



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:38 AM UTC

PDB ID : 3NUT / pdb_00003nut
Title : Crystal structure of the methyltransferase CobJ
Authors : Pickersgill, R.W.; Hutchison, T.; To, T.T.
Deposited on : 2010-07-07
Resolution : 2.22 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

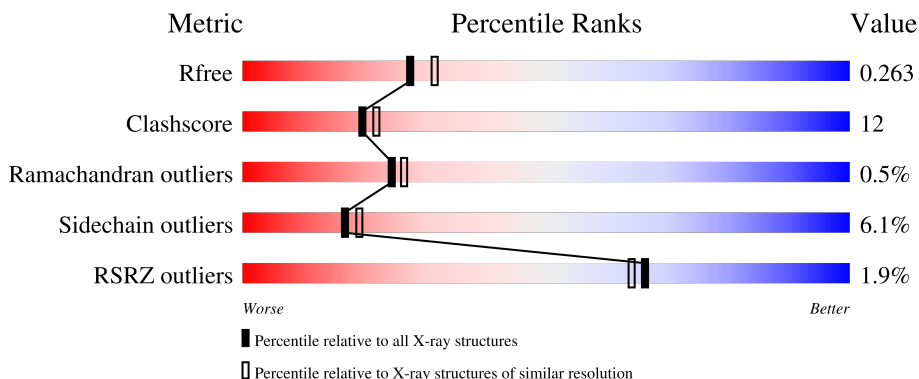
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.22 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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	251	 2% (red), 78% (green), 16% (yellow), 2% (orange), 2% (grey)
1	B	251	 72% (green), 20% (yellow), 2% (orange), 2% (grey)
1	C	251	 2% (red), 80% (green), 16% (yellow), 2% (orange), 2% (grey)
1	D	251	 3% (red), 68% (green), 24% (yellow), 2% (orange), 5% (grey)

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
3	GOL	B	246	-	-	X	-
3	GOL	C	246	-	-	X	-
3	GOL	C	247	-	-	X	-

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 7643 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 Precorrin-3 methylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	240	Total 1824	C 1149	N 332	O 335	S 8	0	0	0
1	B	240	Total 1822	C 1148	N 332	O 334	S 8	0	0	0
1	C	242	Total 1835	C 1155	N 334	O 337	S 9	0	0	0
1	D	239	Total 1813	C 1143	N 328	O 334	S 8	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

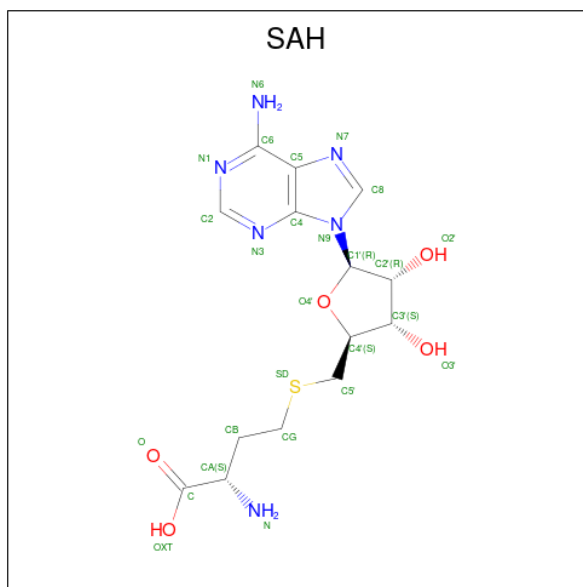
Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	expression tag	UNP O68097
A	-4	HIS	-	expression tag	UNP O68097
A	-3	HIS	-	expression tag	UNP O68097
A	-2	HIS	-	expression tag	UNP O68097
A	-1	HIS	-	expression tag	UNP O68097
A	0	HIS	-	expression tag	UNP O68097
B	-5	HIS	-	expression tag	UNP O68097
B	-4	HIS	-	expression tag	UNP O68097
B	-3	HIS	-	expression tag	UNP O68097
B	-2	HIS	-	expression tag	UNP O68097
B	-1	HIS	-	expression tag	UNP O68097
B	0	HIS	-	expression tag	UNP O68097
C	-5	HIS	-	expression tag	UNP O68097
C	-4	HIS	-	expression tag	UNP O68097
C	-3	HIS	-	expression tag	UNP O68097
C	-2	HIS	-	expression tag	UNP O68097
C	-1	HIS	-	expression tag	UNP O68097
C	0	HIS	-	expression tag	UNP O68097
D	-5	HIS	-	expression tag	UNP O68097
D	-4	HIS	-	expression tag	UNP O68097
D	-3	HIS	-	expression tag	UNP O68097

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	HIS	-	expression tag	UNP O68097
D	-1	HIS	-	expression tag	UNP O68097
D	0	HIS	-	expression tag	UNP O68097

- Molecule 2 is S-ADENOSYL-L-HOMOCYSTEINE (CCD ID: SAH) (formula: C₁₄H₂₀N₆O₅S).



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

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	C	1	Total C O 6 3 3	0	0
3	C	1	Total C O 6 3 3	0	0

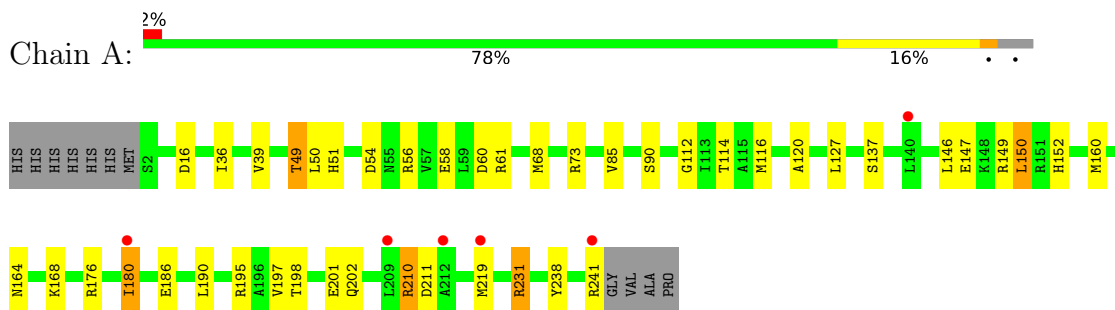
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	46	Total O 46 46	0	0
4	B	54	Total O 54 54	0	0
4	C	59	Total O 59 59	0	0
4	D	38	Total O 38 38	0	0

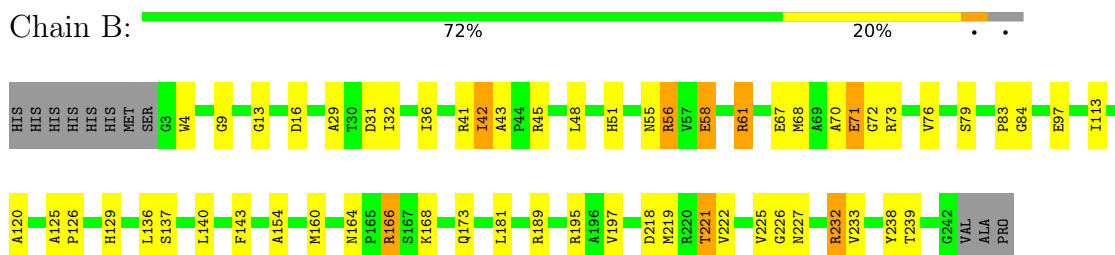
3 Residue-property plots [i](#)

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.

