Full wwPDB X-ray Structure Validation Report ⓘ

Jun 25, 2024 – 01:04 PM EDT

PDB ID	:	5ZU0
Title	:	Proteobacterial origin of protein arginine methylation and regulation of Complex I assembly by MidA
Authors	:	Arold, S.T.; Swaminathan, K.; Hameed, U.F.S.
Deposited on	:	2018-05-05
Resolution	:	2.76 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:

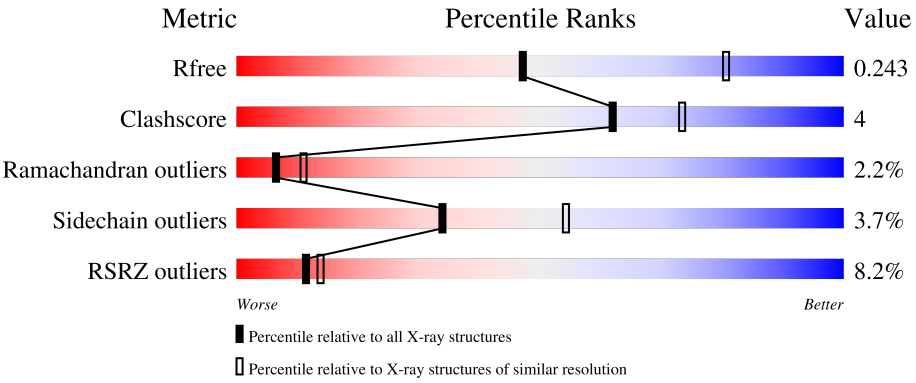
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.76 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	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)

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 <=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	414	6% 88% 8%
1	B	414	10% 87% 8%
1	C	414	7% 81% 12%

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	SAH	A	501	-	-	-	X

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 9832 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.

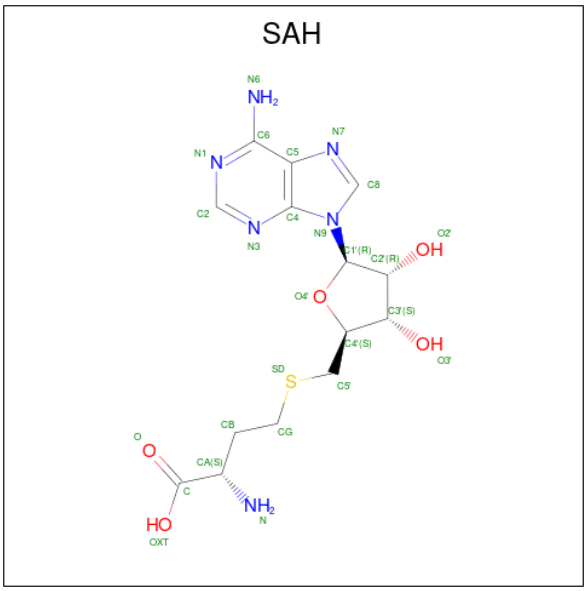
- Molecule 1 is a protein called Protein arginine methyltransferase NDUFAF7 homolog, mitochondrial.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	400	Total	C	N	O	S	Se	0	0	0
			3193	2062	519	599	3	10			
1	B	400	Total	C	N	O	S	Se	0	0	0
			3193	2062	519	599	3	10			
1	C	400	Total	C	N	O	S	Se	0	0	0
			3193	2062	519	599	3	10			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	71	GLY	-	expression tag	UNP Q54S83
A	72	PRO	-	expression tag	UNP Q54S83
A	73	LEU	-	expression tag	UNP Q54S83
A	74	GLY	-	expression tag	UNP Q54S83
A	75	SER	-	expression tag	UNP Q54S83
B	71	GLY	-	expression tag	UNP Q54S83
B	72	PRO	-	expression tag	UNP Q54S83
B	73	LEU	-	expression tag	UNP Q54S83
B	74	GLY	-	expression tag	UNP Q54S83
B	75	SER	-	expression tag	UNP Q54S83
C	71	GLY	-	expression tag	UNP Q54S83
C	72	PRO	-	expression tag	UNP Q54S83
C	73	LEU	-	expression tag	UNP Q54S83
C	74	GLY	-	expression tag	UNP Q54S83
C	75	SER	-	expression tag	UNP Q54S83

- Molecule 2 is S-ADENOSYL-L-HOMOCYSTEINE (three-letter code: SAH) (formula: $C_{14}H_{20}N_6O_5S$).



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

- Molecule 3 is water.

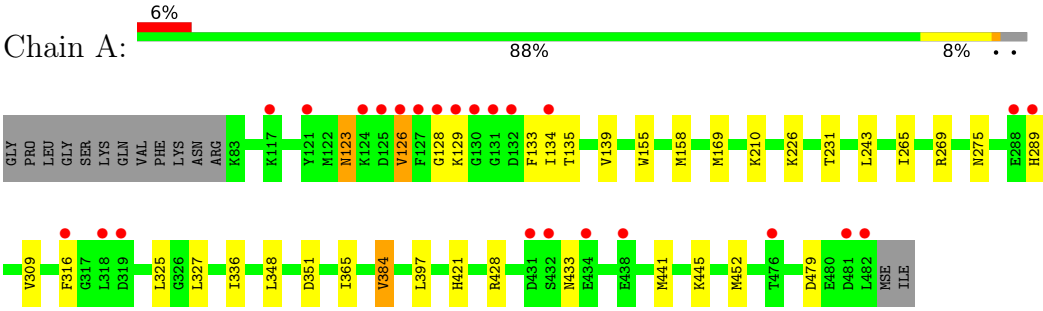
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	48	Total	O	0	0
			48	48		
3	B	37	Total	O	0	0
			37	37		
3	C	90	Total	O	0	0
			90	90		

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.

