



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:26 AM UTC

PDB ID : 7EEW / pdb\_00007eew  
Title : Crystal structure of the intact MTase from *Vibrio vulnificus* YJ016 in complex with the DNA-mimicking Ocr protein and the S-adenosyl-L-homocysteine (SAH)  
Authors : Seo, P.W.; Park, S.Y.; Kim, J.S.  
Deposited on : 2021-03-19  
Resolution : 2.90 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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

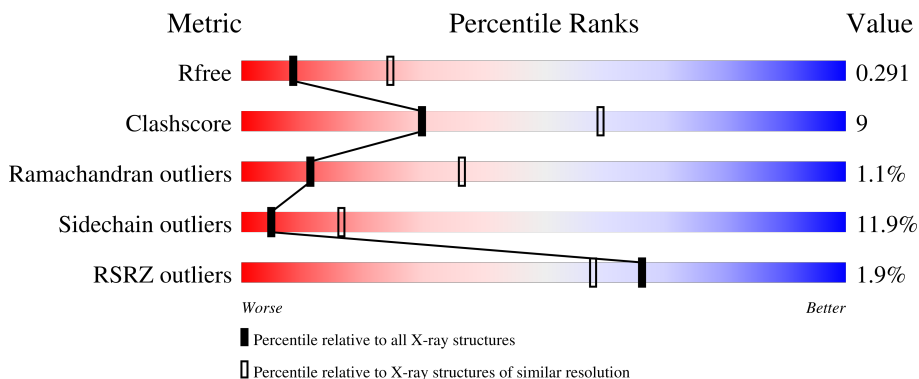
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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  $\geq 3$ , 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	638	
2	B	117	



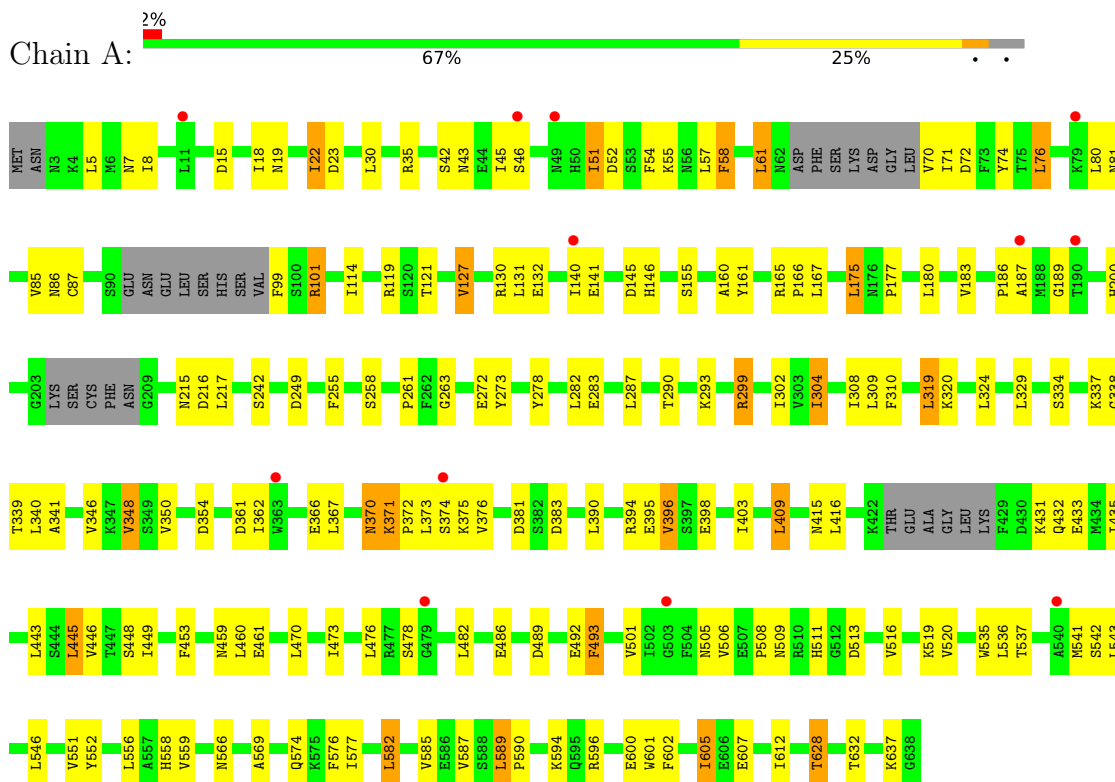
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	54	Total 54	O 54	0	0
4	B	6	Total 6	O 6	0	0

