



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 19, 2022 – 06:00 AM JST

PDB ID : 7VKQ
Title : Crystal structure of D. melanogaster SAMTOR in the SAH bound form
Authors : Tang, X.; Zhang, T.; Ding, J.
Deposited on : 2021-09-30
Resolution : 2.09 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 : 4.02b-467
Mogul : ?? (??), CSD ??CSD?? (????)
Xtriage (Phenix) : 1.13
EDS : 2.29
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) : 2.29

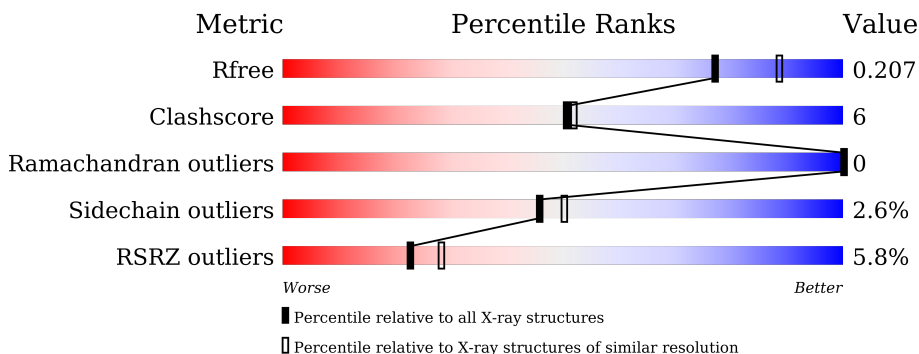
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.09 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 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	304	<div> <div>2%</div> <div> <div></div> <div>61%</div> <div>13%</div> <div>•</div> <div>25%</div> </div> </div>
1	B	304	<div> <div>4%</div> <div> <div></div> <div>61%</div> <div>12%</div> <div>•</div> <div>26%</div> </div> </div>
1	C	304	<div> <div>2%</div> <div> <div></div> <div>60%</div> <div>10%</div> <div></div> <div>31%</div> </div> </div>
1	D	304	<div> <div>9%</div> <div> <div></div> <div>61%</div> <div>8%</div> <div></div> <div>31%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7774 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.

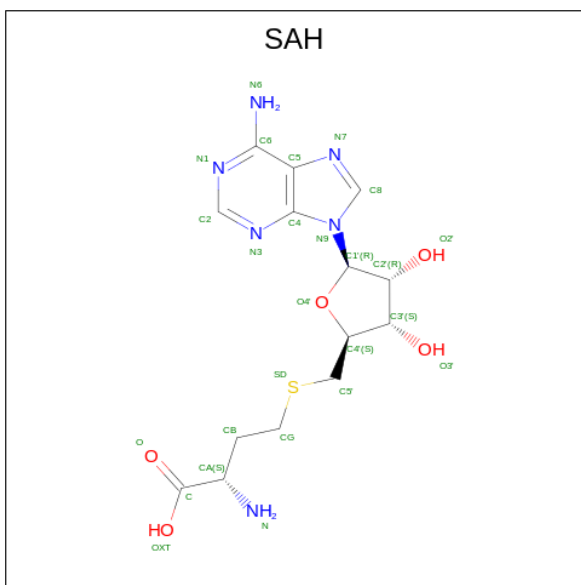
- Molecule 1 is a protein called S-adenosylmethionine sensor upstream of mTORC1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	228	Total	C	N	O	S	0	1	0
			1843	1177	323	330	13			
1	B	225	Total	C	N	O	S	0	1	0
			1819	1164	321	321	13			
1	C	210	Total	C	N	O	S	0	1	0
			1717	1103	297	307	10			
1	D	210	Total	C	N	O	S	0	1	0
			1693	1082	296	305	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q9W138
A	0	SER	-	expression tag	UNP Q9W138
B	-1	GLY	-	expression tag	UNP Q9W138
B	0	SER	-	expression tag	UNP Q9W138
C	-1	GLY	-	expression tag	UNP Q9W138
C	0	SER	-	expression tag	UNP Q9W138
D	-1	GLY	-	expression tag	UNP Q9W138
D	0	SER	-	expression tag	UNP Q9W138

- Molecule 2 is S-ADENOSYL-L-HOMOCYSTEINE (three-letter code: SAH) (formula: $C_{14}H_{20}N_6O_5S$) (labeled as "Ligand of Interest" by depositor).



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

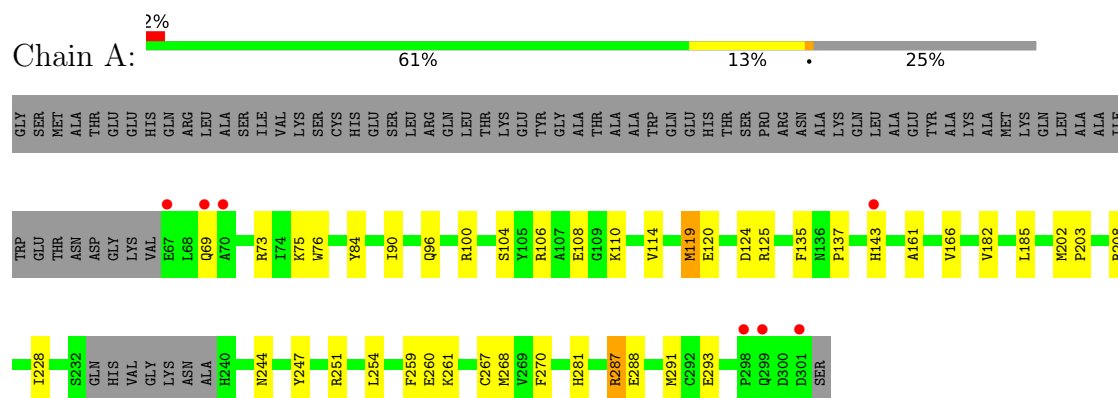
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	159	Total O 159 159	0	0
3	B	145	Total O 145 145	0	0
3	C	164	Total O 164 164	0	0
3	D	130	Total O 130 130	0	0