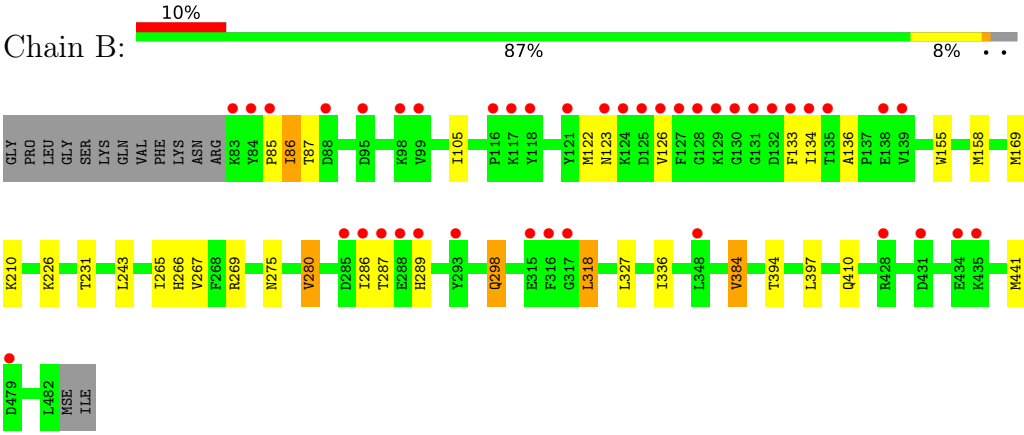
• Molecule 1: Protein arginine methyltransferase NDUFAF7 homolog, mitochondrial

Chain A:



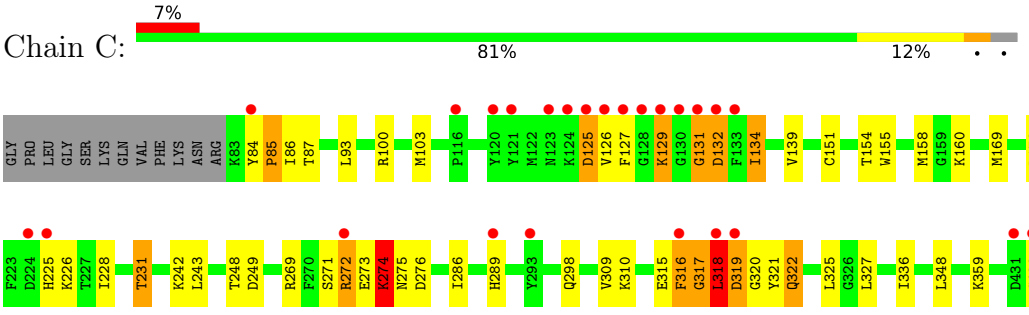
• Molecule 1: Protein arginine methyltransferase NDUFAF7 homolog, mitochondrial

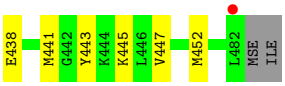
Chain B:



• Molecule 1: Protein arginine methyltransferase NDUFAF7 homolog, mitochondrial

Chain C:




E438
M441
G442
Y443
K444
K445
L446
V447
M452
L482
MSE
ILE

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	72.83Å 105.16Å 201.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.61 – 2.76 29.61 – 2.76	Depositor EDS
% Data completeness (in resolution range)	98.1 (100.61-2.76) 98.2 (29.61-2.76)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.18	Depositor
$< I/\sigma(I) >$ [1]	4.55 (at 2.76Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.200 , 0.236 0.207 , 0.243	Depositor DCC
R_{free} test set	1975 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å^2)	44.2	Xtriage
Anisotropy	0.359	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 40.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	9832	wwPDB-VP
Average B, all atoms (Å^2)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.74% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.54	0/3252	0.72	0/4374
1	B	0.49	0/3252	0.71	0/4374
1	C	0.66	0/3252	0.74	0/4374
All	All	0.57	0/9756	0.72	0/13122

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	3193	0	3225	18	0
1	B	3193	0	3225	20	0
1	C	3193	0	3225	51	0
2	A	26	0	19	0	0
2	B	26	0	19	0	0
2	C	26	0	19	0	0
3	A	48	0	0	0	0
3	B	37	0	0	0	0
3	C	90	0	0	2	0
All	All	9832	0	9732	86	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 4.