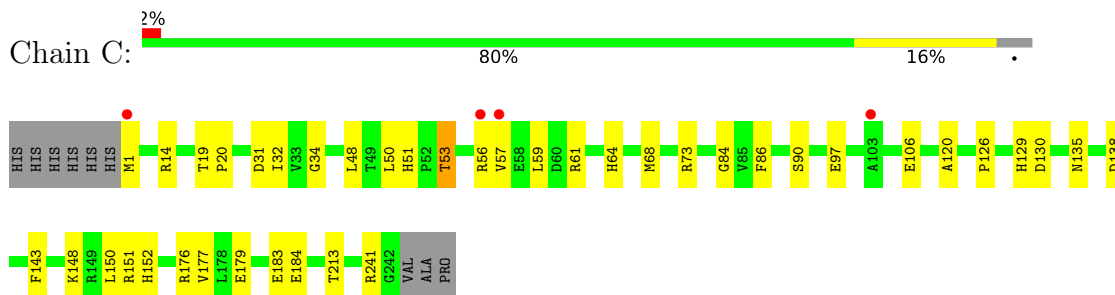
- Molecule 1: Precorrin-3 methylase



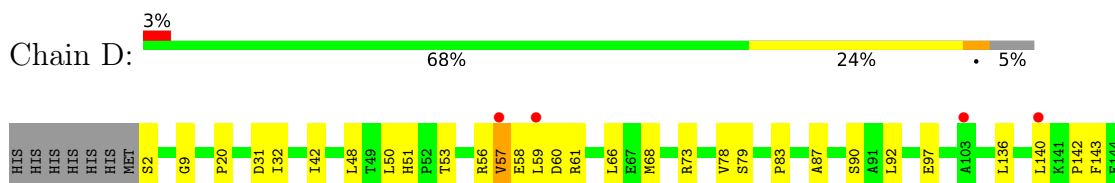
- Molecule 1: Precorrin-3 methylase

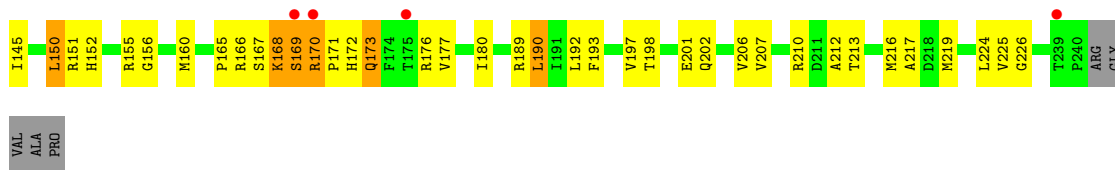


- Molecule 1: Precorrin-3 methylase



- Molecule 1: Precorrin-3 methylase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	85.64Å 110.99Å 116.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	58.56 – 2.22 58.56 – 2.22	Depositor EDS
% Data completeness (in resolution range)	99.1 (58.56-2.22) 99.1 (58.56-2.22)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 2.22Å)	Xtrriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.216 , 0.264 0.215 , 0.263	Depositor DCC
R_{free} test set	2773 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	37.8	Xtrriage
Anisotropy	0.126	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 32.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.000 for -h,l,k	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7643	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.95% 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: GOL, 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	0.82	1/1863 (0.1%)	1.00	3/2539 (0.1%)
1	B	0.92	2/1861 (0.1%)	1.05	4/2536 (0.2%)
1	C	0.82	0/1874	0.99	3/2552 (0.1%)
1	D	0.80	0/1852	0.98	2/2525 (0.1%)
All	All	0.84	3/7450 (0.0%)	1.00	12/10152 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	36	ILE	CA-CB	8.34	1.59	1.53
1	B	154	ALA	CA-CB	-5.32	1.45	1.53
1	A	112	GLY	N-CA	5.19	1.50	1.45

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	4	TRP	N-CA-C	6.11	117.52	108.60
1	B	136	LEU	CA-C-N	-5.36	115.15	122.87
1	B	136	LEU	C-N-CA	-5.36	115.15	122.87
1	C	213	THR	CA-C-N	5.26	125.32	119.32
1	C	213	THR	C-N-CA	5.26	125.32	119.32
1	C	34	GLY	N-CA-C	5.19	118.33	110.75
1	A	231	ARG	N-CA-C	5.17	115.11	108.34
1	A	186	GLU	CA-C-N	5.15	125.19	119.32
1	A	186	GLU	C-N-CA	5.15	125.19	119.32
1	D	170	ARG	CA-C-N	5.12	125.41	119.47
1	D	170	ARG	C-N-CA	5.12	125.41	119.47
1	B	154	ALA	N-CA-C	5.03	116.45	111.07

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	1824	0	1830	31	0
1	B	1822	0	1828	69	0
1	C	1835	0	1841	39	0
1	D	1813	0	1817	42	0
2	A	26	0	19	1	0
2	B	26	0	19	3	0
2	C	26	0	19	1	0
2	D	26	0	19	0	0
3	A	18	0	24	2	0
3	B	18	0	24	14	0
3	C	12	0	16	14	0
4	A	46	0	0	5	0
4	B	54	0	0	2	0
4	C	59	0	0	4	0
4	D	38	0	0	0	0
All	All	7643	0	7456	172	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 12.

