



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 09:24 am BST

PDB ID : 4YNM  
Title : ASH1L wild-type SET domain in complex with S-adenosyl methionine (SAM)  
Authors : Rogawski, D.S.; Ndoj, J.; Cho, H.-J.; Maillard, I.; Grembecka, J.; Cierpicki, T.  
Deposited on : 2015-03-10  
Resolution : 2.19 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](mailto: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.

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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

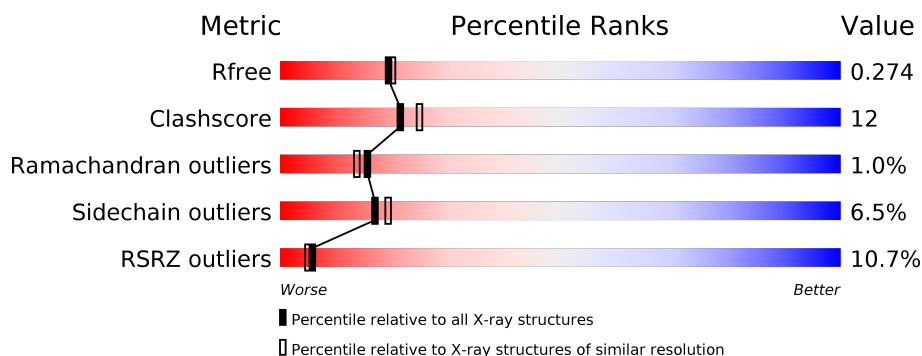
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.19 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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  $\geq 3$ , 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	226	
1	B	226	

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
2	ZN	B	2302	-	-	X	-

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3497 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 Histone-lysine N-methyltransferase ASH1L.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	215	Total	C	N	O	S	0	0	0
			1735	1075	310	328	22			
1	B	196	Total	C	N	O	S	0	0	0
			1598	997	284	297	20			

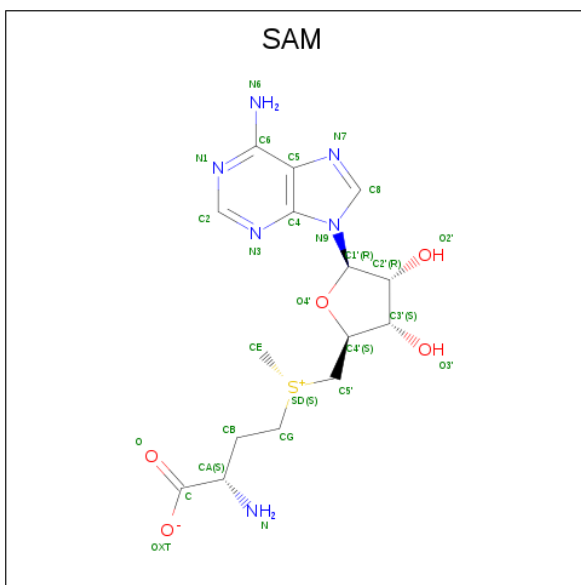
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2063	GLY	-	expression tag	UNP Q9NR48
A	2064	ALA	-	expression tag	UNP Q9NR48
A	2065	MET	-	expression tag	UNP Q9NR48
A	2066	ALA	-	expression tag	UNP Q9NR48
A	2067	GLY	-	expression tag	UNP Q9NR48
A	2068	SER	-	expression tag	UNP Q9NR48
B	2063	GLY	-	expression tag	UNP Q9NR48
B	2064	ALA	-	expression tag	UNP Q9NR48
B	2065	MET	-	expression tag	UNP Q9NR48
B	2066	ALA	-	expression tag	UNP Q9NR48
B	2067	GLY	-	expression tag	UNP Q9NR48
B	2068	SER	-	expression tag	UNP Q9NR48

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is S-ADENOSYLMETHIONINE (three-letter code: SAM) (formula: C<sub>15</sub>H<sub>22</sub>N<sub>6</sub>O<sub>5</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 27	C 15	N 6	O 5	S 1	0	0
3	B	1	Total 27	C 15	N 6	O 5	S 1	0	0