All (86) 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:C:432:SER:HB3	1:C:435:LYS:HG3	1.36	1.04
1:B:410:GLN:HE22	1:B:452:MSE:CE	1.79	0.95
1:C:155:TRP:HA	1:C:158:MSE:HE2	1.51	0.92
1:C:321:TYR:CE1	1:C:359:LYS:HE3	2.05	0.91
1:C:432:SER:CB	1:C:435:LYS:HG3	2.00	0.91
1:C:276:ASP:OD1	1:C:310:LYS:HE3	1.71	0.90
1:C:155:TRP:HA	1:C:158:MSE:CE	2.02	0.89
1:C:445:LYS:HD3	1:C:452:MSE:HE2	1.55	0.87
1:B:410:GLN:HE22	1:B:452:MSE:HE2	1.50	0.76
1:C:151:CYS:O	1:C:154:THR:HG22	1.85	0.76
1:B:267:VAL:HB	1:B:280:VAL:HG13	1.71	0.73
1:C:134:ILE:HD12	1:C:134:ILE:N	2.03	0.73
1:C:317:GLY:O	1:C:318:LEU:HB2	1.92	0.69
1:B:410:GLN:HE22	1:B:452:MSE:HE3	1.59	0.68
1:C:231:THR:HG21	3:C:604:HOH:O	1.94	0.67
1:A:265:ILE:HD13	1:A:384:VAL:HG21	1.78	0.66
1:B:265:ILE:HD13	1:B:384:VAL:HG21	1.77	0.66
1:C:139:VAL:HG21	1:C:452:MSE:HE1	1.77	0.66
1:A:421:HIS:CD2	1:B:397:LEU:HD13	2.32	0.65
1:C:155:TRP:CA	1:C:158:MSE:HE2	2.26	0.64
1:C:272:ARG:O	1:C:272:ARG:HG2	1.97	0.64
1:C:272:ARG:HB2	1:C:319:ASP:HA	1.80	0.63
1:C:131:GLY:O	1:C:132:ASP:HB3	1.99	0.63
1:A:128:GLY:O	1:A:129:LYS:HG3	1.98	0.62
1:C:321:TYR:HE1	1:C:359:LYS:HE3	1.64	0.62
1:A:139:VAL:HG21	1:A:452:MSE:HE1	1.81	0.62
1:C:155:TRP:CH2	1:C:160:LYS:HG2	2.36	0.60
1:C:322:GLN:NE2	3:C:602:HOH:O	2.34	0.60
1:A:479:ASP:N	1:A:479:ASP:OD1	2.33	0.60
1:A:134:ILE:HG21	1:A:445:LYS:HE2	1.84	0.59
1:C:271:SER:C	1:C:273:GLU:H	2.04	0.59
1:C:169:MSE:HE1	1:C:243:LEU:HD22	1.83	0.59
1:C:129:LYS:HG3	1:C:173:ARG:HH12	1.67	0.59
1:C:154:THR:O	1:C:158:MSE:HG3	2.02	0.59
1:B:155:TRP:HA	1:B:158:MSE:HE3	1.85	0.58
1:C:248:THR:O	1:C:249:ASP:HB2	2.04	0.57
1:C:309:VAL:HG11	1:C:327:LEU:HD21	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:PHE:HZ	1:A:365:ILE:HG21	1.70	0.56
1:B:410:GLN:NE2	1:B:452:MSE:HE2	2.20	0.56
1:C:274:LYS:O	1:C:275:ASN:C	2.43	0.56
1:B:169:MSE:HE1	1:B:243:LEU:HD22	1.89	0.54
1:A:126:VAL:HG11	1:C:269:ARG:HH22	1.72	0.54
1:C:155:TRP:HA	1:C:158:MSE:HE3	1.86	0.54
1:C:139:VAL:HG21	1:C:452:MSE:CE	2.39	0.53
1:A:309:VAL:HG11	1:A:327:LEU:HD11	1.91	0.53
1:A:155:TRP:HA	1:A:158:MSE:HE3	1.91	0.52
1:A:169:MSE:HE3	1:A:336:ILE:HD11	1.92	0.52
1:C:134:ILE:HD12	1:C:134:ILE:H	1.72	0.52
1:C:169:MSE:HE3	1:C:336:ILE:HD11	1.92	0.52
1:A:139:VAL:HG21	1:A:452:MSE:CE	2.40	0.51
1:C:273:GLU:C	1:C:275:ASN:H	2.15	0.50
1:A:269:ARG:HD3	1:A:298:GLN:HE21	1.76	0.50
1:B:269:ARG:HD3	1:B:298:GLN:HE21	1.76	0.50
1:B:210:LYS:HE2	1:B:226:LYS:O	2.11	0.50
1:A:169:MSE:HE1	1:A:243:LEU:HD22	1.92	0.50
1:C:309:VAL:HG11	1:C:327:LEU:CD2	2.41	0.50
1:B:169:MSE:HE3	1:B:336:ILE:HD11	1.94	0.49
1:C:432:SER:OG	1:C:435:LYS:HG3	2.12	0.49
1:A:210:LYS:HE2	1:A:226:LYS:O	2.14	0.48
1:B:179:ASP:OD1	1:B:182:ARG:NH1	2.46	0.48
1:C:271:SER:C	1:C:273:GLU:N	2.68	0.48
1:C:271:SER:O	1:C:273:GLU:N	2.47	0.47
1:A:133:PHE:CZ	1:A:365:ILE:HG21	2.50	0.47
1:C:321:TYR:HE1	1:C:359:LYS:CE	2.26	0.47
1:C:321:TYR:CD1	1:C:359:LYS:HE3	2.48	0.46
1:C:125:ASP:HB2	1:C:127:PHE:CE2	2.51	0.46
1:C:316:PHE:CD1	1:C:316:PHE:O	2.69	0.46
1:C:226:LYS:HE3	1:C:228:ILE:O	2.17	0.45
1:C:274:LYS:H	1:C:274:LYS:HG2	1.50	0.45
1:B:267:VAL:HG12	1:B:280:VAL:CG1	2.48	0.44
1:B:266:HIS:CD2	1:B:327:LEU:HD12	2.53	0.44
1:C:321:TYR:CE1	1:C:359:LYS:CE	2.88	0.43
1:C:316:PHE:CD1	1:C:316:PHE:C	2.90	0.43
1:C:286:ILE:O	1:C:286:ILE:HD12	2.18	0.42
1:B:105:ILE:HD11	1:B:280:VAL:HG22	2.01	0.42
1:C:155:TRP:CZ3	1:C:160:LYS:HG2	2.54	0.42
1:C:443:TYR:O	1:C:447:VAL:HG22	2.20	0.42
1:C:316:PHE:O	1:C:316:PHE:HD1	2.03	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:428:ARG:NH1	1:B:394:THR:HG23	2.34	0.42
1:C:93:LEU:HB3	1:C:103:MSE:HE1	2.02	0.41
1:C:100:ARG:HE	1:C:100:ARG:HB3	1.67	0.41
1:A:135:THR:O	1:A:139:VAL:HG23	2.21	0.41
1:B:136:ALA:HB2	1:B:452:MSE:HE1	2.03	0.41
1:B:267:VAL:CG1	1:B:280:VAL:CG1	3.00	0.40
1:B:269:ARG:CD	1:B:298:GLN:HE21	2.33	0.40
1:C:84:TYR:HB2	1:C:85:PRO:CD	2.52	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	398/414 (96%)	378 (95%)	16 (4%)	4 (1%)	15 27
1	B	398/414 (96%)	372 (94%)	16 (4%)	10 (2%)	5 9
1	C	398/414 (96%)	366 (92%)	20 (5%)	12 (3%)	4 6
All	All	1194/1242 (96%)	1116 (94%)	52 (4%)	26 (2%)	6 11

