



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 01:25 PM UTC

PDB ID : 6X9I / pdb_00006x9i
Title : Human DNMT1(729-1600) Bound to Zebularine-Containing 12mer dsDNA
Authors : Pathuri, S.; Horton, J.R.; Cheng, X.
Deposited on : 2020-06-02
Resolution : 2.20 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

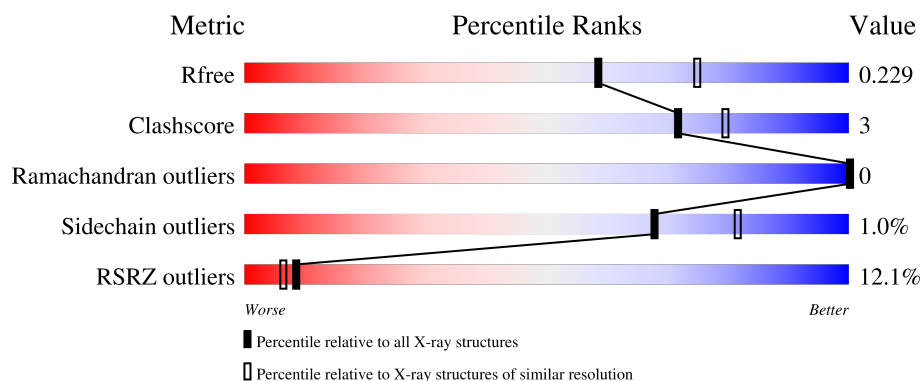
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (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 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	874	
2	C	12	
3	D	12	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6956 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 DNA (cytosine-5)-methyltransferase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	815	Total	C	N	O	S	0	0	0
			6194	3935	1088	1129	42			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	727	HIS	-	expression tag	UNP P26358
A	728	MET	-	expression tag	UNP P26358

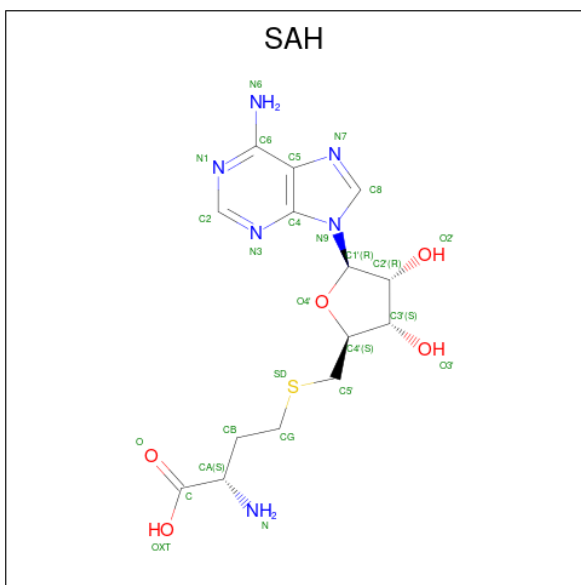
- Molecule 2 is a DNA chain called DNA (5'-D(*GP*AP*GP*GP*CP*(5CM)P*GP*CP*CP*TP*GP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	12	Total	C	N	O	P	0	0	0
			244	116	47	70	11			

- Molecule 3 is a DNA chain called DNA (5'-D(*GP*CP*AP*GP*G)-R(P*(PYO))-D(P*GP*GP*CP*CP*TP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	12	Total	C	N	O	P	0	0	0
			243	115	46	71	11			