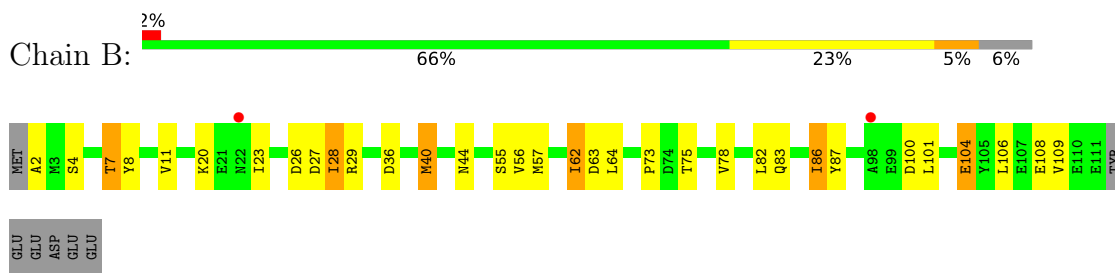
### 3 Residue-property plots [i](#)

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 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: Type I restriction-modification system methyltransferase subunit



- Molecule 2: Overcome classical restriction gp0.3



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	133.28Å 133.28Å 236.81Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	14.96 – 2.90 14.96 – 2.90	Depositor EDS
% Data completeness (in resolution range)	93.9 (14.96-2.90) 93.1 (14.96-2.90)	Depositor EDS
$R_{merge}$	0.19	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.95 (at 2.91Å)	Xtrriage
Refinement program	PHENIX 1.12_2829	Depositor
R, $R_{free}$	0.272 , 0.293 0.272 , 0.291	Depositor DCC
$R_{free}$ test set	1397 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	58.4	Xtrriage
Anisotropy	0.551	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.22 , 0.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.36$ , $\langle L^2 \rangle = 0.19$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	5829	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	63.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.33% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	A	0.10	0/4929	0.29	0/6654
2	B	0.09	0/920	0.27	0/1249
All	All	0.10	0/5849	0.29	0/7903

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

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	4840	0	4861	81	1
2	B	903	0	825	18	0
3	A	26	0	19	6	0
4	A	54	0	0	1	0
4	B	6	0	0	0	0
All	All	5829	0	5705	98	1

