHE:CE2	2.49	0.47
2:B:197:ILE:CD1	2:B:203:MET:HE2	2.45	0.47
2:B:74:LEU:O	2:B:76:PRO:HD3	2.15	0.47
3:N:41:U:H2'	3:N:42:G:O4'	2.15	0.46
2:B:57:VAL:O	2:B:57:VAL:HG13	2.15	0.46
1:A:243:VAL:CG2	1:A:265:PHE:CE2	2.99	0.46
2:B:54:LEU:O	2:B:57:VAL:HG12	2.16	0.46
2:B:31:GLU:HG2	2:B:152:LYS:HE2	1.98	0.46
1:A:83:ARG:CD	1:A:143:LYS:HG2	2.46	0.45
1:A:1:MET:HB2	1:A:14:ASP:OD1	2.16	0.45
2:B:107:THR:HG22	2:B:108:GLN:N	2.32	0.45
3:N:66:U:C4	3:N:67:U:C4	3.04	0.45
1:A:83:ARG:HD3	1:A:143:LYS:HG2	1.98	0.45
2:B:68:VAL:HG12	2:B:69:THR:N	2.32	0.45
3:N:73:G:O2'	3:N:75:C:C5	2.71	0.44
2:B:158:ILE:HD12	2:B:158:ILE:N	2.31	0.44
1:A:37:GLN:CD	1:A:42:VAL:HG22	2.43	0.44
2:B:58:ILE:HG22	2:B:59:GLY:N	2.32	0.44
1:A:1:MET:HB3	1:A:2:SER:CA	2.36	0.44
3:N:7:A:H5'	6:N:232:HOH:O	2.18	0.44
3:N:75:C:O2	3:N:75:C:C2'	2.65	0.44
2:B:111:ILE:CG2	2:B:112:LYS:N	2.80	0.44
2:B:419:GLY:HA3	2:B:461:MET:SD	2.58	0.44
1:A:48:ASP:O	1:A:52:ARG:HD3	2.18	0.43
2:B:207:GLU:HG3	2:B:209:CYS:H	1.83	0.43
2:B:110:ASP:O	2:B:113:ALA:HB3	2.19	0.43
2:B:384:LEU:CD1	2:B:414:LEU:HD11	2.48	0.43
1:A:252:THR:HG22	1:A:253:GLY:N	2.32	0.42
2:B:114:LEU:HD22	2:B:114:LEU:N	2.34	0.42
2:B:184:ARG:HD2	2:B:186:ASP:OD1	2.20	0.42
2:B:108:GLN:O	2:B:111:ILE:HG22	2.19	0.42
1:A:1:MET:HE2	1:A:14:ASP:HA	2.01	0.42
1:A:17:ILE:HD13	1:A:71:PRO:HA	2.02	0.42
2:B:199:ALA:HB1	2:B:224:PHE:HB2	2.02	0.42
3:N:36:U:OP1	6:N:202:HOH:O	2.22	0.42
3:N:59:G:O2'	3:N:60:U:H5''	2.19	0.41
2:B:65:ALA:C	2:B:66:PHE:CD2	2.98	0.41
3:N:17:C:C4	3:N:19:G:O6	2.74	0.41
1:A:12:GLU:CD	1:A:32:ARG:HG3	2.45	0.41
2:B:110:ASP:O	2:B:114:LEU:HD23	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:172:TYR:HB3	2:B:180:ILE:HD13	2.03	0.41
3:N:52:G:O2'	3:N:53:G:H5'	2.21	0.41
3:N:59:G:H1'	3:N:60:U:P	2.60	0.41
2:B:405:GLU:N	2:B:406:PRO:CD	2.83	0.41
3:N:0:G:N7	3:N:74:C:C5	2.89	0.41
2:B:368:ASN:HA	2:B:392:ALA:HB2	2.03	0.41
2:B:405:GLU:HB2	2:B:406:PRO:HD3	2.03	0.41
2:B:384:LEU:HD23	2:B:384:LEU:HA	1.93	0.40
3:N:58[A]:A:H3'	6:N:229:HOH:O	2.20	0.40
2:B:57:VAL:HG23	2:B:66:PHE:CE2	2.57	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)
6:A:409:HOH:O	6:A:409:HOH:O[7_555]	2.00	0.20

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	278/289 (96%)	272 (98%)	6 (2%)	0	100	100
2	B	365/497 (73%)	352 (96%)	13 (4%)	0	100	100
All	All	643/786 (82%)	624 (97%)	19 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

resolution.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	233/238 (98%)	231 (99%)	2 (1%)	70	78
2	B	309/428 (72%)	306 (99%)	3 (1%)	68	75
All	All	542/666 (81%)	537 (99%)	5 (1%)	70	78