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

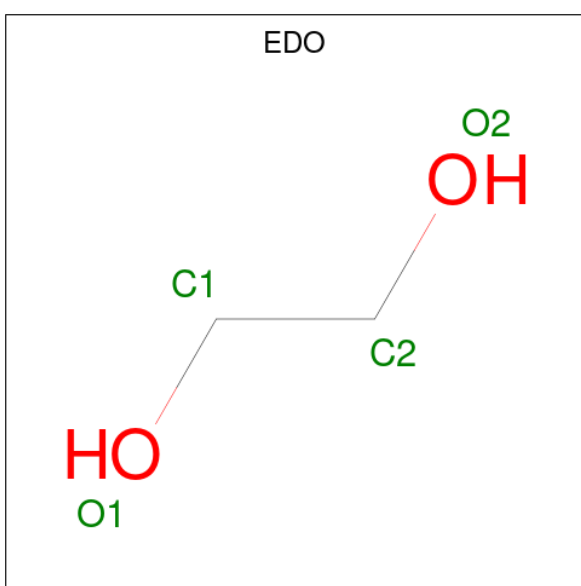


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			26	14	6	5	1		

- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

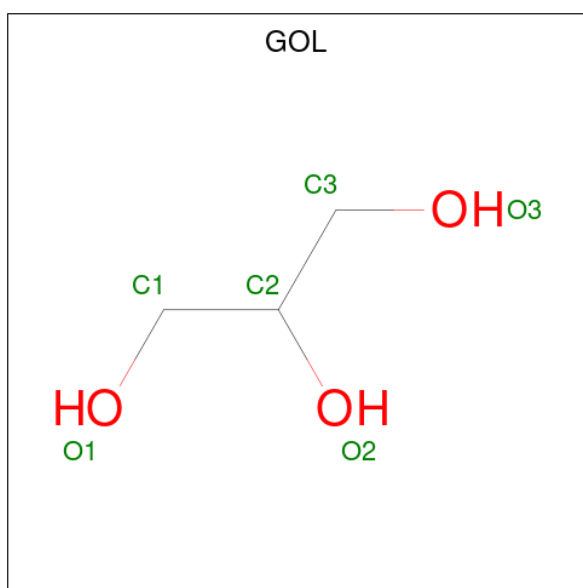
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Zn	0	0
			2	2		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0

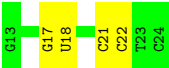
- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C O 6 3 3	0	0

- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	208	Total O 208 208	0	0
8	C	4	Total O 4 4	0	0
8	D	13	Total O 13 13	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	161.89Å 77.70Å 115.41Å 90.00° 125.67° 90.00°	Depositor
Resolution (Å)	45.21 – 2.20 45.21 – 2.20	Depositor EDS
% Data completeness (in resolution range)	93.5 (45.21-2.20) 93.5 (45.21-2.20)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, R_{free}	0.194 , 0.225 0.196 , 0.229	Depositor DCC
R_{free} test set	2000 reflections (3.39%)	wwPDB-VP
Wilson B-factor (Å ²)	45.6	Xtriage
Anisotropy	0.337	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 56.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6956	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.62% 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: 5CM, PYO, ZN, SAH, EDO, GOL

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	0.19	0/6356	0.39	0/8654
2	C	0.26	0/250	0.48	0/382
3	D	0.23	0/250	0.44	0/382
All	All	0.19	0/6856	0.40	0/9418