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 (98) 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:309:LEU:HB2	1:A:416:LEU:HB2	1.67	0.76
1:A:175:LEU:O	1:A:299:ARG:NH2	2.19	0.75
1:A:310:PHE:O	1:A:320:LYS:NZ	2.23	0.71
1:A:51:ILE:HD11	1:A:86:ASN:HD22	1.60	0.66
1:A:340:LEU:HD13	1:A:348:VAL:HG21	1.80	0.64
1:A:215:ASN:HD21	1:A:242:SER:HB3	1.63	0.64
1:A:155:SER:O	4:A:801:HOH:O	2.15	0.63
1:A:216:ASP:OD2	3:A:700:SAH:O3'	2.18	0.61
2:B:26:ASP:O	2:B:28:ILE:N	2.36	0.59
1:A:19:ASN:HD21	1:A:22:ILE:HG13	1.67	0.59
1:A:189:GLY:HA3	3:A:700:SAH:HA	1.85	0.58
1:A:334:SER:HB3	1:A:366:GLU:HA	1.85	0.58
1:A:71:ILE:HG13	1:A:121:THR:HG22	1.84	0.58
1:A:470:LEU:HD12	1:A:585:VAL:HB	1.86	0.57
1:A:371:LYS:HD3	1:A:372:PRO:HD2	1.86	0.57
2:B:23:ILE:HA	2:B:28:ILE:HD11	1.86	0.57
1:A:35:ARG:HD2	1:A:131:LEU:HB3	1.88	0.55
1:A:54:PHE:HB2	1:A:57:LEU:HD12	1.89	0.55
1:A:186:PRO:HD2	1:A:258:SER:HB2	1.88	0.54
1:A:263:GLY:HA2	1:A:282:LEU:H	1.72	0.54
2:B:56:VAL:HG11	2:B:86:ILE:HD12	1.89	0.53
1:A:509:ASN:HB2	1:A:535:TRP:CD1	2.44	0.53
2:B:57:MET:HG3	2:B:82:LEU:HD13	1.90	0.53
1:A:431:LYS:HG2	1:A:432:GLN:H	1.74	0.53
1:A:160:ALA:O	3:A:700:SAH:H5'1	2.09	0.52
1:A:361:ASP:O	1:A:394:ARG:NH2	2.41	0.52
2:B:73:PRO:HB2	2:B:75:THR:HG22	1.91	0.52
2:B:100:ASP:O	2:B:104:GLU:HG3	2.09	0.52
1:A:290:THR:HG21	1:A:302:ILE:HG12	1.92	0.52
1:A:520:VAL:HG23	1:A:576:PHE:HD1	1.75	0.52
1:A:7:ASN:HB3	1:A:130:ARG:HH21	1.76	0.51
1:A:177:PRO:HD2	1:A:200:HIS:CD2	2.46	0.51
2:B:29:ARG:HH22	2:B:109:VAL:HG11	1.76	0.51
1:A:58:PHE:HA	1:A:61:LEU:HD12	1.93	0.50
1:A:324:LEU:HD23	1:A:409:LEU:HD23	1.94	0.50
2:B:11:VAL:HG11	2:B:86:ILE:HD11	1.94	0.50
1:A:71:ILE:HD11	1:A:121:THR:HA	1.94	0.49
1:A:283:GLU:HG3	1:A:308:ILE:HG12	1.93	0.49
1:A:478:SER:HA	1:A:541:MET:HA	1.94	0.49
1:A:556:LEU:HA	1:A:559:VAL:HG12	1.93	0.49
1:A:551:VAL:HG12	1:A:589:LEU:HG	1.95	0.49
1:A:448:SER:HB3	1:A:612:ILE:HD11	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:381:ASP:OD1	1:A:381:ASP:N	2.41	0.48
1:A:299:ARG:HA	1:A:354:ASP:HA	1.95	0.48
1:A:304:ILE:HB	1:A:308:ILE:HD11	1.95	0.48
1:A:367:LEU:HD11	1:A:383:ASP:HB3	1.95	0.48
1:A:374:SER:OG	1:A:375:LYS:N	2.46	0.48