All (172) 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:166:ARG:HH11	1:B:166:ARG:HG2	1.04	1.18
1:B:84:GLY:HA3	3:C:246:GOL:H2	1.10	1.08
1:D:213:THR:H	1:D:216:MET:HE2	1.20	1.03
3:B:246:GOL:H2	1:C:84:GLY:CA	1.90	0.99
3:B:246:GOL:H2	1:C:84:GLY:HA3	0.99	0.98
1:B:84:GLY:HA3	3:C:246:GOL:C2	1.95	0.96
1:B:84:GLY:CA	3:C:246:GOL:H2	1.95	0.95
3:B:246:GOL:C2	1:C:84:GLY:HA3	1.95	0.95
1:B:232:ARG:HH11	1:B:232:ARG:HG2	1.34	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:166:ARG:HH11	1:B:166:ARG:CG	1.81	0.92
1:B:56:ARG:HH11	1:B:56:ARG:HG3	1.32	0.91
1:B:166:ARG:HG2	1:B:166:ARG:NH1	1.86	0.89
1:B:56:ARG:HH11	1:B:56:ARG:CG	1.90	0.85
1:C:120:ALA:HB1	3:C:246:GOL:H32	1.59	0.85
1:D:213:THR:H	1:D:216:MET:CE	1.92	0.82
1:D:212:ALA:HA	1:D:216:MET:HE1	1.61	0.82
1:C:120:ALA:HB1	3:C:246:GOL:C3	2.10	0.82
1:B:120:ALA:HB1	3:B:246:GOL:C3	2.13	0.79
1:B:42:ILE:HD13	1:B:48:LEU:HD21	1.66	0.76
1:B:113:ILE:HD11	1:C:19:THR:HG21	1.67	0.75
1:B:113:ILE:HD11	1:C:19:THR:CG2	2.17	0.75
1:C:135:ASN:HD22	3:C:247:GOL:H11	1.52	0.74
1:B:232:ARG:HG2	1:B:232:ARG:NH1	2.03	0.73
1:B:218:ASP:H	1:B:221:THR:HG23	1.54	0.72
1:B:120:ALA:HB3	3:B:246:GOL:H12	1.71	0.72
1:B:68:MET:HE3	1:B:73:ARG:CZ	2.20	0.72
1:D:213:THR:N	1:D:216:MET:HE2	2.02	0.71
1:C:53:THR:HG22	1:C:61:ARG:HH21	1.55	0.71
1:D:56:ARG:HG3	1:D:57:VAL:H	1.56	0.70
1:B:120:ALA:HB1	3:B:246:GOL:H31	1.74	0.69
1:B:233:VAL:CG2	1:B:238:TYR:HE2	2.06	0.69
1:C:51:HIS:NE2	1:C:68:MET:HE1	2.08	0.69
1:D:68:MET:HE3	1:D:73:ARG:CZ	2.24	0.68
1:C:120:ALA:HB3	3:C:246:GOL:H12	1.76	0.68
1:A:210:ARG:HG2	1:A:211:ASP:N	2.10	0.67
1:B:218:ASP:H	1:B:221:THR:CG2	2.08	0.66
1:B:166:ARG:CG	1:B:166:ARG:NH1	2.49	0.66
1:B:222:VAL:HG23	2:B:1501:SAH:N3	2.11	0.65
3:A:246:GOL:H32	4:A:1637:HOH:O	1.98	0.64
1:A:68:MET:CE	1:A:73:ARG:CZ	2.75	0.63
1:B:129:HIS:HD2	4:C:1672:HOH:O	1.81	0.63
1:B:31:ASP:OD1	1:B:68:MET:HE1	1.99	0.63
1:A:51:HIS:NE2	1:A:68:MET:HE1	2.14	0.62
1:B:120:ALA:HB1	3:B:246:GOL:H32	1.80	0.62
1:B:218:ASP:OD1	1:B:221:THR:HG22	1.99	0.62
1:B:73:ARG:HD3	4:B:1699:HOH:O	1.99	0.62
1:C:143:PHE:CZ	1:C:176:ARG:HG2	2.34	0.62
1:B:197:VAL:HG12	2:B:1501:SAH:N1	2.15	0.62
1:B:233:VAL:CG2	1:B:238:TYR:CE2	2.82	0.62
1:D:32:ILE:HG12	1:D:48:LEU:HD11	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:1508:HOH:O	1:B:51:HIS:HD2	1.81	0.62
1:B:218:ASP:OD1	1:B:221:THR:CG2	2.48	0.62
1:A:146:LEU:O	1:A:150:LEU:HD22	1.99	0.61
1:D:56:ARG:HG3	1:D:57:VAL:N	2.15	0.61
1:D:142:PRO:HG2	1:D:145:ILE:HD12	1.83	0.60
1:A:68:MET:HE2	1:A:73:ARG:CZ	2.32	0.59
1:B:113:ILE:CD1	1:C:19:THR:HG21	2.32	0.59
1:B:233:VAL:HG23	1:B:238:TYR:CE2	2.37	0.59
1:B:120:ALA:CB	3:B:246:GOL:H12	2.32	0.59
1:A:241:ARG:HD3	1:D:140:LEU:HD13	1.85	0.58
1:C:14:ARG:NH2	4:C:1639:HOH:O	2.36	0.58
1:D:58:GLU:O	1:D:60:ASP:N	2.37	0.58
1:A:51:HIS:CE1	1:A:68:MET:HE1	2.39	0.56
1:A:85:VAL:HG21	1:A:114:THR:HG21	1.87	0.56
1:C:32:ILE:HG12	1:C:48:LEU:HD11	1.88	0.56
1:D:160:MET:HB2	1:D:225:VAL:HB	1.88	0.56
1:D:193:PHE:CE2	1:D:216:MET:HE3	2.41	0.56
1:C:86:PHE:HA	3:C:247:GOL:H31	1.88	0.56
1:D:83:PRO:O	1:D:87:ALA:HB3	2.05	0.55
1:B:233:VAL:HG23	1:B:238:TYR:HE2	1.70	0.55
1:B:140:LEU:CD1	1:C:241:ARG:HH11	2.19	0.55
1:C:151:ARG:NH1	1:C:184:GLU:OE2	2.38	0.55
1:B:126:PRO:HD2	3:B:246:GOL:H32	1.89	0.55
1:B:140:LEU:HD13	1:C:241:ARG:HD3	1.88	0.54
1:D:197:VAL:HG12	1:D:198:THR:HG23	1.89	0.54
1:B:143:PHE:HD1	1:B:173:GLN:HE21	1.55	0.54
1:A:58:GLU:HG3	1:A:60:ASP:H	1.73	0.54
1:A:152:HIS:HB3	1:D:152:HIS:O	2.07	0.54
1:A:68:MET:HE3	1:A:73:ARG:NH1	2.22	0.54
1:D:42:ILE:HD11	1:D:78:VAL:HG11	1.90	0.54
1:C:126:PRO:HD2	3:C:246:GOL:H32	1.90	0.54
1:C:31:ASP:OD1	1:C:51:HIS:HE1	1.91	0.53
1:B:58:GLU:OE2	1:B:61:ARG:HD2	2.08	0.53
1:A:68:MET:HB3	1:A:73:ARG:HD3	1.91	0.53
1:A:39:VAL:HA	3:A:247:GOL:H31	1.90	0.52
1:C:53:THR:HG22	1:C:61:ARG:NH2	2.24	0.52
1:B:219:MET:HA	4:B:1531:HOH:O	2.09	0.52
1:C:138:ASP:OD2	3:C:247:GOL:H12	2.10	0.52
1:A:195:ARG:O	1:A:202:GLN:HA	2.09	0.51
1:B:56:ARG:HG3	1:B:56:ARG:NH1	2.12	0.51
1:D:180:ILE:H	1:D:180:ILE:HD12	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:68:MET:HE2	4:C:1609:HOH:O	2.10	0.51
1:B:13:GLY:O	3:B:248:GOL:H2	2.10	0.51
1:C:120:ALA:HB1	3:C:246:GOL:H31	1.88	0.51
1:B:56:ARG:NH1	1:B:58:GLU:OE1	2.44	0.51
1:A:120:ALA:HB2	1:A:127:LEU:HD12	1.92	0.51
1:A:238:TYR:OH	1:D:97:GLU:OE1	2.17	0.51
2:C:1501:SAH:HG1	4:C:1564:HOH:O	2.11	0.51
1:B:120:ALA:CB	3:B:246:GOL:H31	2.41	0.50
1:C:31:ASP:OD2	1:C:73:ARG:HD3	2.11	0.50
1:B:227:ASN:HD21	1:B:239:THR:HG21	1.77	0.50
1:B:31:ASP:OD1	1:B:68:MET:CE	2.59	0.50
1:D:168:LYS:HA	1:D:171:PRO:HA	1.94	0.50
1:A:16:ASP:O	1:D:20:PRO:HD2	2.10	0.50
1:B:51:HIS:CE1	1:B:68:MET:HE1	2.47	0.50
1:A:176:ARG:O	1:A:180:ILE:HG12	2.12	0.49
1:B:232:ARG:HH11	1:B:232:ARG:CG	2.14	0.49
1:C:53:THR:CG2	1:C:61:ARG:HH21	2.23	0.49
1:D:167:SER:C	1:D:169:SER:H	2.19	0.49
1:D:56:ARG:CG	1:D:57:VAL:N	2.73	0.49
1:D:150:LEU:HD21	1:D:177:VAL:HG13	1.95	0.49
1:B:56:ARG:CG	1:B:56:ARG:NH1	2.60	0.49
1:D:9:GLY:HA3	1:D:83:PRO:HD3	1.95	0.49
1:A:197:VAL:HG12	1:A:198:THR:HG23	1.95	0.49
1:D:143:PHE:HD1	1:D:173:GLN:HE21	1.60	0.48
1:B:137:SER:OG	1:B:164:ASN:HB3	2.14	0.48
1:C:120:ALA:CB	3:C:246:GOL:H12	2.44	0.48
1:A:36:ILE:HD12	1:A:50:LEU:HD23	1.95	0.48
1:B:32:ILE:HD12	1:B:42:ILE:HD12	1.96	0.48
1:D:166:ARG:HD2	1:D:217:ALA:O	2.13	0.48
1:C:64:HIS:NE2	1:C:68:MET:HE3	2.29	0.48
1:D:172:HIS:CD2	1:D:173:GLN:H	2.33	0.47
1:C:150:LEU:HD11	1:C:177:VAL:HG22	1.96	0.47
1:C:135:ASN:HD22	3:C:247:GOL:H32	1.80	0.47
1:B:238:TYR:OH	1:C:97:GLU:OE1	2.28	0.47
1:D:192:LEU:HG	1:D:206:VAL:HG22	1.97	0.46
1:C:179:GLU:O	1:C:183:GLU:HG3	2.15	0.46
1:C:148:LYS:O	1:C:152:HIS:HD2	1.97	0.46
1:A:147:GLU:HG3	1:A:180:ILE:HD12	1.97	0.46
1:D:66:LEU:HD21	1:D:92:LEU:HD11	1.98	0.46
1:B:67:GLU:O	1:B:70:ALA:HB3	2.16	0.45
1:B:160:MET:HB2	1:B:225:VAL:HB	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:68:MET:HE3	1:B:73:ARG:NH1	2.31	0.45
1:D:31:ASP:OD2	1:D:51:HIS:HE1	1.98	0.45
1:B:143:PHE:HD1	1:B:173:GLN:NE2	2.15	0.45
1:A:54:ASP:OD1	1:A:54:ASP:C	2.59	0.45
1:D:189:ARG:HD2	1:D:226:GLY:O	2.17	0.45
1:C:32:ILE:HB	1:C:50:LEU:HD23	1.99	0.45
1:B:120:ALA:HB3	3:B:246:GOL:C1	2.43	0.45
1:A:116:MET:HE2	1:A:127:LEU:HD13	1.99	0.44
1:C:53:THR:CG2	1:C:61:ARG:NH2	2.80	0.44
1:B:70:ALA:C	1:B:72:GLY:H	2.25	0.44
1:B:222:VAL:HG22	2:B:1501:SAH:H1'	2.00	0.44
1:D:68:MET:CE	1:D:73:ARG:CZ	2.95	0.43
1:D:192:LEU:HB2	1:D:224:LEU:HB2	2.00	0.43
1:B:29:ALA:HB2	1:B:76:VAL:HG23	2.00	0.43
1:A:137:SER:HB2	1:A:164:ASN:HB3	2.00	0.43
1:B:9:GLY:HA3	1:B:83:PRO:HD3	1.99	0.43
1:B:31:ASP:OD1	1:B:51:HIS:HE1	2.02	0.43
1:B:42:ILE:HG12	1:B:43:ALA:N	2.34	0.43
1:A:149:ARG:NE	1:D:156:GLY:O	2.51	0.43
1:D:192:LEU:HD12	1:D:192:LEU:N	2.33	0.43
1:D:151:ARG:HD3	1:D:155:ARG:HH21	1.84	0.43
1:D:136:LEU:HD12	1:D:165:PRO:HA	2.01	0.42
1:A:164:ASN:HA	1:A:219:MET:HE1	2.01	0.42
1:B:125:ALA:HA	3:B:246:GOL:H32	2.02	0.42
1:B:195:ARG:HB2	1:B:195:ARG:NH1	2.35	0.42
1:B:218:ASP:N	1:B:221:THR:HG23	2.29	0.42
1:A:49:THR:CG2	4:A:1528:HOH:O	2.67	0.42
1:A:114:THR:HG22	2:A:1501:SAH:O	2.20	0.42
1:D:172:HIS:CG	1:D:173:GLN:H	2.37	0.42
1:D:190:LEU:HD12	1:D:190:LEU:HA	1.89	0.42
1:B:16:ASP:O	1:C:20:PRO:HD2	2.20	0.41
1:C:129:HIS:O	1:C:130:ASP:C	2.64	0.41
1:C:138:ASP:CG	3:C:247:GOL:H12	2.44	0.41
1:A:56:ARG:O	1:A:61:ARG:NH1	2.54	0.41
1:B:189:ARG:HD2	1:B:226:GLY:O	2.21	0.41
1:A:68:MET:HE2	1:A:73:ARG:NE	2.36	0.41
4:A:1509:HOH:O	3:B:247:GOL:H12	2.20	0.41
1:D:197:VAL:HA	1:D:202:GLN:HG3	2.03	0.41
1:D:176:ARG:O	1:D:180:ILE:HD12	2.21	0.41
1:A:231:ARG:HB2	4:A:1700:HOH:O	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	238/251 (95%)	228 (96%)	10 (4%)	0	100	100
1	B	238/251 (95%)	225 (94%)	12 (5%)	1 (0%)	30	33
1	C	240/251 (96%)	228 (95%)	12 (5%)	0	100	100
1	D	237/251 (94%)	220 (93%)	13 (6%)	4 (2%)	7	5
All	All	953/1004 (95%)	901 (94%)	47 (5%)	5 (0%)	24	26