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	275	ASN
1	B	85	PRO
1	B	87	THR
1	B	275	ASN
1	B	287	THR
1	C	87	THR
1	C	132	ASP
1	C	318	LEU
1	A	123	ASN

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Mol	Chain	Res	Type
1	B	86	ILE
1	B	122	MSE
1	B	123	ASN
1	C	86	ILE
1	C	125	ASP
1	C	272	ARG
1	A	126	VAL
1	B	126	VAL
1	B	318	LEU
1	C	126	VAL
1	C	317	GLY
1	C	320	GLY
1	A	316	PHE
1	C	85	PRO
1	C	131	GLY
1	C	274	LYS
1	B	134	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	351/352 (100%)	340 (97%)	11 (3%)	40	60
1	B	351/352 (100%)	341 (97%)	10 (3%)	43	63
1	C	351/352 (100%)	333 (95%)	18 (5%)	24	41
All	All	1053/1056 (100%)	1014 (96%)	39 (4%)	34	54

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	ASN
1	A	231	THR
1	A	289	HIS
1	A	298	GLN
1	A	325	LEU

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Mol	Chain	Res	Type
1	A	348	LEU
1	A	351	ASP
1	A	384	VAL
1	A	397	LEU
1	A	433	ASN
1	A	441	MSE
1	B	86	ILE
1	B	133	PHE
1	B	231	THR
1	B	280	VAL
1	B	286	ILE
1	B	289	HIS
1	B	298	GLN
1	B	318	LEU
1	B	384	VAL
1	B	441	MSE
1	C	129	LYS
1	C	134	ILE
1	C	219	LYS
1	C	225	HIS
1	C	231	THR
1	C	242	LYS
1	C	274	LYS
1	C	289	HIS
1	C	298	GLN
1	C	315	GLU
1	C	316	PHE
1	C	318	LEU
1	C	319	ASP
1	C	322	GLN
1	C	325	LEU
1	C	348	LEU
1	C	438	GLU
1	C	441	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	298	GLN
1	A	421	HIS
1	B	298	GLN
1	B	410	GLN

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Mol	Chain	Res	Type
1	C	322	GLN

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

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	$\#\|Z\| > 2$
2	SAH	B	501	-	24,28,28	1.22	4 (16%)	25,40,40	1.49	6 (24%)
2	SAH	C	501	-	24,28,28	1.29	3 (12%)	25,40,40	1.43	5 (20%)
2	SAH	A	501	-	24,28,28	1.21	3 (12%)	25,40,40	1.50	5 (20%)