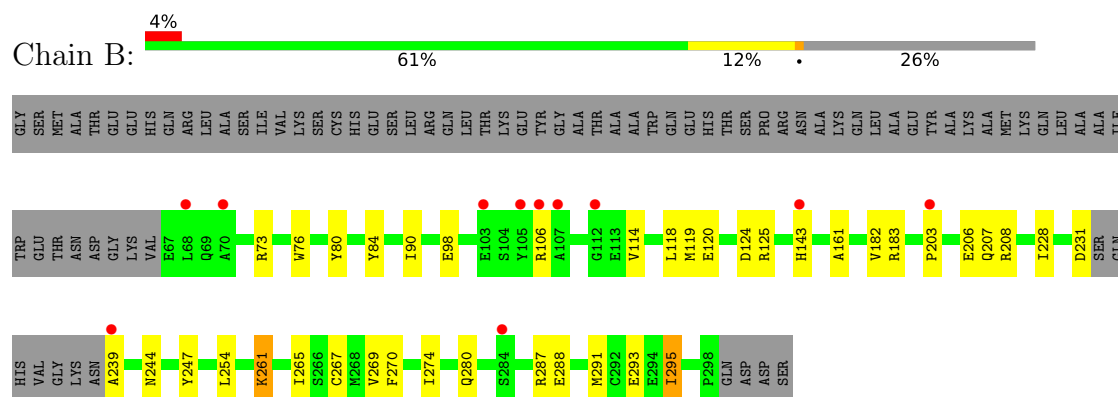
3 Residue-property plots [i](#)

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.

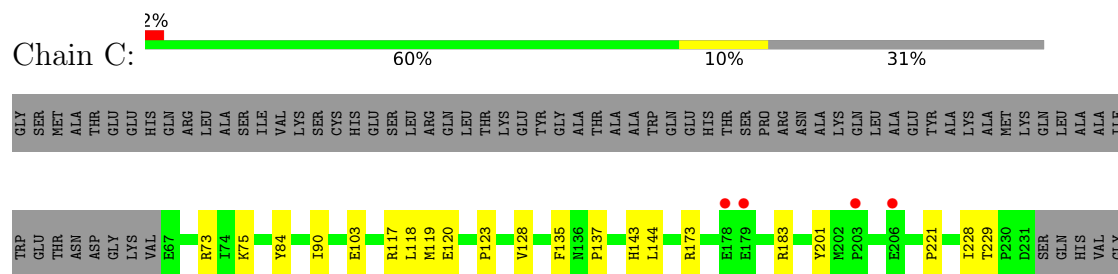
- Molecule 1: S-adenosylmethionine sensor upstream of mTORC1



- Molecule 1: S-adenosylmethionine sensor upstream of mTORC1

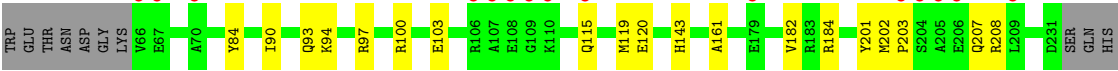
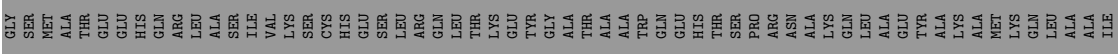


- Molecule 1: S-adenosylmethionine sensor upstream of mTORC1





● Molecule 1: S-adenosylmethionine sensor upstream of mTORC1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	49.93Å 63.99Å 80.27Å 90.26° 93.38° 96.79°	Depositor
Resolution (Å)	41.02 – 2.09 41.51 – 2.09	Depositor EDS
% Data completeness (in resolution range)	94.8 (41.02-2.09) 94.8 (41.51-2.09)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.166 , 0.206 0.166 , 0.207	Depositor DCC
R_{free} test set	1996 reflections (3.63%)	wwPDB-VP
Wilson B-factor (Å ²)	18.8	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 55.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7774	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.24% 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: 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	0.39	0/1883	0.56	0/2548
1	B	0.38	0/1859	0.55	0/2515
1	C	0.42	0/1757	0.54	0/2380
1	D	0.38	0/1729	0.53	0/2342
All	All	0.40	0/7228	0.55	0/9785

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 [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	1843	0	1832	24	0
1	B	1819	0	1822	20	0
1	C	1717	0	1715	18	0
1	D	1693	0	1684	21	0
2	A	26	0	19	0	0
2	B	26	0	19	0	0
2	C	26	0	19	1	0
2	D	26	0	19	2	0
3	A	159	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	145	0	0	3	0
3	C	164	0	0	4	0
3	D	130	0	0	3	0
All	All	7774	0	7129	81	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 6.