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	59	LEU
1	D	57	VAL
1	B	71	GLU
1	D	173	GLN
1	D	168	LYS

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	186/195 (95%)	177 (95%)	9 (5%)	23	28
1	B	185/195 (95%)	170 (92%)	15 (8%)	11	11
1	C	187/195 (96%)	180 (96%)	7 (4%)	30	39
1	D	185/195 (95%)	171 (92%)	14 (8%)	12	13
All	All	743/780 (95%)	698 (94%)	45 (6%)	17	19

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	THR
1	A	90	SER
1	A	150	LEU
1	A	160	MET
1	A	168	LYS
1	A	180	ILE
1	A	190	LEU
1	A	201	GLU
1	A	210	ARG
1	B	41	ARG
1	B	42	ILE
1	B	45	ARG
1	B	55	ASN
1	B	56	ARG
1	B	58	GLU
1	B	61	ARG
1	B	71	GLU
1	B	79	SER
1	B	97	GLU
1	B	166	ARG
1	B	168	LYS
1	B	181	LEU
1	B	221	THR
1	B	232	ARG
1	C	1	MET
1	C	53	THR
1	C	56	ARG
1	C	57	VAL
1	C	59	LEU
1	C	90	SER
1	C	106	GLU
1	D	2	SER
1	D	50	LEU
1	D	53	THR
1	D	61	ARG
1	D	79	SER
1	D	90	SER
1	D	150	LEU
1	D	169	SER
1	D	170	ARG
1	D	190	LEU
1	D	201	GLU

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Mol	Chain	Res	Type
1	D	207	VAL
1	D	210	ARG
1	D	219	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	139	ASN
1	A	227	ASN
1	B	51	HIS
1	B	129	HIS
1	B	173	GLN
1	C	51	HIS
1	C	129	HIS
1	C	172	HIS
1	C	173	GLN
1	D	51	HIS
1	D	139	ASN
1	D	164	ASN
1	D	173	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

12 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
3	GOL	A	248	-	5,5,5	0.39	0	5,5,5	0.80	0
3	GOL	B	246	-	5,5,5	0.85	0	5,5,5	1.10	0
3	GOL	B	247	-	5,5,5	0.30	0	5,5,5	0.37	0
3	GOL	A	247	-	5,5,5	0.42	0	5,5,5	0.34	0
2	SAH	D	1501	-	27,28,28	1.40	4 (14%)	36,40,40	1.93	6 (16%)
3	GOL	B	248	-	5,5,5	0.40	0	5,5,5	0.33	0
3	GOL	A	246	-	5,5,5	0.39	0	5,5,5	0.40	0
2	SAH	B	1501	-	27,28,28	1.48	3 (11%)	36,40,40	2.07	11 (30%)
2	SAH	C	1501	-	27,28,28	1.54	4 (14%)	36,40,40	2.16	11 (30%)
2	SAH	A	1501	-	27,28,28	1.46	3 (11%)	36,40,40	2.07	9 (25%)
3	GOL	C	247	-	5,5,5	0.32	0	5,5,5	1.01	0
3	GOL	C	246	-	5,5,5	0.77	0	5,5,5	1.07	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
3	GOL	A	248	-	-	3/4/4/4	-
3	GOL	B	246	-	-	2/4/4/4	-
3	GOL	B	247	-	-	2/4/4/4	-
3	GOL	A	247	-	-	2/4/4/4	-
2	SAH	D	1501	-	-	0/15/31/31	0/3/3/3
3	GOL	B	248	-	-	2/4/4/4	-
3	GOL	A	246	-	-	0/4/4/4	-
2	SAH	B	1501	-	-	0/15/31/31	0/3/3/3
2	SAH	C	1501	-	-	2/15/31/31	0/3/3/3
2	SAH	A	1501	-	-	1/15/31/31	0/3/3/3
3	GOL	C	247	-	-	1/4/4/4	-
3	GOL	C	246	-	-	2/4/4/4	-

