



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:53 AM UTC

PDB ID : 5U0B / pdb_00005u0b
Title : Structure of full-length Zika virus NS5
Authors : Zhao, B.; Du, F.
Deposited on : 2016-11-23
Resolution : 3.00 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

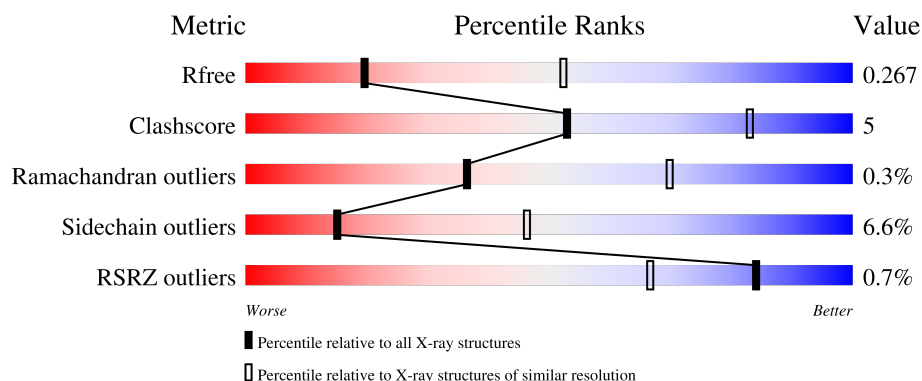
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	903	
1	B	903	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 14282 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

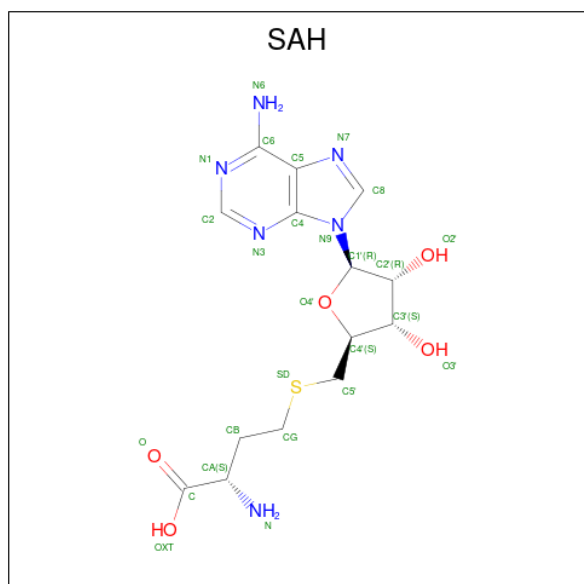
- Molecule 1 is a protein called Genome polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	883	Total	C	N	O	S	0	0	0
			7098	4459	1290	1301	48			
1	B	883	Total	C	N	O	S	0	0	0
			7098	4459	1290	1301	48			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	62	LEU	ILE	engineered mutation	UNP Q32ZE1
A	212	SER	CYS	engineered mutation	UNP Q32ZE1
B	62	LEU	ILE	engineered mutation	UNP Q32ZE1
B	212	SER	CYS	engineered mutation	UNP Q32ZE1

- Molecule 2 is S-ADENOSYL-L-HOMOCYSTEINE (CCD ID: 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		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		
3	B	2	Total	Zn	0	0
			2	2		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

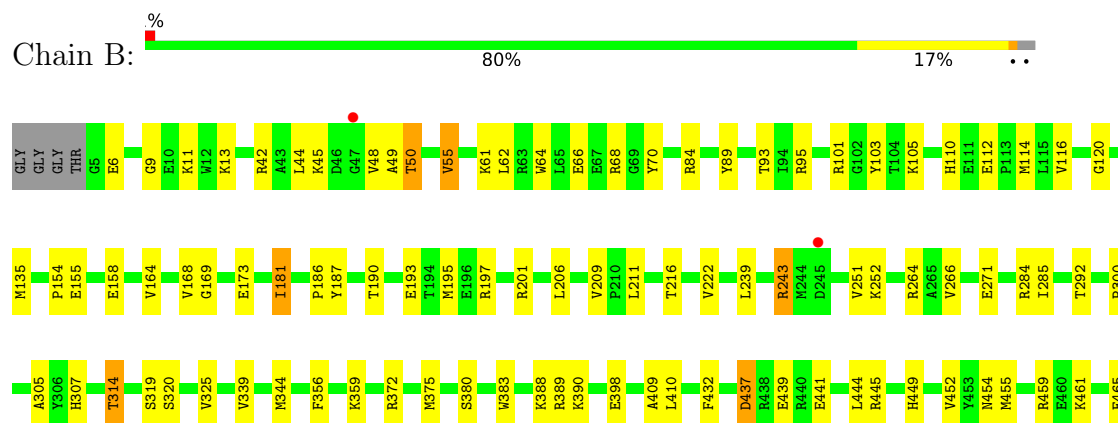
3 Residue-property plots [i](#)