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	SAH	B	501	-	-	1/11/31/31	0/3/3/3
2	SAH	C	501	-	-	1/11/31/31	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SAH	A	501	-	-	1/11/31/31	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	SAH	C5-C4	2.87	1.48	1.40
2	A	501	SAH	C5-C4	2.82	1.48	1.40
2	B	501	SAH	C5-C4	2.79	1.48	1.40
2	B	501	SAH	C2-N3	2.57	1.36	1.32
2	A	501	SAH	OXT-C	-2.54	1.22	1.30
2	A	501	SAH	C2-N3	2.44	1.36	1.32
2	C	501	SAH	C2-N3	2.20	1.35	1.32
2	C	501	SAH	O4'-C1'	2.19	1.44	1.41
2	B	501	SAH	O4'-C1'	2.05	1.43	1.41
2	B	501	SAH	OXT-C	-2.05	1.23	1.30

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	501	SAH	N3-C2-N1	-3.91	122.57	128.68
2	A	501	SAH	N3-C2-N1	-3.80	122.74	128.68
2	C	501	SAH	N3-C2-N1	-3.72	122.87	128.68
2	A	501	SAH	OXT-C-O	-2.82	117.69	124.09
2	B	501	SAH	CB-CG-SD	-2.76	107.11	113.31
2	B	501	SAH	OXT-C-O	-2.75	117.84	124.09
2	C	501	SAH	CB-CG-SD	-2.66	107.33	113.31
2	A	501	SAH	CB-CG-SD	-2.58	107.52	113.31
2	C	501	SAH	OXT-C-O	-2.57	118.26	124.09
2	A	501	SAH	OXT-C-CA	2.37	121.45	113.38
2	B	501	SAH	C4-C5-N7	-2.29	107.02	109.40
2	C	501	SAH	C4-C5-N7	-2.24	107.06	109.40
2	A	501	SAH	C4-C5-N7	-2.23	107.08	109.40
2	C	501	SAH	C2-N1-C6	2.05	122.27	118.75
2	B	501	SAH	OXT-C-CA	2.04	120.35	113.38
2	B	501	SAH	C2-N1-C6	2.02	122.21	118.75

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	501	SAH	CB-CG-SD-C5'

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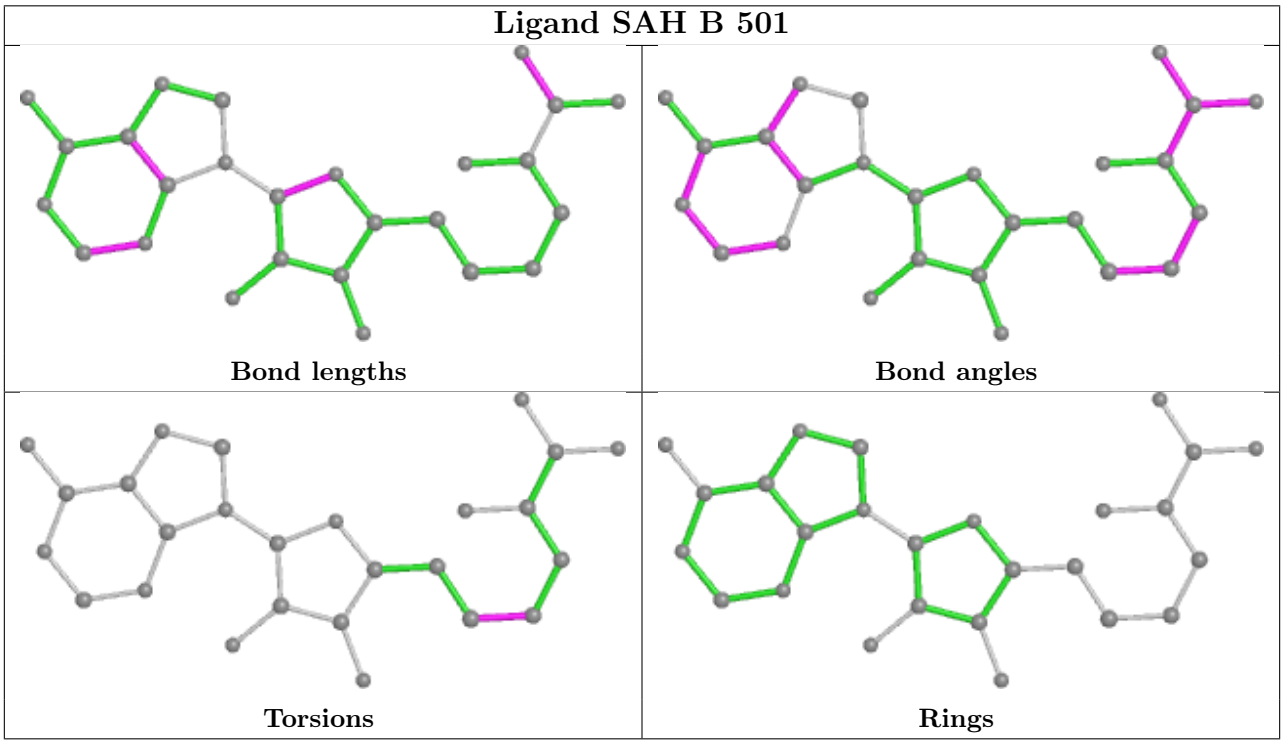
Mol	Chain	Res	Type	Atoms
2	C	501	SAH	CB-CG-SD-C5'
2	A	501	SAH	CB-CG-SD-C5'