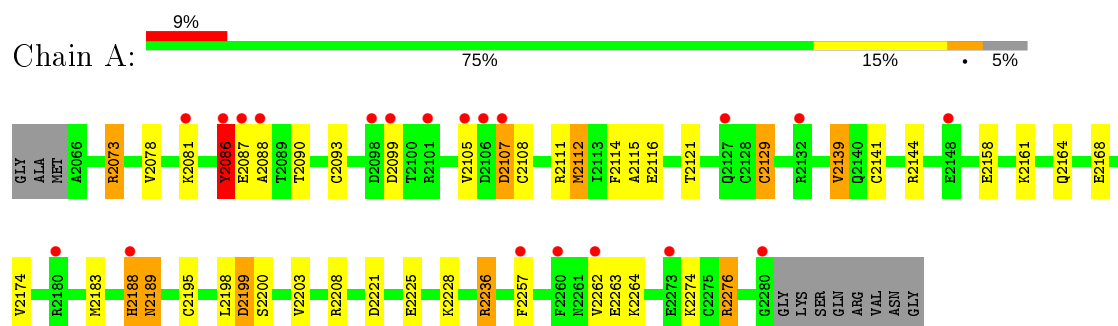
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	51	Total O 51 51	0	0
4	B	53	Total O 53 53	0	0

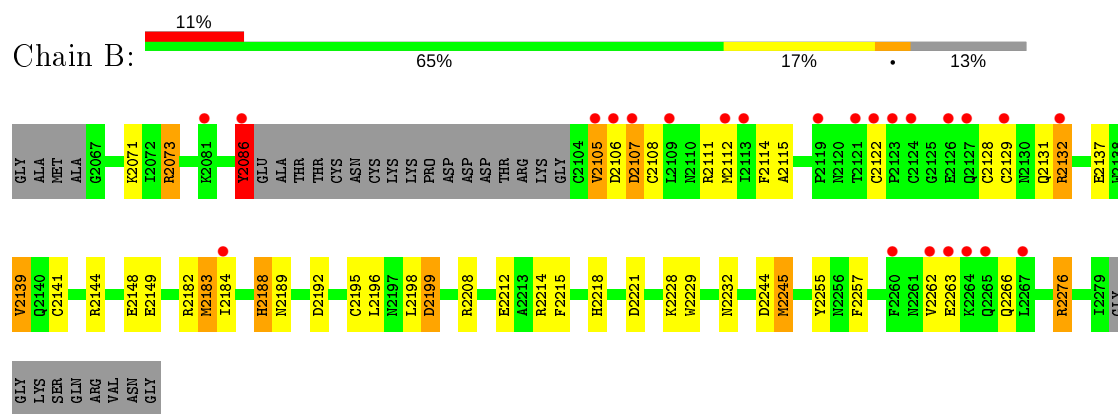
### 3 Residue-property plots [i](#)

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: Histone-lysine N-methyltransferase ASH1L



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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	59.14Å 59.14Å 230.98Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.82 – 2.19 46.82 – 2.19	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.82-2.19) 99.9 (46.82-2.19)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.43 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.242 , 0.278 0.236 , 0.274	Depositor DCC
$R_{free}$ test set	1271 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.9	Xtriage
Anisotropy	0.023	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 36.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.044 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3497	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.13% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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: ZN, SAM

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	0.80	0/1772	0.92	9/2383 (0.4%)
1	B	0.87	3/1633 (0.2%)	1.02	13/2195 (0.6%)
All	All	0.84	3/3405 (0.1%)	0.97	22/4578 (0.5%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	2195	CYS	CB-SG	-5.85	1.72	1.81
1	B	2199	ASP	CB-CG	-5.15	1.41	1.51
1	B	2129	CYS	CB-SG	-5.04	1.73	1.81