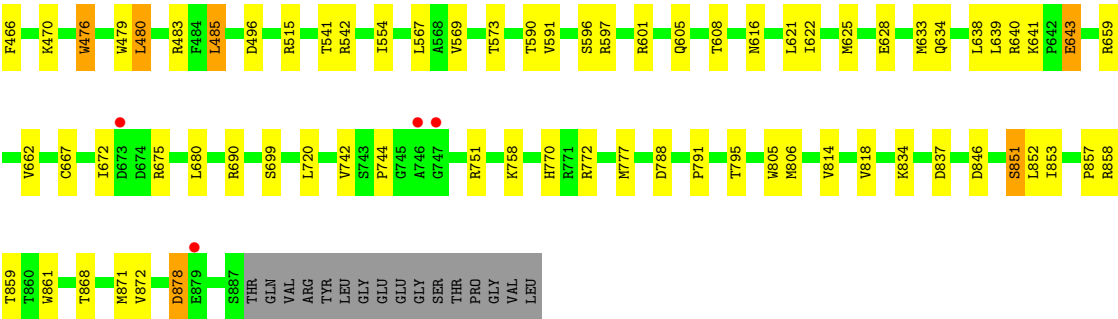
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Genome polyprotein



• Molecule 1: Genome polyprotein





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	136.50Å 197.00Å 95.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	85.78 – 3.00 85.78 – 3.00	Depositor EDS
% Data completeness (in resolution range)	96.7 (85.78-3.00) 86.0 (85.78-3.00)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 3.01Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155)	Depositor
R, R_{free}	0.231 , 0.268 0.231 , 0.267	Depositor DCC
R_{free} test set	2529 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	37.9	Xtriage
Anisotropy	0.543	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 26.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	14282	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.98 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.2224e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.11	0/7263	0.33	0/9812
1	B	0.12	0/7263	0.34	0/9812
All	All	0.11	0/14526	0.33	0/19624

