



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2023 – 11:16 AM EDT

PDB ID : 2PX5
Title : Crystal structure of the Murray Valley Encephalitis Virus NS5 2'-O Methyltransferase domain in complex with SAH (Orthorhombic crystal form)
Authors : Assenberg, R.; Ren, J.; Verma, A.; Walter, T.S.; Alderton, D.; Hurrelbrink, R.J.; Fuller, S.D.; Owens, R.J.; Stuart, D.I.; Grimes, J.M.; Oxford Protein Production Facility (OPPF)
Deposited on : 2007-05-14
Resolution : 2.30 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](#) ①) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.35
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) : 2.35

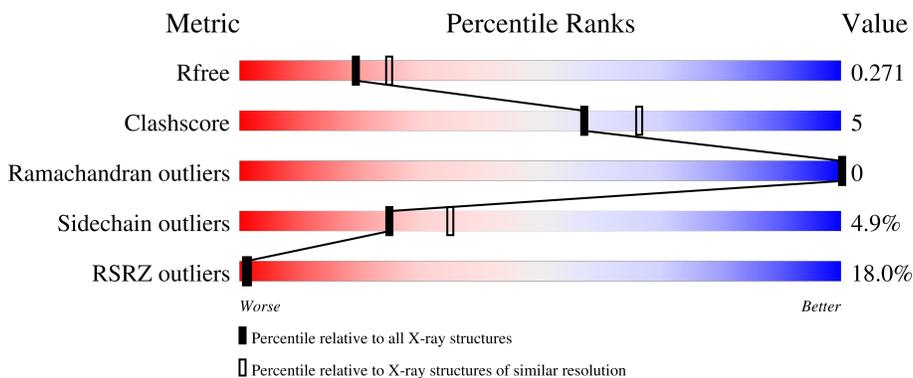
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	269	
1	B	269	

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	SO4	B	910	-	-	-	X

2 Entry composition [i](#)

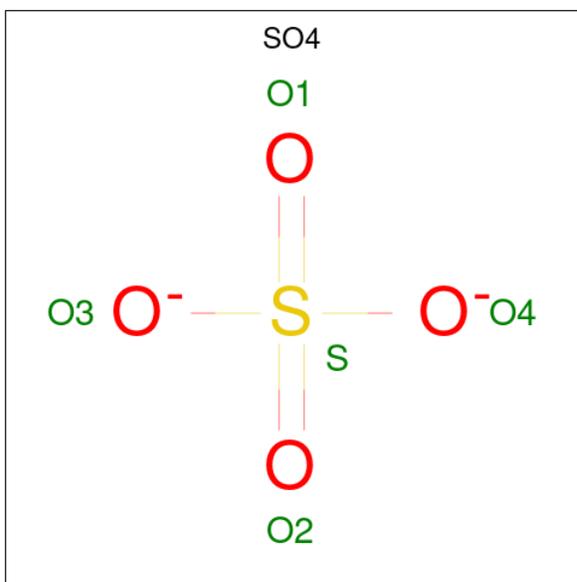
There are 4 unique types of molecules in this entry. The entry contains 4613 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Genome polyprotein [Contains: Capsid protein C (Core protein); Envelope protein M (Matrix protein); Major envelope protein E; Non-structural protein 1 (NS1); Non-structural protein 2A (NS2A); Flavivirin protease NS2B regulatory subunit; Flavivirin protease NS3 catalytic subunit; Non-structural protein 4A (NS4A); Non-structural protein 4B (NS4B); RNA-directed RNA polymerase (EC 2.7.7.48) (NS5)].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	265	Total 2095	C 1317	N 381	O 383	S 14	0	0	0
1	B	263	Total 2086	C 1312	N 379	O 381	S 14	0	0	0