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	6194	0	5780	41	0
2	C	244	0	137	2	0
3	D	243	0	133	2	0
4	A	26	0	19	0	0
5	A	2	0	0	0	0
6	A	16	0	24	0	0
7	A	6	0	8	0	0
8	A	208	0	0	1	0
8	C	4	0	0	0	0
8	D	13	0	0	0	0
All	All	6956	0	6101	42	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (42) 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:1232:MET:HG3	3:D:17:DG:H2'	1.58	0.83
1:A:1015:LYS:NZ	1:A:1052:ASP:OD1	2.22	0.72
1:A:995:ARG:NH2	1:A:1062:VAL:O	2.35	0.60
1:A:749:LYS:HG3	1:A:758:GLU:HA	1.85	0.57
1:A:1464:ASN:O	1:A:1473:ARG:NH1	2.36	0.57
1:A:793:HIS:ND1	1:A:822:ASP:OD1	2.39	0.56
1:A:1006:ASN:OD1	1:A:1007:GLY:N	2.38	0.56
1:A:1203:THR:HG23	1:A:1204:THR:HG23	1.90	0.54
1:A:790:GLN:HB3	1:A:825:LEU:HD12	1.90	0.52
1:A:1154:GLY:HA3	1:A:1585:ALA:HB3	1.90	0.52
1:A:1535:LYS:HE3	2:C:7:DG:O6	2.10	0.51
1:A:1443:GLU:HB3	1:A:1451:MET:HE3	1.92	0.50
1:A:1455:LEU:HB2	1:A:1494:THR:OG1	2.12	0.49
1:A:866:GLN:HG3	1:A:867:LEU:HG	1.95	0.49
1:A:1281:LYS:NZ	8:A:1817:HOH:O	2.47	0.48
1:A:1160:ILE:HD12	1:A:1219:MET:HE1	1.95	0.47
1:A:1514:TYR:HE2	1:A:1533:MET:HG3	1.79	0.47
1:A:1392:PRO:HB3	1:A:1401:ARG:HD3	1.97	0.47
1:A:941:LEU:HB2	1:A:992:ARG:HB2	1.97	0.46
1:A:1195:LEU:HG	1:A:1199:MET:HE2	1.97	0.46
1:A:1187:PHE:HB3	1:A:1189:GLU:HG3	1.97	0.46
1:A:1446:LEU:HD21	1:A:1452:ALA:HB2	1.99	0.45
1:A:1500:LEU:HB2	1:A:1501:PRO:HD3	1.98	0.45
1:A:971:GLU:OE1	1:A:974:ARG:HD2	2.17	0.44
1:A:1414:CYS:SG	1:A:1546:ARG:HD2	2.58	0.44
1:A:1155:PHE:HZ	1:A:1264:LEU:HD13	1.82	0.44
1:A:1476:CYS:SG	1:A:1477:SER:N	2.92	0.43
1:A:837:LYS:HG2	1:A:864:PHE:HB2	2.00	0.43
1:A:775:TYR:HB3	1:A:796:TRP:CZ3	2.54	0.42
1:A:913:GLN:HG2	1:A:923:TYR:CE2	2.54	0.42
1:A:1419:ALA:HB1	1:A:1446:LEU:HB3	2.01	0.42
1:A:748:LYS:C	1:A:749:LYS:HD2	2.44	0.42
1:A:777:ALA:HB2	1:A:796:TRP:CE3	2.55	0.42
1:A:898:ARG:H	1:A:898:ARG:HG2	1.68	0.41
1:A:1415:LYS:HB3	1:A:1417:MET:HE3	2.03	0.41
1:A:1269:ARG:O	1:A:1272:VAL:HG22	2.21	0.41
3:D:21:DC:H2''	3:D:22:DC:C6	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1415:LYS:HE2	1:A:1417:MET:HE1	2.02	0.40
1:A:1263:PHE:HB3	1:A:1317:ALA:HB3	2.03	0.40
1:A:1533:MET:HE1	2:C:4:DG:H2'	2.04	0.40
1:A:1002:PRO:HD2	1:A:1013:ASP:O	2.21	0.40
1:A:1441:ASN:O	1:A:1454:LYS:HE3	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	805/874 (92%)	782 (97%)	23 (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	621/753 (82%)	615 (99%)	6 (1%)	68	81

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	826	SER

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Mol	Chain	Res	Type
1	A	1045	SER
1	A	1240	TYR
1	A	1417	MET
1	A	1533	MET
1	A	1536	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1019	ASN
1	A	1181	ASN
1	A	1245	ASN
1	A	1505	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	PYO	D	18	3	16,20,21	1.20	2 (12%)	21,28,31	0.71	0
2	5CM	C	6	2,3	18,21,22	1.69	6 (33%)	24,30,33	1.69	7 (29%)

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	PYO	D	18	3	-	4/7/25/26	0/2/2/2
2	5CM	C	6	2,3	-	5/7/21/22	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	6	5CM	O2-C2	-3.60	1.17	1.23
2	C	6	5CM	C2-N1	3.26	1.46	1.40
3	D	18	PYO	C2-N1	3.09	1.46	1.40
2	C	6	5CM	C6-C5	2.98	1.39	1.34
2	C	6	5CM	C4-N3	2.25	1.37	1.34
2	C	6	5CM	C5-C4	-2.23	1.42	1.44
3	D	18	PYO	O2-C2	-2.21	1.19	1.23
2	C	6	5CM	C6-N1	2.11	1.41	1.38

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	6	5CM	O4'-C1'-N1	3.75	114.52	107.86
2	C	6	5CM	C6-N1-C2	-3.07	116.86	120.95
2	C	6	5CM	C1'-N1-C2	2.89	122.78	117.83
2	C	6	5CM	C5A-C5-C6	-2.73	119.16	122.85
2	C	6	5CM	C4-N3-C2	-2.25	117.68	120.81
2	C	6	5CM	C5-C6-N1	-2.23	120.89	123.31
2	C	6	5CM	N1-C2-N3	2.01	122.30	118.80

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	18	PYO	C2'-C1'-N1-C6
3	D	18	PYO	C2'-C1'-N1-C2
2	C	6	5CM	O4'-C1'-N1-C6
2	C	6	5CM	C2'-C1'-N1-C6
3	D	18	PYO	O4'-C1'-N1-C6
2	C	6	5CM	C3'-C4'-C5'-O5'
3	D	18	PYO	O4'-C1'-N1-C2
2	C	6	5CM	C2'-C1'-N1-C2
2	C	6	5CM	O4'-C1'-N1-C2

