



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 11:28 PM UTC

PDB ID : 1FJX / pdb_00001fjx
Title : STRUCTURE OF TERNARY COMPLEX OF HHAI METHYLTRANSFERASE MUTANT (T250G) IN COMPLEX WITH DNA AND ADOHCY
Authors : Vilkaitis, G.; Dong, A.; Weinhold, E.; Cheng, X.; Klimasauskas, S.
Deposited on : 2000-08-08
Resolution : 2.26 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

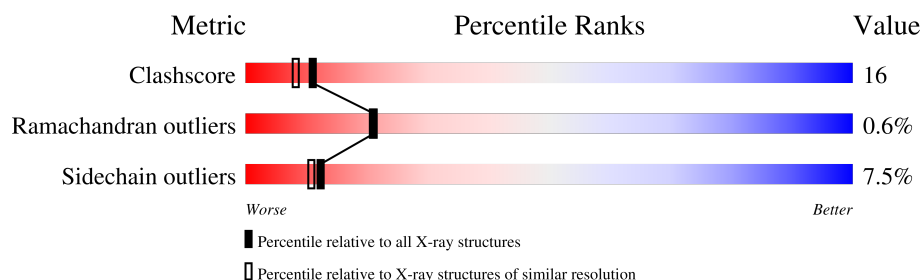
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.26 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))
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)

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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	C	13	
1	D	13	
2	A	327	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4517 atoms, of which 1121 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 DNA (5'-D(*TP*GP*AP*TP*AP*GP*CP*GP*CP*TP*AP*TP*C)-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	C	13	Total	C	H	N	O	P	0	0	0
			290	127	27	47	77	12			
1	D	13	Total	C	H	N	O	P	0	0	0
			290	127	27	47	77	12			

- Molecule 2 is a protein called HHAI DNA METHYLTRANSFERASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	A	327	Total	C	H	N	O	S	0	0	0
			3190	1658	591	443	485	13			

There is a discrepancy between the modelled and reference sequences:

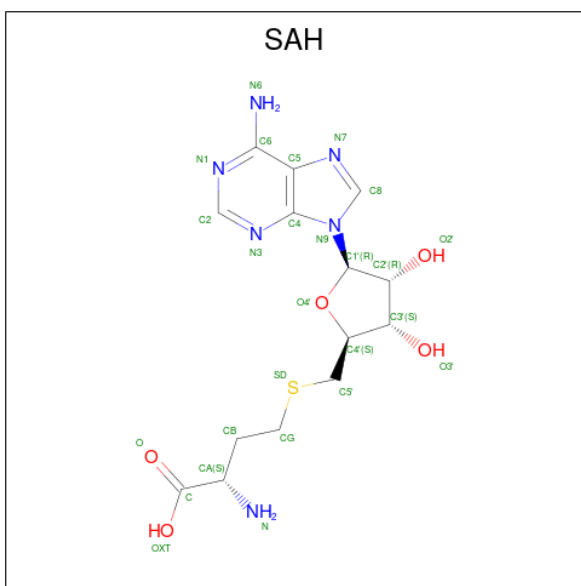
Chain	Residue	Modelled	Actual	Comment	Reference
A	250	GLY	THR	engineered mutation	UNP P05102

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is S-ADENOSYL-L-HOMOCYSTEINE (CCD ID: SAH) (formula: $C_{14}H_{20}N_6O_5S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	H	N	O	S	0	0
			32	14	6	6	5	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	17	Total 51	H 34	O 17	0	0
5	D	24	Total 72	H 48	O 24	0	0
5	A	194	Total 582	H 388	O 194	0	0

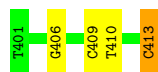
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

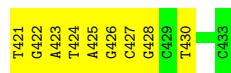
- Molecule 1: DNA (5'-D(*TP*GP*AP*TP*AP*GP*CP*GP*CP*TP*AP*TP*C)-3')

Chain C: 