All (81) 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:B:119:MET:SD	3:B:612:HOH:O	2.35	0.84
1:B:73:ARG:HB2	1:B:228:ILE:HD13	1.69	0.74
1:A:124:ASP:O	1:A:143:HIS:HB2	1.95	0.67
1:D:97[A]:ARG:NH2	3:D:501:HOH:O	2.24	0.65
1:C:243:LYS:HD3	1:C:245[B]:TRP:HB2	1.78	0.65
1:C:247:TYR:CE2	1:C:251:ARG:HD2	2.32	0.65
1:D:97[B]:ARG:HG2	1:D:100:ARG:NH1	2.12	0.65
1:B:280:GLN:NE2	3:B:501:HOH:O	2.21	0.62
1:D:184:ARG:NH2	3:D:502:HOH:O	2.33	0.62
1:B:119:MET:HG2	1:B:120:GLU:O	2.00	0.61
1:C:201:TYR:CD1	2:C:401:SAH:H8	2.35	0.61
1:A:73:ARG:HB2	1:A:228:ILE:HD13	1.83	0.61
1:A:106:ARG:HH21	1:A:114:VAL:H	1.51	0.59
1:B:161:ALA:HB2	1:B:182:VAL:HB	1.86	0.58
1:A:119:MET:HG2	1:A:120:GLU:O	2.04	0.58
1:C:258:ARG:NH1	1:C:260:GLU:OE1	2.37	0.57
1:B:239:ALA:N	3:B:504:HOH:O	2.38	0.57
1:C:73:ARG:HB2	1:C:228:ILE:HD13	1.86	0.56
1:A:288:GLU:O	1:A:291:MET:HG2	2.06	0.56
1:C:247:TYR:CZ	1:C:251:ARG:HD2	2.40	0.55
1:C:229:THR:HG21	1:C:245[A]:TRP:CZ3	2.42	0.55
1:A:96:GLN:OE1	1:A:100[B]:ARG:NH2	2.40	0.54
1:D:161:ALA:HB2	1:D:182:VAL:HB	1.88	0.54
1:B:288:GLU:O	1:B:291:MET:HG2	2.08	0.53
1:A:202:MET:O	1:A:208:ARG:NH1	2.42	0.52
1:B:106:ARG:HH21	1:B:114:VAL:H	1.57	0.51
1:A:161:ALA:HB2	1:A:182:VAL:HB	1.92	0.51
1:A:96:GLN:HE22	1:D:258:ARG:HH12	1.58	0.50
1:B:125:ARG:HD2	1:B:143:HIS:HA	1.94	0.50
1:D:201:TYR:CD1	2:D:401:SAH:H8	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:TYR:HA	1:A:90:ILE:HB	1.93	0.50
1:B:247:TYR:CE1	1:B:287:ARG:HB2	2.47	0.50
1:D:84:TYR:HA	1:D:90:ILE:HB	1.94	0.49
1:D:119:MET:HG2	1:D:120:GLU:O	2.12	0.49
1:C:183:ARG:NH2	3:C:502:HOH:O	2.22	0.49
1:C:243:LYS:HD2	3:C:623:HOH:O	2.13	0.49
1:D:97[B]:ARG:HG2	1:D:100:ARG:HH11	1.78	0.48
1:D:93:GLN:O	1:D:97[B]:ARG:HG3	2.13	0.48
1:A:76:TRP:HE1	1:A:260:GLU:CD	2.17	0.47
1:B:76:TRP:CH2	1:B:267:CYS:HB3	2.50	0.47
1:B:208:ARG:HB3	1:B:295:ILE:HG12	1.96	0.47
1:C:229:THR:HG21	1:C:245[A]:TRP:HZ3	1.79	0.47
1:D:247:TYR:CE2	1:D:251:ARG:HD2	2.49	0.47
1:A:119:MET:HB3	1:A:119:MET:HE3	1.77	0.47
1:C:119:MET:HG2	1:C:120:GLU:O	2.14	0.46
1:A:135:PHE:CE2	1:A:137:PRO:HB3	2.50	0.46
1:A:125:ARG:HD2	1:A:143:HIS:HA	1.98	0.46
1:A:259:PHE:CD2	1:A:268:MET:HG2	2.49	0.46
1:C:135:PHE:CE2	1:C:137:PRO:HB3	2.51	0.46
1:C:143:HIS:HB2	3:C:634:HOH:O	2.16	0.46
1:B:84:TYR:HA	1:B:90:ILE:HB	1.98	0.45
1:B:203:PRO:HD2	1:B:207:GLN:OE1	2.17	0.45
1:D:143:HIS:HB2	3:D:591:HOH:O	2.16	0.45
1:D:203:PRO:HD2	1:D:207:GLN:OE1	2.17	0.44
1:C:84:TYR:HA	1:C:90:ILE:HB	2.00	0.44
1:B:124:ASP:O	1:B:143:HIS:HB3	2.19	0.43
1:A:69:GLN:O	1:A:75:LYS:HG3	2.18	0.43
1:A:247:TYR:CZ	1:A:287:ARG:HG3	2.53	0.43
1:B:261:LYS:HE2	1:B:265:ILE:O	2.19	0.43
1:A:254:LEU:HB3	1:A:270:PHE:HB3	2.00	0.43
1:A:251:ARG:NH2	3:A:504:HOH:O	2.42	0.43
1:B:183[A]:ARG:HA	1:B:183[A]:ARG:HD2	1.75	0.42
1:B:80:TYR:CE2	1:B:269:VAL:HG11	2.54	0.42
1:A:76:TRP:CH2	1:A:267:CYS:HB3	2.54	0.42
1:C:117:ARG:NH1	1:C:221:PRO:HG3	2.34	0.42
1:D:246:ARG:HE	1:D:246:ARG:HB3	1.67	0.42
1:A:110:LYS:HB3	1:A:281:HIS:CE1	2.54	0.42
1:A:96:GLN:NE2	1:D:258:ARG:HH12	2.17	0.42
1:C:128:VAL:HG21	1:C:144:LEU:HD22	2.02	0.42
1:D:202:MET:HA	1:D:203:PRO:HD3	1.96	0.42
1:A:104:SER:O	1:A:108:GLU:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:201:TYR:CG	2:D:401:SAH:H8	2.55	0.41
1:D:115:GLN:HG3	1:D:115:GLN:O	2.20	0.41
1:C:277:GLU:OE2	3:C:501:HOH:O	2.21	0.41
1:D:94:LYS:HZ3	1:D:97[B]:ARG:NH2	2.18	0.41
1:B:98:GLU:OE1	1:B:274:ILE:HG23	2.21	0.41
1:B:254:LEU:HB3	1:B:270:PHE:HB3	2.03	0.41
1:D:202:MET:HB2	1:D:208:ARG:HG2	2.03	0.41
1:D:249:LEU:HB3	1:D:254:LEU:HB2	2.02	0.40
1:A:166:VAL:HG11	1:A:185:LEU:HD12	2.02	0.40
1:C:135:PHE:CZ	1:C:137:PRO:HB3	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	225/304 (74%)	220 (98%)	5 (2%)	0	100	100
1	B	222/304 (73%)	219 (99%)	3 (1%)	0	100	100
1	C	207/304 (68%)	197 (95%)	10 (5%)	0	100	100
1	D	207/304 (68%)	202 (98%)	5 (2%)	0	100	100
All	All	861/1216 (71%)	838 (97%)	23 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	
1	A	199/266 (75%)	193 (97%)	6 (3%)	41	44
1	B	196/266 (74%)	189 (96%)	7 (4%)	35	36
1	C	187/266 (70%)	181 (97%)	6 (3%)	39	41
1	D	183/266 (69%)	182 (100%)	1 (0%)	88	92
All	All	765/1064 (72%)	745 (97%)	20 (3%)	46	50