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



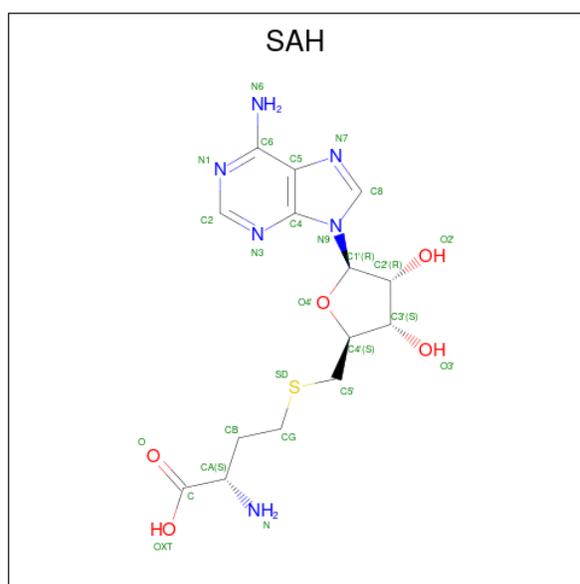
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	Total 5	O 4	S 1	0	0
2	A	1	Total 5	O 4	S 1	0	0

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

- Molecule 3 is S-ADENOSYL-L-HOMOCYSTEINE (three-letter code: SAH) (formula: $C_{14}H_{20}N_6O_5S$).



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

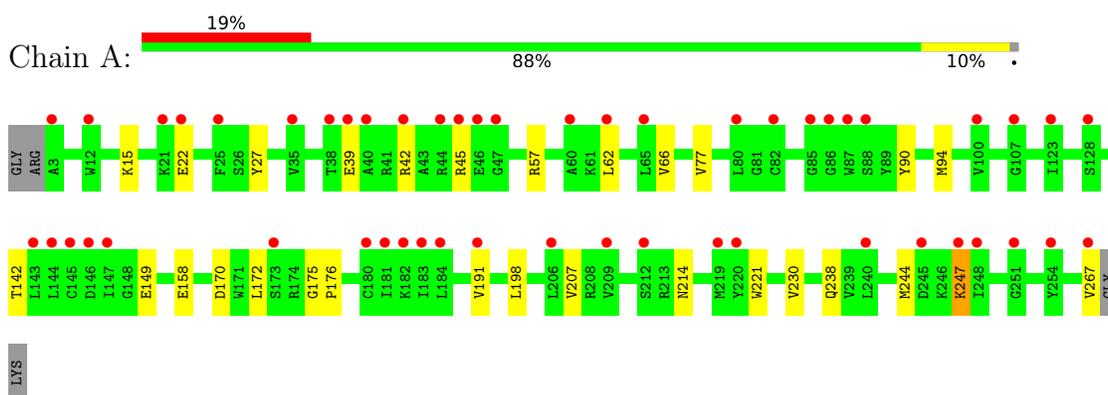
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	175	Total 175	O 175	0	0
4	B	160	Total 160	O 160	0	0

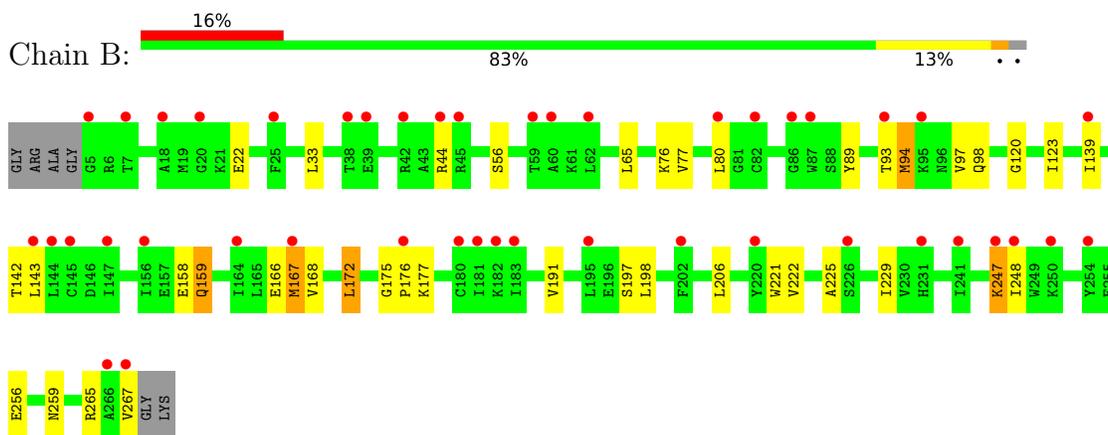
3 Residue-property plots [i](#)