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	2132	ARG	NE-CZ-NH2	-10.81	114.89	120.30
1	B	2132	ARG	NE-CZ-NH1	9.54	125.07	120.30
1	A	2236	ARG	NE-CZ-NH1	8.55	124.58	120.30
1	A	2199	ASP	CB-CG-OD2	-8.36	110.78	118.30
1	B	2086	TYR	CA-CB-CG	6.53	125.81	113.40
1	B	2086	TYR	CB-CG-CD1	-6.37	117.18	121.00
1	B	2086	TYR	CB-CG-CD2	6.35	124.81	121.00
1	A	2199	ASP	CB-CG-OD1	6.11	123.80	118.30
1	A	2144	ARG	NE-CZ-NH2	6.06	123.33	120.30
1	B	2245	MET	CG-SD-CE	-6.05	90.52	100.20
1	A	2086	TYR	CA-CB-CG	5.89	124.59	113.40
1	B	2244	ASP	CB-CG-OD2	5.87	123.58	118.30
1	B	2221	ASP	CB-CG-OD1	5.66	123.39	118.30
1	B	2132	ARG	CG-CD-NE	-5.59	100.05	111.80
1	A	2236	ARG	NE-CZ-NH2	-5.49	117.55	120.30
1	B	2073	ARG	NE-CZ-NH2	5.42	123.01	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed( <sup>o</sup> )	Ideal( <sup>o</sup> )
1	A	2189	ASN	N-CA-CB	-5.30	101.06	110.60
1	B	2196	LEU	CA-CB-CG	5.20	127.26	115.30
1	A	2221	ASP	CB-CG-OD2	5.08	122.87	118.30
1	B	2182	ARG	NE-CZ-NH2	-5.07	117.77	120.30
1	B	2192	ASP	CB-CG-OD2	5.04	122.84	118.30
1	A	2225	GLU	OE1-CD-OE2	-5.01	117.29	123.30

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	1735	0	1642	53	2
1	B	1598	0	1513	35	2
2	A	3	0	0	0	0
2	B	3	0	0	2	0
3	A	27	0	22	0	0
3	B	27	0	22	2	0
4	A	51	0	0	3	0
4	B	53	0	0	0	0
All	All	3497	0	3199	79	2

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 12.