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, 2 are monoatomic - leaving 6 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
6	EDO	A	1708	-	3,3,3	0.43	0	2,2,2	0.38	0
7	GOL	A	1705	-	5,5,5	0.92	0	5,5,5	1.05	0
6	EDO	A	1706	-	3,3,3	0.43	0	2,2,2	0.37	0
6	EDO	A	1704	-	3,3,3	0.46	0	2,2,2	0.40	0
4	SAH	A	1701	-	27,28,28	1.13	2 (7%)	36,40,40	0.86	1 (2%)
6	EDO	A	1707	-	3,3,3	0.46	0	2,2,2	0.32	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
6	EDO	A	1708	-	-	1/1/1/1	-
7	GOL	A	1705	-	-	2/4/4/4	-
6	EDO	A	1706	-	-	0/1/1/1	-
6	EDO	A	1704	-	-	0/1/1/1	-
4	SAH	A	1701	-	-	1/15/31/31	0/3/3/3
6	EDO	A	1707	-	-	1/1/1/1	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1701	SAH	C8-N9	3.32	1.43	1.37
4	A	1701	SAH	C5-C4	2.07	1.42	1.39

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1701	SAH	O4'-C1'-N9	-2.11	104.03	108.09

There are no chirality outliers.

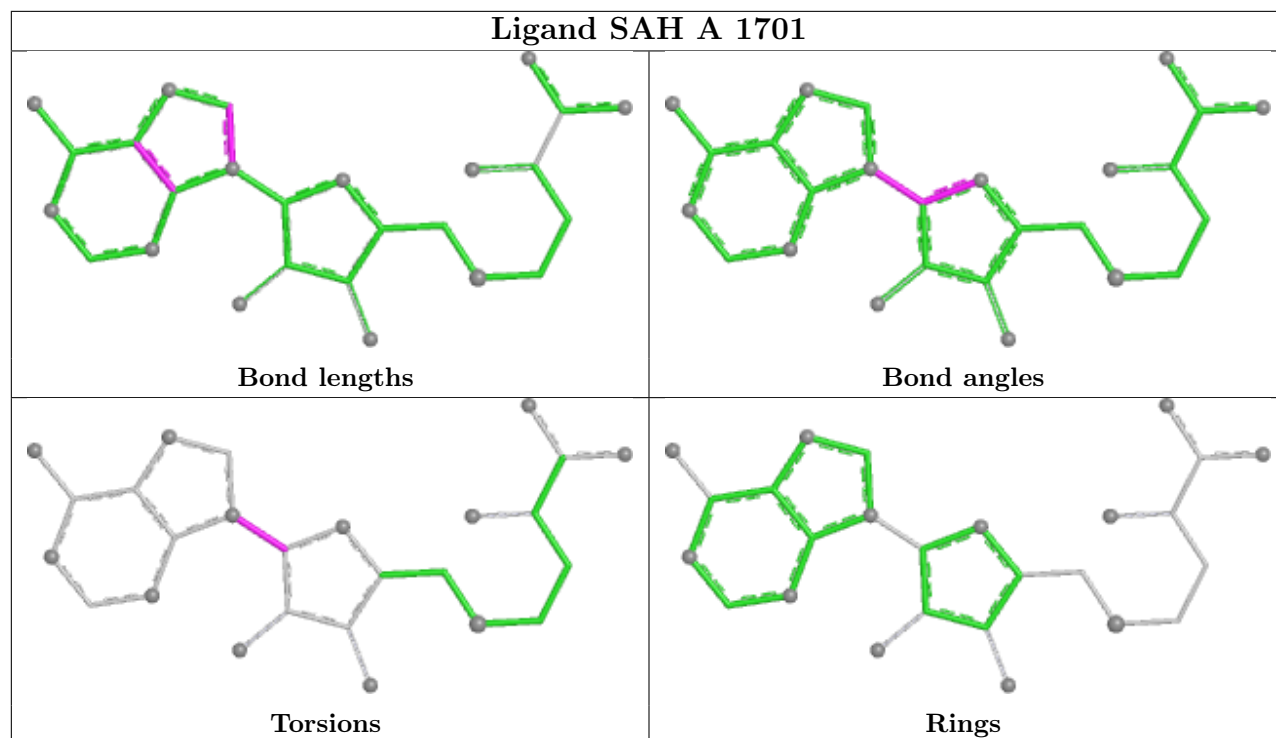
All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	1705	GOL	O1-C1-C2-C3
7	A	1705	GOL	O1-C1-C2-O2
4	A	1701	SAH	C2'-C1'-N9-C8
6	A	1707	EDO	O1-C1-C2-O2
6	A	1708	EDO	O1-C1-C2-O2