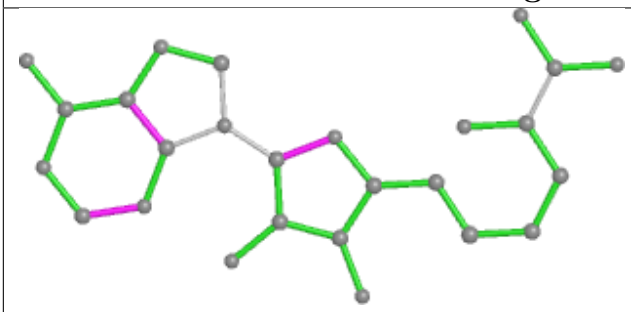
There are no ring outliers.

No monomer is involved in short contacts.

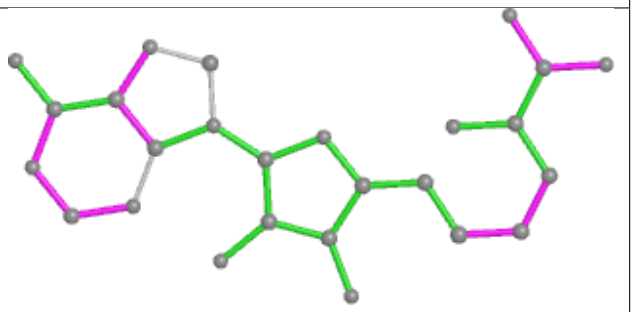
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