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1501	SAH	C5-C4	4.95	1.47	1.39
2	C	1501	SAH	C5-C4	4.67	1.47	1.39
2	D	1501	SAH	C5-C4	4.57	1.47	1.39
2	A	1501	SAH	C5-C4	4.52	1.47	1.39
2	A	1501	SAH	C5-C6	3.23	1.50	1.41
2	C	1501	SAH	C8-N7	3.21	1.37	1.31
2	A	1501	SAH	C8-N7	2.93	1.37	1.31
2	C	1501	SAH	C5-C6	2.92	1.49	1.41
2	B	1501	SAH	C8-N7	2.79	1.37	1.31
2	B	1501	SAH	C5-N7	-2.38	1.34	1.39
2	D	1501	SAH	C4-N9	-2.36	1.32	1.37
2	D	1501	SAH	C5-C6	2.29	1.47	1.41
2	D	1501	SAH	C8-N7	2.16	1.35	1.31
2	C	1501	SAH	O4'-C1'	2.13	1.46	1.42

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1501	SAH	C5-C4-N3	-6.30	118.05	126.72
2	B	1501	SAH	C5-C4-N3	-6.12	118.28	126.72
2	A	1501	SAH	C5-C4-N3	-6.05	118.38	126.72
2	D	1501	SAH	C5-C4-N3	-5.43	119.24	126.72
2	B	1501	SAH	N3-C4-N9	5.30	136.18	127.17
2	D	1501	SAH	N3-C4-N9	4.71	135.18	127.17
2	A	1501	SAH	C4-C5-N7	-4.47	105.47	110.58
2	C	1501	SAH	C4-C5-N7	-4.34	105.62	110.58
2	B	1501	SAH	N3-C2-N1	-4.33	122.02	128.58
2	C	1501	SAH	N3-C4-N9	4.28	134.45	127.17
2	A	1501	SAH	N3-C4-N9	4.23	134.37	127.17
2	A	1501	SAH	C2-N3-C4	4.11	121.86	111.83
2	C	1501	SAH	C2-N3-C4	4.10	121.83	111.83
2	D	1501	SAH	C2-N3-C4	3.88	121.30	111.83
2	B	1501	SAH	C2-N3-C4	3.88	121.30	111.83
2	C	1501	SAH	N3-C2-N1	-3.86	122.74	128.58
2	D	1501	SAH	N3-C2-N1	-3.66	123.04	128.58
2	A	1501	SAH	N3-C2-N1	-3.56	123.19	128.58
2	D	1501	SAH	C4-N9-C8	3.35	109.25	105.74
2	C	1501	SAH	C5-N7-C8	3.25	108.56	103.45
2	A	1501	SAH	C5-N7-C8	3.23	108.53	103.45
2	A	1501	SAH	C6-C5-N7	2.87	137.62	132.09
2	A	1501	SAH	C4-N9-C8	2.73	108.60	105.74
2	A	1501	SAH	N9-C8-N7	-2.67	110.14	113.94
2	D	1501	SAH	C4-C5-N7	-2.63	107.58	110.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1501	SAH	N6-C6-N1	2.57	124.10	118.38
2	B	1501	SAH	C4-N9-C8	2.55	108.42	105.74
2	C	1501	SAH	O4'-C1'-N9	2.49	112.87	108.09
2	B	1501	SAH	C4-C5-N7	-2.42	107.81	110.58
2	C	1501	SAH	C6-C5-N7	2.30	136.52	132.09
2	B	1501	SAH	C5-C6-N6	-2.25	117.73	123.29
2	C	1501	SAH	N9-C8-N7	-2.22	110.78	113.94
2	C	1501	SAH	CB-CG-SD	-2.20	108.55	113.45
2	C	1501	SAH	C2-N1-C6	2.20	122.34	118.73
2	B	1501	SAH	C4'-C5'-SD	-2.14	106.16	113.78
2	B	1501	SAH	C2-N1-C6	2.02	122.06	118.73
2	B	1501	SAH	C5-N7-C8	2.01	106.61	103.45

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	247	GOL	C1-C2-C3-O3
3	A	247	GOL	O2-C2-C3-O3
3	B	246	GOL	C1-C2-C3-O3
3	B	247	GOL	C1-C2-C3-O3
3	B	248	GOL	C1-C2-C3-O3
3	C	246	GOL	C1-C2-C3-O3
3	C	247	GOL	C1-C2-C3-O3
3	B	246	GOL	O2-C2-C3-O3
3	B	247	GOL	O2-C2-C3-O3
3	C	246	GOL	O2-C2-C3-O3
3	B	248	GOL	O2-C2-C3-O3
3	A	248	GOL	O1-C1-C2-C3
3	A	248	GOL	O2-C2-C3-O3
2	A	1501	SAH	C3'-C4'-C5'-SD
2	C	1501	SAH	CB-CG-SD-C5'
3	A	248	GOL	O1-C1-C2-O2
2	C	1501	SAH	C-CA-CB-CG

There are no ring outliers.

10 monomers are involved in 35 short contacts:

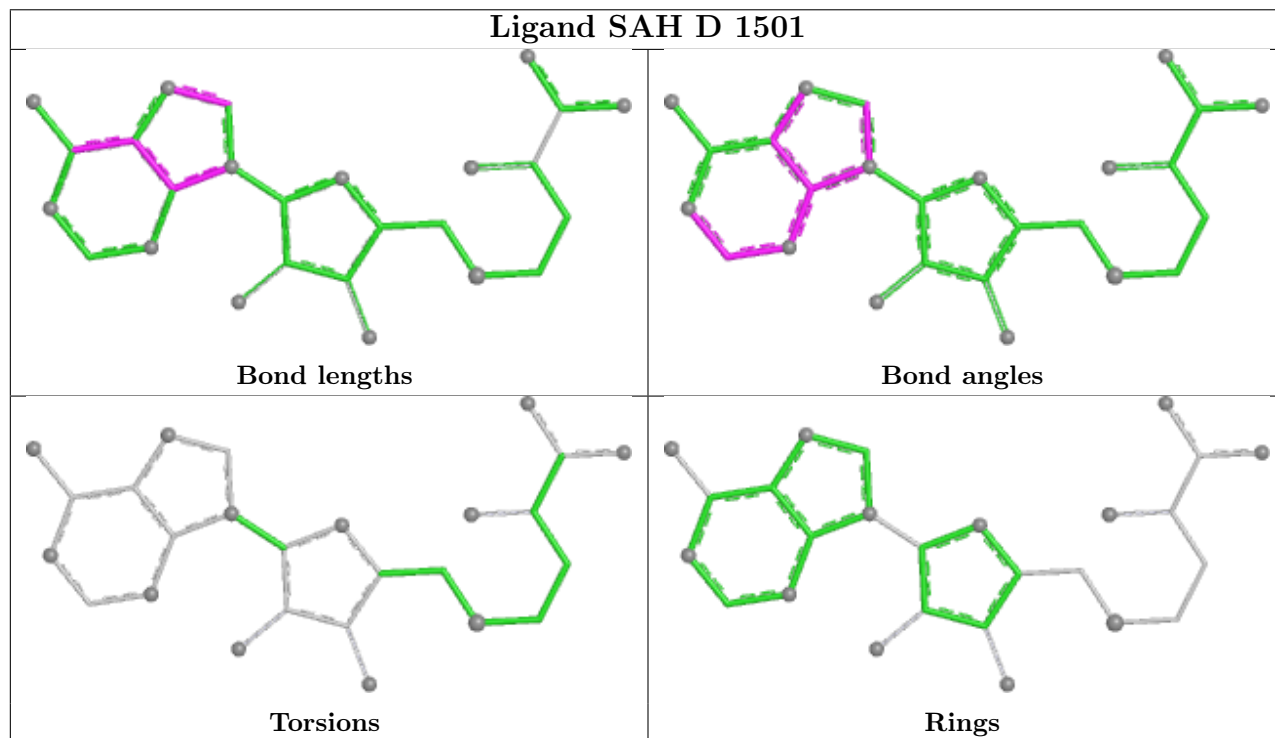
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	246	GOL	12	0
3	B	247	GOL	1	0