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	815/874 (93%)	0.67	101 (12%) 8 6	30, 66, 126, 158	0
2	C	11/12 (91%)	0.65	0 100 100	68, 87, 126, 142	0
3	D	11/12 (91%)	0.27	0 100 100	62, 78, 127, 142	0
All	All	837/898 (93%)	0.67	101 (12%) 8 6	30, 66, 126, 158	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1477	SER	5.0
1	A	754	ALA	4.3
1	A	1007	GLY	3.9
1	A	1456	ARG	3.9
1	A	890	PHE	3.9
1	A	884	PRO	3.7
1	A	787	SER	3.7
1	A	753	ASP	3.7
1	A	1486	ASP	3.6
1	A	976	TYR	3.5
1	A	1005	SER	3.5
1	A	950	ILE	3.4
1	A	1441	ASN	3.4
1	A	963	PRO	3.4
1	A	1489	ALA	3.4
1	A	1600	ALA	3.4
1	A	1240	TYR	3.3
1	A	1492	PHE	3.3
1	A	1450	THR	3.3
1	A	731	ILE	3.3
1	A	1481	ALA	3.2
1	A	738	VAL	3.2
1	A	1457	TYR	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	889	LYS	3.2
1	A	1465	GLY	3.2
1	A	1502	HIS	3.1
1	A	1239	THR	3.1
1	A	1453	ARG	3.1
1	A	756	THR	3.1
1	A	1448	ASP	3.1
1	A	732	SER	3.0
1	A	755	GLU	3.0
1	A	1452	ALA	3.0
1	A	1464	ASN	3.0
1	A	1472	LEU	2.9
1	A	1475	VAL	2.9
1	A	1462	ARG	2.9
1	A	1455	LEU	2.9
1	A	1485	CYS	2.9
1	A	751	CYS	2.9
1	A	967	ASP	2.8
1	A	1443	GLU	2.8
1	A	783	TRP	2.7
1	A	1471	ALA	2.7
1	A	740	THR	2.7
1	A	745	SER	2.7
1	A	1478	CYS	2.7
1	A	1498	TRP	2.7
1	A	979	TYR	2.7
1	A	1460	HIS	2.6
1	A	1488	ALA	2.6
1	A	892	PHE	2.6
1	A	730	ARG	2.6
1	A	850	ASP	2.6
1	A	750	VAL	2.6
1	A	779	VAL	2.6
1	A	1103	ALA	2.5
1	A	741	ASP	2.5
1	A	1006	ASN	2.5
1	A	1466	ARG	2.4
1	A	1479	VAL	2.4
1	A	789	GLY	2.4
1	A	835	ILE	2.4
1	A	918	ASP	2.4
1	A	968	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	1458	THR	2.4
1	A	964	VAL	2.4
1	A	769	ASP	2.3
1	A	737	ALA	2.3
1	A	1442	ILE	2.3
1	A	917	LEU	2.3
1	A	775	TYR	2.3
1	A	948	PHE	2.2
1	A	916	ASP	2.2
1	A	978	ASP	2.2
1	A	1497	PRO	2.2
1	A	1491	GLN	2.2
1	A	1496	ILE	2.2
1	A	1467	SER	2.2
1	A	788	ASN	2.2
1	A	888	ASN	2.2
1	A	1104	ARG	2.2
1	A	1490	ARG	2.2
1	A	734	VAL	2.2
1	A	965	ASP	2.2
1	A	1444	VAL	2.2
1	A	1531	GLU	2.2
1	A	1459	HIS	2.2
1	A	786	SER	2.1
1	A	973	TYR	2.1
1	A	1136	LEU	2.1
1	A	1487	PRO	2.1
1	A	796	TRP	2.1
1	A	1449	GLY	2.1
1	A	772	LYS	2.0
1	A	891	LYS	2.0
1	A	899	LEU	2.0
1	A	1505	ASN	2.0
1	A	1004	LYS	2.0
1	A	1348	LYS	2.0
1	A	1463	LYS	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column

labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	5CM	C	6	20/21	0.94	0.12	78,81,92,103	0
3	PYO	D	18	19/20	0.96	0.09	39,50,59,60	0

