



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:16 PM UTC

PDB ID : 5LSU / pdb_00005lsu
Title : Structure of the Epigenetic Oncogene MMSET and inhibition by N-Alkyl Sine-fungin Derivatives
Authors : Tisi, D.; Pathuri, P.; Heightman, T.
Deposited on : 2016-09-05
Resolution : 2.14 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

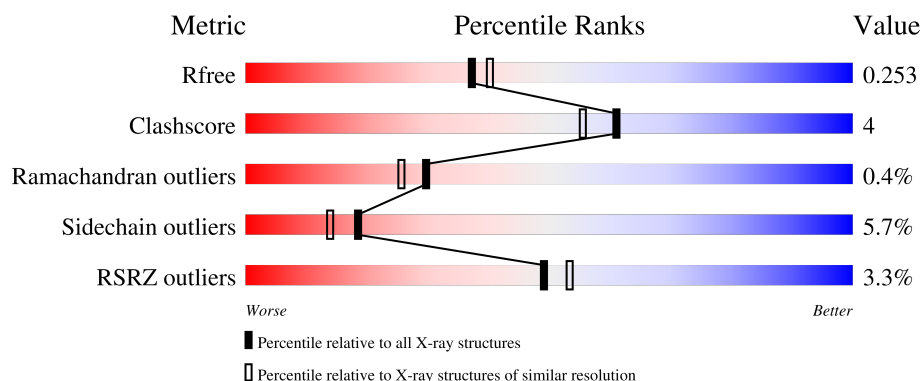
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

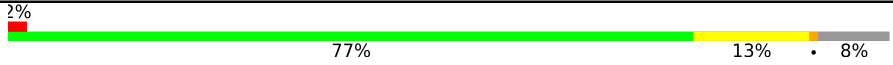
The reported resolution of this entry is 2.14 Å.

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	3689 (2.16-2.12)
Clashscore	190562	3812 (2.16-2.12)
Ramachandran outliers	187476	3773 (2.16-2.12)
Sidechain outliers	187428	3772 (2.16-2.12)
RSRZ outliers	180081	3691 (2.16-2.12)

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	252	
1	B	252	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4212 atoms, of which 46 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 Histone-lysine N-methyltransferase NSD2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	231	Total	C	N	O	S	0	1	0
			1825	1140	321	343	21			
1	B	222	Total	C	N	O	S	0	2	0
			1753	1100	310	321	22			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	952	MET	-	initiating methionine	UNP O96028
A	953	GLY	-	expression tag	UNP O96028
A	954	SER	-	expression tag	UNP O96028
A	955	SER	-	expression tag	UNP O96028
A	956	HIS	-	expression tag	UNP O96028
A	957	HIS	-	expression tag	UNP O96028
A	958	HIS	-	expression tag	UNP O96028
A	959	HIS	-	expression tag	UNP O96028
A	960	HIS	-	expression tag	UNP O96028
A	961	HIS	-	expression tag	UNP O96028
A	962	SER	-	expression tag	UNP O96028
A	963	SER	-	expression tag	UNP O96028
A	964	GLY	-	expression tag	UNP O96028
A	965	LEU	-	expression tag	UNP O96028
A	966	VAL	-	expression tag	UNP O96028
A	967	PRO	-	expression tag	UNP O96028
A	968	ARG	-	expression tag	UNP O96028
A	969	GLY	-	expression tag	UNP O96028
A	970	SER	-	expression tag	UNP O96028
A	971	HIS	-	expression tag	UNP O96028
A	972	MET	-	expression tag	UNP O96028
A	975	LEU	GLN	engineered mutation	UNP O96028
A	978	LEU	ALA	engineered mutation	UNP O96028
A	1071	LEU	ASP	engineered mutation	UNP O96028
A	1072	GLN	GLY	engineered mutation	UNP O96028

