



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 15, 2020 – 12:03 am BST

PDB ID : 3G5T
Title : Crystal structure of trans-aconitate 3-methyltransferase from yeast
Authors : Burgie, E.S.; Bingman, C.A.; Wesenberg, G.E.; Phillips Jr., G.N.; Center for Eukaryotic Structural Genomics (CESG)
Deposited on : 2009-02-05
Resolution : 1.12 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

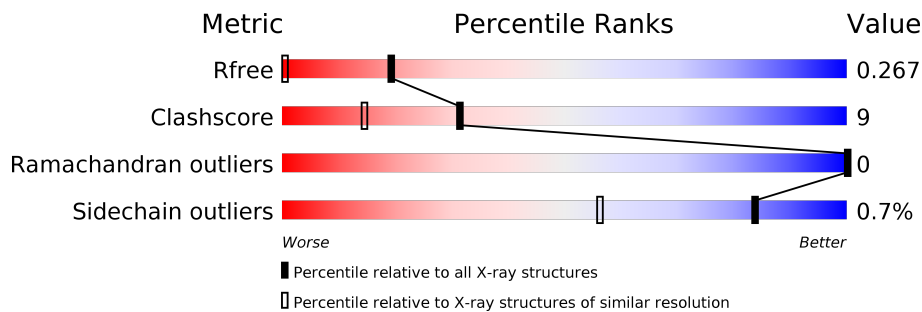
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 1.12 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	1168 (1.14-1.10)
Clashscore	141614	1205 (1.14-1.10)
Ramachandran outliers	138981	1168 (1.14-1.10)
Sidechain outliers	138945	1165 (1.14-1.10)

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	299	

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	EDO	A	304	-	-	X	-
5	DMS	A	310	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5507 atoms, of which 2264 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.

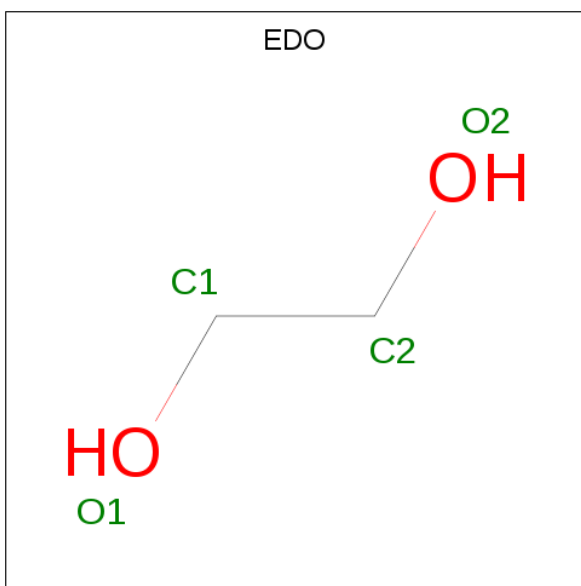
- Molecule 1 is a protein called Trans-aconitate 3-methyltransferase.

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S	Se			
1	A	299	4891	1669	2264	436	511	3	8	0	42	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	SER	-	EXPRESSION TAG	UNP P32643

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



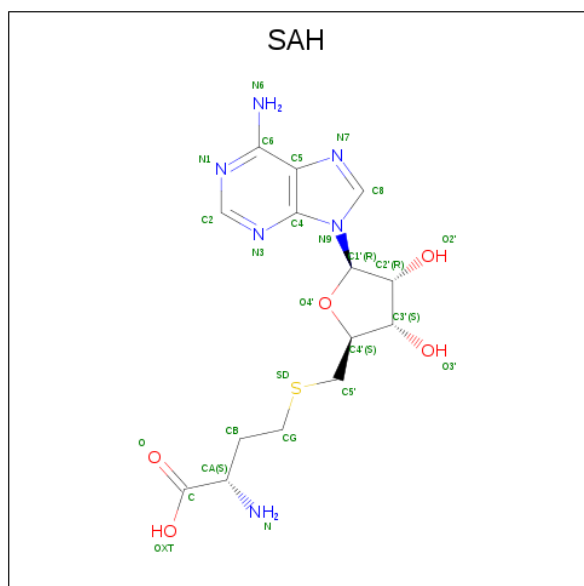
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	4	2	2	0	0
2	A	1	4	2	2	0	0
2	A	1	4	2	2	0	0