1:A:187:ALA:O	3:A:700:SAH:N	2.46	0.48
1:A:5:LEU:HD21	1:A:140:ILE:HD11	1.96	0.47
1:A:390:LEU:HD13	1:A:396:VAL:HG21	1.95	0.47
2:B:7:THR:OG1	2:B:8:TYR:N	2.47	0.47
1:A:337:LYS:HA	1:A:338:GLY:HA2	1.53	0.47
1:A:30:LEU:HD22	1:A:114:ILE:HG22	1.97	0.47
1:A:35:ARG:HE	1:A:132:GLU:HA	1.80	0.46
1:A:145:ASP:OD1	1:A:146:HIS:N	2.49	0.46
1:A:161:TYR:HA	3:A:700:SAH:H5'1	1.97	0.46
1:A:19:ASN:O	1:A:23:ASP:HB2	2.15	0.46
1:A:278:TYR:HE2	1:A:319:LEU:HD12	1.80	0.46
1:A:601:TRP:O	1:A:605:ILE:HG22	2.15	0.46
1:A:45:ILE:HG21	1:A:87:CYS:HA	1.99	0.45
2:B:20:LYS:HG2	2:B:101:LEU:HD21	1.98	0.45
1:A:8:ILE:HG23	1:A:127:VAL:HG23	1.98	0.45
1:A:536:LEU:HD22	1:A:542:SER:HB3	1.98	0.45
1:A:446:VAL:HA	1:A:449:ILE:HG12	2.00	0.44
1:A:101:ARG:N	1:A:101:ARG:HD2	2.32	0.44
1:A:519:LYS:NZ	1:A:541:MET:SD	2.89	0.43
1:A:58:PHE:HB2	1:A:76:LEU:HB2	2.00	0.43
1:A:370:ASN:N	1:A:370:ASN:OD1	2.51	0.43
2:B:104:GLU:HG3	2:B:104:GLU:H	1.67	0.43
1:A:263:GLY:CA	1:A:282:LEU:H	2.32	0.43
1:A:505:ASN:OD1	1:A:506:VAL:N	2.51	0.43
1:A:175:LEU:HD12	1:A:255:PHE:CE2	2.53	0.43
1:A:183:VAL:HA	1:A:255:PHE:HB2	2.01	0.43
1:A:582:LEU:HD23	1:A:582:LEU:HA	1.87	0.42
1:A:558:HIS:CD2	1:A:602:PHE:HB3	2.55	0.42
1:A:569:ALA:HB2	1:A:577:ILE:HG13	2.01	0.42
1:A:19:ASN:ND2	1:A:22:ILE:HG13	2.33	0.42
1:A:628:THR:O	1:A:632:THR:HG23	2.20	0.42
1:A:81:ASN:O	1:A:85:VAL:HG23	2.19	0.42
2:B:57:MET:HE2	2:B:62:ILE:HD13	2.02	0.42
1:A:493:PHE:HB3	1:A:505:ASN:O	2.19	0.42
1:A:493:PHE:CE1	1:A:508:PRO:HB3	2.55	0.42
1:A:42:SER:O	1:A:46:SER:HB2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:83:GLN:HA	2:B:86:ILE:HG22	2.02	0.41
1:A:459:ASN:O	1:A:459:ASN:ND2	2.53	0.41
1:A:552:TYR:CE2	1:A:596:ARG:HG2	2.56	0.41
1:A:15:ASP:OD2	1:A:119:ARG:HG3	2.19	0.41
1:A:165:ARG:N	1:A:166:PRO:HD2	2.36	0.41
1:A:476:LEU:HD21	1:A:543:LEU:HD12	2.02	0.41
1:A:574:GLN:HG3	2:B:87:TYR:CD2	2.55	0.41
2:B:40:MET:O	2:B:44:ASN:ND2	2.53	0.41
1:A:445:LEU:HD13	1:A:449:ILE:HD13	2.02	0.41
2:B:29:ARG:HG3	2:B:106:LEU:HD11	2.03	0.41
1:A:15:ASP:CG	1:A:119:ARG:HG3	2.46	0.40
2:B:4:SER:HB3	2:B:55:SER:OG	2.21	0.40
1:A:161:TYR:C	3:A:700:SAH:HG2	2.46	0.40
1:A:329:LEU:O	1:A:362:ILE:HG12	2.22	0.40
2:B:2:ALA:O	2:B:4:SER:N	2.53	0.40