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	7098	0	6990	70	0
1	B	7098	0	6990	73	0
2	A	26	0	19	1	0
2	B	26	0	19	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	15	0	0	0	0
4	B	15	0	0	0	0
All	All	14282	0	14018	142	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (142) 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:99:GLU:OE2	1:A:101:ARG:NH1	2.18	0.77
1:B:9:GLY:HA3	1:B:187:TYR:HB2	1.67	0.76
1:B:70:TYR:HB3	1:B:222:VAL:HG11	1.70	0.74
1:A:551:GLU:OE2	1:A:601:ARG:NH2	2.25	0.67
1:B:465:GLU:HG3	1:B:470:LYS:HG3	1.76	0.66
1:A:82:CYS:SG	1:A:104:THR:OG1	2.54	0.65
1:B:168:VAL:HG21	1:B:181:ILE:HD12	1.79	0.64
1:A:787:VAL:HG11	1:A:883:MET:HB2	1.80	0.63
1:A:149:GLU:HG2	1:A:160:ARG:NH1	2.14	0.63
1:B:120:GLY:HA2	1:B:264:ARG:HB2	1.80	0.63
1:B:193:GLU:HG2	1:B:197:ARG:HH11	1.64	0.62
1:A:197:ARG:HG2	1:A:200:ARG:HH12	1.64	0.62
1:A:718:LEU:HD21	1:A:841:LEU:HD23	1.80	0.62
1:B:372:ARG:HA	1:B:375:MET:HE2	1.82	0.61
1:B:878:ASP:OD2	1:B:878:ASP:N	2.34	0.60
1:B:13:LYS:HG2	1:B:154:PRO:HG3	1.86	0.57
1:B:834:LYS:NZ	1:B:837:ASP:OD2	2.30	0.57
1:A:826:MET:O	1:A:829:LYS:NZ	2.37	0.56
1:A:405:ARG:HH12	1:A:794:ARG:HH12	1.51	0.56
1:A:569:VAL:O	1:A:573:THR:HB	2.05	0.56
1:B:44:LEU:HD21	1:B:55:VAL:HA	1.86	0.56
1:B:569:VAL:O	1:B:573:THR:HB	2.06	0.55
1:A:476:TRP:CD1	1:A:476:TRP:N	2.75	0.54
1:B:320:SER:HB2	1:B:344:MET:HB2	1.88	0.54
1:A:84:ARG:HD2	1:A:114:MET:HE2	1.89	0.54
1:B:319:SER:O	1:B:459:ARG:NH1	2.41	0.53
1:B:476:TRP:N	1:B:476:TRP:CD1	2.77	0.53
1:A:105:LYS:HD2	1:A:110:HIS:CD2	2.44	0.53
1:B:852:LEU:O	1:B:858:ARG:HG2	2.08	0.52
1:B:193:GLU:HG2	1:B:197:ARG:NH1	2.24	0.52
1:A:168:VAL:HG11	1:A:181:ILE:HG13	1.91	0.52
1:A:622:ILE:HA	1:A:625:MET:HE3	1.92	0.52
1:B:788:ASP:OD1	1:B:788:ASP:N	2.42	0.52
1:B:50:THR:O	1:B:50:THR:OG1	2.26	0.52
1:A:730:CYS:HB2	1:A:770:HIS:HE1	1.75	0.51
1:A:29:LYS:NZ	1:A:34:GLU:OE2	2.43	0.51
1:A:193:GLU:HA	1:A:196:GLU:HB2	1.92	0.51
1:A:262:GLY:HA3	1:A:586:GLU:HG2	1.93	0.50
1:B:758:LYS:HD3	1:B:791:PRO:HG3	1.94	0.50
1:B:858:ARG:HA	1:B:861:TRP:HB3	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:857:PRO:O	1:B:859:THR:N	2.45	0.49
1:A:9:GLY:HA3	1:A:187:TYR:HB2	1.95	0.49
1:B:805:TRP:HE1	1:B:806:MET:HE2	1.78	0.49
1:B:662:VAL:HG22	1:B:667:CYS:HB2	1.95	0.49
1:A:101:ARG:HH21	1:A:125:ARG:HB2	1.78	0.48
1:B:305:ALA:HB3	1:B:596:SER:HB3	1.95	0.48
1:A:120:GLY:HA2	1:A:264:ARG:HG2	1.94	0.48
1:B:616:ASN:HD21	1:B:667:CYS:HB3	1.79	0.48
1:A:731:ARG:HD3	1:A:736:LEU:HD21	1.95	0.48
1:A:868:THR:O	1:A:872:VAL:HG23	2.14	0.48
1:B:659:ARG:HH12	1:B:672:ILE:HA	1.78	0.48
1:B:112:GLU:OE2	1:B:466:PHE:HA	2.14	0.47
1:A:50:THR:HG21	1:A:588:GLY:HA2	1.97	0.47
1:B:307:HIS:NE2	1:B:596:SER:HB2	2.30	0.47
1:B:284:ARG:NH2	1:B:449:HIS:O	2.44	0.47
1:A:805:TRP:HE1	1:A:806:MET:HE2	1.79	0.47
1:A:640:ARG:HH22	1:B:11:LYS:NZ	2.13	0.47
1:A:724:ARG:HB3	1:A:826:MET:HE1	1.97	0.46
1:B:49:ALA:HB1	1:B:116:VAL:HA	1.97	0.46
1:A:245:ASP:HA	1:A:246:GLY:HA2	1.64	0.46
1:A:877:GLY:O	1:A:882:TYR:OH	2.33	0.46
1:A:158:GLU:HG3	1:A:191:MET:HG2	1.98	0.46
1:A:329:LEU:HD22	1:A:780:ALA:HB1	1.98	0.46
1:B:195:MET:HE2	1:B:206:LEU:HD21	1.98	0.46
1:B:805:TRP:NE1	1:B:806:MET:HE2	2.30	0.46
1:A:103:TYR:CE2	1:A:135:MET:HE1	2.51	0.45
1:A:748:TRP:HZ3	1:A:756:LEU:HD22	1.81	0.45
1:B:690:ARG:NH2	1:B:699:SER:OG	2.45	0.45
1:B:772:ARG:HB2	1:B:846:ASP:OD2	2.16	0.45
1:A:77:VAL:HG22	1:A:142:THR:HB	1.98	0.45
