

Full wwPDB X-ray Structure Validation Report (i)

May 25, 2020 - 07:34 am BST

PDB ID : 1M0E

Title : ZEBULARINE: A NOVEL DNA METHYLATION INHIBITOR THAT

FORMS A COVALENT COMPLEX WITH DNA METHYLTRANSFERASE

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Deposited on : 2002-06-12

Resolution : 2.50 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity: 4.02b-467

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : NOT EXECUTED

EDS : NOT EXECUTED

buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

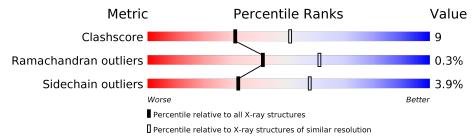
Validation Pipeline (wwPDB-VP) : 2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$

The reported resolution of this entry is 2.50 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\mathring{\rm A})}) \end{array}$
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)

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 >=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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality	y of chain	
1	С	12	42%	42%	17%
2	D	12	50%	50%	
3	A	327	81%		17% •



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 3379 atoms, of which 26 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 DNA chain called 5'-D(P*CP*CP*AP*TP*GP*CP*GP*CP*TP*GP*AP*C)-3'.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	С	12	Total 243	C 115	N 44	O 72	P 12	0	0	0

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

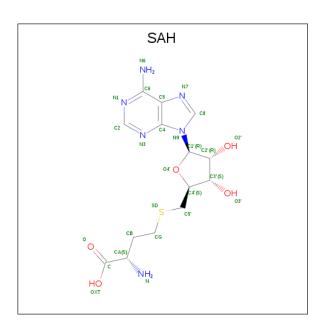
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	D	12	Total 248	C 117	N 47	O 72	P 12	0	0	0

• Molecule 3 is a protein called Modification methylase HhaI.

Mol	Chain	Residues	${f Atoms}$					ZeroOcc	AltConf	Trace	
3	A	327	Total 2617	C 1651	H 26	N 441	O 486	S 13	0	0	0

• Molecule 4 is S-ADENOSYL-L-HOMOCYSTEINE (three-letter code: SAH) (formula: $C_{14}H_{20}N_6O_5S$).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
1	Α	1	Total	С	N	О	S	0	0
4	A	1	26	14	6	5	1	U	0

• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	С	15	Total O 15 15	0	0
5	D	15	Total O 15 15	0	0
5	A	215	Total O 215 215	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. 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. 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.

Note EDS was not executed.



Chain C: 42% 42% 17%



• Molecule 2: 5'-D(P*GP*TP*CP*AP*GP*(Z)P*GP*CP*AP*TP*GP*G)-3'

Chain D: 50% 50%



• Molecule 3: Modification methylase HhaI

Chain A: 17% • 17%



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants	96.19Å 96.19Å 315.75Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.70 - 2.50	Depositor
% Data completeness	98.1 (39.70-2.50)	Depositor
(in resolution range)	,	Берозног
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.177 , 0.243	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3379	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP



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, Z

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		
MIOI	Chain	RMSZ	# Z >5	RMSZ	# Z > 5	
1	С	0.70	0/271	0.94	1/415~(0.2%)	
2	D	0.75	0/257	0.83	0/393	
3	A	0.54	0/2646	0.71	1/3569 (0.0%)	
All	All	0.58	0/3174	0.74	$2/4377 \ (0.0\%)$	

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	С	0	1
2	D	0	2
All	All	0	3

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^o)$
1	С	409	DC	C1'-O4'-C4'	-7.28	102.82	110.10
3	A	78	GLY	N-CA-C	-5.36	99.71	113.10

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	С	413	DC	Sidechain
2	D	422	DG	Sidechain
2	D	424	DC	Sidechain



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	С	243	0	135	6	0
2	D	248	0	133	4	0
3	A	2591	26	2549	47	0
4	A	26	0	19	0	0
5	A	215	0	0	4	0
5	С	15	0	0	0	0
5	D	15	0	0	0	0
All	All	3353	26	2836	54	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 9.