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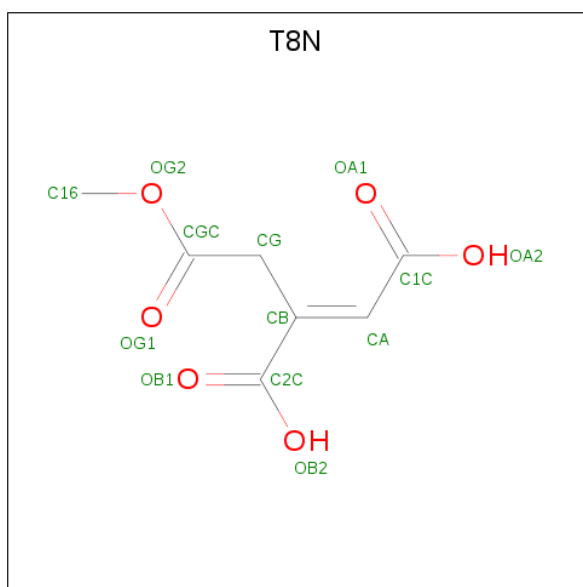
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		

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



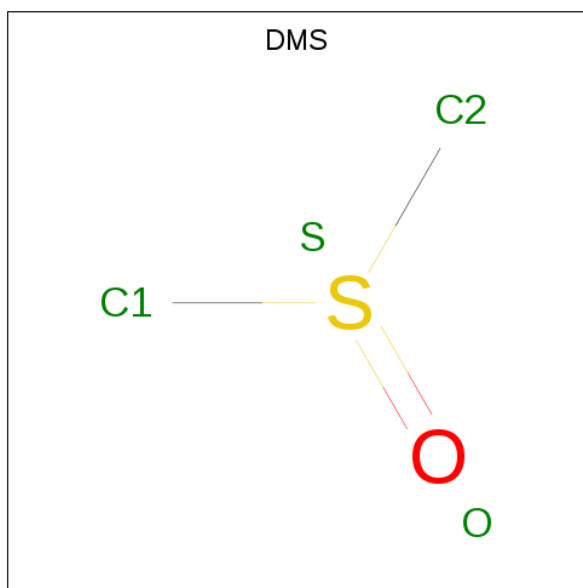
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			26	14	6	5	1		

- Molecule 4 is (2E)-2-(2-methoxy-2-oxoethyl)but-2-enedioic acid (three-letter code: T8N) (formula: $C_7H_8O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	7	6		

- Molecule 5 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C₂H₆OS).



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

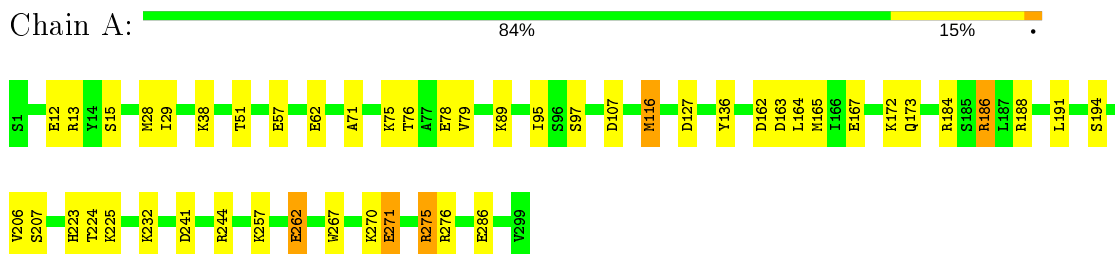
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	522	Total 541	O 541	0	19

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.

- Molecule 1: Trans-aconitate 3-methyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	36.83Å 91.89Å 104.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.47 – 1.12 28.74 – 1.12	Depositor EDS
% Data completeness (in resolution range)	98.2 (34.47-1.12) 98.2 (28.74-1.12)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.99 (at 1.12Å)	Xtrriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.120 , 0.137 0.256 , 0.267	Depositor DCC
R_{free} test set	6751 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	8.8	Xtrriage
Anisotropy	0.478	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 39.2	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	5507	wwPDB-VP
Average B, all atoms (Å ²)	12.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.01% 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: T8N, SAH, DMS, EDO

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	1.34	34/2908 (1.2%)	1.28	19/3904 (0.5%)