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

Mol	Chain	Res	Type
1	A	4	VAL
1	A	136	PHE
2	B	31	GLU
2	B	207	GLU
2	B	384	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	56	ASN
2	B	60	HIS

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	N	75/77 (97%)	14 (18%)	1 (1%)

All (14) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	N	4	C
3	N	6	G
3	N	15	G
3	N	19	G
3	N	23	A
3	N	30	G
3	N	36	U
3	N	37	A

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Mol	Chain	Res	Type
3	N	50	A
3	N	54	U
3	N	57	A
3	N	60	U
3	N	73	G
3	N	76	A

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
3	N	59	G

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](#)

Of 2 ligands modelled in this entry, 1 is 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
4	SAH	A	301	-	27,28,28	1.03	3 (11%)	36,40,40	1.81	9 (25%)

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
4	SAH	A	301	-	-	0/15/31/31	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	301	SAH	C2-N1	2.41	1.38	1.33
4	A	301	SAH	C2-N3	2.19	1.37	1.33
4	A	301	SAH	C8-N7	2.08	1.35	1.31

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	301	SAH	N3-C2-N1	-5.54	120.20	128.58
4	A	301	SAH	N9-C8-N7	-3.94	108.34	113.94
4	A	301	SAH	C5-C4-N3	-3.81	121.47	126.72
4	A	301	SAH	C2-N3-C4	3.05	119.28	111.83
4	A	301	SAH	C5-N7-C8	3.05	108.24	103.45
4	A	301	SAH	N3-C4-N9	2.41	131.26	127.17
4	A	301	SAH	C4-N9-C8	2.15	108.00	105.74
4	A	301	SAH	C4-C5-N7	-2.13	108.14	110.58
4	A	301	SAH	C5'-SD-CG	-2.06	96.16	102.26

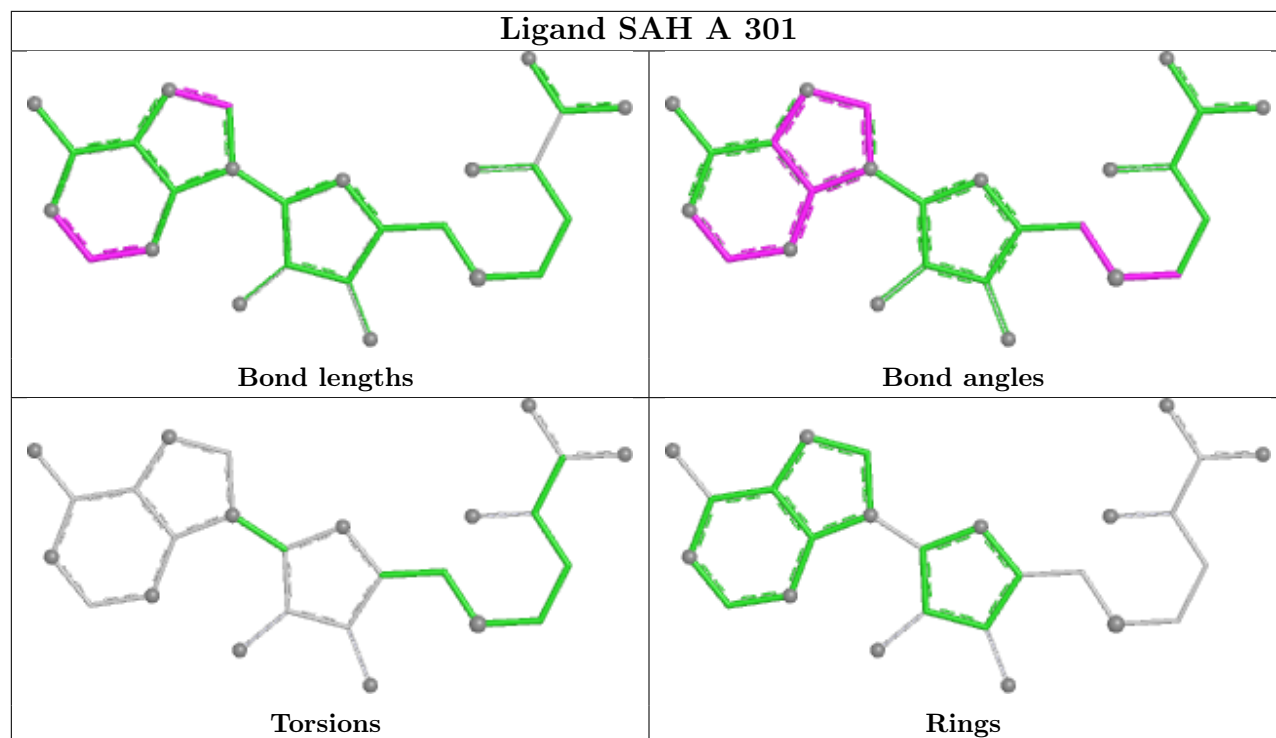
There are no chirality outliers.