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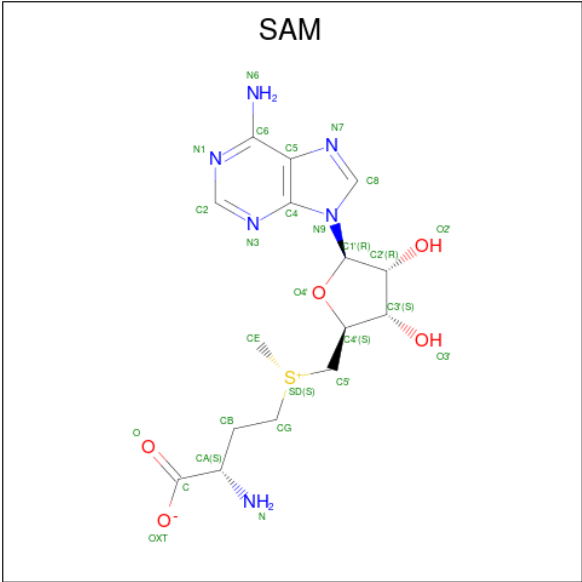
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Chain	Residue	Modelled	Actual	Comment	Reference
A	1073	ARG	LYS	engineered mutation	UNP O96028
B	952	MET	-	initiating methionine	UNP O96028
B	953	GLY	-	expression tag	UNP O96028
B	954	SER	-	expression tag	UNP O96028
B	955	SER	-	expression tag	UNP O96028
B	956	HIS	-	expression tag	UNP O96028
B	957	HIS	-	expression tag	UNP O96028
B	958	HIS	-	expression tag	UNP O96028
B	959	HIS	-	expression tag	UNP O96028
B	960	HIS	-	expression tag	UNP O96028
B	961	HIS	-	expression tag	UNP O96028
B	962	SER	-	expression tag	UNP O96028
B	963	SER	-	expression tag	UNP O96028
B	964	GLY	-	expression tag	UNP O96028
B	965	LEU	-	expression tag	UNP O96028
B	966	VAL	-	expression tag	UNP O96028
B	967	PRO	-	expression tag	UNP O96028
B	968	ARG	-	expression tag	UNP O96028
B	969	GLY	-	expression tag	UNP O96028
B	970	SER	-	expression tag	UNP O96028
B	971	HIS	-	expression tag	UNP O96028
B	972	MET	-	expression tag	UNP O96028
B	975	LEU	GLN	engineered mutation	UNP O96028
B	978	LEU	ALA	engineered mutation	UNP O96028
B	1071	LEU	ASP	engineered mutation	UNP O96028
B	1072	GLN	GLY	engineered mutation	UNP O96028
B	1073	ARG	LYS	engineered mutation	UNP O96028

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	3	Total Zn 3 3	0	0
2	B	3	Total Zn 3 3	0	0

- Molecule 3 is S-ADENOSYLMETHIONINE (CCD ID: SAM) (formula: C₁₅H₂₂N₆O₅S).

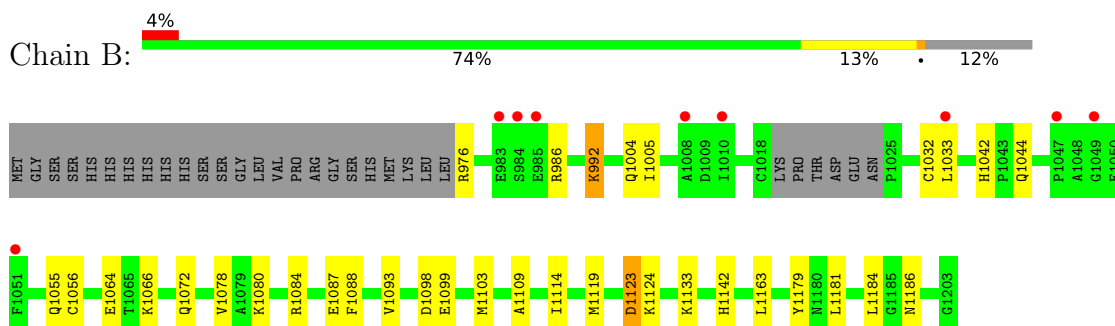


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	S	0	0
			50	15	23	6	5	1		
3	B	1	Total	C	H	N	O	S	0	0
			50	15	23	6	5	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	280	Total	O	0	0
			280	280		
4	B	248	Total	O	0	0
			248	248		

- Molecule 1: Histone-lysine N-methyltransferase NSD2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	94.58Å 63.11Å 81.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.81 – 2.14 49.81 – 2.14	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.81-2.14) 99.9 (49.81-2.14)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 2.14Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.171 , 0.238 0.184 , 0.253	Depositor DCC
R_{free} test set	1391 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	35.0	Xtriage
Anisotropy	0.678	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 78.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4212	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.52% 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: SAM, ZN

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	0.99	0/1868	1.24	1/2521 (0.0%)
1	B	0.98	1/1797 (0.1%)	1.27	10/2421 (0.4%)
All	All	0.98	1/3665 (0.0%)	1.25	11/4942 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1005	ILE	CG1-CD1	5.52	1.73	1.51