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: Genome polyprotein [Contains: Capsid protein C (Core protein); Envelope protein M (Matrix protein); Major envelope protein E; Non-structural protein 1 (NS1); Non-structural protein 2A (NS2A); Flavivirin protease NS2B regulatory subunit; Flavivirin protease NS3 catalytic subunit; Non-structural protein 4A (NS4A); Non-structural protein 4B (NS4B); RNA-directed RNA polymerase (EC 2.7.7.48) (NS5)]



- Molecule 1: Genome polyprotein [Contains: Capsid protein C (Core protein); Envelope protein M (Matrix protein); Major envelope protein E; Non-structural protein 1 (NS1); Non-structural protein 2A (NS2A); Flavivirin protease NS2B regulatory subunit; Flavivirin protease NS3 catalytic subunit; Non-structural protein 4A (NS4A); Non-structural protein 4B (NS4B); RNA-directed RNA polymerase (EC 2.7.7.48) (NS5)]



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.90Å 89.10Å 101.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.12 – 2.30 28.51 – 2.15	Depositor EDS
% Data completeness (in resolution range)	96.1 (20.12-2.30) 91.3 (28.51-2.15)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.16Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.187 , 0.260 0.201 , 0.271	Depositor DCC
R_{free} test set	1551 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	29.5	Xtrriage
Anisotropy	0.788	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 65.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4613	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.61% 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

5.1 Standard geometry

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

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	0.48	0/2138	0.59	0/2874
1	B	0.47	0/2129	0.59	0/2862
All	All	0.48	0/4267	0.59	0/5736

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	2095	0	2103	16	0
1	B	2086	0	2095	24	0
2	A	25	0	0	0	0
2	B	20	0	0	1	0
3	A	26	0	19	1	0
3	B	26	0	19	1	0
4	A	175	0	0	4	0
4	B	160	0	0	2	0
All	All	4613	0	4236	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 5.

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:158:GLU:OE2	4:A:655:HOH:O	2.06	0.73
3:A:500:SAH:N	4:A:520:HOH:O	2.29	0.65
1:B:247:LYS:C	1:B:248:ILE:HD12	2.19	0.63
1:B:123:ILE:HG22	1:B:267:VAL:HG21	1.79	0.63
1:B:80:LEU:CD1	1:B:167:MET:HE1	2.37	0.54
2:B:906:SO4:O2	4:B:577:HOH:O	2.17	0.53
1:B:259:ASN:ND2	4:B:861:HOH:O	2.42	0.53
1:A:90:TYR:CE2	1:A:94:MET:HE1	2.44	0.52
1:B:247:LYS:O	1:B:248:ILE:HD12	2.11	0.51
1:B:56:SER:OG	3:B:500:SAH:OXT	2.25	0.50
1:A:15:LYS:HD3	1:A:244:MET:SD	2.52	0.49
1:B:80:LEU:HD13	1:B:167:MET:CE	2.42	0.49
1:B:80:LEU:HD13	1:B:167:MET:HE1	1.95	0.49
1:A:57:ARG:HD2	4:A:760:HOH:O	2.14	0.48
1:B:143:LEU:HD23	1:B:172:LEU:HD13	1.97	0.46
1:B:158:GLU:HG3	1:B:191:VAL:HG23	1.98	0.46
1:B:65:LEU:HD22	1:B:222:VAL:HG11	1.98	0.46
1:A:175:GLY:N	1:A:176:PRO:CD	2.80	0.45
1:B:175:GLY:N	1:B:176:PRO:CD	2.78	0.45
1:B:77:VAL:HG22	1:B:142:THR:HB	2.00	0.44
1:A:175:GLY:N	1:A:176:PRO:HD2	2.32	0.44
1:A:238:GLN:NE2	1:A:238:GLN:HA	2.32	0.44
1:B:94:MET:HB2	1:B:97:VAL:HG23	2.01	0.43
1:A:90:TYR:CE2	1:A:94:MET:CE	3.02	0.43
1:B:206:LEU:HD22	1:B:229:ILE:HG23	2.01	0.43
1:B:222:VAL:HG23	1:B:225:ALA:CB	2.49	0.43
1:A:27:TYR:CZ	1:A:214:ASN:HB3	2.54	0.43
1:A:39:GLU:O	1:A:42:ARG:HG3	2.19	0.43
1:A:230:VAL:HG23	4:A:751:HOH:O	2.18	0.43
1:B:98:GLN:HE21	1:B:98:GLN:HB2	1.64	0.43
1:B:120:GLY:HA2	1:B:265:ARG:HG2	1.99	0.42
1:B:80:LEU:CD1	1:B:167:MET:CE	2.97	0.42
1:A:198:LEU:HB3	1:A:221:TRP:CH2	2.54	0.42
1:B:198:LEU:HB3	1:B:221:TRP:CH2	2.54	0.42
1:A:247:LYS:N	1:A:247:LYS:HD2	2.35	0.42
1:B:167:MET:HE3	1:B:168:VAL:HA	2.02	0.42
1:B:167:MET:HE3	1:B:167:MET:C	2.40	0.41
1:B:159:GLN:HE21	1:B:159:GLN:HB3	1.67	0.41
1:A:77:VAL:HG22	1:A:142:THR:HB	2.03	0.41
1:B:89:TYR:O	1:B:93:THR:HG23	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:GLU:HG3	1:A:191:VAL:HG23	2.03	0.41
1:A:62:LEU:O	1:A:66:VAL:HG23	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	263/269 (98%)	261 (99%)	2 (1%)	0	100	100
1	B	261/269 (97%)	260 (100%)	1 (0%)	0	100	100
All	All	524/538 (97%)	521 (99%)	3 (1%)	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	223/225 (99%)	215 (96%)	8 (4%)	35	49
1	B	223/225 (99%)	209 (94%)	14 (6%)	18	24
All	All	446/450 (99%)	424 (95%)	22 (5%)	25	35