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	A	0	1

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	97[A]	SER	CB-OG	-9.54	1.29	1.42
1	A	97[B]	SER	CB-OG	-9.54	1.29	1.42
1	A	223	HIS	C-O	-9.14	1.05	1.23
1	A	257	LYS	CB-CG	-8.90	1.28	1.52
1	A	186[A]	ARG	CZ-NH1	8.36	1.44	1.33
1	A	186[B]	ARG	CZ-NH1	8.36	1.44	1.33
1	A	275	ARG	NE-CZ	-8.29	1.22	1.33
1	A	232	LYS	CE-NZ	-7.62	1.30	1.49
1	A	241	ASP	CB-CG	7.50	1.67	1.51
1	A	57	GLU	CG-CD	-7.31	1.41	1.51
1	A	116[A]	MSE	SE-CE	-7.25	1.52	1.95
1	A	116[B]	MSE	SE-CE	-7.25	1.52	1.95
1	A	286	GLU	CG-CD	-7.00	1.41	1.51
1	A	15[A]	SER	CB-OG	-6.85	1.33	1.42
1	A	15[B]	SER	CB-OG	-6.85	1.33	1.42
1	A	271[A]	GLU	CB-CG	-6.67	1.39	1.52
1	A	271[B]	GLU	CB-CG	-6.67	1.39	1.52
1	A	275	ARG	CG-CD	-6.63	1.35	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	262[A]	GLU	CG-CD	-6.24	1.42	1.51
1	A	262[B]	GLU	CG-CD	-6.24	1.42	1.51
1	A	167[A]	GLU	CG-CD	6.14	1.61	1.51
1	A	167[B]	GLU	CG-CD	6.14	1.61	1.51
1	A	172	LYS	CE-NZ	6.13	1.64	1.49
1	A	62[A]	GLU	CD-OE2	6.03	1.32	1.25
1	A	62[B]	GLU	CD-OE2	6.03	1.32	1.25
1	A	136	TYR	CE2-CZ	-5.70	1.31	1.38
1	A	165	MSE	SE-CE	-5.66	1.62	1.95
1	A	167[A]	GLU	CD-OE2	5.23	1.31	1.25
1	A	167[B]	GLU	CD-OE2	5.23	1.31	1.25
1	A	271[A]	GLU	CD-OE1	5.10	1.31	1.25
1	A	271[B]	GLU	CD-OE1	5.10	1.31	1.25
1	A	257	LYS	CE-NZ	5.03	1.61	1.49
1	A	194[A]	SER	CA-CB	-5.01	1.45	1.52
1	A	194[B]	SER	CA-CB	-5.01	1.45	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	275	ARG	NE-CZ-NH2	-31.83	104.39	120.30
1	A	275	ARG	NH1-CZ-NH2	14.32	135.16	119.40
1	A	184	ARG	NE-CZ-NH2	-12.52	114.04	120.30
1	A	13	ARG	NE-CZ-NH2	-11.54	114.53	120.30
1	A	13	ARG	NE-CZ-NH1	10.58	125.59	120.30
1	A	244[A]	ARG	NE-CZ-NH1	-8.98	115.81	120.30
1	A	244[B]	ARG	NE-CZ-NH1	-8.98	115.81	120.30
1	A	186[A]	ARG	NE-CZ-NH2	-8.09	116.26	120.30
1	A	186[B]	ARG	NE-CZ-NH2	-8.09	116.26	120.30
1	A	184	ARG	NE-CZ-NH1	8.01	124.30	120.30
1	A	276	ARG	NE-CZ-NH1	6.74	123.67	120.30
1	A	275	ARG	CD-NE-CZ	6.48	132.67	123.60
1	A	271[A]	GLU	OE1-CD-OE2	6.23	130.77	123.30
1	A	271[B]	GLU	OE1-CD-OE2	6.23	130.77	123.30
1	A	116[A]	MSE	CG-SE-CE	-5.90	85.93	98.90
1	A	116[B]	MSE	CG-SE-CE	-5.90	85.93	98.90
1	A	276	ARG	NE-CZ-NH2	-5.47	117.56	120.30
1	A	127	ASP	CB-CG-OD1	5.39	123.15	118.30
1	A	162	ASP	CB-CG-OD2	-5.14	113.68	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	224	THR	Peptide