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1098	ASP	CA-CB-CG	7.10	119.70	112.60
1	A	1093	VAL	CB-CA-C	6.30	119.97	110.90
1	B	1123	ASP	CA-CB-CG	6.06	118.66	112.60
1	B	1055	GLN	CA-C-N	5.93	128.55	120.54
1	B	1055	GLN	C-N-CA	5.93	128.55	120.54
1	B	1186	ASN	CA-C-N	5.34	127.44	120.28
1	B	1186	ASN	C-N-CA	5.34	127.44	120.28
1	B	1032	CYS	CA-C-N	5.06	127.06	120.28
1	B	1032	CYS	C-N-CA	5.06	127.06	120.28
1	B	976	ARG	CA-C-N	5.05	127.46	120.29
1	B	976	ARG	C-N-CA	5.05	127.46	120.29

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	1825	0	1746	20	0
1	B	1753	0	1692	13	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	27	23	22	3	0
3	B	27	23	22	0	0
4	A	280	0	0	0	0
4	B	248	0	0	1	0
All	All	4166	46	3482	30	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 4.

All (30) 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:999:PRO:HB3	1:A:1005[A]:ILE:HD11	1.65	0.79
1:A:1071:LEU:HD13	1:B:1124:LYS:HG2	1.74	0.69
1:A:1071:LEU:CD1	1:B:1124:LYS:HG2	2.25	0.67
1:A:1141:ASN:HD22	3:A:1304:SAM:C8	2.17	0.58
1:A:1042:HIS:HD2	1:A:1044:GLN:H	1.52	0.56
1:A:973:LYS:HG3	1:A:976:ARG:HH21	1.70	0.55
1:A:1084:ARG:O	1:A:1087:GLU:HB2	2.10	0.51
1:A:1141:ASN:HD22	3:A:1304:SAM:H8	1.73	0.51
1:A:1096:LEU:HD21	1:A:1122:ILE:HD12	1.94	0.49
1:A:1057:PHE:HB3	1:A:1160:ARG:NH2	2.29	0.47
1:A:1141:ASN:HD21	3:A:1304:SAM:HN2	1.62	0.47
1:A:1099:GLU:O	1:A:1103:MET:HG3	2.16	0.46
1:B:1099:GLU:O	1:B:1103:MET:HG3	2.15	0.46
1:A:1192:ARG:HA	1:B:1004:GLN:HE22	1.81	0.46
1:A:1042:HIS:CD2	1:A:1044:GLN:H	2.32	0.45
1:A:1109:ALA:HA	1:A:1114:ILE:HD12	2.00	0.44
1:A:1057:PHE:HB3	1:A:1160:ARG:HH22	1.83	0.44
1:B:1119[A]:MET:HE2	4:B:1547:HOH:O	2.18	0.44
1:A:1026:CYS:HA	1:A:1032:CYS:HB2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1091:GLU:OE2	1:A:1160:ARG:HD2	2.18	0.44
1:B:992:LYS:HB3	1:B:1133[A]:LYS:HG2	1.98	0.44
1:B:1109:ALA:HA	1:B:1114:ILE:HD12	2.00	0.44
1:A:1064:GLU:HB3	1:A:1080:LYS:HD3	1.99	0.43
1:B:1142:HIS:HB2	1:B:1179:TYR:CG	2.54	0.43
1:B:1119[B]:MET:HB2	1:B:1119[B]:MET:HE3	1.56	0.43
1:B:1042:HIS:HD2	1:B:1044:GLN:H	1.66	0.42
1:B:1056:CYS:HB3	1:B:1088:PHE:CD2	2.55	0.41
1:B:1064:GLU:HB3	1:B:1080:LYS:HD3	2.00	0.41
1:A:1010:ILE:H	1:A:1010:ILE:HD12	1.85	0.41
1:B:1084:ARG:O	1:B:1087:GLU:HB2	2.21	0.41

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	230/252 (91%)	222 (96%)	6 (3%)	2 (1%)	14	8
1	B	220/252 (87%)	211 (96%)	9 (4%)	0	100	100
All	All	450/504 (89%)	433 (96%)	15 (3%)	2 (0%)	30	26

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	982	GLN
1	A	1010	ILE

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	200/224 (89%)	189 (94%)	11 (6%)	19	15
1	B	191/224 (85%)	180 (94%)	11 (6%)	18	13
All	All	391/448 (87%)	369 (94%)	22 (6%)	18	14