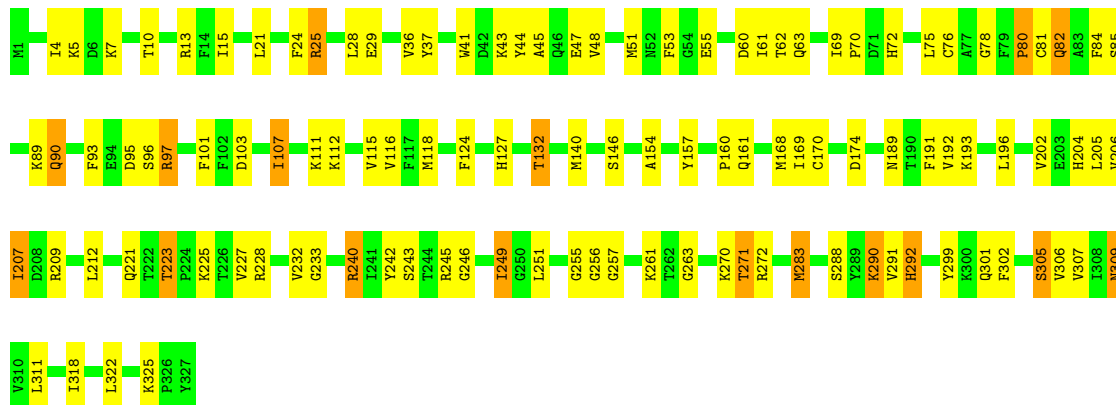
- Molecule 1: DNA (5'-D(*TP*GP*AP*TP*AP*GP*CP*GP*CP*TP*AP*TP*C)-3')

Chain D: 



- Molecule 2: HHAI DNA METHYLTRANSFERASE

Chain A: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	98.46 Å 98.46 Å 323.22 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.26	Depositor
% Data completeness (in resolution range)	82.9 (20.00-2.26)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.207 , 0.283	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4517	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, SAH

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	C	0.46	0/294	0.92	0/452
1	D	0.48	0/294	0.87	0/452
2	A	0.74	1/2654 (0.0%)	1.09	17/3576 (0.5%)
All	All	0.70	1/3242 (0.0%)	1.05	17/4480 (0.4%)

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	C	0	1
2	A	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	283	MET	SD-CE	-7.47	1.60	1.79

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	318	ILE	N-CA-C	-8.17	102.85	110.53
2	A	249	ILE	N-CA-C	-7.66	101.55	110.21
2	A	305	SER	N-CA-C	7.41	120.93	109.60
2	A	10	THR	N-CA-C	6.91	122.78	111.37
2	A	41	TRP	N-CA-C	6.59	121.49	113.38
2	A	204	HIS	N-CA-C	6.33	120.46	112.87
2	A	45	ALA	N-CA-C	-6.23	104.42	111.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	124	PHE	N-CA-C	-5.88	105.00	111.82
2	A	292	HIS	CA-C-N	5.81	125.48	119.56
2	A	292	HIS	C-N-CA	5.81	125.48	119.56
2	A	257	GLY	N-CA-C	-5.65	104.46	111.70
2	A	103	ASP	N-CA-C	-5.59	104.68	112.45
2	A	78	GLY	N-CA-C	-5.47	100.22	113.18
2	A	146	SER	N-CA-C	-5.41	102.89	110.35
2	A	107	ILE	N-CA-C	-5.32	104.31	111.44
2	A	5	LYS	N-CA-C	5.19	118.77	112.23
2	A	85	SER	N-CA-C	5.03	117.73	110.28

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	A	157	TYR	Sidechain
1	C	413	DC	Sidechain

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	C	263	27	149	4	0
1	D	263	27	149	14	0
2	A	2599	591	2577	86	0
3	A	10	0	0	0	0
4	A	26	6	19	1	0
5	A	194	388	0	9	0
5	C	17	34	0	0	0
5	D	24	48	0	4	0
All	All	3396	1121	2894	99	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 16.