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	2627	2264	2582	47	0
2	A	32	0	48	8	0
3	A	26	0	19	0	0
4	A	13	0	6	0	0
5	A	4	0	6	4	0
6	A	541	0	0	17	0
All	All	3243	2264	2661	49	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 (49) 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:75[C]:LYS:HE2	6:A:718:HOH:O	1.04	1.21
1:A:191[B]:LEU:HD23	2:A:304:EDO:H11	1.30	1.11
1:A:271[B]:GLU:OE1	6:A:672:HOH:O	1.85	0.93
1:A:38[C]:LYS:HE3	6:A:787:HOH:O	1.69	0.91
1:A:191[B]:LEU:HD23	2:A:304:EDO:C1	2.11	0.80
1:A:186[B]:ARG:NE	6:A:857:HOH:O	1.93	0.76
1:A:163[B]:ASP:OD1	1:A:275:ARG:NH1	2.22	0.73
1:A:191[B]:LEU:CD2	2:A:304:EDO:H11	2.16	0.72
1:A:186[A]:ARG:HG2	2:A:304:EDO:H12	1.75	0.69
1:A:225:LYS:HD3	1:A:225:LYS:O	1.93	0.67
1:A:71:ALA:HA	1:A:95[B]:ILE:HD11	1.77	0.65
1:A:207[B]:SER:HB2	1:A:225:LYS:HD2	1.77	0.65
1:A:38[B]:LYS:HD2	6:A:787:HOH:O	2.02	0.60
1:A:75[C]:LYS:HD2	1:A:76:THR:N	2.16	0.60
1:A:38[C]:LYS:CE	6:A:787:HOH:O	2.38	0.60
1:A:78[B]:GLU:OE2	6:A:465:HOH:O	2.14	0.59
1:A:206[B]:VAL:HG13	6:A:775:HOH:O	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:186[B]:ARG:HG2	2:A:304:EDO:H12	1.85	0.58
1:A:164:LEU:HD21	1:A:271[A]:GLU:HG2	1.85	0.57
1:A:188:ARG:HB3	5:A:310:DMS:H23	1.87	0.56
1:A:262[A]:GLU:OE2	1:A:270[A]:LYS:NZ	2.36	0.56
1:A:186[B]:ARG:NH2	6:A:857:HOH:O	2.38	0.55
1:A:75[A]:LYS:HD3	6:A:719:HOH:O	2.07	0.54
1:A:267:TRP:CZ2	1:A:271[B]:GLU:HG3	2.43	0.54
1:A:75[C]:LYS:NZ	6:A:413:HOH:O	2.41	0.53
1:A:163[B]:ASP:OD2	1:A:275:ARG:NH1	2.42	0.51
1:A:267:TRP:CE2	1:A:271[B]:GLU:HG3	2.45	0.51
1:A:163[B]:ASP:CG	1:A:275:ARG:NH1	2.63	0.51
1:A:173[A]:GLN:HG3	6:A:601:HOH:O	2.12	0.50
1:A:12[A]:GLU:CG	6:A:413:HOH:O	2.60	0.49
1:A:186[B]:ARG:CG	2:A:304:EDO:H12	2.42	0.49
1:A:186[A]:ARG:CG	2:A:304:EDO:H12	2.43	0.49
1:A:89[A]:LYS:HE2	6:A:354:HOH:O	2.14	0.48
5:A:310:DMS:C2	6:A:483:HOH:O	2.61	0.47
1:A:225:LYS:CD	1:A:225:LYS:O	2.62	0.47
1:A:163[B]:ASP:OD2	1:A:275:ARG:CZ	2.62	0.47
5:A:310:DMS:H22	6:A:483:HOH:O	2.13	0.47
1:A:28[B]:MSE:SE	2:A:305:EDO:H22	2.65	0.47
1:A:188:ARG:HB3	5:A:310:DMS:C2	2.47	0.45
1:A:186[B]:ARG:CZ	6:A:857:HOH:O	2.48	0.44
1:A:163[B]:ASP:OD2	1:A:275:ARG:NH2	2.50	0.44
1:A:75[C]:LYS:HD3	1:A:79:VAL:HG23	1.99	0.43
1:A:163[B]:ASP:OD1	1:A:275:ARG:CZ	2.66	0.42
1:A:29[A]:ILE:HD13	1:A:29[A]:ILE:HG21	1.83	0.41

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	345/299 (115%)	339 (98%)	6 (2%)	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	317/263 (120%)	315 (99%)	2 (1%)	86	61