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	119	MET
1	A	203	PRO
1	A	244	ASN
1	A	261	LYS
1	A	287	ARG
1	A	293	GLU
1	B	118	LEU
1	B	206	GLU
1	B	231	ASP
1	B	244	ASN
1	B	261	LYS
1	B	293	GLU
1	B	295	ILE
1	C	75	LYS
1	C	103	GLU
1	C	118	LEU
1	C	123	PRO
1	C	173	ARG
1	C	261	LYS
1	D	103	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	127	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 ligands are modelled in this entry.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

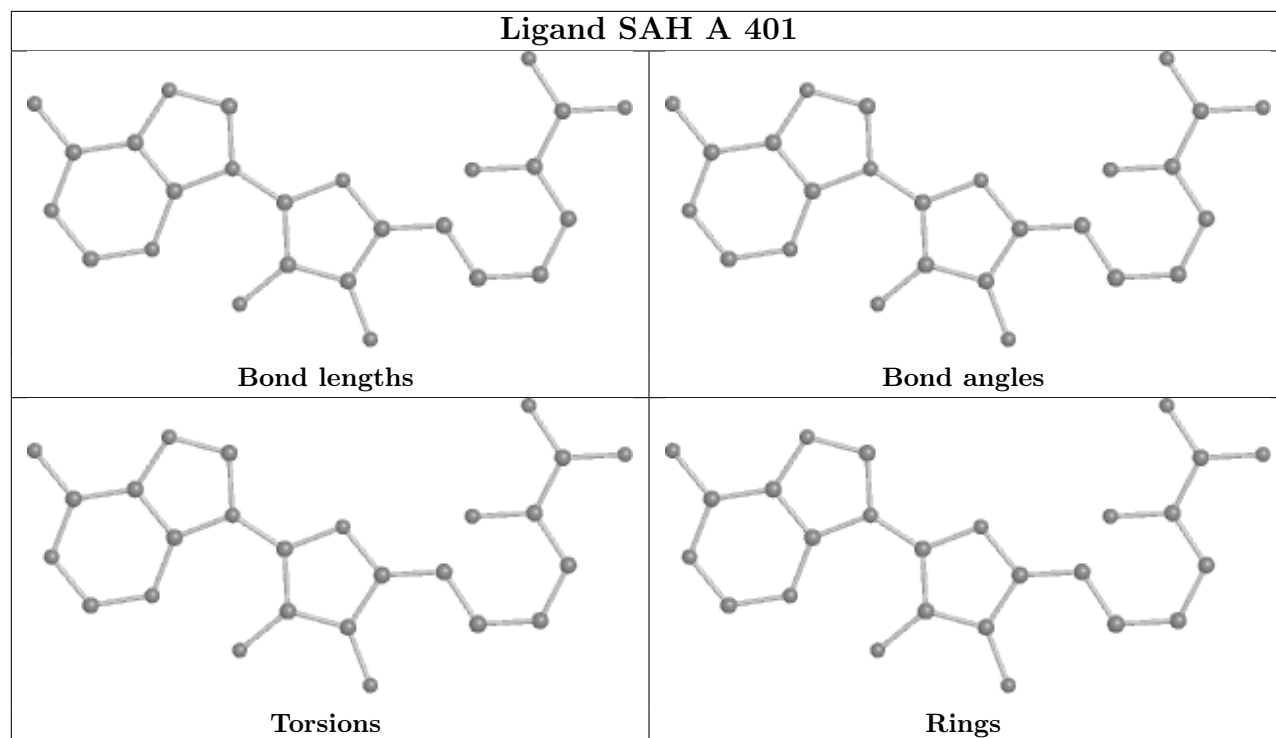
There are no torsion outliers.