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	45	ARG
1	A	149	GLU
1	A	170	ASP
1	A	172	LEU
1	A	207	VAL
1	A	247	LYS
1	A	267	VAL
1	B	22	GLU
1	B	33	LEU
1	B	44	ARG
1	B	76	LYS
1	B	94	MET
1	B	139	ILE
1	B	159	GLN
1	B	166	GLU
1	B	167	MET
1	B	172	LEU
1	B	177	LYS
1	B	197	SER
1	B	247	LYS
1	B	256	GLU

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	11	GLN
1	A	238	GLN
1	B	98	GLN
1	B	159	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

11 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
2	SO4	B	910	-	4,4,4	0.14	0	6,6,6	0.17	0
2	SO4	A	911	-	4,4,4	0.17	0	6,6,6	0.39	0
2	SO4	B	908	-	4,4,4	0.15	0	6,6,6	0.30	0
2	SO4	B	907	-	4,4,4	0.22	0	6,6,6	0.36	0
3	SAH	A	500	-	24,28,28	1.24	4 (16%)	25,40,40	1.73	6 (24%)
2	SO4	A	903	-	4,4,4	0.15	0	6,6,6	0.24	0
2	SO4	A	902	-	4,4,4	0.19	0	6,6,6	0.36	0
2	SO4	A	905	-	4,4,4	0.11	0	6,6,6	0.43	0
2	SO4	A	904	-	4,4,4	0.15	0	6,6,6	0.38	0
2	SO4	B	906	-	4,4,4	0.17	0	6,6,6	0.31	0
3	SAH	B	500	-	24,28,28	1.15	1 (4%)	25,40,40	1.76	4 (16%)

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	B	500	-	-	0/11/31/31	0/3/3/3
3	SAH	A	500	-	-	2/11/31/31	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	500	SAH	C2-N3	3.54	1.37	1.32
3	A	500	SAH	C2-N3	3.51	1.37	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	SAH	C2-N1	2.31	1.38	1.33
3	A	500	SAH	OXT-C	-2.13	1.23	1.30
3	A	500	SAH	C2'-C1'	-2.12	1.50	1.53