All (99) 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:D:427:DC:C6	2:A:81:CYS:SG	2.51	1.03
2:A:227:VAL:HG23	2:A:243:SER:HB3	1.54	0.90
2:A:309:ASN:HD22	2:A:309:ASN:H	1.19	0.86
2:A:202:VAL:HG13	2:A:205:LEU:HD12	1.58	0.85
2:A:69:ILE:O	2:A:111:LYS:HE3	1.77	0.84
2:A:261:LYS:HD3	2:A:299:TYR:HB3	1.59	0.83
1:D:421:DT:H2'	1:D:421:DT:O2	1.82	0.80
1:D:424:DT:H2''	1:D:425:DA:C8	2.19	0.77
2:A:25:ARG:HD2	2:A:53:PHE:CE2	2.22	0.75
2:A:44:TYR:O	2:A:48:VAL:HG23	1.87	0.73
2:A:82:GLN:H	2:A:82:GLN:HE21	1.36	0.71
1:D:427:DC:C5	2:A:81:CYS:SG	2.81	0.70
2:A:15:ILE:HG22	2:A:37:TYR:HB3	1.74	0.69
2:A:43:LYS:O	2:A:47:GLU:HG3	1.92	0.69
2:A:202:VAL:HG11	2:A:272:ARG:NH1	2.08	0.69
2:A:202:VAL:HG11	2:A:272:ARG:CZ	2.23	0.67
2:A:209:ARG:HD3	2:A:232:VAL:HG21	1.78	0.66
2:A:227:VAL:CG2	2:A:243:SER:HB3	2.25	0.65
2:A:101:PHE:HZ	2:A:140:MET:CE	2.09	0.65
2:A:309:ASN:HD22	2:A:309:ASN:N	1.94	0.65
2:A:95:ASP:OD1	2:A:97:ARG:HD3	1.97	0.65
1:C:409:DC:H2''	1:C:410:DT:H5'	1.79	0.64
1:D:426:DG:H8	5:D:454:HOH:O	1.81	0.63
2:A:80:PRO:HB3	2:A:82:GLN:HE22	1.64	0.62
2:A:82:GLN:HE21	2:A:82:GLN:N	1.98	0.61
2:A:93:PHE:CZ	2:A:132:THR:HG22	2.36	0.61
2:A:202:VAL:HG12	2:A:202:VAL:O	1.98	0.61
2:A:223:THR:HG21	2:A:225:LYS:HE2	1.83	0.61
2:A:82:GLN:H	2:A:82:GLN:NE2	2.00	0.58
2:A:207:ILE:O	2:A:271:THR:HG22	2.04	0.57
2:A:140:MET:HG2	2:A:170:CYS:SG	2.46	0.56
1:D:424:DT:H2''	1:D:425:DA:H8	1.69	0.56
2:A:89:LYS:O	2:A:90:GLN:HB2	2.05	0.55
2:A:13:ARG:HE	2:A:36:VAL:HG11	1.70	0.55
2:A:90:GLN:O	2:A:127:HIS:HE1	1.91	0.53
2:A:309:ASN:H	2:A:309:ASN:ND2	1.96	0.53
1:C:409:DC:H2''	1:C:410:DT:C5'	2.39	0.52
2:A:160:PRO:HB3	2:A:191:PHE:HA	1.90	0.52
2:A:306:VAL:HG23	5:A:352:HOH:O	2.09	0.52
2:A:193:LYS:HA	2:A:196:LEU:HD12	1.91	0.52
2:A:101:PHE:CZ	2:A:140:MET:HE1	2.45	0.52
2:A:270:LYS:HB2	5:A:480:HOH:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:425:DA:H2''	5:D:454:HOH:O	2.11	0.50
2:A:24:PHE:CD2	2:A:76:CYS:HB3	2.47	0.49
2:A:48:VAL:HG21	2:A:301:GLN:OE1	2.13	0.49
2:A:228:ARG:NH2	2:A:240:ARG:HD2	2.28	0.49
2:A:48:VAL:HG22	2:A:292:HIS:CD2	2.48	0.49
2:A:80:PRO:HA	5:A:472:HOH:O	2.12	0.49
2:A:263:GLY:HA2	5:A:433:HOH:O	2.12	0.49
2:A:101:PHE:HZ	2:A:140:MET:HE1	1.75	0.49