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	THR
1	A	107	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates 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	EDO	A	306	-	3,3,3	0.41	0	2,2,2	0.99	0
3	SAH	A	308	-	21,28,28	1.25	2 (9%)	20,40,40	1.41	2 (10%)
2	EDO	A	307	-	3,3,3	0.56	0	2,2,2	0.49	0
2	EDO	A	302	-	3,3,3	0.65	0	2,2,2	0.57	0
2	EDO	A	303	-	3,3,3	1.40	1 (33%)	2,2,2	0.06	0
2	EDO	A	304	-	3,3,3	1.03	0	2,2,2	1.34	0
2	EDO	A	300	-	3,3,3	0.33	0	2,2,2	0.23	0
5	DMS	A	310	-	3,3,3	2.58	2 (66%)	3,3,3	0.53	0
4	T8N	A	309	-	6,12,12	1.01	1 (16%)	6,15,15	1.04	0
2	EDO	A	305	-	3,3,3	0.60	0	2,2,2	0.12	0
2	EDO	A	301	-	3,3,3	0.28	0	2,2,2	0.14	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
2	EDO	A	306	-	-	0/1/1/1	-
3	SAH	A	308	-	-	2/7/31/31	0/3/3/3
2	EDO	A	307	-	-	1/1/1/1	-
2	EDO	A	302	-	-	0/1/1/1	-
2	EDO	A	303	-	-	0/1/1/1	-
2	EDO	A	304	-	-	0/1/1/1	-
2	EDO	A	300	-	-	0/1/1/1	-
4	T8N	A	309	-	-	1/8/14/14	-
2	EDO	A	305	-	-	0/1/1/1	-
2	EDO	A	301	-	-	0/1/1/1	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	310	DMS	O-S	3.29	1.72	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	310	DMS	C2-S	-3.02	1.53	1.75
3	A	308	SAH	O4'-C1'	2.76	1.44	1.41
3	A	308	SAH	CG-SD	-2.44	1.72	1.81
2	A	303	EDO	O2-C2	2.36	1.54	1.42
4	A	309	T8N	CG-CGC	-2.14	1.47	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	308	SAH	O4'-C1'-C2'	-3.23	102.21	106.93
3	A	308	SAH	CB-CG-SD	-2.72	107.20	113.31

There are no chirality outliers.

All (4) torsion outliers are listed below:

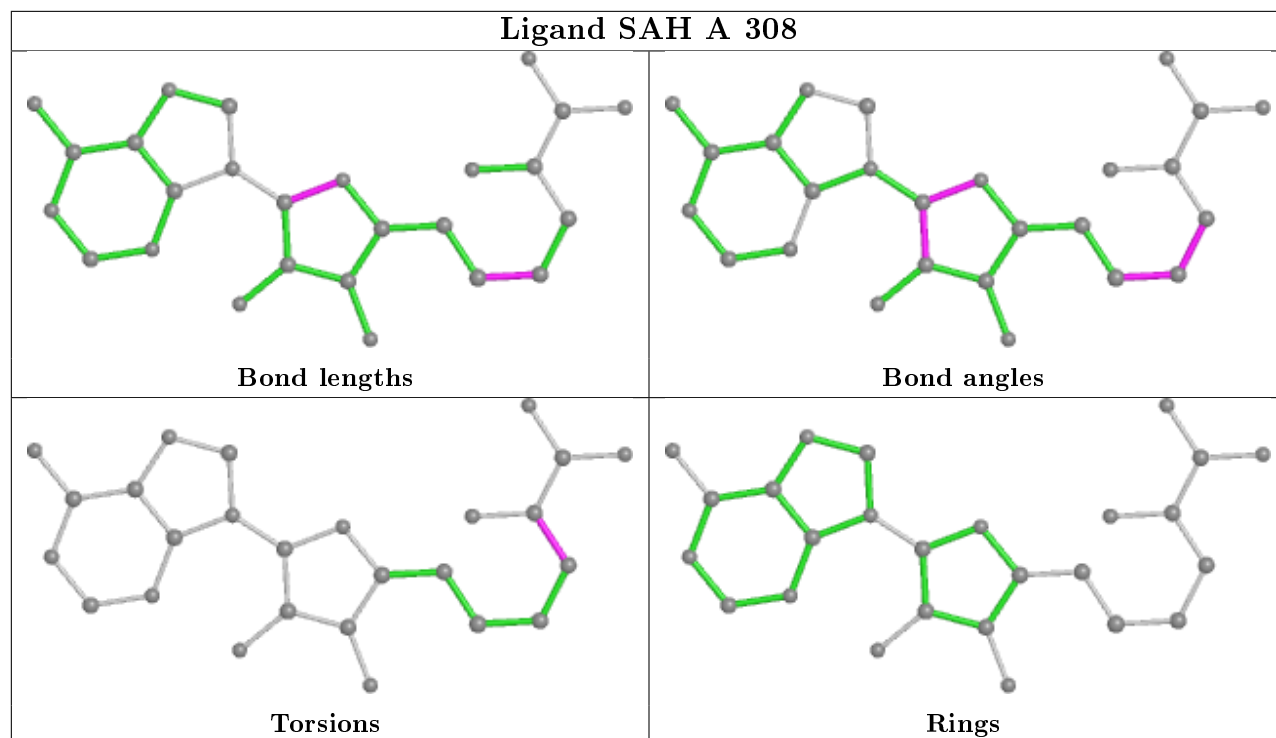
Mol	Chain	Res	Type	Atoms
3	A	308	SAH	N-CA-CB-CG
3	A	308	SAH	C-CA-CB-CG
2	A	307	EDO	O1-C1-C2-O2
4	A	309	T8N	CB-CG-CGC-OG2