1:A:286:ARG:HG3	1:A:293:TRP:CE2	2.52	0.45
1:B:659:ARG:NH1	1:B:672:ILE:HG22	2.31	0.45
1:B:383:TRP:CE3	1:B:554:ILE:HD13	2.52	0.45
1:B:239:LEU:HD11	1:B:243:ARG:NH1	2.32	0.44
1:A:769:PHE:O	1:A:775:ARG:NH1	2.44	0.44
1:B:101:ARG:HA	1:B:101:ARG:HD3	1.78	0.44
1:B:390:LYS:HE3	1:B:496:ASP:O	2.17	0.44
1:A:123:ILE:HD13	1:A:264:ARG:HD3	2.00	0.44
1:B:437:ASP:O	1:B:441:GLU:HG3	2.18	0.44
1:A:197:ARG:HG2	1:A:200:ARG:NH1	2.31	0.44
1:A:636:LEU:HD11	1:A:681:ARG:HB3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:GLN:HE22	1:A:356:PHE:HB2	1.82	0.44
1:A:660:MET:HG2	1:A:669:VAL:HG22	2.00	0.44
1:A:690:ARG:NH2	1:A:699:SER:HB3	2.33	0.44
1:A:853:ILE:H	1:A:853:ILE:HG13	1.44	0.44
1:B:454:ASN:O	1:B:476:TRP:HA	2.17	0.44
1:B:770:HIS:CD2	1:B:770:HIS:H	2.35	0.44
1:B:314:THR:HB	1:B:590:THR:HG21	2.00	0.44
1:B:441:GLU:O	1:B:445:ARG:HG3	2.17	0.44
1:A:458:LYS:HD3	1:A:460:GLU:OE2	2.17	0.43
1:A:383:TRP:CE3	1:A:554:ILE:HD13	2.53	0.43
1:A:157:GLU:HB3	1:A:185:CYS:HB2	2.00	0.43
1:B:359:LYS:O	1:B:541:THR:HG21	2.19	0.43
1:A:55:VAL:HB	1:A:114:MET:HE3	2.00	0.43
1:A:491:GLY:O	1:A:495:GLU:HG2	2.18	0.43
1:A:367:PRO:HD2	1:A:546:PHE:CD1	2.53	0.43
1:B:169:GLY:O	1:B:173:GLU:HG2	2.19	0.43
1:B:409:ALA:O	1:B:483:ARG:NH2	2.31	0.43
1:A:541:THR:HA	1:A:600:GLN:HG2	2.01	0.43
1:A:573:THR:HG22	1:A:574:TYR:CD1	2.53	0.43
1:A:454:ASN:O	1:A:476:TRP:HA	2.18	0.42
1:B:89:TYR:OH	1:B:114:MET:O	2.32	0.42
1:A:70:TYR:HB3	1:A:222:VAL:HG21	2.00	0.42
1:B:105:LYS:HD2	1:B:110:HIS:CE1	2.54	0.42
1:B:158:GLU:OE2	1:B:190:THR:HB	2.19	0.42
1:B:64:TRP:NE1	1:B:68:ARG:NH1	2.67	0.42
1:B:62:LEU:HA	1:B:62:LEU:HD23	1.77	0.42
1:B:209:VAL:HG12	1:B:211:LEU:H	1.84	0.42
1:B:628:GLU:HA	1:B:675:ARG:NH2	2.34	0.42
1:B:777:MET:HG3	1:B:861:TRP:HZ2	1.85	0.42
1:B:846:ASP:O	1:B:851:SER:HB2	2.19	0.42
1:A:146:ASP:OD2	2:A:1001:SAH:HG1	2.20	0.42
1:A:90:TYR:O	1:A:93:THR:OG1	2.36	0.42
1:B:640:ARG:HD2	1:B:641:LYS:HD3	2.02	0.42
1:B:164:VAL:O	1:B:168:VAL:HG23	2.19	0.41
1:B:622:ILE:HA	1:B:625:MET:HE3	2.02	0.41
1:B:356:PHE:CD2	1:B:455:MET:HE1	2.55	0.41
1:B:432:PHE:CZ	1:B:480:LEU:HD12	2.55	0.41
1:B:485:LEU:HD12	1:B:485:LEU:HA	1.91	0.41
1:A:749:SER:OG	1:A:751:ARG:HG2	2.20	0.41
1:A:765:GLN:OE1	1:A:811:MET:HG3	2.20	0.41
1:B:643:GLU:H	1:B:643:GLU:HG3	1.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:CYS:HG	1:A:104:THR:HG1	1.58	0.41
1:A:278:ILE:O	1:A:282:ILE:HG12	2.20	0.41
1:A:436:VAL:HG11	1:A:484:PHE:CD2	2.55	0.41
1:B:344:MET:CE	1:B:461:LYS:HE2	2.51	0.41
1:B:605:GLN:HB2	1:B:608:THR:OG1	2.20	0.41
1:A:352:GLN:NE2	1:A:356:PHE:HB2	2.35	0.41
1:A:437:ASP:OD1	1:A:440:ARG:NH1	2.54	0.41
1:B:868:THR:O	1:B:872:VAL:HG23	2.20	0.41
1:A:63:ARG:O	1:A:67:GLU:HG3	2.21	0.41
1:A:834:LYS:HE3	1:A:837:ASP:OD2	2.21	0.41
1:A:113:PRO:HB3	1:A:126:LEU:HD13	2.03	0.40
1:A:239:LEU:HD11	1:A:243:ARG:NH1	2.36	0.40
1:A:476:TRP:N	1:A:476:TRP:HD1	2.17	0.40
1:A:580:LYS:HG2	1:A:592:MET:SD	2.62	0.40
1:A:633:MET:HE2	1:A:681:ARG:NH1	2.37	0.40
1:B:84:ARG:HG2	1:B:114:MET:HE3	2.03	0.40
1:B:285:ILE:HD11	1:B:452:VAL:HG11	2.03	0.40
1:B:103:TYR:CE2	1:B:135:MET:HE1	2.56	0.40
1:B:398:GLU:H	1:B:398:GLU:CD	2.29	0.40
1:B:61:LYS:HG2	1:B:209:VAL:HG11	2.02	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	881/903 (98%)	834 (95%)	44 (5%)	3 (0%)	36	70
1	B	881/903 (98%)	832 (94%)	46 (5%)	3 (0%)	36	70
All	All	1762/1806 (98%)	1666 (95%)	90 (5%)	6 (0%)	36	70