There are no ring outliers.

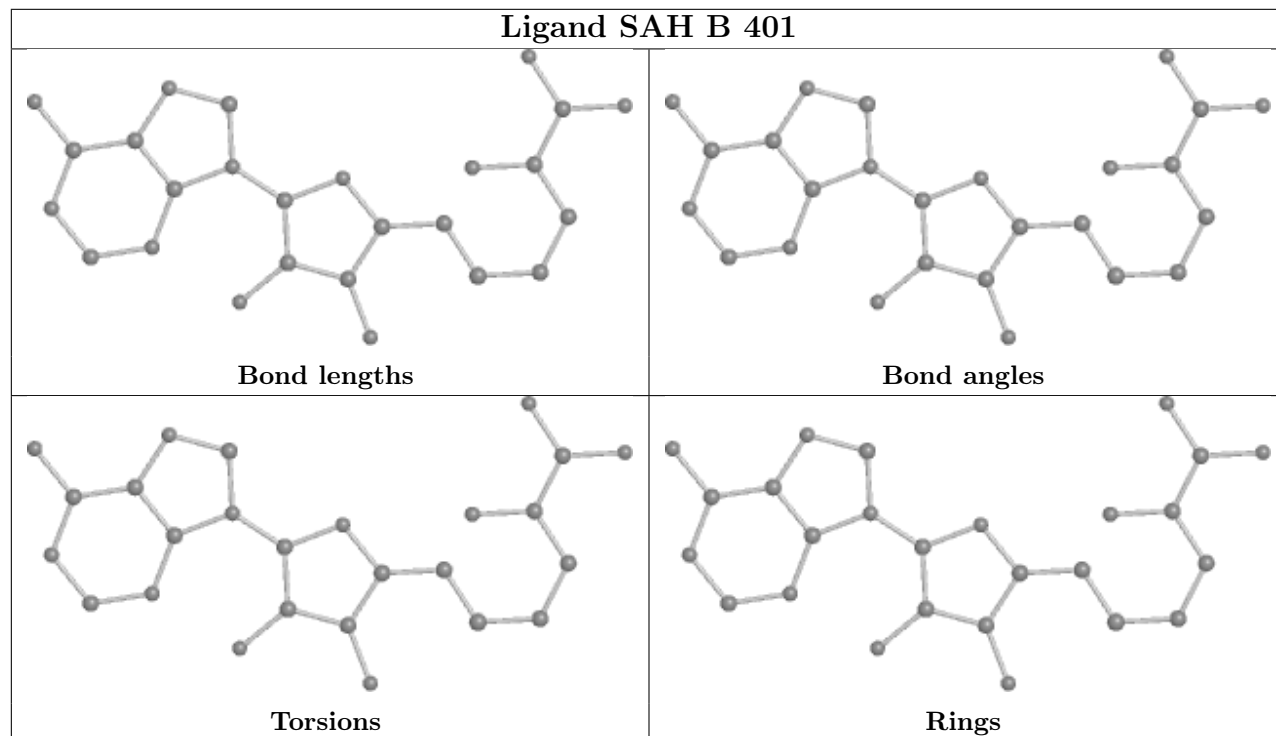
No monomer is involved in short contacts.