All (54) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} \textbf{Interatomic} \\ \textbf{distance} \ (\text{\r{A}}) \end{array}$	Clash overlap (Å)
3:A:172:ARG:HD2	3:A:174:ASP:OD1	1.85	0.76
1:C:406:DG:O6	3:A:256:GLY:HA3	1.86	0.75
3:A:209:ARG:HD3	3:A:258:ILE:HD12	1.72	0.71
3:A:82:GLN:HE21	3:A:82:GLN:H	1.38	0.70
3:A:309:ASN:HD22	3:A:309:ASN:H	1.38	0.69
3:A:156:ASP:HB3	3:A:185:PRO:HB3	1.79	0.64
3:A:89:LYS:HD3	5:A:540:HOH:O	1.96	0.63
3:A:137:LYS:NZ	3:A:141:ASN:HD21	2.00	0.59
3:A:131:ASN:O	3:A:135:VAL:HG23	2.04	0.58
2:D:429:DC:H4'	3:A:89:LYS:HE2	1.85	0.56
1:C:409:DC:H2'	1:C:410:DT:H72	1.88	0.54
3:A:187:GLU:HG2	5:A:459:HOH:O	2.08	0.54
3:A:48:VAL:HG22	3:A:292:HIS:CD2	2.43	0.53
3:A:25:ARG:HD2	3:A:53:PHE:CD2	2.44	0.53
3:A:103:ASP:O	3:A:107:ILE:HG13	2.10	0.52
3:A:309:ASN:HD22	3:A:309:ASN:N	2.07	0.52
3:A:210:LYS:O	3:A:210:LYS:HG2	2.11	0.51
2:D:432:DG:H2'	2:D:433:DG:C8	2.46	0.51
3:A:127:HIS:HD2	3:A:132:THR:OG1	1.94	0.50
3:A:220:GLU:HA	3:A:245:ARG:HH12	1.77	0.49

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A		Interatomic	Clash
Atom-1	Atom-2	${\rm distance}({\rm \AA})$	overlap (Å)
3:A:220:GLU:HA	3:A:245:ARG:NH1	2.27	0.49
1:C:405:DT:H2"	1:C:406:DG:C8	2.47	0.48
3:A:169:ILE:HG13	3:A:180:PHE:CZ	2.48	0.48
3:A:15:ILE:HG22	3:A:37:TYR:HB3	1.96	0.48
3:A:165:ARG:HH22	3:A:250:THR:HG21	1.80	0.47
3:A:169:ILE:HG13	3:A:180:PHE:CE2	2.49	0.47
3:A:48:VAL:HG22	3:A:292:HIS:HD2	1.79	0.47
3:A:66:GLU:HG3	3:A:66:GLU:H	1.47	0.46
3:A:125:ALA:HB2	3:A:166:ILE:HD12	1.97	0.45
3:A:80:PRO:HB3	3:A:82:GLN:HE22	1.82	0.45
3:A:66:GLU:HG2	3:A:106:ARG:HG2	1.99	0.44
2:D:425:DA:O3'	3:A:162:LYS:HE3	2.18	0.44
3:A:219:ILE:HD13	3:A:227:VAL:HG21	2.00	0.43
3:A:139:THR:HA	5:A:500:HOH:O	2.19	0.43
1:C:412:DA:H2"	1:C:413:DC:O4'	2.19	0.42
3:A:79:PHE:HB2	3:A:80:PRO:HD2	2.01	0.42
2:D:432:DG:H2"	2:D:433:DG:H5'	2.01	0.42
1:C:409:DC:H2'	1:C:410:DT:C7	2.49	0.42
3:A:156:ASP:HB3	3:A:185:PRO:CB	2.48	0.42
3:A:19:ALA:HB3	3:A:40:GLU:HB2	2.00	0.42
3:A:15:ILE:HD12	3:A:17:LEU:HD21	2.01	0.42
3:A:74:ILE:HA	3:A:115:VAL:O	2.20	0.42
3:A:155:LEU:HD12	3:A:155:LEU:C	2.40	0.42
3:A:90:GLN:O	3:A:127:HIS:HE1	2.03	0.42
3:A:258:ILE:HG12	3:A:258:ILE:O	2.20	0.41
3:A:79:PHE:HB2	3:A:80:PRO:CD	2.51	0.41
3:A:209:ARG:HB3	3:A:211:ASP:OD1	2.21	0.41
3:A:309:ASN:N	3:A:309:ASN:ND2	2.68	0.41
3:A:1:MET:HE1	5:A:524:HOH:O	2.21	0.41
1:C:408:DG:H1'	1:C:409:DC:C6	2.56	0.41
3:A:309:ASN:H	3:A:309:ASN:ND2	2.12	0.40
3:A:127:HIS:CD2	3:A:132:THR:OG1	2.74	0.40
3:A:59:GLY:O	3:A:60:ASP:C	2.59	0.40
3:A:271:THR:HG22	3:A:271:THR:O	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	$f{Analysed} ig f{Favoured} f{Aller}$		Outliers	Percentiles
3	A	$325/327 \ (99\%)$	309 (95%)	15 (5%)	1 (0%)	41 61