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	640	ARG
1	B	744	PRO
1	A	300	PRO
1	B	300	PRO
1	B	186	PRO
1	A	48	VAL

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	753/767 (98%)	709 (94%)	44 (6%)	18	51
1	B	753/767 (98%)	697 (93%)	56 (7%)	13	42
All	All	1506/1534 (98%)	1406 (93%)	100 (7%)	15	46

All (100) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	THR
1	A	35	VAL
1	A	62	LEU
1	A	101	ARG
1	A	104	THR
1	A	125	ARG
1	A	151	SER
1	A	168	VAL
1	A	181	ILE
1	A	198	LEU
1	A	229	ILE
1	A	243	ARG
1	A	245	ASP
1	A	314	THR
1	A	319	SER
1	A	325	VAL
1	A	337	THR
1	A	339	VAL

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Mol	Chain	Res	Type
1	A	375	MET
1	A	444	LEU
1	A	476	TRP
1	A	479	TRP
1	A	483	ARG
1	A	485	LEU
1	A	515	ARG
1	A	559	GLU
1	A	577	LYS
1	A	591	VAL
1	A	601	ARG
1	A	621	LEU
1	A	639	LEU
1	A	650	GLN
1	A	681	ARG
1	A	683	LEU
1	A	742	VAL
1	A	766	LEU
1	A	775	ARG
1	A	790	VAL
1	A	792	THR
1	A	812	LEU
1	A	846	ASP
1	A	853	ILE
1	A	876	ILE
1	A	883	MET
1	B	6	GLU
1	B	42	ARG
1	B	45	LYS
1	B	48	VAL
1	B	50	THR
1	B	55	VAL
1	B	66	GLU
1	B	93	THR
1	B	95	ARG
1	B	155	GLU
1	B	181	ILE
1	B	201	ARG
1	B	216	THR
1	B	243	ARG
1	B	251	VAL
1	B	252	LYS

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Mol	Chain	Res	Type
1	B	266	VAL
1	B	271	GLU
1	B	292	THR
1	B	314	THR
1	B	325	VAL
1	B	339	VAL
1	B	380	SER
1	B	388	LYS
1	B	389	ARG
1	B	410	LEU
1	B	437	ASP
1	B	439	GLU
1	B	444	LEU
1	B	476	TRP
1	B	479	TRP
1	B	480	LEU
1	B	485	LEU
1	B	515	ARG
1	B	542	ARG
1	B	567	LEU
1	B	591	VAL
1	B	597	ARG
1	B	601	ARG
1	B	621	LEU
1	B	633	MET
1	B	634	GLN
1	B	638	LEU
1	B	639	LEU
1	B	643	GLU
1	B	680	LEU
1	B	720	LEU
1	B	742	VAL
1	B	751	ARG
1	B	795	THR
1	B	814	VAL
1	B	818	VAL
1	B	851	SER
1	B	853	ILE
1	B	871	MET
1	B	878	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	110	HIS
1	A	352	GLN
1	A	402	ASN
1	A	454	ASN
1	A	605	GLN
1	A	620	GLN
1	A	825	HIS
1	B	238	GLN
1	B	287	ASN
1	B	315	GLN
1	B	352	GLN
1	B	497	HIS
1	B	616	ASN
1	B	716	ASN
1	B	762	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](#)

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 for Mogul analysis.