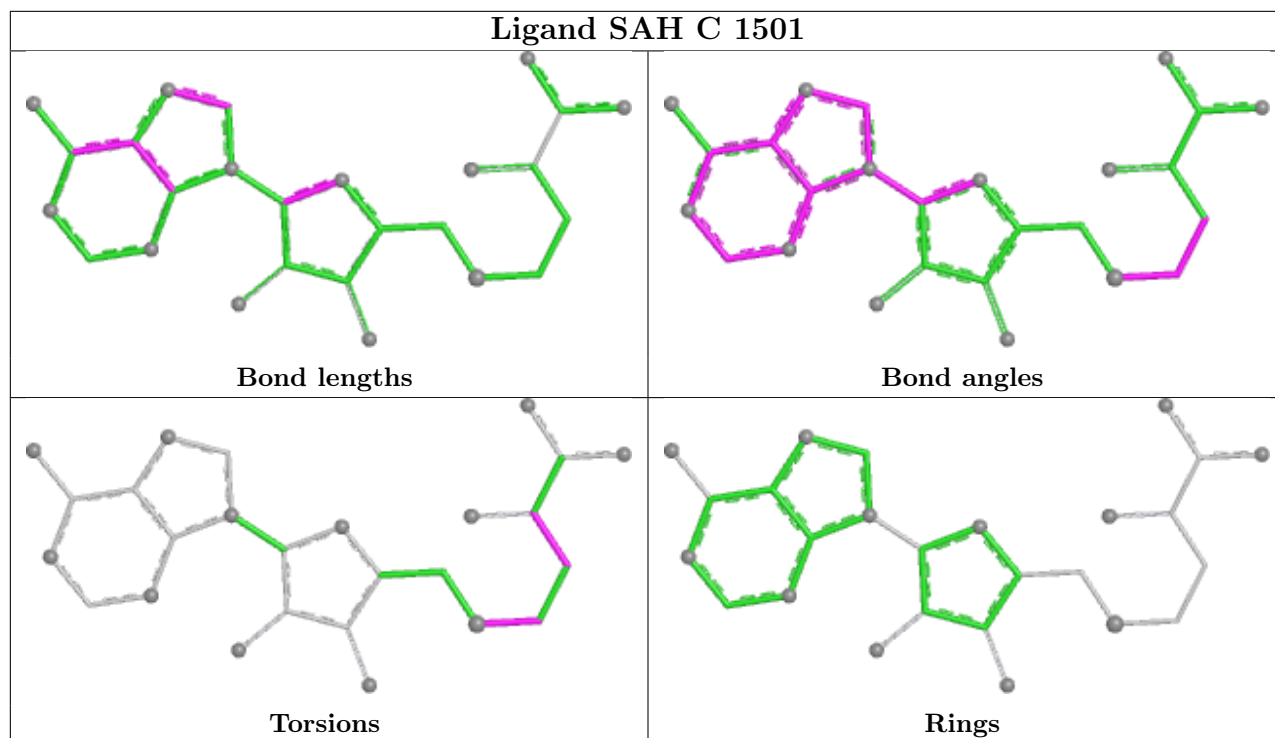
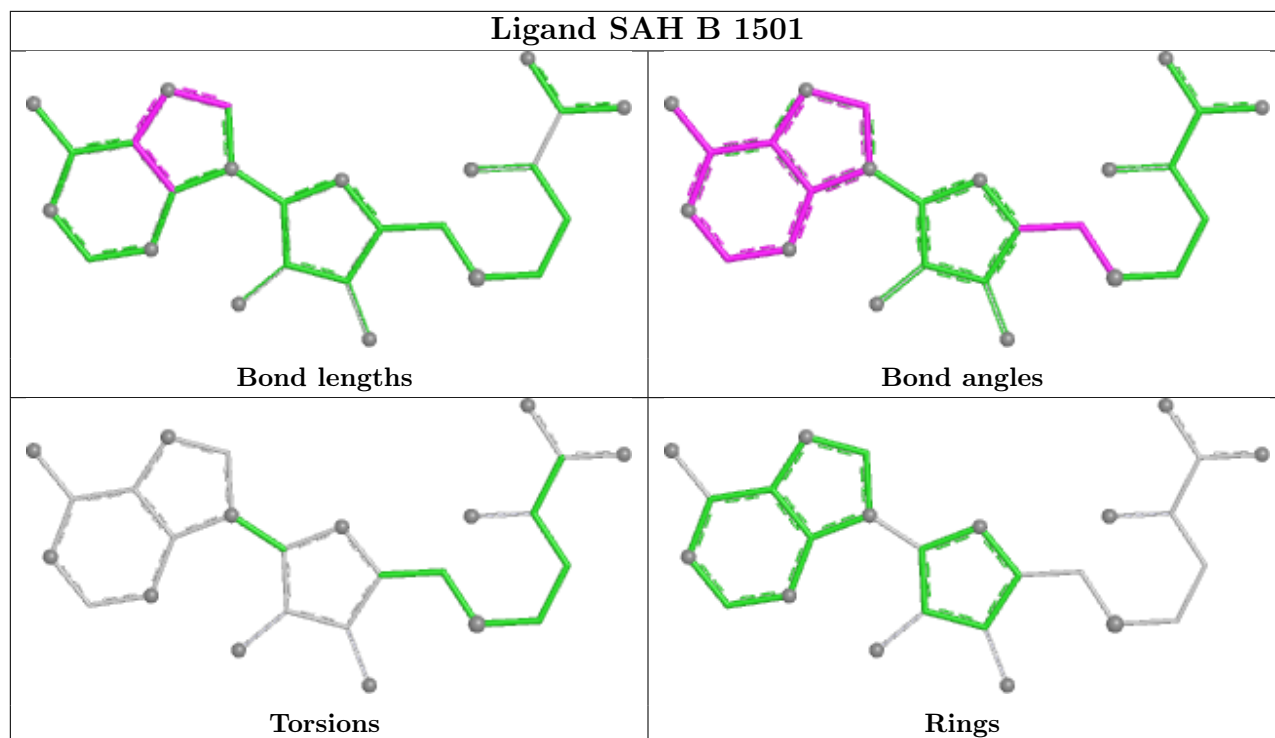
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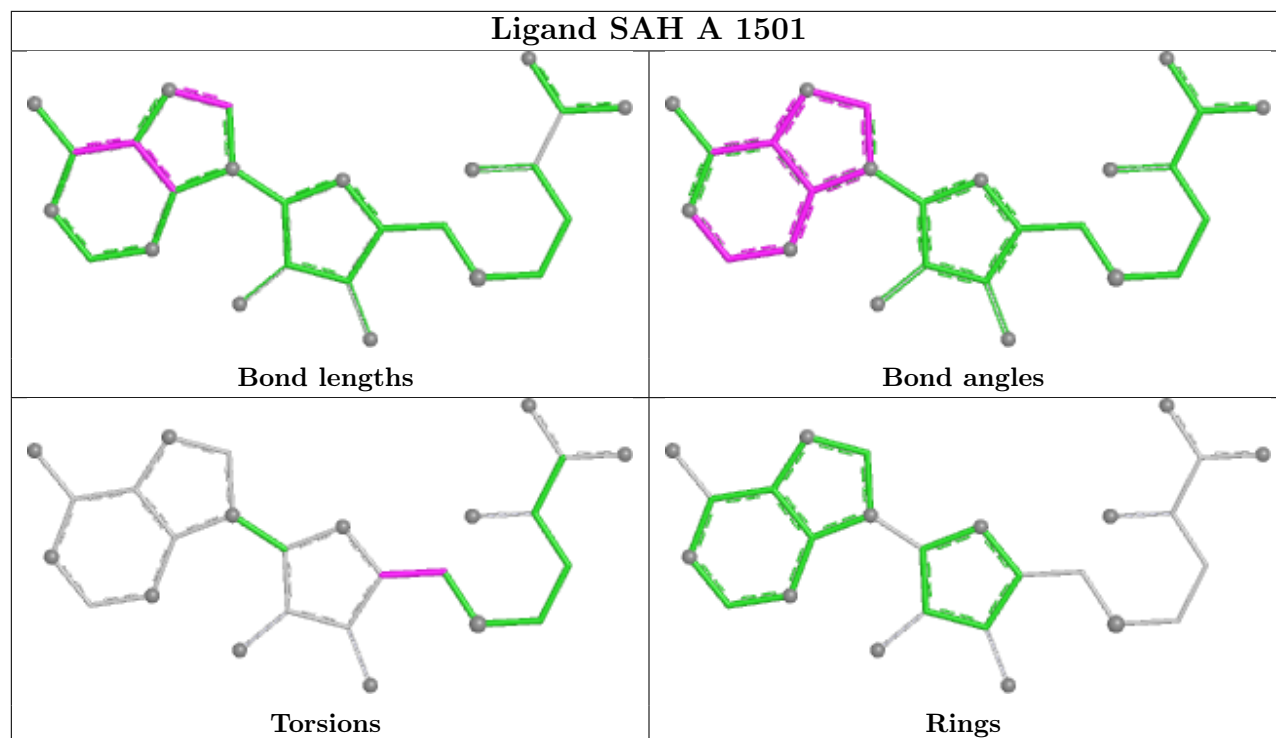
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	247	GOL	1	0
3	B	248	GOL	1	0
3	A	246	GOL	1	0
2	B	1501	SAH	3	0
2	C	1501	SAH	1	0
2	A	1501	SAH	1	0
3	C	247	GOL	5	0
3	C	246	GOL	9	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	240/251 (95%)	0.17	6 (2%) 58 56	20, 38, 62, 72	0
1	B	240/251 (95%)	-0.06	0 100 100	22, 33, 47, 52	0
1	C	242/251 (96%)	0.04	4 (1%) 69 67	22, 37, 49, 61	0
1	D	239/251 (95%)	0.39	8 (3%) 49 46	22, 44, 70, 79	0
All	All	961/1004 (95%)	0.14	18 (1%) 66 64	20, 37, 61, 79	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	57	VAL	6.0
1	D	59	LEU	3.6
1	A	140	LEU	3.3
1	C	56	ARG	3.0
1	D	140	LEU	3.0
1	C	1	MET	2.9
1	D	103	ALA	2.8
1	C	57	VAL	2.6
1	A	219	MET	2.5
1	D	175	THR	2.3
1	C	103	ALA	2.3
1	A	180	ILE	2.3
1	A	241	ARG	2.2
1	A	209	LEU	2.2
1	D	239	THR	2.0
1	A	212	ALA	2.0
1	D	170	ARG	2.0
1	D	169	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