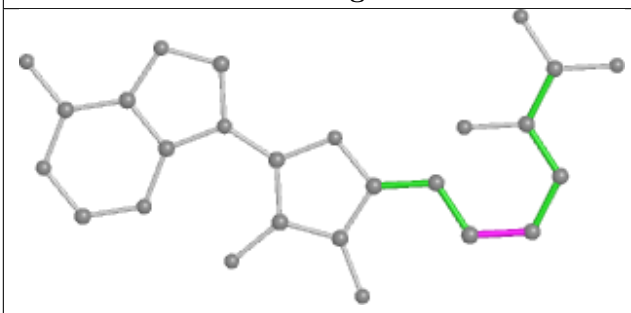
Ligand SAH C 501



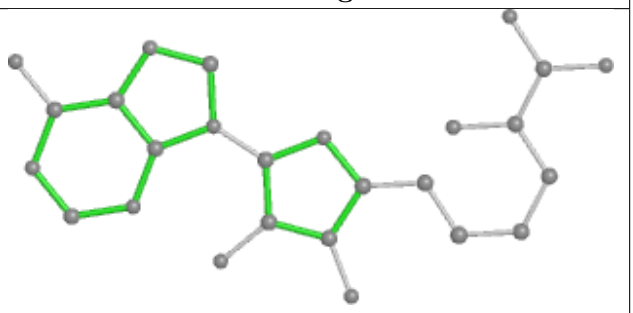
Bond lengths



Bond angles

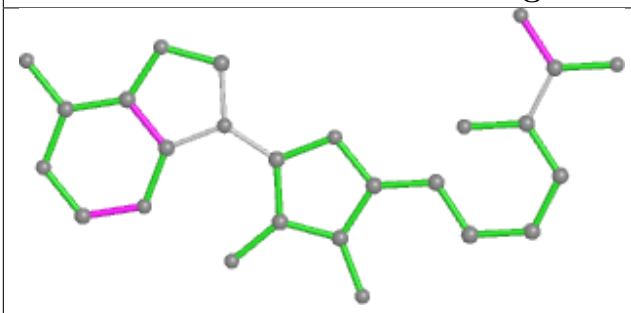


Torsions

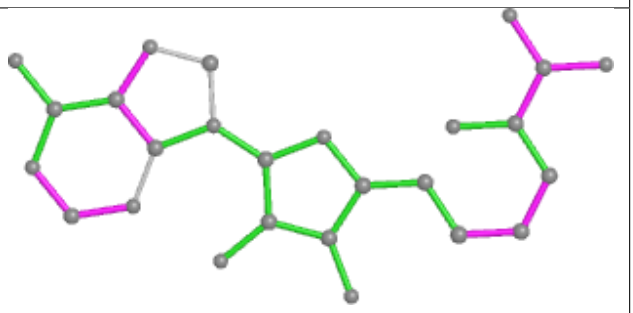


Rings

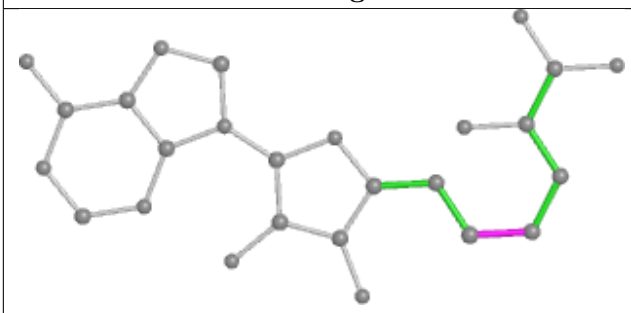
Ligand SAH A 501



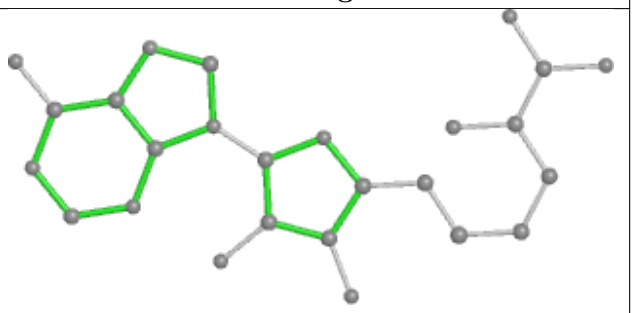
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers

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, 95^{th} 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	390/414 (94%)	0.17	25 (6%)	19	23	30, 51, 88, 162	0
1	B	390/414 (94%)	0.48	42 (10%)	5	6	30, 52, 123, 171	0
1	C	390/414 (94%)	0.30	29 (7%)	14	17	24, 44, 108, 182	0
All	All	1170/1242 (94%)	0.32	96 (8%)	11	14	24, 50, 116, 182	0

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	128	GLY	15.5
1	B	131	GLY	15.2
1	C	131	GLY	14.5
1	B	132	ASP	13.1
1	C	128	GLY	10.6
1	C	130	GLY	10.0
1	C	126	VAL	9.8
1	C	127	PHE	9.3
1	C	129	LYS	9.3
1	A	129	LYS	9.1
1	B	130	GLY	8.9
1	A	127	PHE	7.5
1	B	127	PHE	7.3
1	A	316	PHE	6.3
1	C	133	PHE	6.2
1	A	125	ASP	6.0
1	C	124	LYS	5.9
1	C	132	ASP	5.9
1	C	125	ASP	5.7
1	B	133	PHE	5.5
1	A	130	GLY	5.4
1	B	286	ILE	5.3
1	C	431	ASP	5.3

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Mol	Chain	Res	Type	RSRZ
1	B	118	TYR	5.2
1	B	287	THR	5.1
1	B	431	ASP	4.9
1	A	126	VAL	4.8
1	B	85	PRO	4.7
1	C	272	ARG	4.7
1	B	125	ASP	4.5
1	B	316	PHE	4.4
1	A	289	HIS	4.4
1	B	129	LYS	4.3
1	A	134	ILE	4.3
1	B	84	TYR	4.1
1	C	319	ASP	4.1
1	A	431	ASP	4.0
1	A	132	ASP	4.0
1	B	134	ILE	4.0
1	B	121	TYR	3.9
1	A	131	GLY	3.9
1	B	434	GLU	3.9
1	B	95	ASP	3.9
1	C	225	HIS	3.8
1	B	138	GLU	3.6
1	A	128	GLY	3.6
1	C	121	TYR	3.4
1	A	300	LYS	3.4
1	C	120	TYR	3.4
1	B	99	VAL	3.4
1	C	222	ASN	3.3
1	B	288	GLU	3.3
1	A	124	LYS	3.2
1	A	288	GLU	3.1
1	A	482	LEU	3.1
1	A	434	GLU	2.9
1	A	438	GLU	2.9
1	B	289	HIS	2.9
1	C	316	PHE	2.9
1	B	348	LEU	2.9
1	C	432	SER	2.8
1	B	479	ASP	2.8
1	C	289	HIS	2.8
1	B	139	VAL	2.8
1	C	482	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	476	THR	2.7
1	B	123	ASN	2.7
1	A	432	SER	2.6
1	A	318	LEU	2.6
1	B	126	VAL	2.6
1	B	315	GLU	2.5
1	B	428	ARG	2.5
1	B	124	LYS	2.5
1	B	116	PRO	2.5
1	C	84	TYR	2.5
1	B	293	TYR	2.4
1	C	318	LEU	2.4
1	B	285	ASP	2.4
1	C	293	TYR	2.4
1	B	135	THR	2.4
1	A	319	ASP	2.4
1	B	98	LYS	2.4
1	C	116	PRO	2.4
1	B	117	LYS	2.3
1	C	434	GLU	2.3
1	B	435	LYS	2.3
1	C	123	ASN	2.3
1	A	121	TYR	2.3
1	B	88	ASP	2.2
1	B	317	GLY	2.2
1	A	117	LYS	2.2
1	C	433	ASN	2.2
1	A	481	ASP	2.1
1	B	450	LYS	2.1
1	C	224	ASP	2.0
1	B	83	LYS	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)

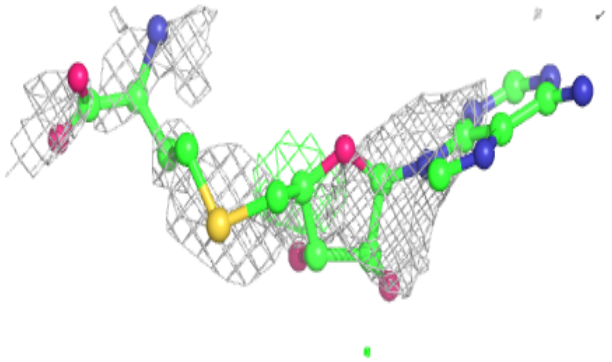
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