There are no ring outliers.

3 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	304	EDO	7	0
5	A	310	DMS	4	0
2	A	305	EDO	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

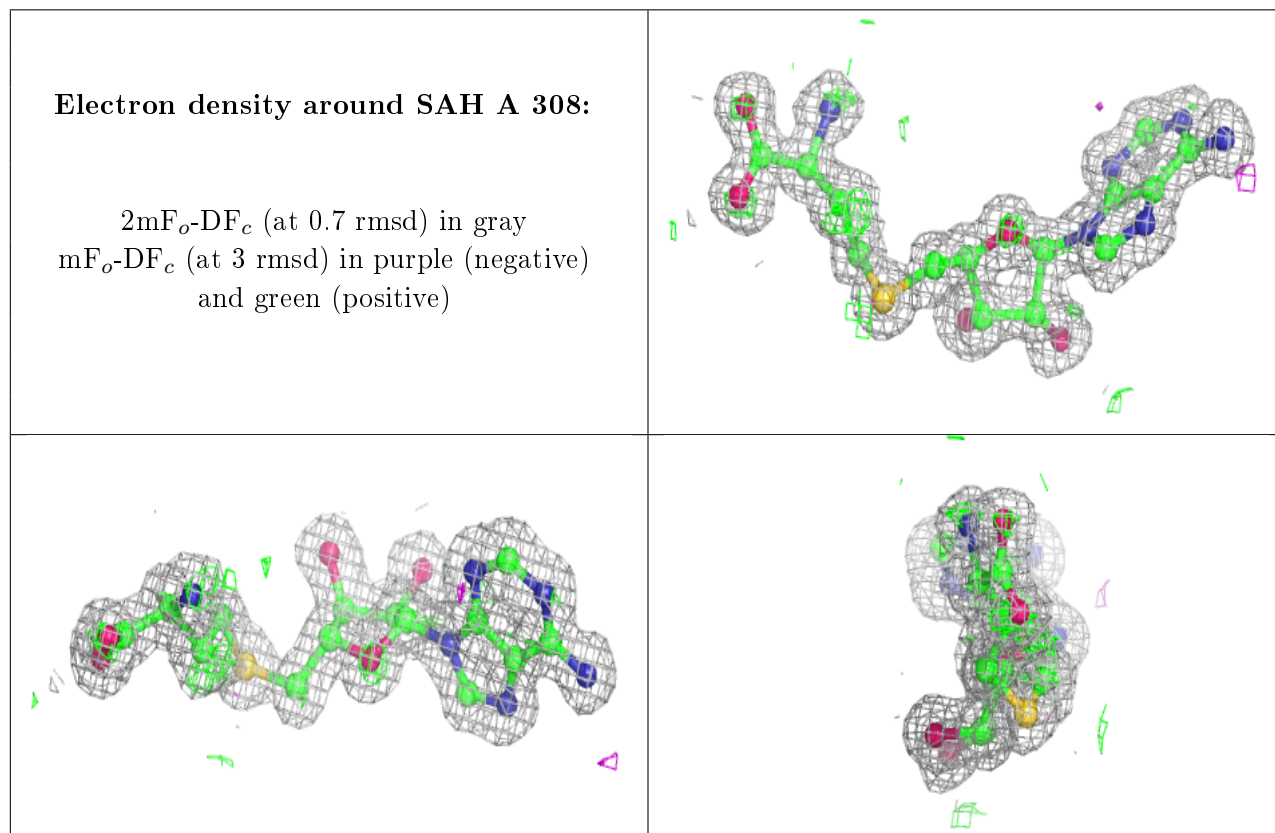
6.3 Carbohydrates [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.