2:A:270:LYS:HD2	5:A:480:HOH:O	2.12	0.48
2:A:61:ILE:HD11	2:A:107:ILE:HD11	1.94	0.48
2:A:115:VAL:HG12	2:A:116:VAL:N	2.28	0.48
2:A:242:TYR:CZ	2:A:249:ILE:HG21	2.49	0.48
2:A:261:LYS:HD3	2:A:299:TYR:CB	2.39	0.48
1:D:426:DG:N7	2:A:240:ARG:NH2	2.62	0.48
2:A:15:ILE:HD11	2:A:75:LEU:HG	1.97	0.47
2:A:325:LYS:HE2	5:A:549:HOH:O	2.14	0.47
1:D:421:DT:H3'	1:D:422:DG:C8	2.50	0.47
2:A:84:PHE:CE2	2:A:132:THR:HG21	2.50	0.47
2:A:189:ASN:HB2	5:A:560:HOH:O	2.15	0.47
2:A:25:ARG:HH22	2:A:55:GLU:CD	2.23	0.46
2:A:48:VAL:HG22	2:A:292:HIS:HD2	1.79	0.46
2:A:168:MET:C	2:A:169:ILE:HD13	2.40	0.46
2:A:306:VAL:HG22	2:A:307:VAL:N	2.31	0.46
1:C:406:DG:O6	2:A:256:GLY:HA3	2.15	0.46
1:D:430:DT:H5'	5:D:444:HOH:O	2.14	0.46
2:A:193:LYS:HA	2:A:196:LEU:CD1	2.46	0.46
2:A:21:LEU:HD21	2:A:301:GLN:HB3	1.97	0.46
2:A:232:VAL:O	2:A:233:GLY:C	2.59	0.45
2:A:227:VAL:HG12	5:A:451:HOH:O	2.15	0.45
2:A:25:ARG:HD3	2:A:29:GLU:OE2	2.17	0.45
2:A:4:ILE:O	2:A:7:LYS:HE2	2.17	0.44
2:A:140:MET:CA	2:A:140:MET:HE2	2.48	0.43
2:A:192:VAL:HG23	2:A:246:GLY:O	2.18	0.43
2:A:283:MET:HE2	2:A:302:PHE:O	2.18	0.43
2:A:206:VAL:HA	2:A:272:ARG:HB3	1.99	0.43
2:A:81:CYS:O	2:A:82:GLN:C	2.61	0.43
2:A:306:VAL:HG21	2:A:311:LEU:CD1	2.48	0.43
2:A:101:PHE:CZ	2:A:140:MET:CE	2.96	0.43
2:A:228:ARG:CZ	2:A:240:ARG:HD2	2.49	0.43
2:A:51:MET:HE3	2:A:51:MET:HB2	1.93	0.43
2:A:51:MET:HG2	2:A:290:LYS:HD3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:428:DG:N7	2:A:255:GLY:HA2	2.35	0.42
2:A:70:PRO:O	2:A:72:HIS:HD2	2.03	0.42
2:A:25:ARG:HD2	2:A:53:PHE:CD2	2.54	0.42
2:A:101:PHE:HE1	2:A:118:MET:SD	2.42	0.42
2:A:62:THR:HG22	2:A:63:GLN:HE21	1.85	0.41
2:A:111:LYS:O	2:A:112:LYS:C	2.63	0.41
2:A:154:ALA:HB3	2:A:161:GLN:O	2.20	0.41
2:A:69:ILE:HA	2:A:70:PRO:HD3	1.90	0.41
1:D:426:DG:OP1	5:D:437:HOH:O	2.22	0.41
1:D:423:DA:H2''	1:D:424:DT:OP2	2.20	0.41
2:A:202:VAL:CG1	2:A:205:LEU:HD12	2.39	0.41
4:A:328:SAH:H5'1	5:A:376:HOH:O	2.21	0.41
1:C:413:DC:C4	1:D:421:DT:H72	2.55	0.40
2:A:322:LEU:O	2:A:325:LYS:HE3	2.21	0.40
2:A:309:ASN:N	2:A:309:ASN:ND2	2.61	0.40
2:A:240:ARG:HG3	2:A:242:TYR:OH	2.22	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
2	A	325/327 (99%)	300 (92%)	23 (7%)	2 (1%)	21 21