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	973	LYS
1	A	979	ARG
1	A	996	VAL
1	A	1004	GLN
1	A	1033	LEU
1	A	1066	LYS
1	A	1078	VAL
1	A	1093	VAL
1	A	1163	LEU
1	A	1181	LEU
1	A	1184	LEU
1	B	986	ARG
1	B	992	LYS
1	B	1033	LEU
1	B	1066	LYS
1	B	1072	GLN
1	B	1078	VAL
1	B	1093	VAL
1	B	1123	ASP
1	B	1163	LEU
1	B	1181	LEU
1	B	1184	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1042	HIS
1	A	1044	GLN
1	A	1061	GLN
1	A	1110	HIS
1	A	1141	ASN
1	B	1004	GLN
1	B	1042	HIS
1	B	1061	GLN
1	B	1072	GLN
1	B	1110	HIS

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](#)

Of 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 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	$\# Z > 2$
3	SAM	A	1304	-	27,29,29	0.62	0	34,42,42	0.56	0
3	SAM	B	1304	-	27,29,29	0.55	0	34,42,42	0.68	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	SAM	A	1304	-	-	3/17/33/33	0/3/3/3
3	SAM	B	1304	-	-	2/17/33/33	0/3/3/3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

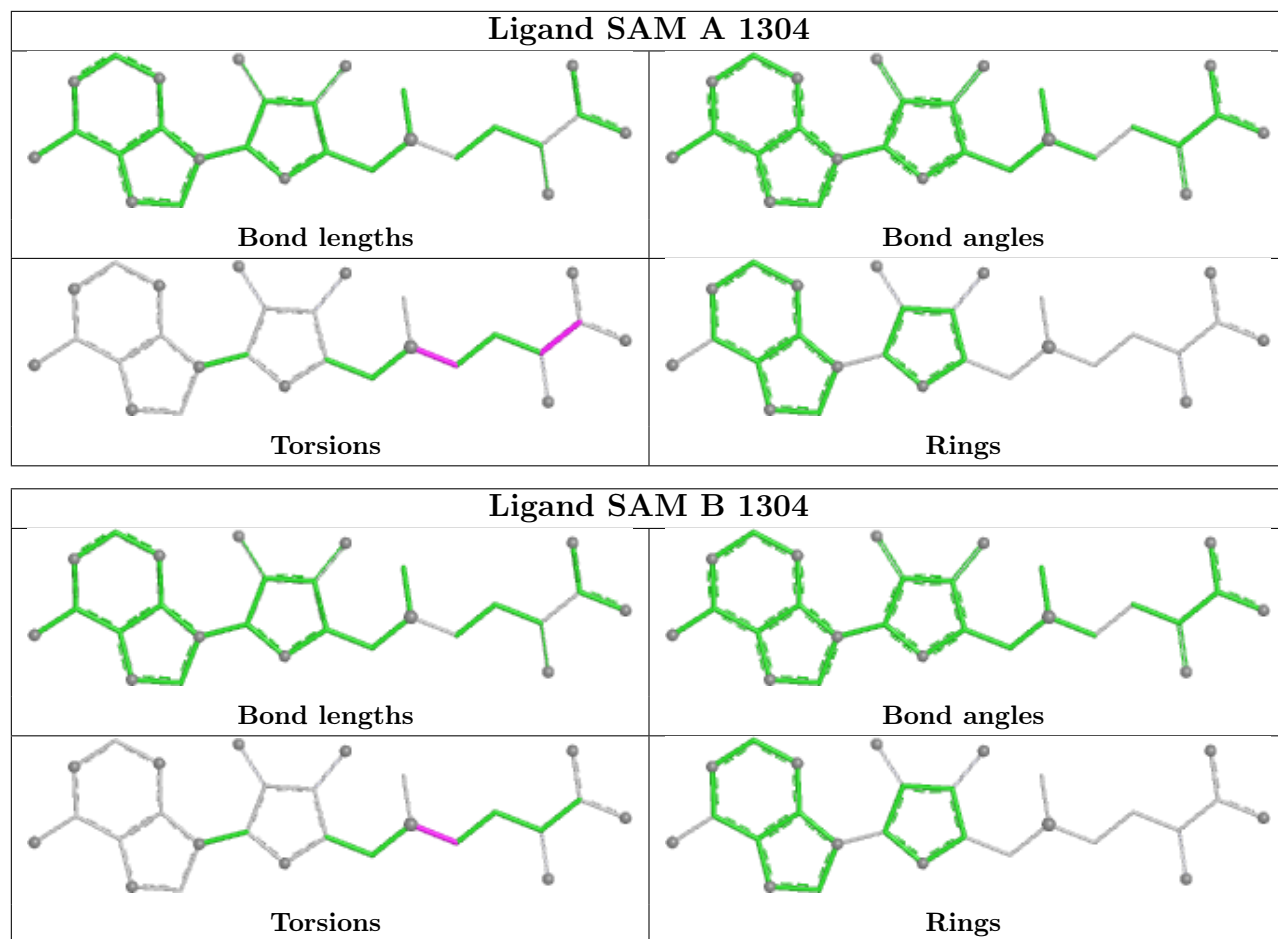
Mol	Chain	Res	Type	Atoms
3	A	1304	SAM	CB-CG-SD-CE
3	A	1304	SAM	CB-CG-SD-C5'
3	B	1304	SAM	CB-CG-SD-CE
3	B	1304	SAM	CB-CG-SD-C5'
3	A	1304	SAM	O-C-CA-N