There are no torsion outliers.

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 [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	282/289 (97%)	0.41	26 (9%) 14 13	32, 45, 82, 117	0
2	B	371/497 (74%)	1.19	93 (25%) 1 1	36, 61, 122, 156	0
3	N	77/77 (100%)	0.53	12 (15%) 5 4	26, 59, 133, 159	1 (1%)
All	All	730/863 (84%)	0.82	131 (17%) 3 3	26, 56, 115, 159	1 (0%)

All (131) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	10.5
2	B	120	LYS	7.6
2	B	52	PHE	6.8
2	B	463	ASN	6.5
2	B	343	ILE	6.4
2	B	74	LEU	6.4
2	B	119	ILE	5.9
2	B	79	LYS	5.7
2	B	340	LYS	5.7
2	B	341	ASP	5.6
3	N	18	G	5.6
2	B	51	TRP	5.5
2	B	18	HIS	5.2
2	B	77	LYS	5.2
2	B	121	GLY	5.1
1	A	62	ARG	5.0
2	B	80	ARG	5.0
1	A	61	GLY	4.8
2	B	107	THR	4.8
2	B	54	LEU	4.7
2	B	93	ASP	4.7
2	B	76	PRO	4.7
1	A	289	GLY	4.6

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Mol	Chain	Res	Type	RSRZ
2	B	342	TYR	4.4
2	B	111	ILE	4.3
1	A	23	GLY	4.3
2	B	58	ILE	4.2
2	B	53	TYR	4.2
2	B	45	VAL	4.1
1	A	22	HIS	4.1
1	A	253	GLY	4.0
2	B	105	LYS	3.9
2	B	57	VAL	3.9
2	B	106	LEU	3.9
1	A	63	GLY	3.8
1	A	261	ASP	3.8
2	B	43	LYS	3.7
2	B	71	GLY	3.7
2	B	46	THR	3.7
1	A	275	VAL	3.7
2	B	49	LYS	3.7
2	B	66	PHE	3.6
2	B	75	GLN	3.6
2	B	114	LEU	3.6
3	N	19	G	3.5
2	B	73	SER	3.4
2	B	56	ASN	3.4
1	A	64	GLY	3.4
2	B	21	ARG	3.4
2	B	78	LYS	3.3
2	B	100	ASP	3.3
2	B	72	GLY	3.3
1	A	42	VAL	3.3
2	B	108	GLN	3.3
2	B	206	MET	3.3
2	B	351	GLU	3.2
3	N	58[A]	A	3.2
1	A	252	THR	3.2
2	B	60	HIS	3.1
3	N	20	U	3.1
2	B	40	GLN	3.1
2	B	55	ASP	3.0
2	B	122	GLU	3.0
2	B	69	THR	3.0
2	B	88	LYS	3.0

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Mol	Chain	Res	Type	RSRZ
2	B	346	LYS	3.0
2	B	20	ILE	3.0
2	B	429	TRP	3.0
2	B	19	ARG	2.9
2	B	449	GLY	2.9
2	B	349	ARG	2.9
3	N	0	G	2.9
2	B	70	SER	2.9
2	B	41	ARG	2.8
3	N	59	G	2.8
2	B	103	SER	2.8
2	B	123	GLU	2.8
1	A	39	ARG	2.8
2	B	415	ARG	2.8
2	B	446	MET	2.7
1	A	41	GLY	2.7
2	B	98	VAL	2.7
3	N	17	C	2.7
2	B	110	ASP	2.7
2	B	234	GLY	2.7
2	B	124	ILE	2.7
2	B	271	SER	2.7
2	B	270	PHE	2.7
2	B	59	GLY	2.7
1	A	60	CYS	2.6
3	N	15	G	2.6
3	N	76	A	2.6
2	B	42	ARG	2.5
2	B	266	LEU	2.5
2	B	102	LYS	2.5
1	A	241	ARG	2.5
2	B	22	ASP	2.5
2	B	353	GLN	2.4
2	B	68	VAL	2.4
2	B	133	THR	2.4
2	B	358	LEU	2.4
2	B	67	GLU	2.4
2	B	157	ILE	2.4
2	B	269	THR	2.4
3	N	74	C	2.4
2	B	350	GLN	2.3
1	A	28	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	264	SER	2.3
1	A	264	PRO	2.3
2	B	101	GLY	2.3
2	B	436	LEU	2.3
1	A	66	VAL	2.3
1	A	242	THR	2.3
2	B	23	GLY	2.3
2	B	384	LEU	2.3
2	B	352	GLU	2.2
2	B	92	THR	2.2
2	B	156	ALA	2.2
2	B	38	GLN	2.2
2	B	61	SER	2.2
1	A	65	TRP	2.2
3	N	60	U	2.2
2	B	356	ARG	2.2
3	N	75	C	2.2
1	A	47	VAL	2.1
1	A	244	SER	2.1
2	B	233	GLY	2.1
1	A	234	LEU	2.1
1	A	32	ARG	2.1
2	B	50	GLN	2.0
2	B	112	LYS	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](#)