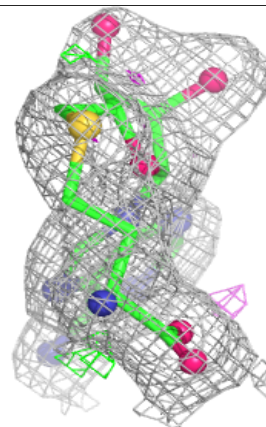
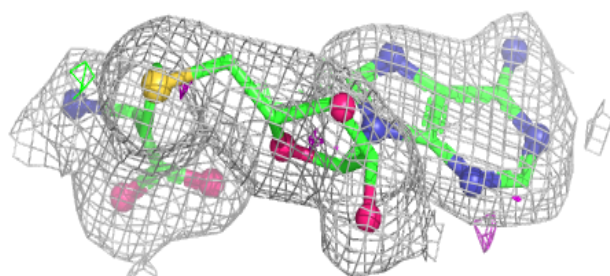
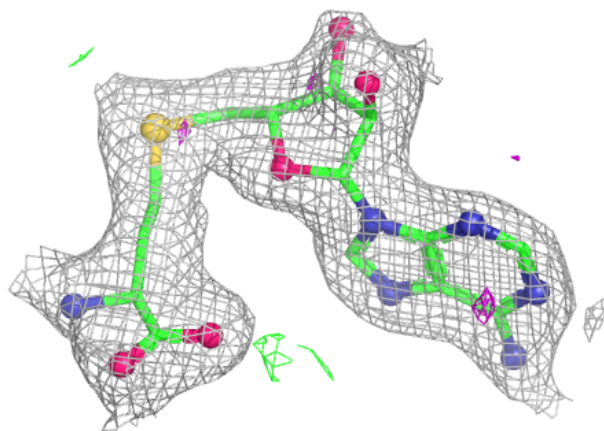
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GOL	B	247	6/6	0.77	0.19	60,62,63,65	0
3	GOL	A	247	6/6	0.82	0.19	62,63,64,64	0
3	GOL	A	248	6/6	0.86	0.12	53,57,58,59	0
3	GOL	B	248	6/6	0.86	0.16	59,66,67,67	0
3	GOL	C	246	6/6	0.86	0.22	41,43,44,45	0
3	GOL	B	246	6/6	0.87	0.20	32,37,39,39	0
3	GOL	C	247	6/6	0.91	0.14	47,48,49,50	0
2	SAH	C	1501	26/26	0.92	0.09	30,35,39,43	0
2	SAH	D	1501	26/26	0.93	0.08	31,38,41,45	0
2	SAH	A	1501	26/26	0.94	0.08	35,38,45,46	0
3	GOL	A	246	6/6	0.95	0.08	34,46,48,51	0
2	SAH	B	1501	26/26	0.95	0.07	27,31,35,39	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.