All (79) 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:2183:MET:CE	1:A:2262:VAL:HG11	1.91	1.00
1:A:2114:PHE:CE2	1:B:2073:ARG:NH1	2.29	0.98
1:B:2199:ASP:OD2	1:B:2228:LYS:NZ	1.97	0.96
1:A:2073:ARG:HH11	1:A:2073:ARG:HG3	1.29	0.96
1:A:2276:ARG:HH11	1:A:2276:ARG:HG3	1.31	0.94
1:A:2183:MET:HE1	1:A:2262:VAL:HG11	1.52	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2195:CYS:SG	1:A:2203:VAL:CG1	2.61	0.88
1:A:2114:PHE:CZ	1:B:2073:ARG:NH1	2.42	0.87
1:B:2183:MET:HE2	1:B:2262:VAL:HG21	1.60	0.82
1:A:2195:CYS:SG	1:A:2203:VAL:HG13	2.19	0.81
1:A:2183:MET:CE	1:A:2262:VAL:CG1	2.58	0.80
1:A:2183:MET:HE2	1:A:2262:VAL:HG11	1.64	0.78
1:B:2276:ARG:HH11	1:B:2276:ARG:HG3	1.48	0.77
1:A:2198:LEU:O	1:A:2199:ASP:HB2	1.85	0.76
1:A:2199:ASP:OD1	1:A:2228:LYS:NZ	2.19	0.76
1:A:2183:MET:HE2	1:A:2262:VAL:CG1	2.17	0.75
1:A:2114:PHE:HE2	1:B:2073:ARG:HH12	1.33	0.74
1:B:2128:CYS:SG	2:B:2302:ZN:ZN	1.75	0.73
1:B:2122:CYS:SG	2:B:2302:ZN:ZN	1.76	0.73
1:A:2183:MET:HE1	1:A:2262:VAL:CG1	2.18	0.71
1:A:2195:CYS:SG	1:A:2203:VAL:HG11	2.33	0.68
1:A:2073:ARG:NH1	1:A:2073:ARG:HG3	2.00	0.68
1:A:2168:GLU:OE2	1:A:2236:ARG:HD2	1.94	0.67
1:B:2144:ARG:HD2	1:B:2215:PHE:CD2	2.30	0.67
1:A:2073:ARG:CG	1:A:2073:ARG:HH11	2.04	0.66
1:A:2262:VAL:O	1:A:2263:GLU:HB2	1.99	0.62
1:B:2132:ARG:HG2	1:B:2137:GLU:OE1	2.01	0.60
1:A:2086:TYR:HE1	1:A:2115:ALA:HB2	1.68	0.59
1:A:2276:ARG:NH1	1:A:2276:ARG:HG3	2.10	0.59
1:A:2088:ALA:CB	1:A:2116:GLU:O	2.51	0.59
1:A:2112:MET:HE2	1:B:2073:ARG:HB2	1.84	0.58
1:A:2276:ARG:CG	1:A:2276:ARG:HH11	2.10	0.58
1:B:2144:ARG:HD3	1:B:2212:GLU:OE2	2.05	0.56
1:A:2161:LYS:O	1:A:2164:GLN:HB3	2.05	0.56
1:B:2108:CYS:O	1:B:2112:MET:HG2	2.05	0.56
1:B:2128:CYS:SG	1:B:2131:GLN:NE2	2.79	0.56
1:A:2188:HIS:ND1	1:A:2189:ASN:N	2.54	0.56
1:B:2139:VAL:HG13	1:B:2141:CYS:SG	2.46	0.56
1:A:2112:MET:CE	1:B:2073:ARG:CB	2.85	0.55
1:A:2141:CYS:CB	1:A:2158:GLU:HG3	2.37	0.55
1:A:2139:VAL:HG13	1:A:2141:CYS:SG	2.47	0.54
1:A:2108:CYS:O	1:A:2112:MET:HG2	2.07	0.54
1:A:2168:GLU:OE2	1:A:2236:ARG:CD	2.56	0.53
1:A:2105:VAL:HG22	1:A:2129:CYS:SG	2.48	0.53
1:A:2093:CYS:SG	1:A:2108:CYS:HA	2.49	0.52
1:B:2105:VAL:O	1:B:2111:ARG:HD2	2.09	0.52
1:B:2183:MET:CE	1:B:2262:VAL:HG11	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2112:MET:CE	1:B:2073:ARG:HB3	2.40	0.51
1:A:2112:MET:HE2	1:B:2073:ARG:CB	2.41	0.51
1:A:2112:MET:CE	1:B:2073:ARG:HB2	2.41	0.50
1:A:2262:VAL:O	1:A:2263:GLU:CB	2.58	0.50
1:B:2218:HIS:HB2	1:B:2255:TYR:CD2	2.47	0.49
1:A:2105:VAL:O	1:A:2111:ARG:HD2	2.13	0.48
1:B:2105:VAL:HG12	1:B:2106:ASP:N	2.29	0.48
1:B:2148:GLU:O	1:B:2149:GLU:HB2	2.13	0.48
1:A:2199:ASP:CG	1:A:2228:LYS:NZ	2.67	0.47
1:B:2214:ARG:O	3:B:2304:SAM:HE1	2.14	0.47
1:A:2274:LYS:HD3	4:A:2450:HOH:O	2.15	0.46
1:B:2276:ARG:NH1	1:B:2276:ARG:HG3	2.25	0.46
1:A:2141:CYS:HB2	1:A:2158:GLU:HG3	1.98	0.46
1:A:2112:MET:CE	1:A:2112:MET:HA	2.47	0.45
1:A:2189:ASN:HB3	4:A:2423:HOH:O	2.15	0.45
1:A:2188:HIS:ND1	1:A:2188:HIS:C	2.70	0.45
1:B:2183:MET:HE2	1:B:2262:VAL:CG2	2.38	0.44
1:A:2112:MET:HE1	1:B:2073:ARG:HB3	1.99	0.43
1:B:2114:PHE:HD2	1:B:2229:TRP:CZ2	2.36	0.43
1:B:2128:CYS:HB3	1:B:2131:GLN:NE2	2.33	0.43
1:B:2188:HIS:ND1	1:B:2189:ASN:N	2.66	0.42
1:A:2183:MET:CE	1:A:2262:VAL:HG13	2.46	0.42
1:A:2121:THR:OG1	4:A:2401:HOH:O	2.19	0.42
1:A:2262:VAL:HG12	1:A:2262:VAL:O	2.20	0.42
1:B:2183:MET:HE2	1:B:2262:VAL:HG11	2.00	0.42
1:B:2232:ASN:CG	1:B:2232:ASN:O	2.58	0.42
1:A:2198:LEU:O	1:A:2199:ASP:CB	2.54	0.42
1:B:2266:GLN:HE22	3:B:2304:SAM:C3'	2.29	0.41
1:A:2088:ALA:HB1	1:A:2116:GLU:O	2.19	0.41
1:A:2078:VAL:HB	1:A:2174:VAL:HG12	2.03	0.41
1:A:2188:HIS:ND1	1:A:2189:ASN:HB2	2.36	0.40
1:B:2086:TYR:HE1	1:B:2115:ALA:HB2	1.87	0.40