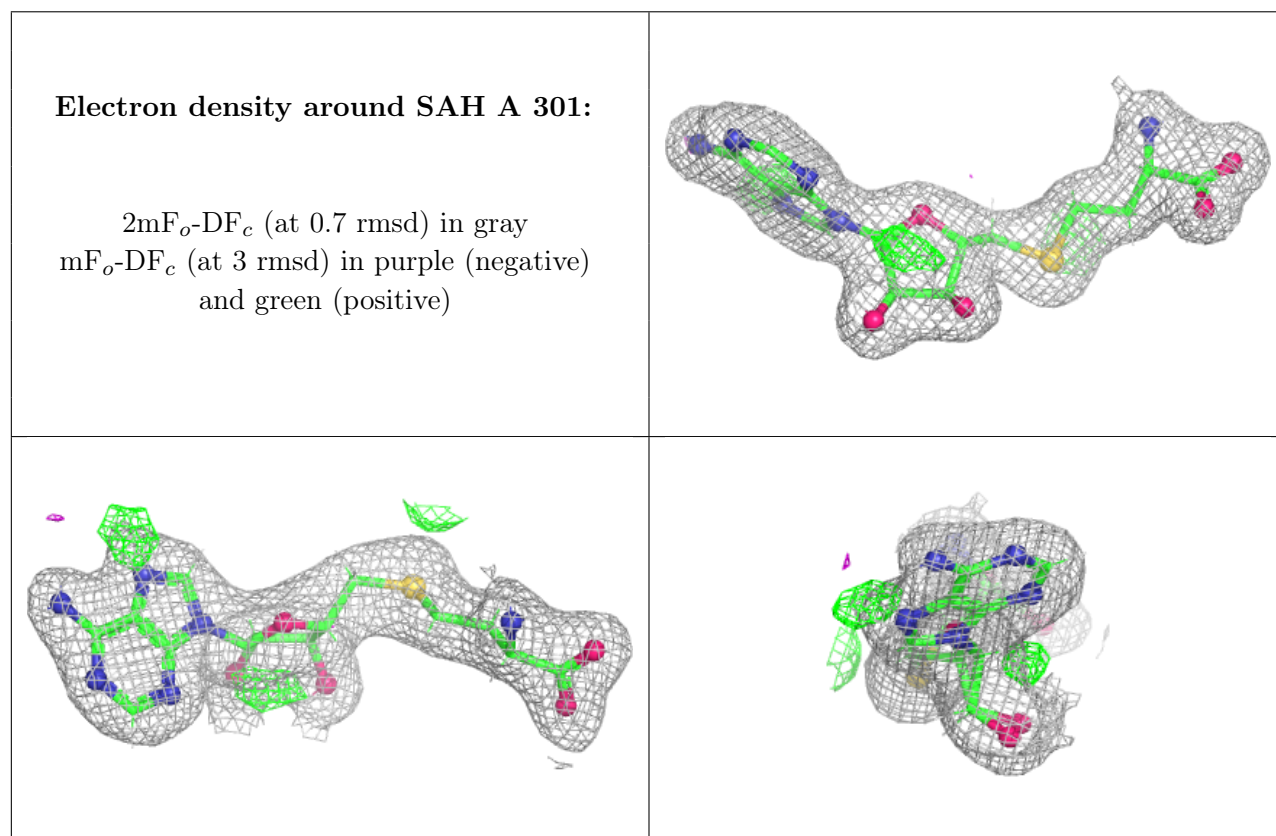
There are no oligosaccharides in this entry.

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(\AA^2)	Q<0.9
4	SAH	A	301	26/26	0.98	0.07	29,36,43,47	0
5	NA	N	101	1/1	0.99	0.04	42,42,42,42	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 [i](#)

There are no such residues in this entry.