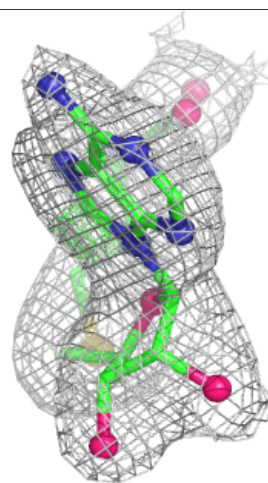
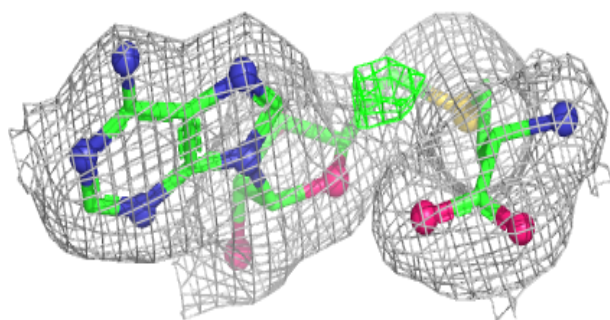
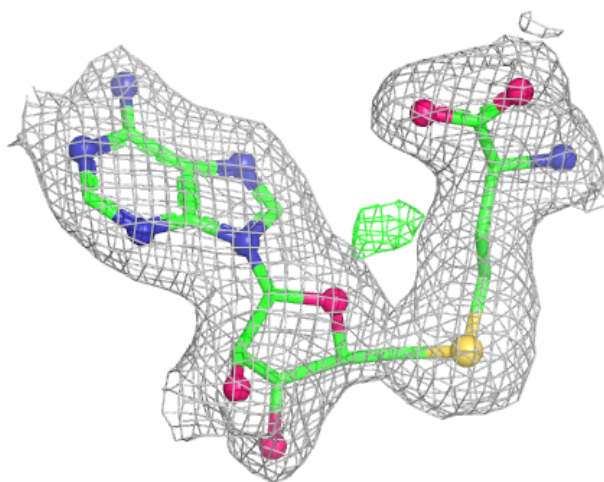
Electron density around SAH C 1501:

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



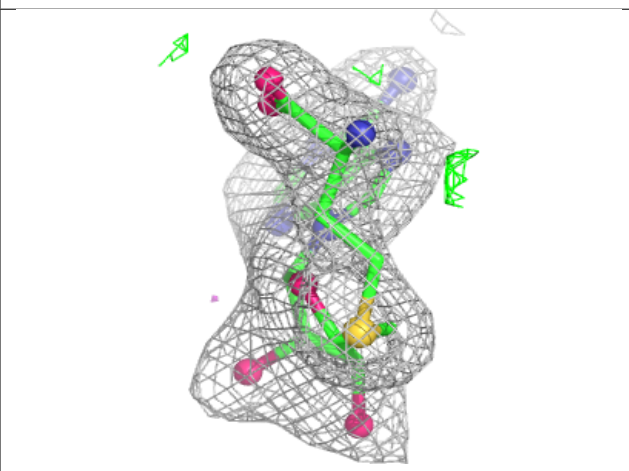
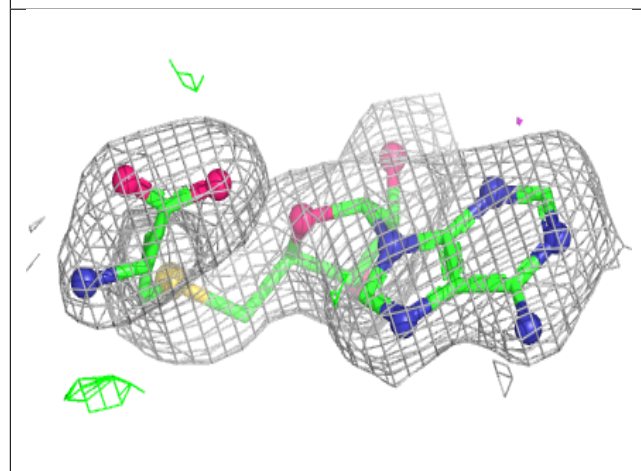
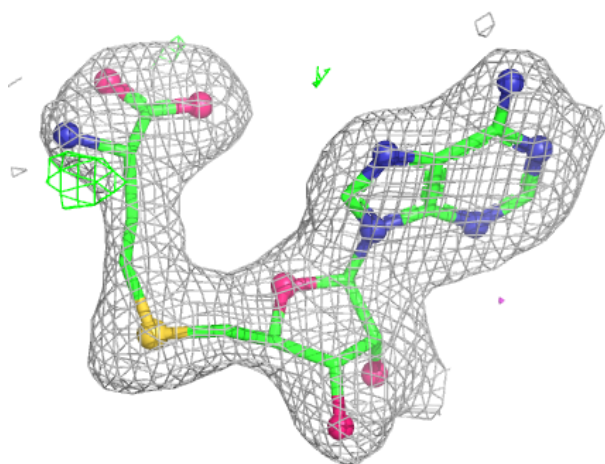
Electron density around SAH D 1501:

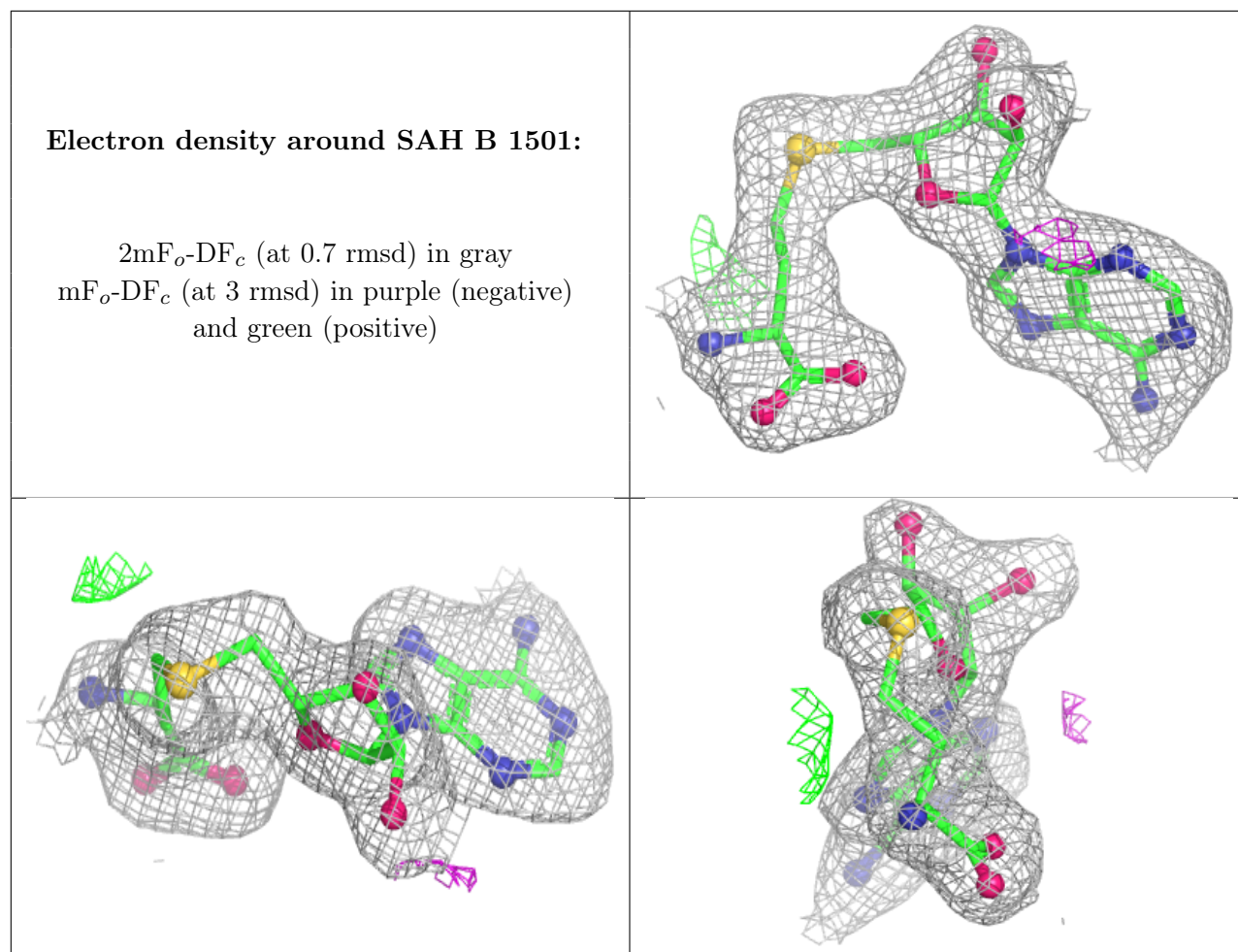
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around SAH A 1501:

$2mF_o-DF_c$ (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.