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	500	SAH	N3-C2-N1	-5.81	119.61	128.68
3	A	500	SAH	N3-C2-N1	-5.58	119.95	128.68
3	B	500	SAH	OXT-C-O	-3.11	117.03	124.09
3	B	500	SAH	OXT-C-CA	2.95	123.44	113.38
3	A	500	SAH	C5'-SD-CG	-2.73	94.09	102.27
3	A	500	SAH	C4-C5-N7	-2.64	106.65	109.40
3	B	500	SAH	C5'-SD-CG	-2.61	94.42	102.27
3	A	500	SAH	OXT-C-O	-2.40	118.64	124.09
3	A	500	SAH	C1'-N9-C4	-2.24	122.70	126.64
3	A	500	SAH	OXT-C-CA	2.06	120.41	113.38

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	500	SAH	C-CA-CB-CG
3	A	500	SAH	N-CA-CB-CG

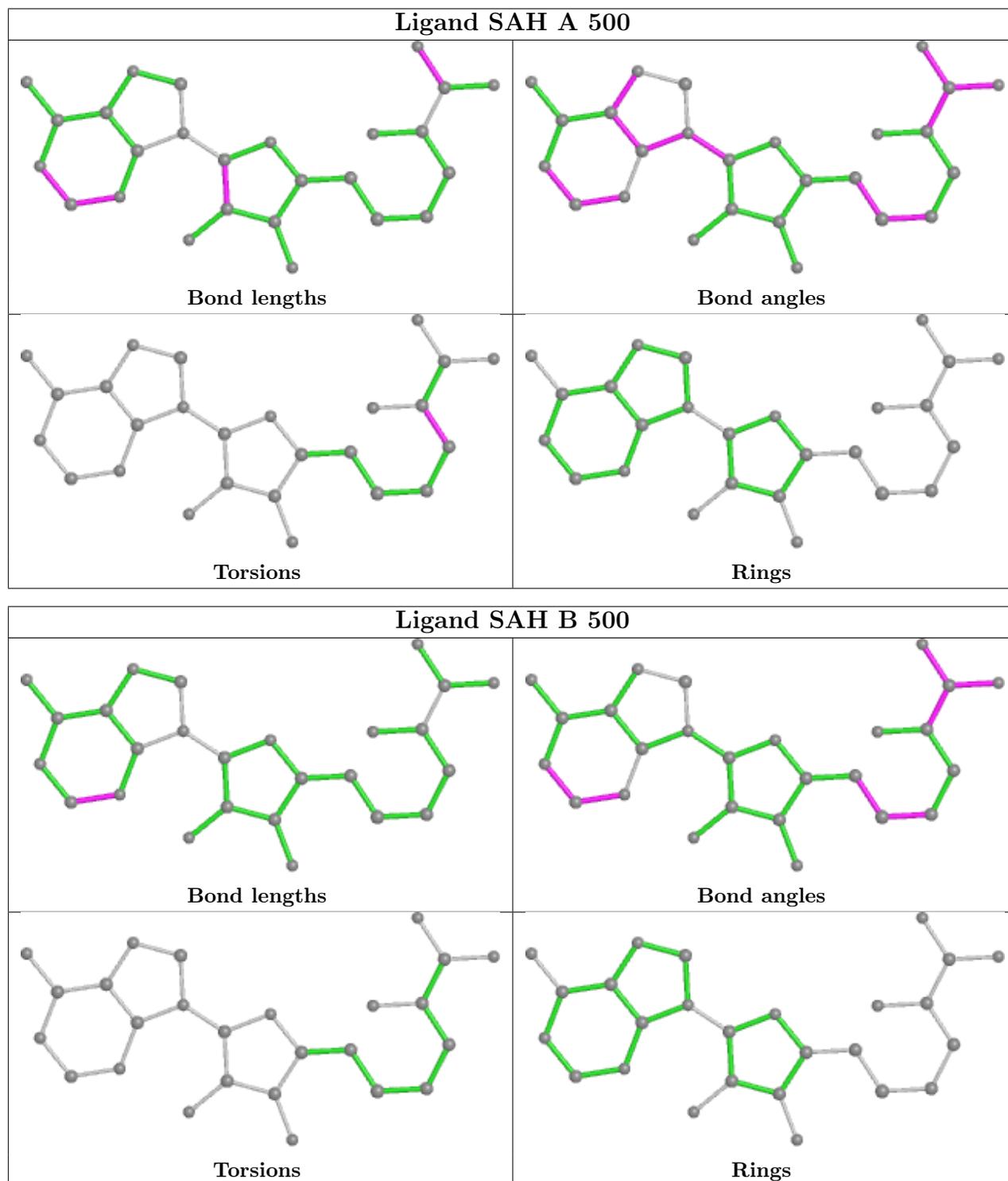
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	500	SAH	1	0
2	B	906	SO4	1	0
3	B	500	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