All (2) 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)
1:A:2081:LYS:CE	1:B:2184:ILE:CG1[4_455]	1.93	0.27
1:A:2208:ARG:NH1	1:B:2257:PHE:CD2[1_545]	2.17	0.03

## 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	213/226 (94%)	197 (92%)	14 (7%)	2 (1%)	17	16
1	B	192/226 (85%)	181 (94%)	9 (5%)	2 (1%)	15	14
All	All	405/452 (90%)	378 (93%)	23 (6%)	4 (1%)	15	14

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2107	ASP
1	B	2107	ASP
1	A	2090	THR
1	B	2105	VAL

### 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	191/198 (96%)	178 (93%)	13 (7%)	16	17
1	B	176/198 (89%)	165 (94%)	11 (6%)	18	20
All	All	367/396 (93%)	343 (94%)	24 (6%)	17	19

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2073	ARG
1	A	2086	TYR

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Mol	Chain	Res	Type
1	A	2087	GLU
1	A	2099	ASP
1	A	2107	ASP
1	A	2112	MET
1	A	2129	CYS
1	A	2139	VAL
1	A	2188	HIS
1	A	2200	SER
1	A	2257	PHE
1	A	2264	LYS
1	A	2276	ARG
1	B	2071	LYS
1	B	2086	TYR
1	B	2107	ASP
1	B	2139	VAL
1	B	2183	MET
1	B	2188	HIS
1	B	2198	LEU
1	B	2208	ARG
1	B	2245	MET
1	B	2263	GLU
1	B	2276	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	2266	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

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	B	2304	-	21,29,29	1.36	2 (9%)	18,42,42	1.57	1 (5%)
3	SAM	A	2304	-	21,29,29	1.33	2 (9%)	18,42,42	1.58	1 (5%)

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	B	2304	-	-	3/8/33/33	0/3/3/3
3	SAM	A	2304	-	-	2/8/33/33	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	2304	SAM	C2-N3	4.50	1.39	1.32
3	B	2304	SAM	C2-N3	4.30	1.39	1.32
3	B	2304	SAM	C2-N1	2.80	1.39	1.33
3	A	2304	SAM	C2-N1	2.49	1.38	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	2304	SAM	N3-C2-N1	-5.65	119.84	128.68
3	B	2304	SAM	N3-C2-N1	-5.30	120.39	128.68

There are no chirality outliers.