All (1) 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:273:TYR:OH	1:A:489:ASP:OD2[6_554]	2.15	0.05

## 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	600/638 (94%)	551 (92%)	43 (7%)	6 (1%)	12 39
2	B	108/117 (92%)	105 (97%)	1 (1%)	2 (2%)	6 23
All	All	708/755 (94%)	656 (93%)	44 (6%)	8 (1%)	11 36

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	473	ILE
1	A	590	PRO
2	B	27	ASP
1	A	341	ALA
1	A	398	GLU
1	A	261	PRO
2	B	28	ILE
1	A	18	ILE

### 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	548/573 (96%)	481 (88%)	67 (12%)	5	16
2	B	98/105 (93%)	88 (90%)	10 (10%)	7	23
All	All	646/678 (95%)	569 (88%)	77 (12%)	5	16

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ILE
1	A	43	ASN
1	A	51	ILE
1	A	52	ASP
1	A	55	LYS
1	A	58	PHE
1	A	61	LEU
1	A	70	VAL
1	A	72	ASP
1	A	74	TYR
1	A	76	LEU
1	A	80	LEU
1	A	99	PHE
1	A	101	ARG
1	A	127	VAL
1	A	141	GLU

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	167	LEU
1	A	175	LEU
1	A	180	LEU
1	A	217	LEU
1	A	249	ASP
1	A	272	GLU
1	A	287	LEU
1	A	293	LYS
1	A	299	ARG
1	A	304	ILE
1	A	319	LEU
1	A	339	THR
1	A	346	VAL
1	A	348	VAL
1	A	350	VAL
1	A	370	ASN
1	A	371	LYS
1	A	373	LEU
1	A	376	VAL
1	A	395	GLU
1	A	396	VAL
1	A	403	ILE
1	A	409	LEU
1	A	415	ASN
1	A	433	GLU
1	A	435	ILE
1	A	443	LEU
1	A	445	LEU
1	A	453	PHE
1	A	460	LEU
1	A	461	GLU
1	A	482	LEU
1	A	486	GLU
1	A	492	GLU
1	A	493	PHE
1	A	501	VAL
1	A	511	HIS
1	A	513	ASP
1	A	516	VAL
1	A	537	THR
1	A	546	LEU
1	A	566	ASN

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Mol	Chain	Res	Type
1	A	582	LEU
1	A	587	VAL
1	A	589	LEU
1	A	594	LYS
1	A	600	GLU
1	A	605	ILE
1	A	607	GLU
1	A	628	THR
1	A	637	LYS
2	B	7	THR
2	B	36	ASP
2	B	40	MET
2	B	62	ILE
2	B	63	ASP
2	B	64	LEU
2	B	78	VAL
2	B	86	ILE
2	B	104	GLU
2	B	108	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	176	ASN
1	A	232	ASN
1	A	322	GLN
1	A	326	GLN
1	A	399	ASN
1	A	451	ASN
2	B	103	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](#)

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).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	SAH	A	700	-	27,28,28	1.67	5 (18%)	36,40,40	2.43	13 (36%)

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	SAH	A	700	-	-	7/15/31/31	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	700	SAH	C5-N7	-3.74	1.32	1.39
3	A	700	SAH	C4-N9	-3.15	1.31	1.37
3	A	700	SAH	C5-C4	3.11	1.44	1.39
3	A	700	SAH	C8-N9	-2.65	1.33	1.37
3	A	700	SAH	OXT-C	-2.65	1.22	1.30

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	700	SAH	O4'-C1'-N9	7.52	122.53	108.09
3	A	700	SAH	C5-C4-N3	-5.08	119.73	126.72
3	A	700	SAH	C4-C5-N7	-4.05	105.96	110.58
3	A	700	SAH	C2-N3-C4	3.41	120.17	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	700	SAH	N3-C2-N1	-3.26	123.65	128.58
3	A	700	SAH	N3-C4-N9	3.05	132.36	127.17
3	A	700	SAH	C6-C5-N7	3.02	137.91	132.09
3	A	700	SAH	O4'-C1'-C2'	-2.87	100.48	106.62
3	A	700	SAH	C5'-C4'-C3'	-2.83	107.99	115.06
3	A	700	SAH	O3'-C3'-C4'	-2.58	103.67	111.08
3	A	700	SAH	C5-N7-C8	2.56	107.48	103.45
3	A	700	SAH	C4-N9-C1'	-2.24	121.39	126.63
3	A	700	SAH	C2-N1-C6	2.08	122.14	118.73

There are no chirality outliers.