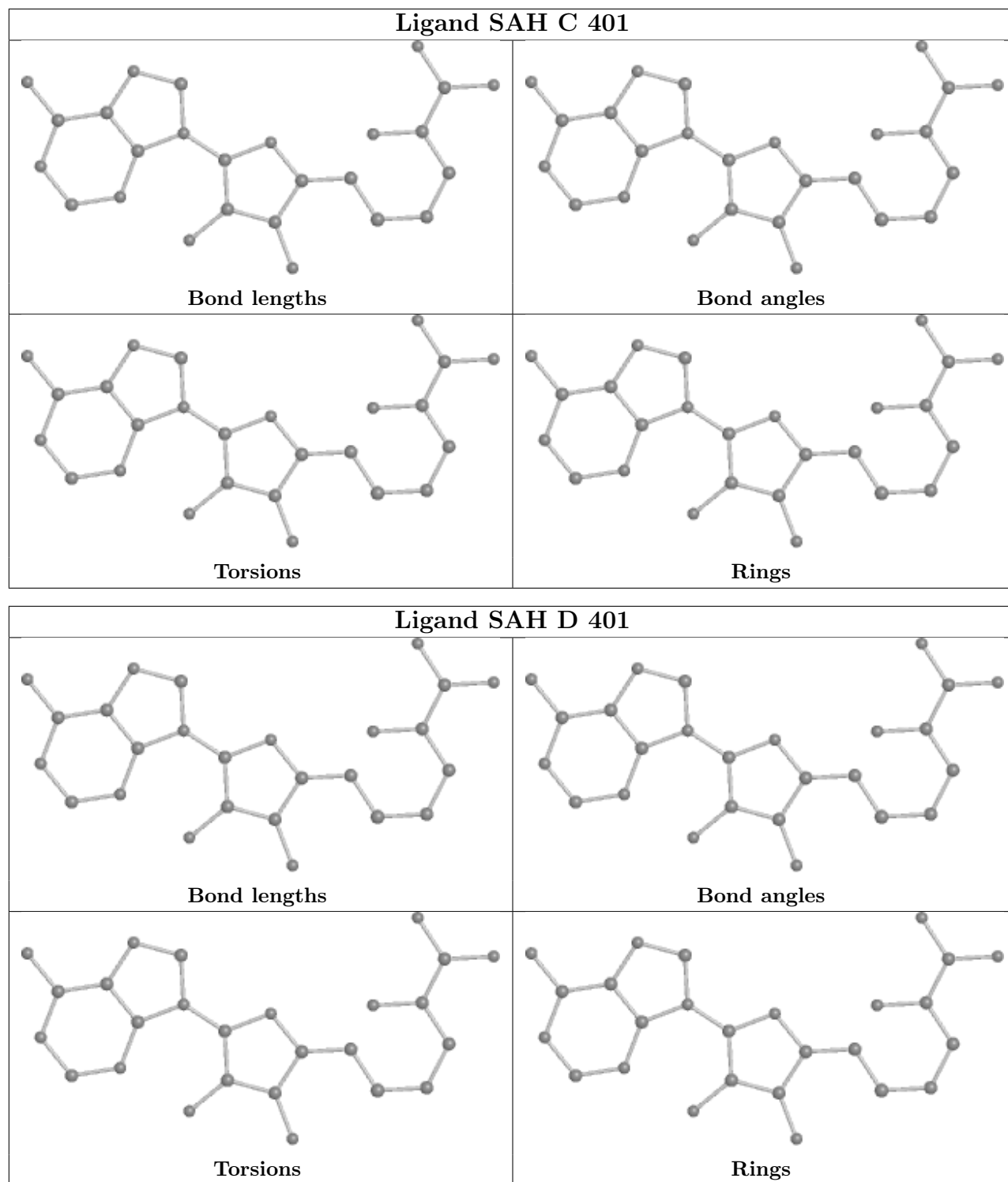
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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.

Ligand SAH A 401



Ligand SAH B 401





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	228/304 (75%)	0.15	7 (3%) 49 55	10, 23, 46, 65	0
1	B	225/304 (74%)	-0.03	11 (4%) 29 35	9, 23, 51, 69	0
1	C	210/304 (69%)	0.06	6 (2%) 51 57	9, 21, 44, 57	0
1	D	210/304 (69%)	0.35	27 (12%) 3 4	12, 26, 56, 66	0
All	All	873/1216 (71%)	0.13	51 (5%) 23 28	9, 23, 51, 69	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	301	ASP	7.5
1	D	107	ALA	5.6
1	D	203	PRO	5.3
1	D	245	TRP	5.0
1	D	109	GLY	4.6
1	D	66	VAL	4.4
1	C	245[A]	TRP	4.3
1	D	285	ILE	4.1
1	D	70	ALA	3.8
1	A	70	ALA	3.6
1	D	284	SER	3.5
1	D	244	ASN	3.5
1	D	281	HIS	3.4
1	A	143	HIS	3.4
1	D	251	ARG	3.3
1	C	203	PRO	3.3
1	D	110	LYS	3.3
1	C	179	GLU	3.3
1	D	115	GLN	3.1
1	D	286	HIS	3.1
1	D	287	ARG	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	299	GLN	3.0
1	A	298	PRO	2.9
1	B	70	ALA	2.9
1	D	108	GLU	2.8
1	D	106	ARG	2.8
1	C	244	ASN	2.8
1	D	204	SER	2.8
1	B	103	GLU	2.8
1	D	283	ALA	2.7
1	D	206	GLU	2.7
1	B	107	ALA	2.7
1	B	143	HIS	2.5
1	D	247	TYR	2.5
1	B	203	PRO	2.5
1	D	179	GLU	2.4
1	D	67	GLU	2.3
1	B	112	GLY	2.3
1	D	252	ILE	2.3
1	B	68	LEU	2.3
1	B	239	ALA	2.2
1	C	206	GLU	2.2
1	D	209	LEU	2.2
1	D	282	TRP	2.1
1	A	69	GLN	2.1
1	B	106	ARG	2.1
1	B	105	TYR	2.1
1	B	284	SER	2.0
1	A	67	GLU	2.0
1	C	178	GLU	2.0
1	D	205	ALA	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