All (5) torsion outliers are listed below:

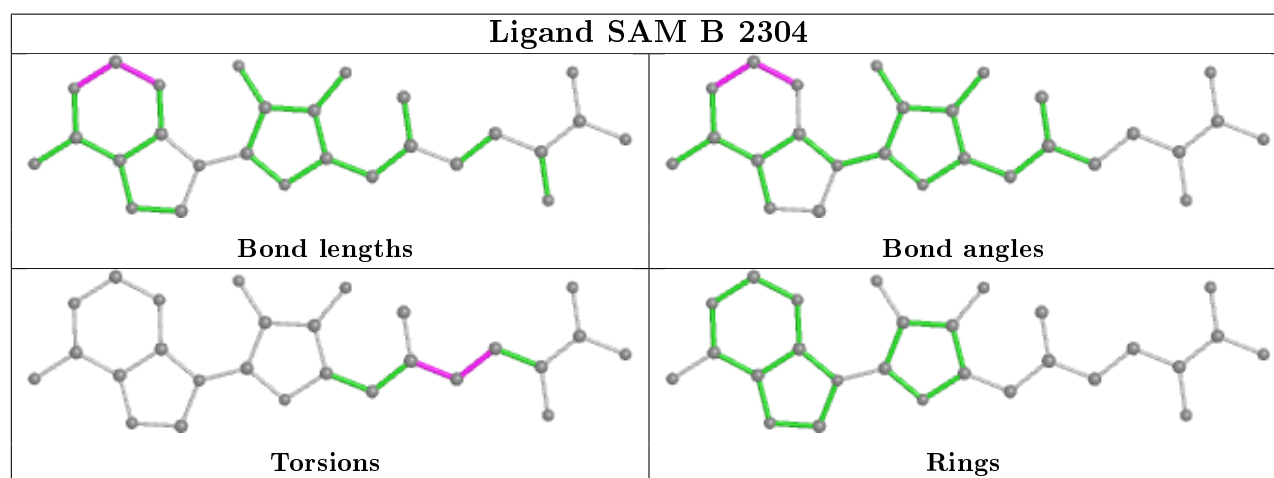
Mol	Chain	Res	Type	Atoms
3	A	2304	SAM	CB-CG-SD-CE
3	A	2304	SAM	CB-CG-SD-C5'
3	B	2304	SAM	CB-CG-SD-CE
3	B	2304	SAM	CB-CG-SD-C5'
3	B	2304	SAM	CA-CB-CG-SD