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	Α	60	ASP

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	Analysed Rotameric (Percentiles	
3	A	279/283 (99%)	268 (96%)	11 (4%)	32 57	

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	25	ARG
3	A	28	LEU
3	A	41	TRP
3	A	66	GLU
3	A	82	GLN
3	A	96	SER
3	A	111	LYS
3	A	162	LYS
3	A	251	LEU
3	A	258	ILE

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Mol	Chain	Res	Type
3	A	276	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	72	HIS
3	A	82	GLN
3	A	127	HIS
3	A	141	ASN
3	A	268	ASN
3	A	309	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	Bo	nd leng	$ ag{ths}$	В	ond ang	les
MIOI	туре	Chain	res	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	Z	D	427	3,2	14,19,20	1.14	1 (7%)	16,26,29	3.13	3 (18%)

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	\mathbf{Rings}
2	Z	D	427	3,2	-	1/4/21/22	0/2/2/2

All (1) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	$Ideal(\AA)$
2	D	427	Z	C6-N1	3.57	1.40	1.35

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
2	D	427	Z	C4-N3-C2	11.74	120.21	114.42
2	D	427	Z	C5-C6-N1	-2.75	117.05	119.49
2	D	427	Z	C6-N1-C2	-2.06	117.93	121.20

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	427	Z	O4'-C1'-N1-C6

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

1 ligand is 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).

1/4	[ol	Type	Chain	Res	Link	Bo	nd leng	ths	Bond angles		
101	101	туре	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
	4	SAH	Α	328	_	21,28,28	0.62	0	20,40,40	0.87	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
4	SAH	A	328	_	_	1/7/31/31	0/3/3/3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

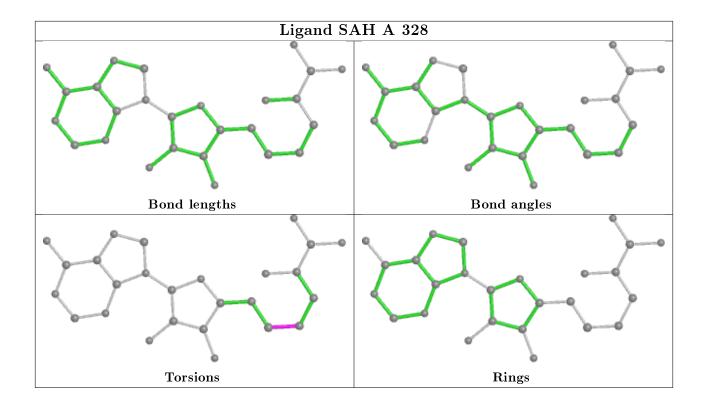
Mol	Chain	Res	Type	Atoms
4	A	328	SAH	CB-CG-SD-C5'

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 then 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)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