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	265/269 (98%)	1.17	51 (19%) 1 1	49, 58, 74, 85	0
1	B	263/269 (97%)	1.13	44 (16%) 1 2	49, 57, 71, 97	0
All	All	528/538 (98%)	1.15	95 (17%) 1 1	49, 57, 72, 97	0

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	267	VAL	8.0
1	B	5	GLY	6.8
1	B	250	LYS	5.9
1	A	44	ARG	4.5
1	A	144	LEU	4.5
1	A	248	ILE	4.5
1	B	44	ARG	4.4
1	A	107	GLY	4.4
1	A	247	LYS	4.4
1	A	25	PHE	4.3
1	A	145	CYS	4.3
1	A	181	ILE	4.2
1	A	267	VAL	4.0
1	B	145	CYS	3.9
1	B	87	TRP	3.9
1	B	226	SER	3.8
1	A	245	ASP	3.8
1	A	21	LYS	3.8
1	A	100	VAL	3.7
1	B	266	ALA	3.7
1	A	42	ARG	3.6
1	A	80	LEU	3.5
1	B	144	LEU	3.5
1	B	38	THR	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	147	ILE	3.5
1	A	38	THR	3.4
1	A	184	LEU	3.4
1	B	247	LYS	3.4
1	B	42	ARG	3.3
1	A	22	GLU	3.3
1	A	183	ILE	3.2
1	A	182	LYS	3.2
1	A	87	TRP	3.2
1	B	183	ILE	3.1
1	A	209	VAL	3.1
1	B	45	ARG	3.1
1	A	254	TYR	3.0
1	A	220	TYR	3.0
1	B	241	ILE	2.9
1	B	20	GLY	2.8
1	B	7	THR	2.8
1	B	25	PHE	2.8
1	A	40	ALA	2.8
1	A	46	GLU	2.7
1	A	35	VAL	2.7
1	B	220	TYR	2.7
1	B	139	ILE	2.7
1	B	62	LEU	2.7
1	A	47	GLY	2.7
1	B	147	ILE	2.7
1	B	82	CYS	2.6
1	A	86	GLY	2.6
1	A	3	ALA	2.6
1	B	95	LYS	2.6
1	A	88	SER	2.6
1	A	251	GLY	2.6
1	B	254	TYR	2.5
1	A	45	ARG	2.5
1	A	82	CYS	2.5
1	A	62	LEU	2.5
1	A	65	LEU	2.5
1	A	206	LEU	2.5
1	A	180	CYS	2.5
1	B	86	GLY	2.5
1	A	143	LEU	2.5
1	B	202	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	176	PRO	2.5
1	B	180	CYS	2.5
1	B	181	ILE	2.5
1	B	59	THR	2.4
1	B	182	LYS	2.4
1	A	12	TRP	2.4
1	B	195	LEU	2.4
1	B	156	ILE	2.3
1	A	60	ALA	2.3
1	A	219	MET	2.3
1	B	248	ILE	2.3
1	B	231	HIS	2.2
1	A	123	ILE	2.2
1	A	191	VAL	2.2
1	A	128	SER	2.2
1	B	18	ALA	2.2
1	B	93	THR	2.2
1	B	80	LEU	2.2
1	A	240	LEU	2.2
1	B	39	GLU	2.2
1	A	146	ASP	2.2
1	B	164	ILE	2.1
1	A	39	GLU	2.1
1	A	212	SER	2.1
1	A	85	GLY	2.1
1	B	143	LEU	2.1
1	B	167	MET	2.1
1	B	60	ALA	2.1
1	A	173	SER	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](#)