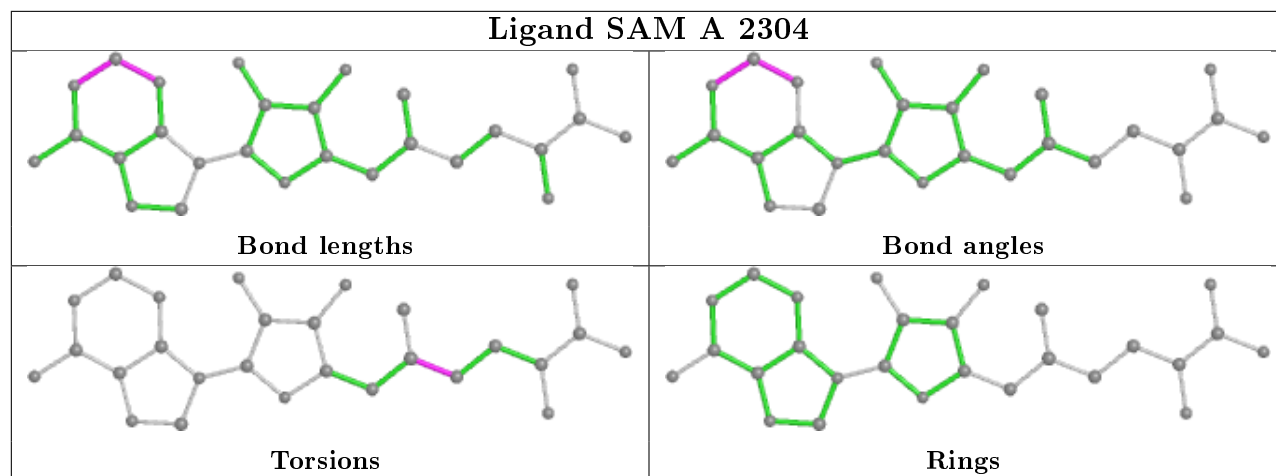
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	2304	SAM	2	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	215/226 (95%)	0.57	20 (9%) <b>8</b> <b>7</b>	23, 40, 77, 90	0
1	B	196/226 (86%)	0.77	24 (12%) <b>4</b> <b>3</b>	23, 39, 67, 83	0
All	All	411/452 (90%)	0.66	44 (10%) <b>6</b> <b>5</b>	23, 40, 73, 90	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	2109	LEU	8.7
1	A	2088	ALA	6.7
1	B	2264	LYS	6.7
1	A	2106	ASP	6.1
1	A	2257	PHE	6.1
1	A	2101	ARG	5.7
1	B	2086	TYR	5.7
1	A	2262	VAL	5.3
1	B	2107	ASP	4.9
1	B	2123	PRO	4.2
1	B	2106	ASP	4.1
1	A	2148	GLU	3.8
1	B	2127	GLN	3.8
1	A	2107	ASP	3.6
1	B	2124	CYS	3.5
1	B	2129	CYS	3.4
1	B	2260	PHE	3.2
1	A	2280	GLY	3.2
1	B	2262	VAL	3.2
1	B	2105	VAL	3.0
1	B	2113	ILE	2.9
1	A	2132	ARG	2.9
1	A	2099	ASP	2.8
1	B	2263	GLU	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	2087	GLU	2.8
1	B	2121	THR	2.7
1	A	2180	ARG	2.7
1	A	2188	HIS	2.6
1	A	2081	LYS	2.6
1	B	2081	LYS	2.6
1	A	2273	GLU	2.6
1	A	2086	TYR	2.5
1	B	2265	GLN	2.5
1	B	2119	PRO	2.4
1	B	2112	MET	2.3
1	A	2127	GLN	2.3
1	A	2260	PHE	2.2
1	B	2122	CYS	2.1
1	A	2098	ASP	2.1
1	B	2126	GLU	2.1
1	B	2132	ARG	2.1
1	B	2184	ILE	2.0
1	B	2267	LEU	2.0
1	A	2105	VAL	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 carbohydrates 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, 95<sup>th</sup> 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
2	ZN	B	2303	1/1	0.91	0.13	79,79,79,79	1
3	SAM	B	2304	27/27	0.94	0.12	32,36,40,48	0
3	SAM	A	2304	27/27	0.96	0.11	28,31,37,49	0