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
4	SO4	A	1005	-	4,4,4	0.24	0	6,6,6	0.06	0
4	SO4	B	1005	-	4,4,4	0.26	0	6,6,6	0.07	0
4	SO4	B	1004	-	4,4,4	0.23	0	6,6,6	0.09	0
2	SAH	B	1001	-	27,28,28	1.08	4 (14%)	36,40,40	2.02	9 (25%)
4	SO4	B	1006	-	4,4,4	0.24	0	6,6,6	0.08	0
4	SO4	A	1006	-	4,4,4	0.24	0	6,6,6	0.10	0
4	SO4	A	1004	-	4,4,4	0.24	0	6,6,6	0.04	0
2	SAH	A	1001	-	27,28,28	1.09	4 (14%)	36,40,40	2.03	10 (27%)

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	1001	-	-	7/15/31/31	0/3/3/3
2	SAH	A	1001	-	-	7/15/31/31	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	SAH	C2-N3	2.53	1.38	1.33
2	B	1001	SAH	C2-N1	2.51	1.38	1.33
2	A	1001	SAH	C2-N1	2.50	1.38	1.33
2	B	1001	SAH	C2-N3	2.49	1.38	1.33
2	A	1001	SAH	OXT-C	-2.22	1.23	1.30
2	B	1001	SAH	C8-N7	2.20	1.35	1.31
2	B	1001	SAH	OXT-C	-2.20	1.23	1.30
2	A	1001	SAH	C8-N7	2.19	1.35	1.31

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1001	SAH	N3-C2-N1	-5.62	120.08	128.58
2	A	1001	SAH	N3-C2-N1	-5.58	120.14	128.58
2	A	1001	SAH	C5-C4-N3	-4.56	120.43	126.72
2	B	1001	SAH	C5-C4-N3	-4.33	120.76	126.72
2	B	1001	SAH	N9-C8-N7	-3.97	108.31	113.94
2	A	1001	SAH	N9-C8-N7	-3.84	108.49	113.94
2	B	1001	SAH	C5'-SD-CG	-3.50	91.86	102.26
2	A	1001	SAH	C2-N3-C4	3.50	120.39	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	SAH	C5'-SD-CG	-3.45	92.01	102.26
2	B	1001	SAH	C2-N3-C4	3.45	120.25	111.83
2	B	1001	SAH	C5-N7-C8	3.37	108.74	103.45
2	A	1001	SAH	C5-N7-C8	3.34	108.70	103.45
2	A	1001	SAH	N3-C4-N9	2.94	132.17	127.17
2	B	1001	SAH	N3-C4-N9	2.73	131.81	127.17
2	B	1001	SAH	OXT-C-O	-2.72	117.91	124.08
2	A	1001	SAH	OXT-C-O	-2.64	118.09	124.08
2	B	1001	SAH	C4-C5-N7	-2.42	107.81	110.58
2	A	1001	SAH	C4-C5-N7	-2.38	107.86	110.58
2	A	1001	SAH	O4'-C1'-N9	2.11	112.14	108.09

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1001	SAH	N-CA-CB-CG
2	B	1001	SAH	N-CA-CB-CG
2	A	1001	SAH	C-CA-CB-CG
2	B	1001	SAH	C-CA-CB-CG
2	A	1001	SAH	OXT-C-CA-CB
2	A	1001	SAH	C2'-C1'-N9-C8
2	B	1001	SAH	C2'-C1'-N9-C8
2	B	1001	SAH	O-C-CA-CB
2	B	1001	SAH	OXT-C-CA-CB
2	A	1001	SAH	O-C-CA-CB
2	A	1001	SAH	CB-CG-SD-C5'
2	B	1001	SAH	CB-CG-SD-C5'
2	A	1001	SAH	O4'-C1'-N9-C8
2	B	1001	SAH	O4'-C1'-N9-C8