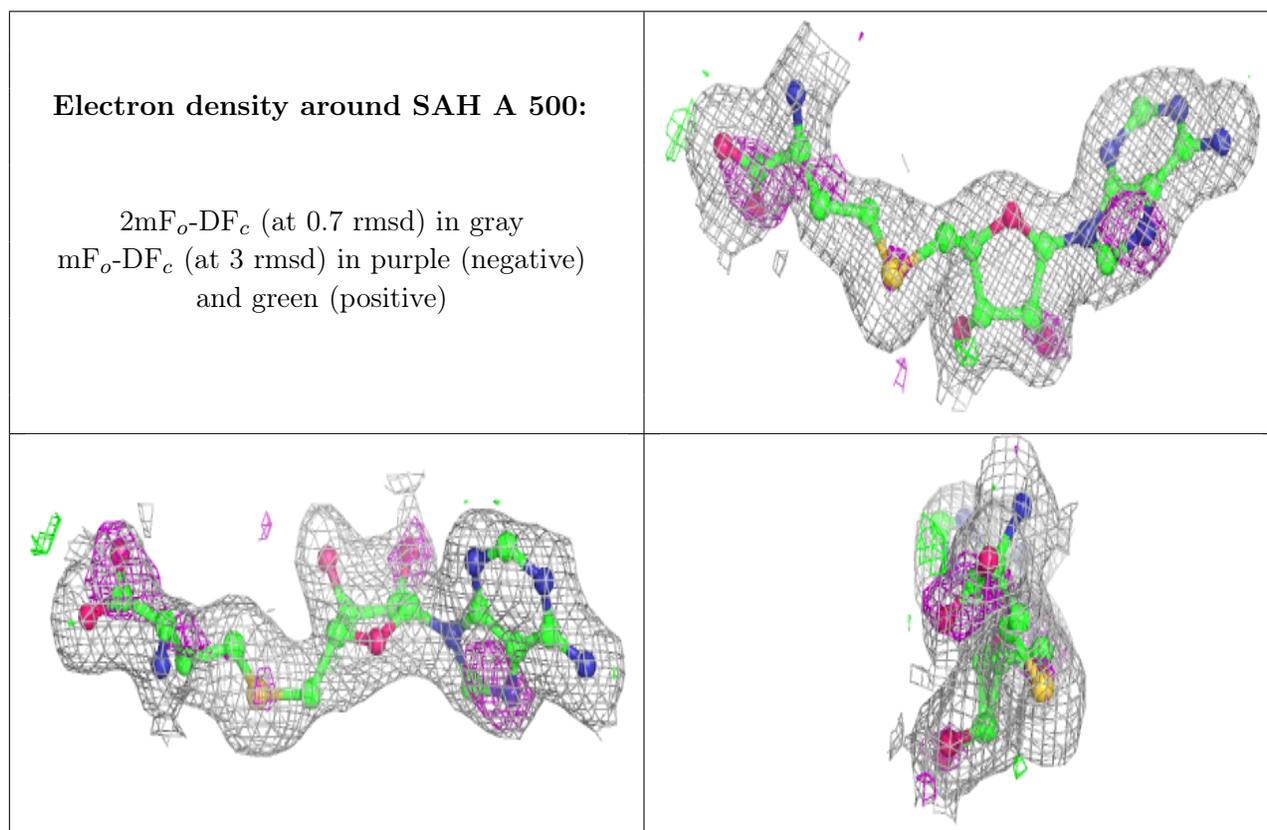
There are no monosaccharides in this entry.