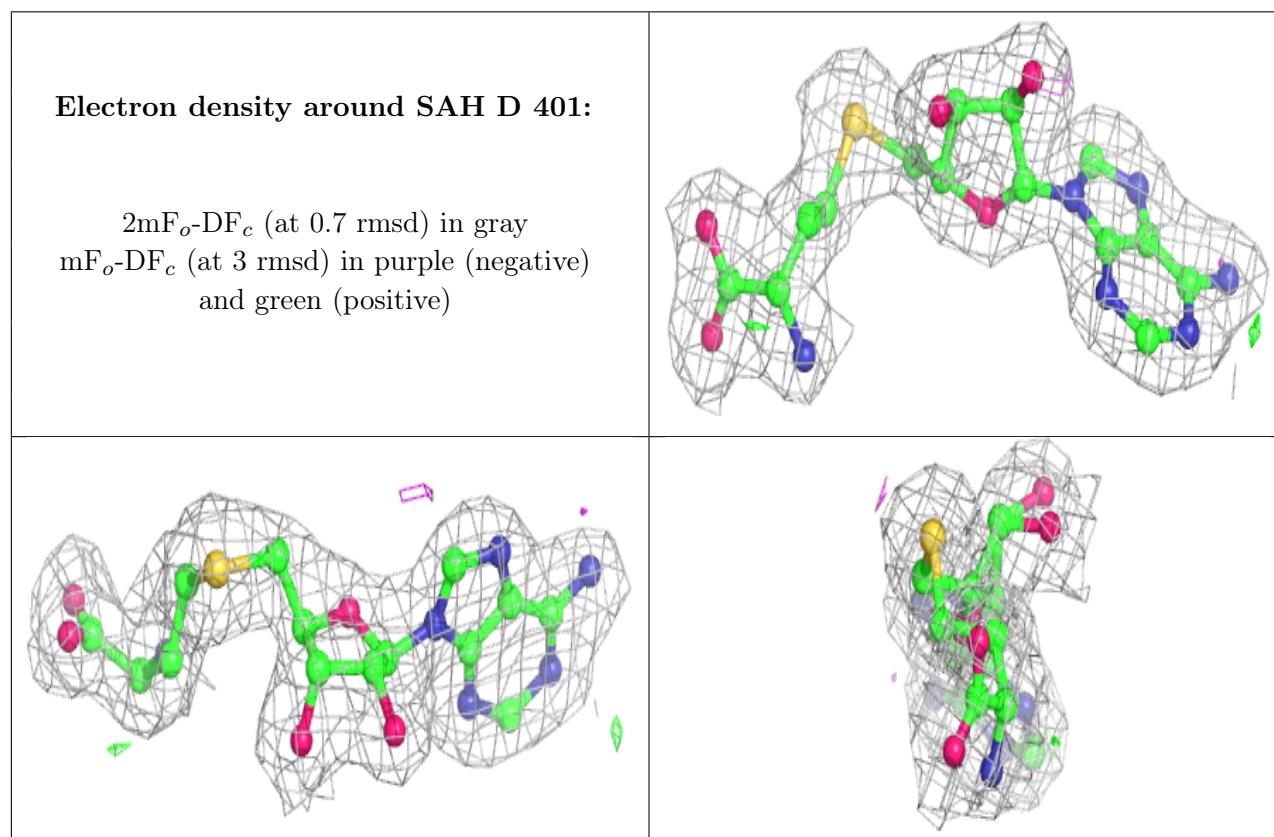
There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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.

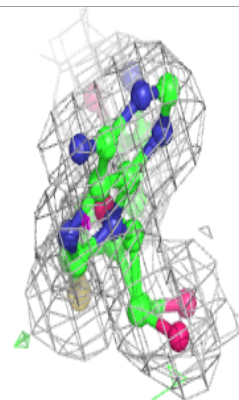
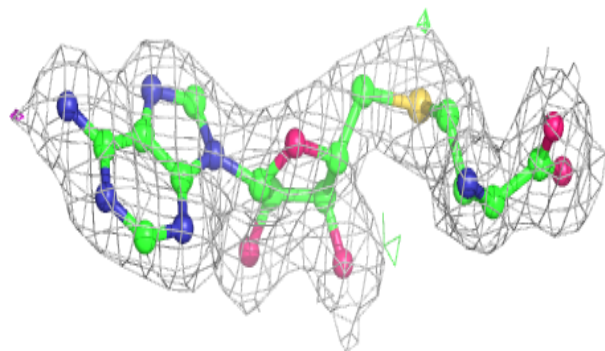
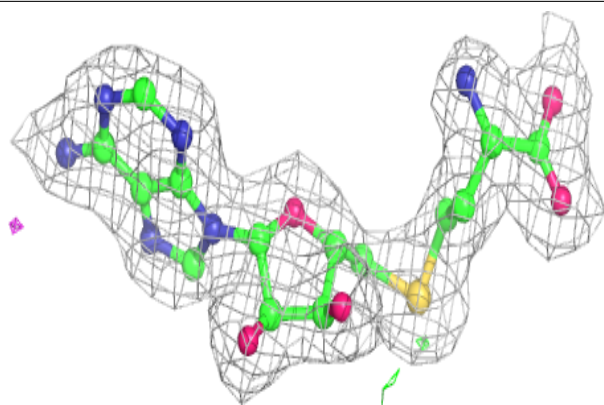
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SAH	D	401	26/26	0.96	0.09	16,20,23,25	0
2	SAH	C	401	26/26	0.97	0.10	10,16,20,22	0
2	SAH	A	401	26/26	0.97	0.10	8,14,20,22	0
2	SAH	B	401	26/26	0.98	0.09	7,13,18,24	0

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.