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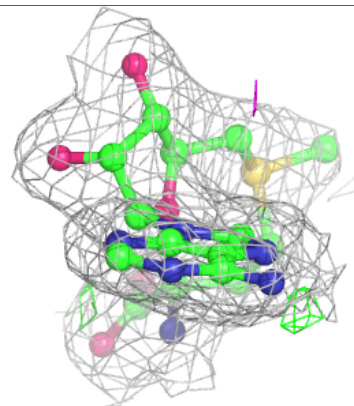
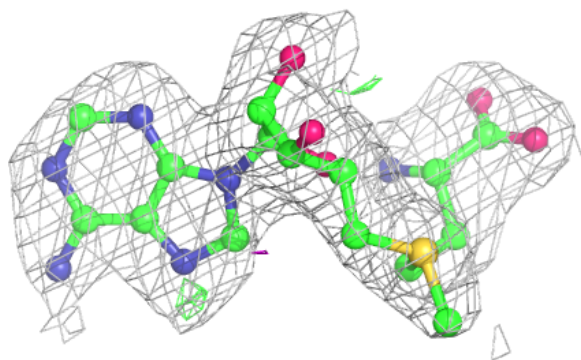
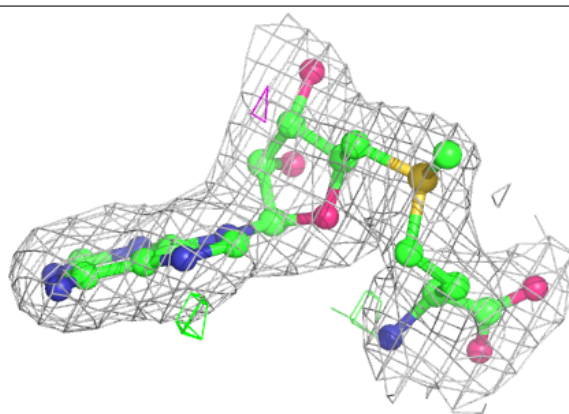
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ZN	B	2302	1/1	0.96	0.17	53,53,53,53	1
2	ZN	B	2301	1/1	0.98	0.16	28,28,28,28	1
2	ZN	A	2303	1/1	0.98	0.11	37,37,37,37	1
2	ZN	A	2301	1/1	0.99	0.13	30,30,30,30	1
2	ZN	A	2302	1/1	0.99	0.10	33,33,33,33	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.

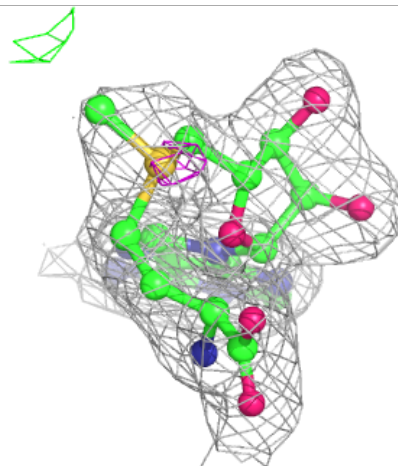
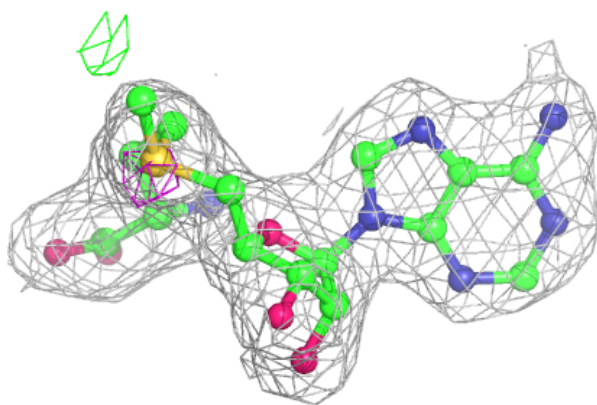
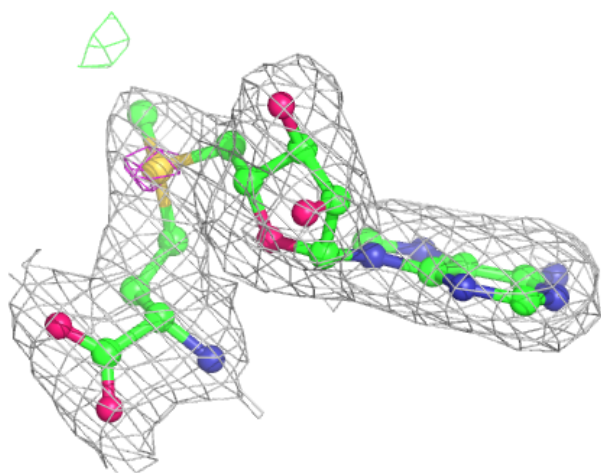
**Electron density around SAM B 2304:**

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around SAM A 2304:**

$2mF_o-DF_c$  (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.