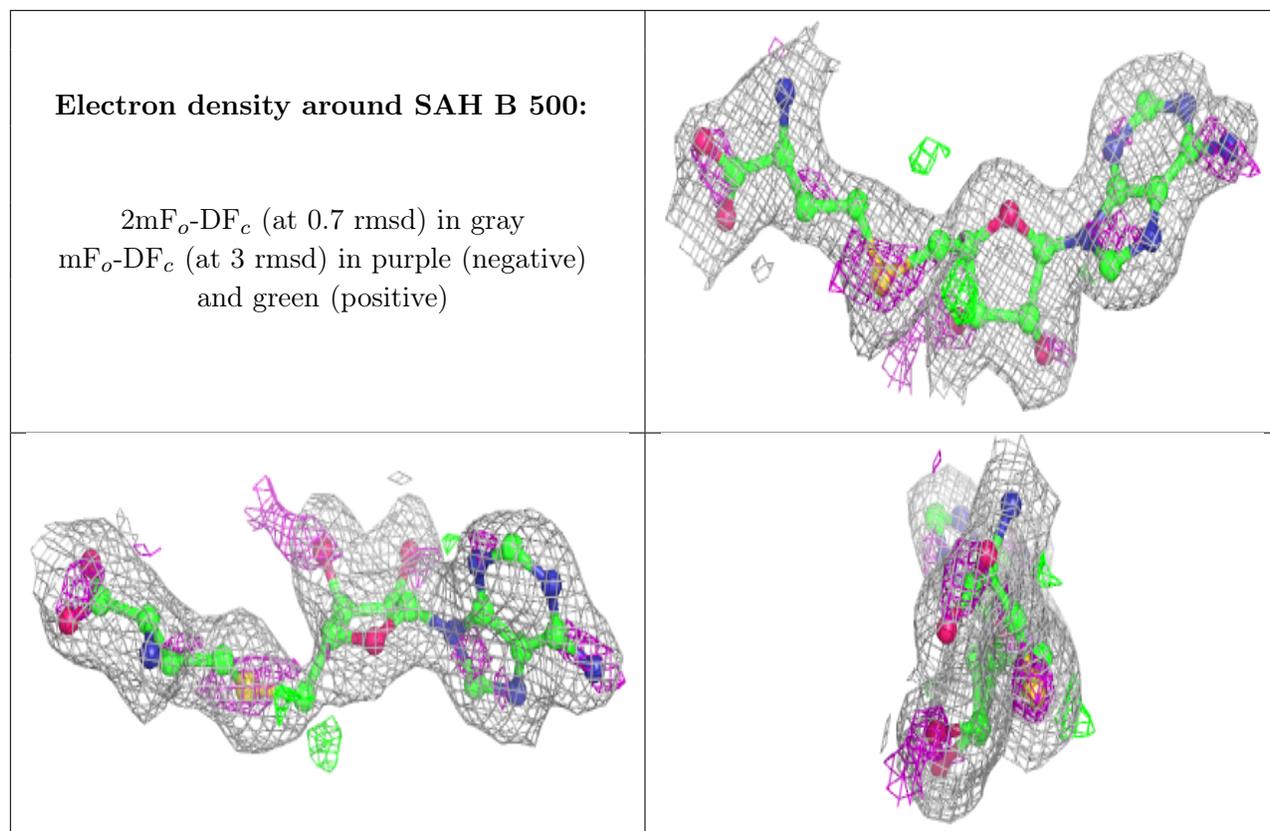
6.4 Ligands [i](#)

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(Å ²)	Q<0.9
2	SO4	B	910	5/5	0.79	0.43	89,113,116,117	0
2	SO4	A	902	5/5	0.80	0.36	87,110,116,120	0
2	SO4	A	905	5/5	0.81	0.38	65,94,104,108	0
2	SO4	A	911	5/5	0.88	0.52	71,73,92,92	0
3	SAH	A	500	26/26	0.88	0.16	39,51,58,61	0
3	SAH	B	500	26/26	0.90	0.15	38,49,57,65	0
2	SO4	B	906	5/5	0.94	0.26	60,72,77,80	0
2	SO4	B	908	5/5	0.95	0.25	79,79,89,91	0
2	SO4	A	903	5/5	0.98	0.38	64,66,68,71	0
2	SO4	B	907	5/5	0.98	0.14	47,53,69,75	0
2	SO4	A	904	5/5	0.98	0.15	55,56,65,66	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.