All (7) torsion outliers are listed below:

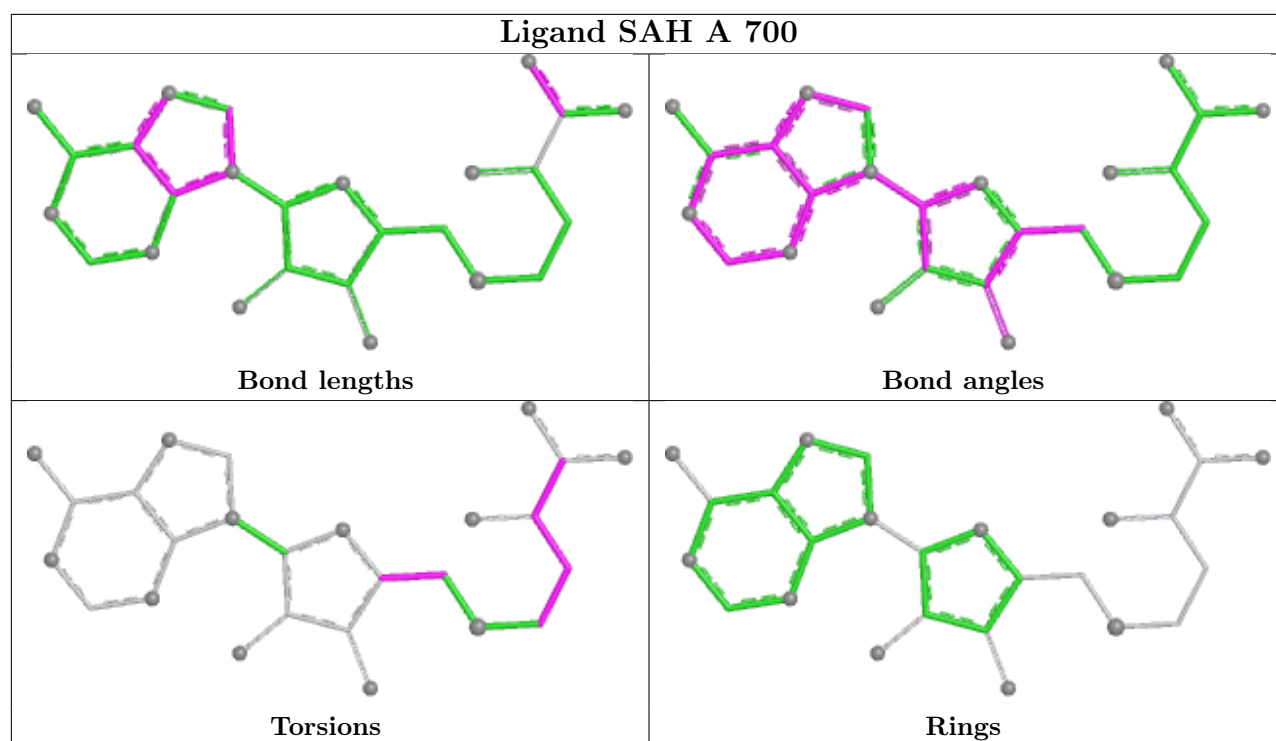
Mol	Chain	Res	Type	Atoms
3	A	700	SAH	O4'-C4'-C5'-SD
3	A	700	SAH	C3'-C4'-C5'-SD
3	A	700	SAH	CA-CB-CG-SD
3	A	700	SAH	C-CA-CB-CG
3	A	700	SAH	O-C-CA-CB
3	A	700	SAH	OXT-C-CA-CB
3	A	700	SAH	N-CA-CB-CG

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	700	SAH	6	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, 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	610/638 (95%)	0.01	12 (1%) 65 56	27, 61, 100, 144	0
2	B	110/117 (94%)	-0.04	2 (1%) 67 59	25, 48, 117, 148	0
All	All	720/755 (95%)	0.00	14 (1%) 66 58	25, 60, 101, 148	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	11	LEU	4.4
1	A	140	ILE	4.4
2	B	98	ALA	4.0
1	A	479	GLY	4.0
1	A	540	ALA	2.7
1	A	363	TRP	2.5
1	A	46	SER	2.4
1	A	49	ASN	2.3
2	B	22	ASN	2.3
1	A	190	THR	2.3
1	A	374	SER	2.2
1	A	79	LYS	2.2
1	A	503	GLY	2.1
1	A	187	ALA	2.1