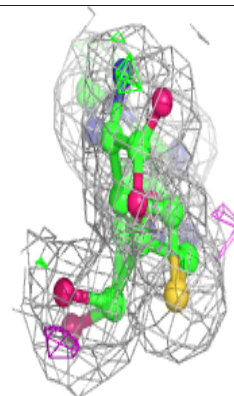
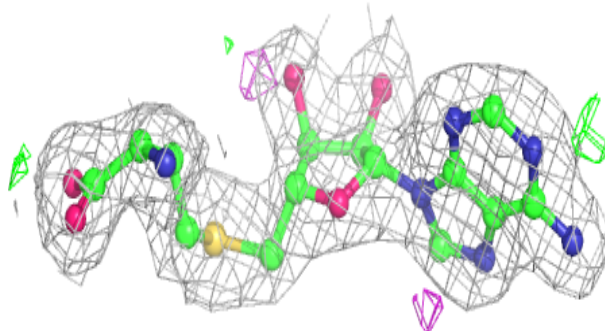
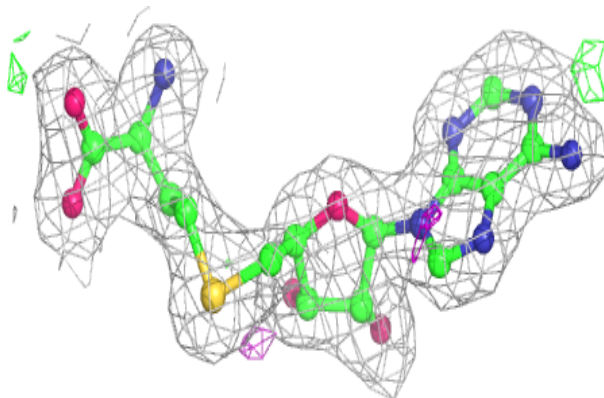


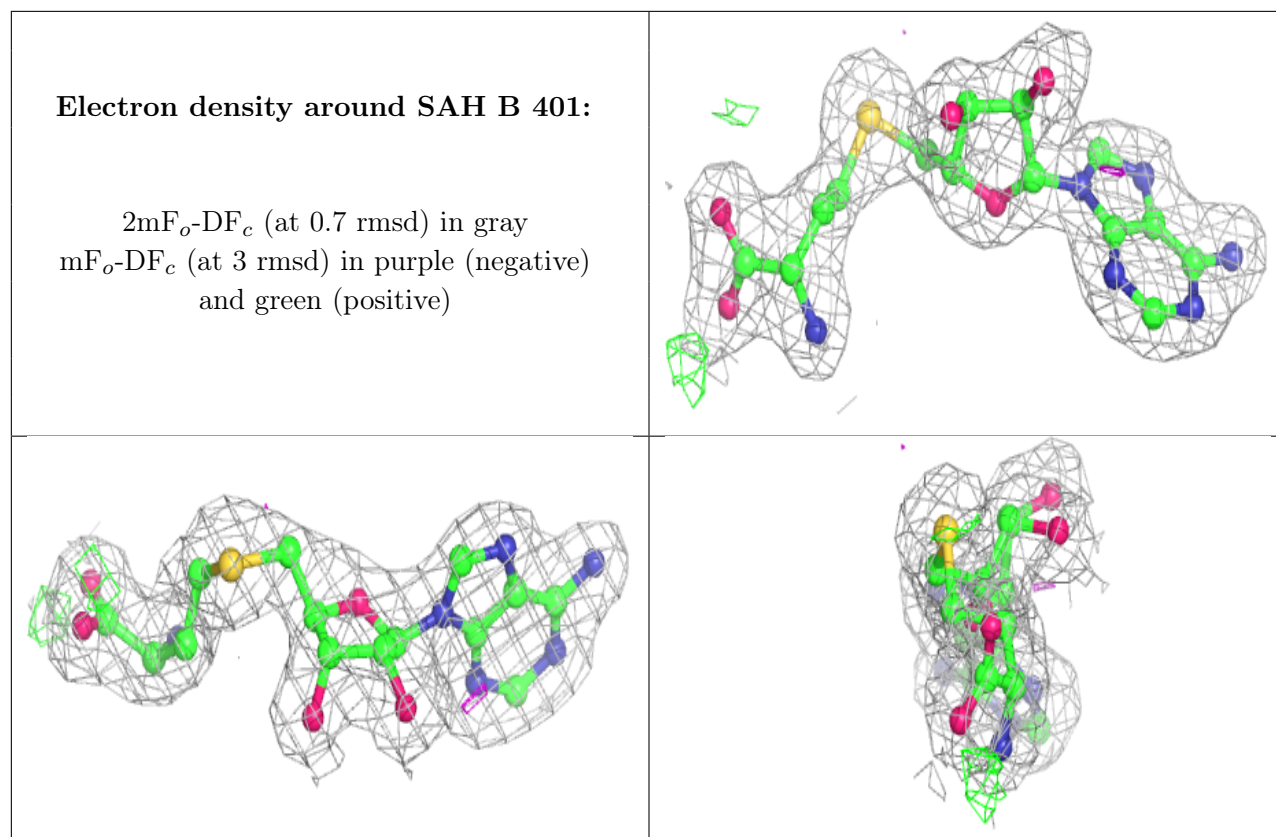
Electron density around SAH C 401:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around SAH A 401:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.