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1304	SAM	3	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	231/252 (91%)	0.04	6 (2%) 57 61	22, 48, 94, 128	1 (0%)
1	B	222/252 (88%)	0.10	9 (4%) 41 46	17, 46, 105, 116	2 (0%)
All	All	453/504 (89%)	0.07	15 (3%) 49 53	17, 48, 104, 128	3 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1008	ALA	3.9
1	A	977	GLU	3.4
1	B	1033	LEU	3.1
1	B	1051	PHE	2.9
1	A	981	THR	2.8
1	B	984	SER	2.8
1	A	974	LEU	2.8
1	B	1010	ILE	2.3
1	A	978	LEU	2.3
1	B	983	GLU	2.2
1	B	985	GLU	2.1
1	B	1047	PRO	2.1
1	A	982	GLN	2.1
1	B	1049	GLY	2.0
1	A	1010	ILE	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](#)

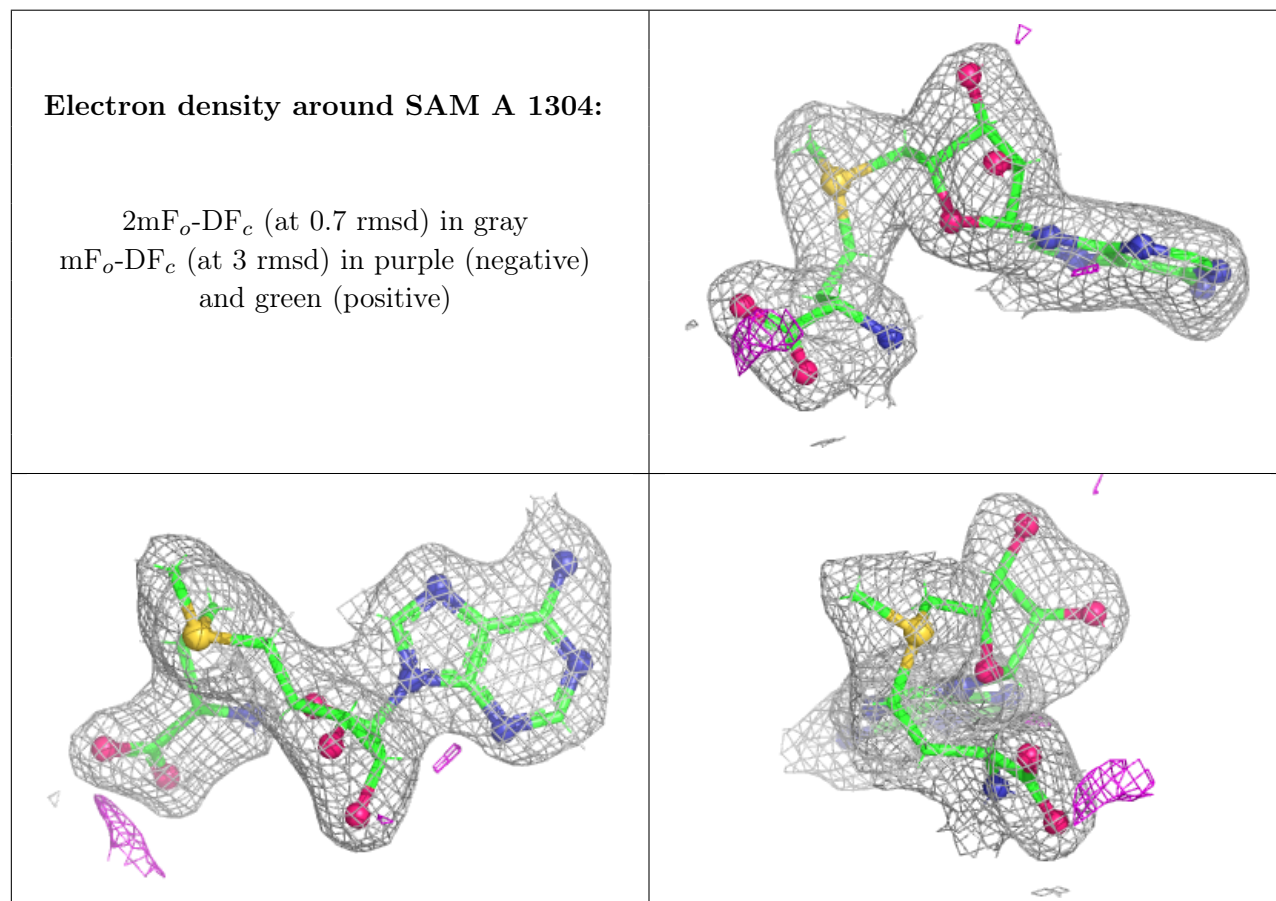
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