### 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](#)

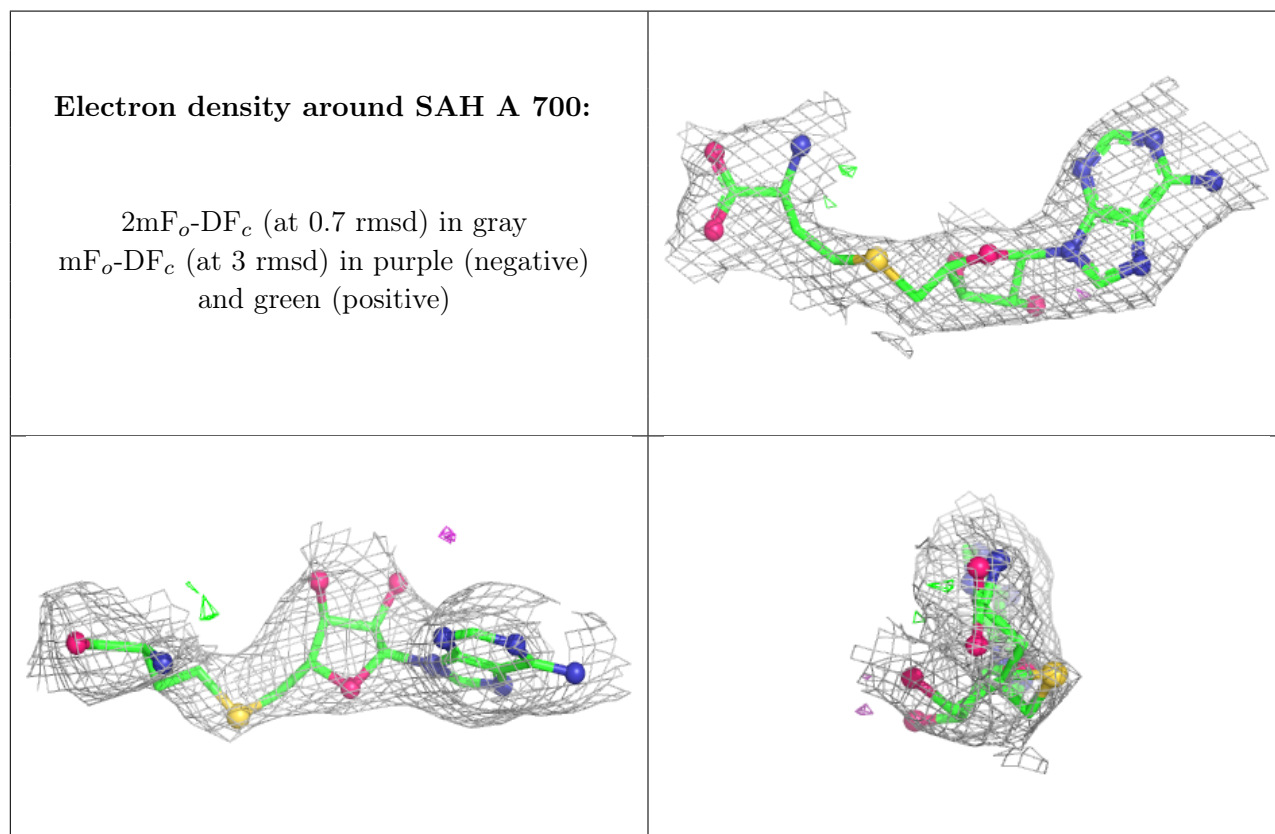
There are no oligosaccharides in this entry.

## 6.4 Ligands [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(Å <sup>2</sup> )	Q<0.9
3	SAH	A	700	26/26	0.79	0.09	43,51,58,65	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.