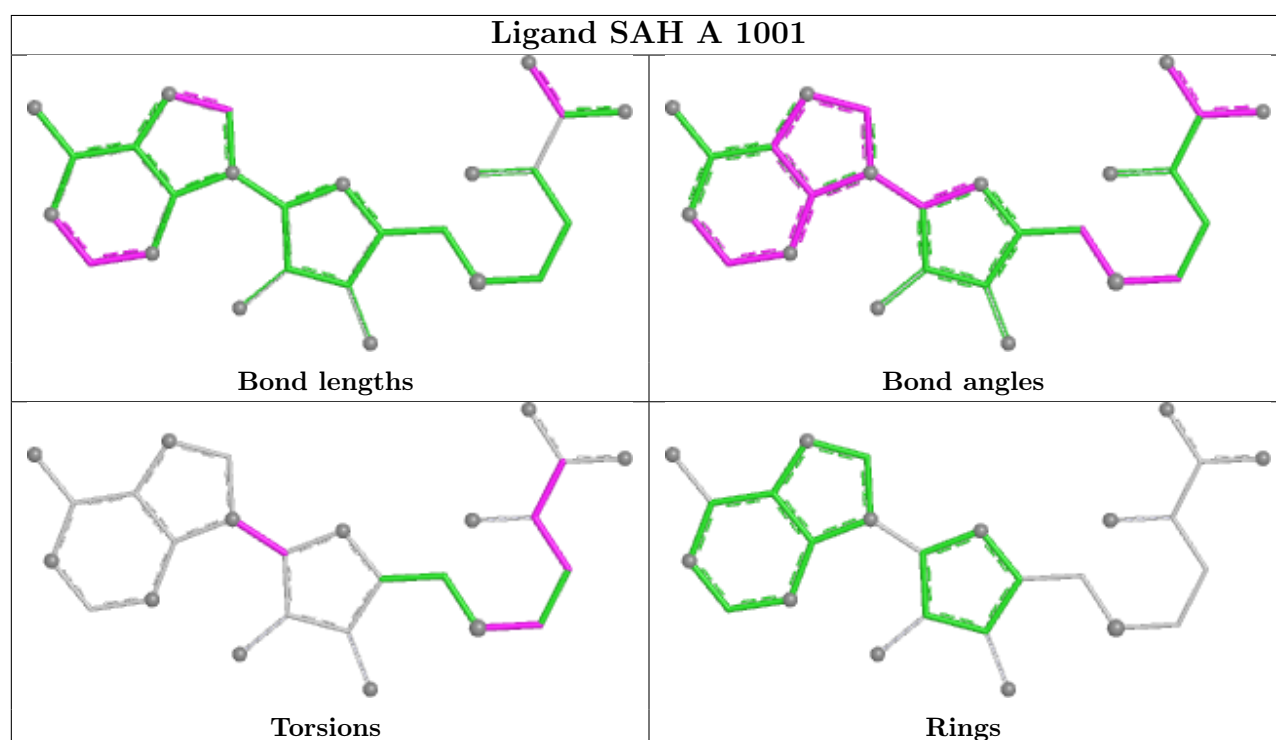
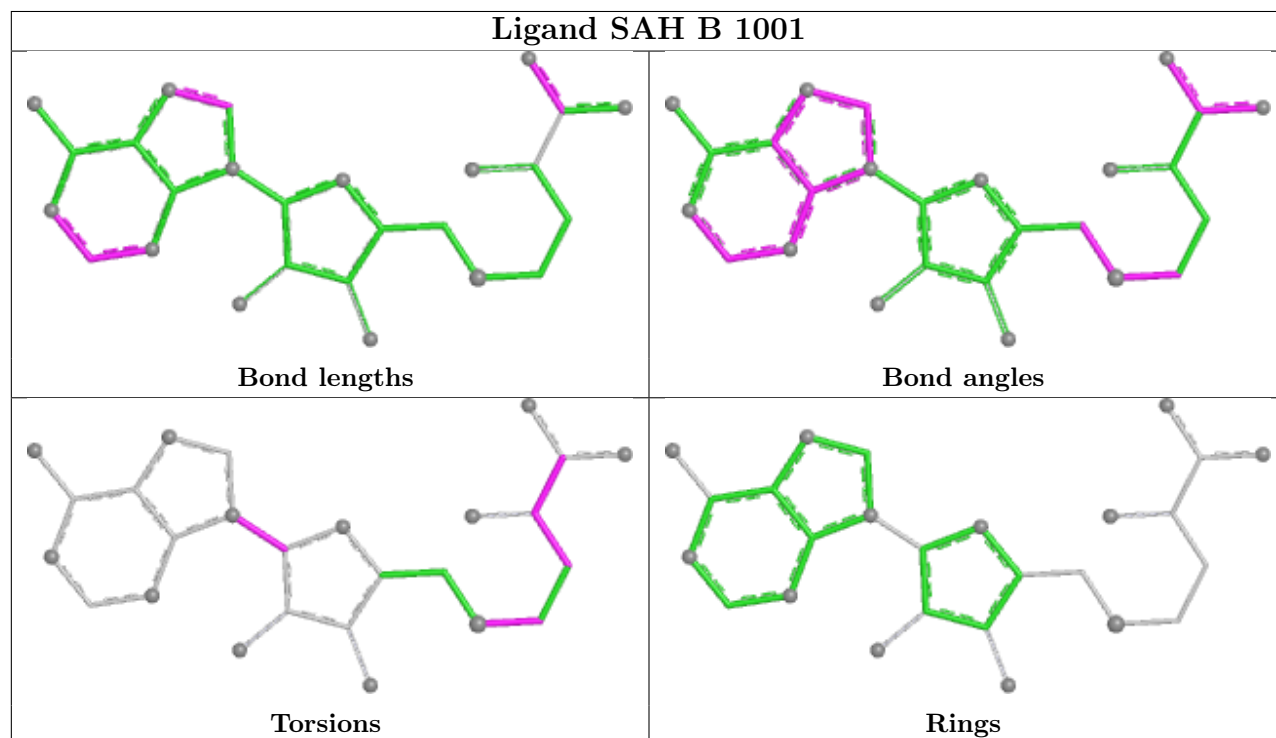
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1001	SAH	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier.

Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	883/903 (97%)	0.30	6 (0%) 84 66	30, 42, 56, 68	0
1	B	883/903 (97%)	0.22	6 (0%) 84 66	24, 34, 48, 60	0
All	All	1766/1806 (97%)	0.26	12 (0%) 84 66	24, 39, 53, 68	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	634	GLN	4.9
1	A	148	GLY	3.0
1	A	39	GLU	2.7
1	A	640	ARG	2.6
1	A	746	ALA	2.4
1	B	746	ALA	2.4
1	B	747	GLY	2.3
1	B	879	GLU	2.2
1	A	522	GLU	2.1
1	B	245	ASP	2.1
1	B	673	ASP	2.0
1	B	47	GLY	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

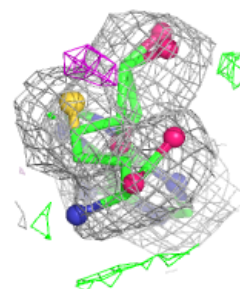
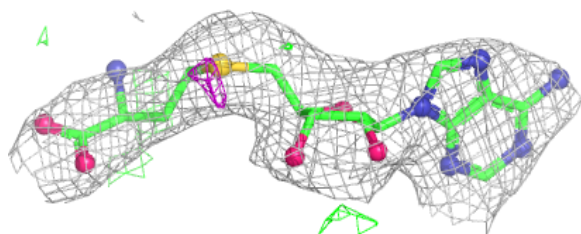
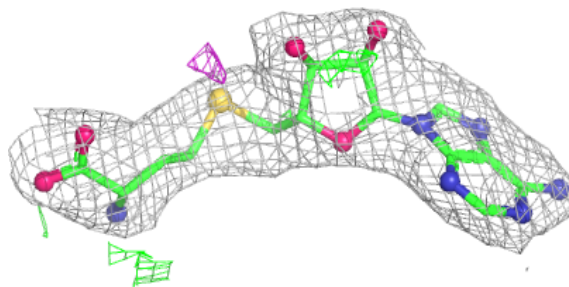
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	SO4	A	1005	5/5	0.69	0.20	49,62,68,77	0
2	SAH	B	1001	26/26	0.90	0.12	23,38,57,60	0
4	SO4	B	1005	5/5	0.90	0.14	44,48,56,65	0
4	SO4	B	1004	5/5	0.91	0.11	27,44,46,48	0
4	SO4	A	1006	5/5	0.91	0.11	45,52,56,59	0
4	SO4	B	1006	5/5	0.91	0.12	35,51,55,57	0
2	SAH	A	1001	26/26	0.93	0.12	28,43,62,65	0
4	SO4	A	1004	5/5	0.94	0.09	36,46,49,51	0
3	ZN	A	1002	1/1	0.97	0.06	30,30,30,30	0
3	ZN	A	1003	1/1	0.98	0.04	44,44,44,44	0
3	ZN	B	1002	1/1	0.99	0.03	27,27,27,27	0
3	ZN	B	1003	1/1	0.99	0.03	29,29,29,29	0

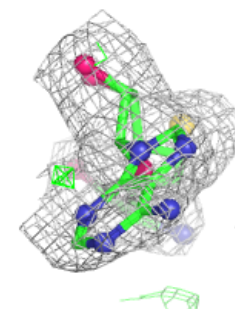
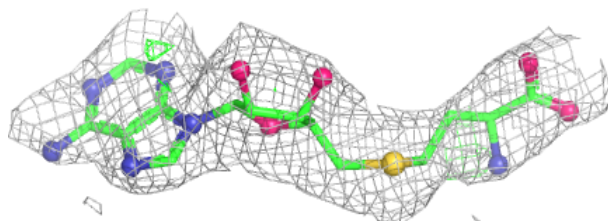
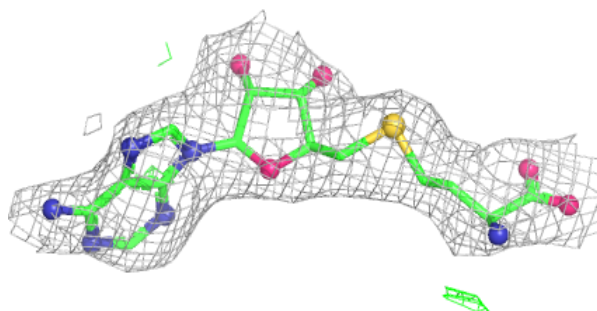
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around SAH B 1001:

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

**Electron density around SAH A 1001:**

$2mF_o-DF_c$ (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.