6.3 Carbohydrates [i](#)

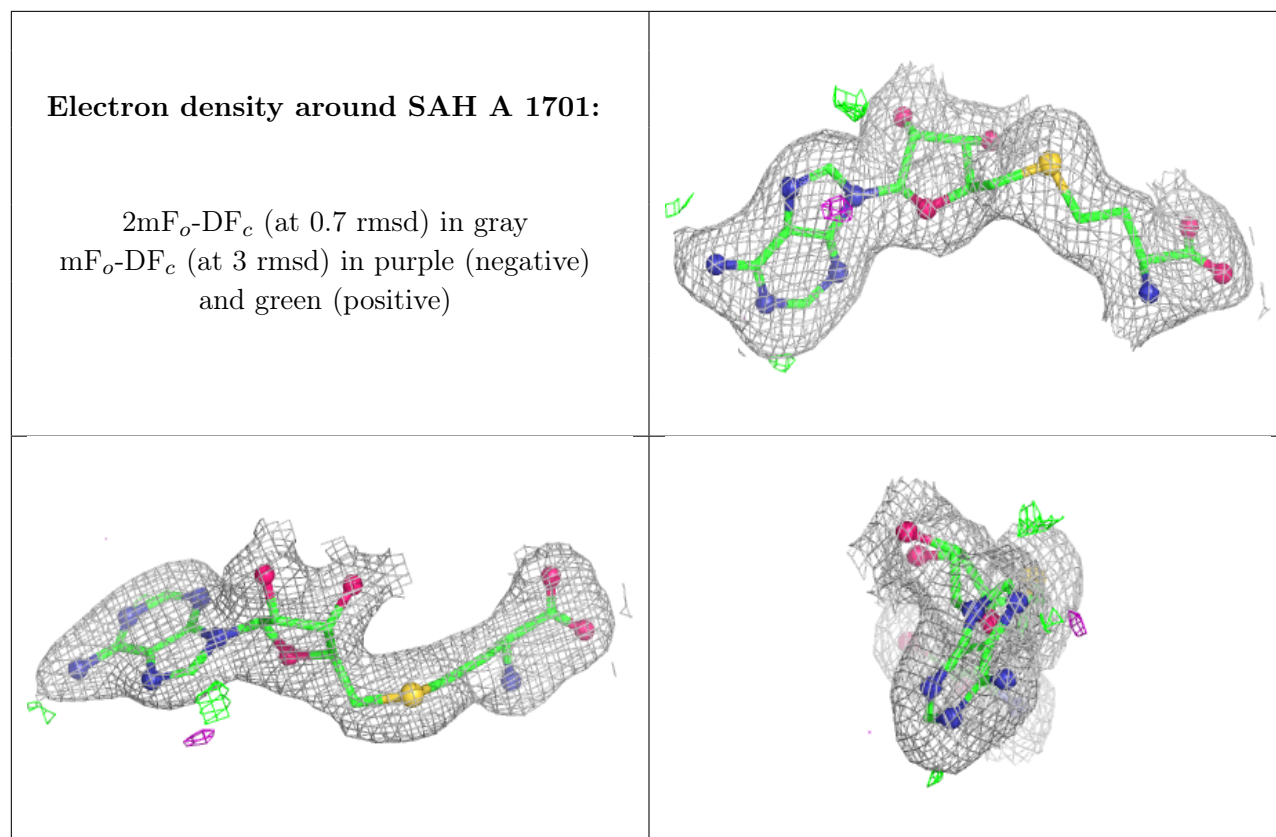
There are no oligosaccharides in this entry.

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
6	EDO	A	1704	4/4	0.82	0.16	63,71,71,76	0
7	GOL	A	1705	6/6	0.82	0.22	65,68,79,80	0
6	EDO	A	1708	4/4	0.84	0.28	51,56,57,63	0
6	EDO	A	1706	4/4	0.88	0.14	78,80,80,81	0
6	EDO	A	1707	4/4	0.92	0.13	61,64,68,71	0
5	ZN	A	1702	1/1	0.95	0.14	89,89,89,89	0
4	SAH	A	1701	26/26	0.96	0.07	33,39,48,50	0
5	ZN	A	1703	1/1	0.99	0.06	78,78,78,78	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.