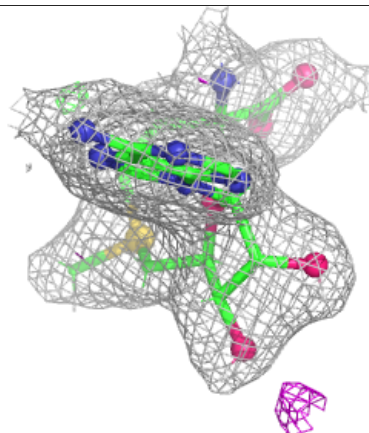
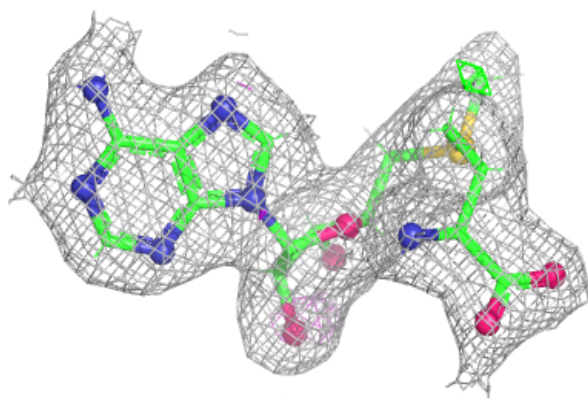
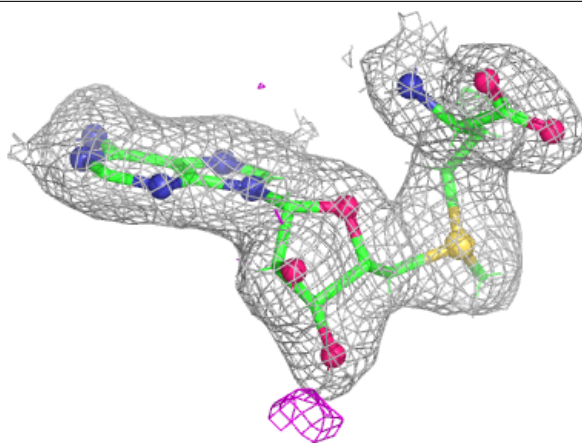
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	SAM	A	1304	27/27	0.95	0.07	28,44,51,54	0
2	ZN	B	1303	1/1	0.97	0.05	85,85,85,85	0
3	SAM	B	1304	27/27	0.97	0.05	24,31,39,42	0
2	ZN	B	1302	1/1	0.99	0.04	81,81,81,81	0
2	ZN	A	1301	1/1	0.99	0.05	42,42,42,42	0
2	ZN	A	1303	1/1	1.00	0.02	48,48,48,48	0
2	ZN	B	1301	1/1	1.00	0.04	32,32,32,32	0
2	ZN	A	1302	1/1	1.00	0.03	54,54,54,54	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.



Electron density around SAM B 1304:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.