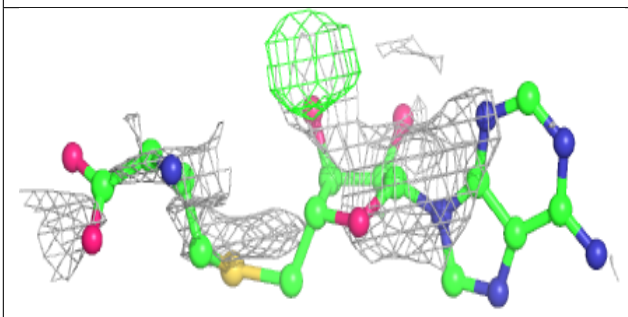
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	SAH	A	501	26/26	0.66	0.42	105,154,161,162	0
2	SAH	C	501	26/26	0.77	0.32	99,110,121,123	0
2	SAH	B	501	26/26	0.82	0.30	75,96,107,109	0

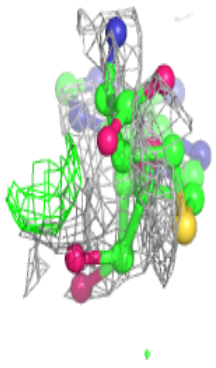
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around SAH A 501:

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

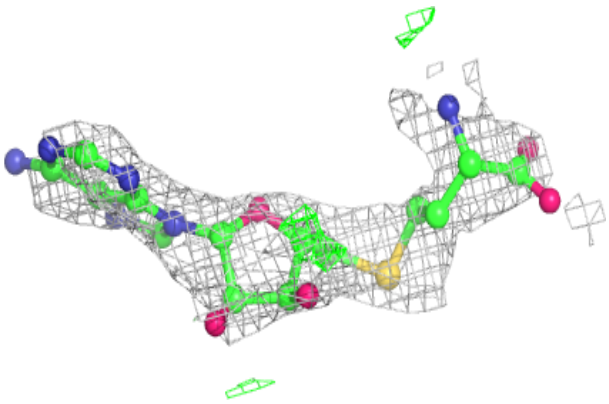


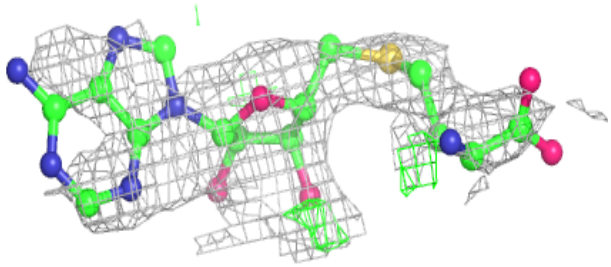


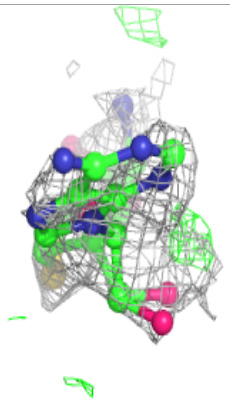


Electron density around SAH C 501:

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

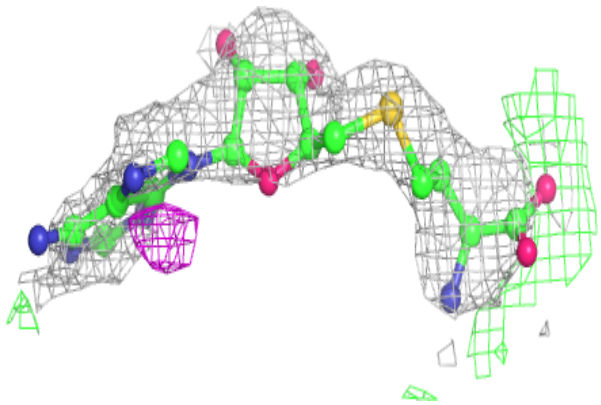


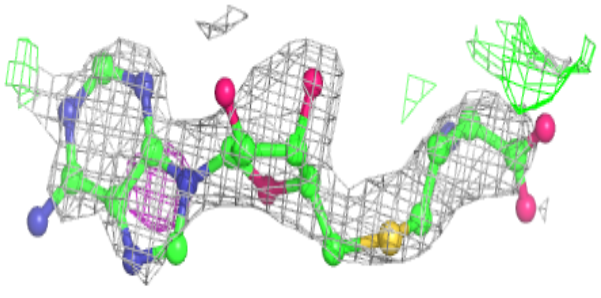


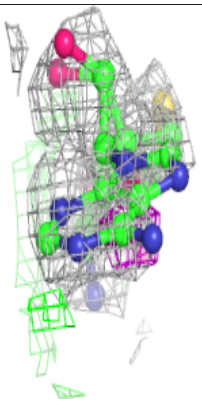


Electron density around SAH B 501:

$2mF_o$-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.