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	90	GLN
2	A	60	ASP

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	
2	A	281/282 (100%)	260 (92%)	21 (8%)	12	11

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	25	ARG
2	A	28	LEU
2	A	80	PRO
2	A	82	GLN
2	A	96	SER
2	A	97	ARG
2	A	132	THR
2	A	174	ASP
2	A	207	ILE
2	A	212	LEU
2	A	221	GLN
2	A	223	THR
2	A	240	ARG
2	A	245	ARG
2	A	251	LEU
2	A	271	THR
2	A	288	SER
2	A	290	LYS
2	A	291	VAL
2	A	305	SER
2	A	309	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	63	GLN
2	A	72	HIS
2	A	82	GLN
2	A	127	HIS

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Mol	Chain	Res	Type
2	A	216	ASN
2	A	221	GLN
2	A	268	ASN
2	A	297	GLN
2	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](#)

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](#)

3 ligands 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	SO4	A	600	-	4,4,4	0.50	0	6,6,6	0.67	0
3	SO4	A	601	-	4,4,4	0.44	0	6,6,6	0.51	0
4	SAH	A	328	-	27,28,28	0.58	1 (3%)	36,40,40	1.09	2 (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
4	SAH	A	328	-	-	4/15/31/31	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	328	SAH	O-C	2.38	1.29	1.22

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	328	SAH	C5'-SD-CG	4.31	115.06	102.26
4	A	328	SAH	C4'-C5'-SD	-3.94	99.71	113.78

There are no chirality outliers.

All (4) torsion outliers are listed below:

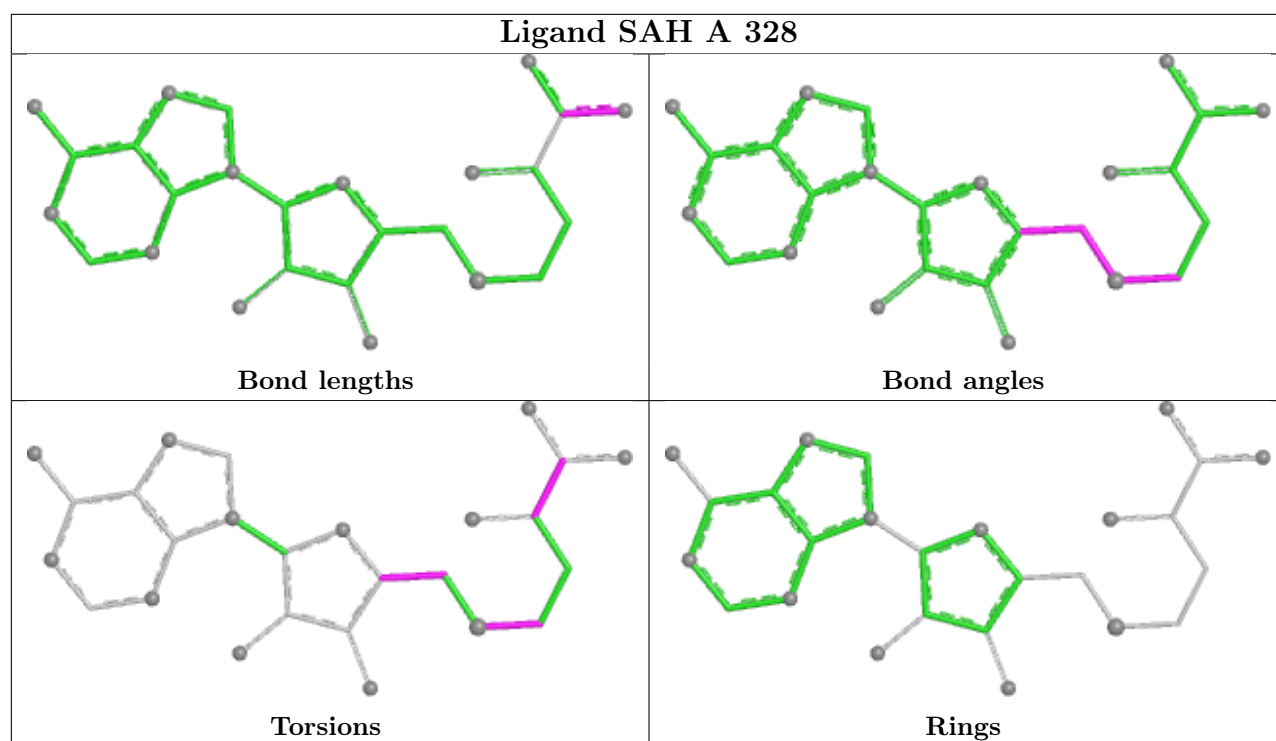
Mol	Chain	Res	Type	Atoms
4	A	328	SAH	OXT-C-CA-CB
4	A	328	SAH	O4'-C4'-C5'-SD
4	A	328	SAH	CB-CG-SD-C5'
4	A	328	SAH	O-C-CA-CB

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	328	SAH	1	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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.