



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 02:16 AM UTC

PDB ID : 1WY7 / pdb_00001wy7
Title : crystal structure of a putative RNA methyltransferase PH1948 from *Pyrococcus horikoshii*
Authors : Gao, Y.G.; Yao, M.; Tanaka, I.
Deposited on : 2005-02-08
Resolution : 2.20 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

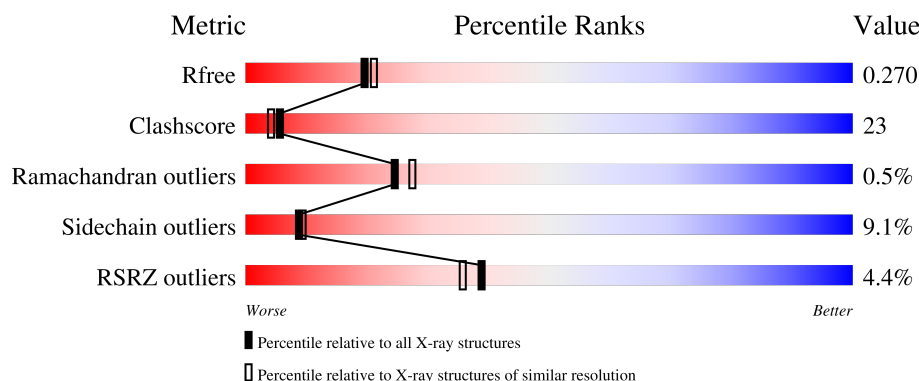
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	207	3% 63% 26% 5% • 5%
1	B	207	3% 64% 26% 5% •
1	C	207	6% 63% 27% 6% • •
1	D	207	5% 57% 35% 5% • •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6823 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.

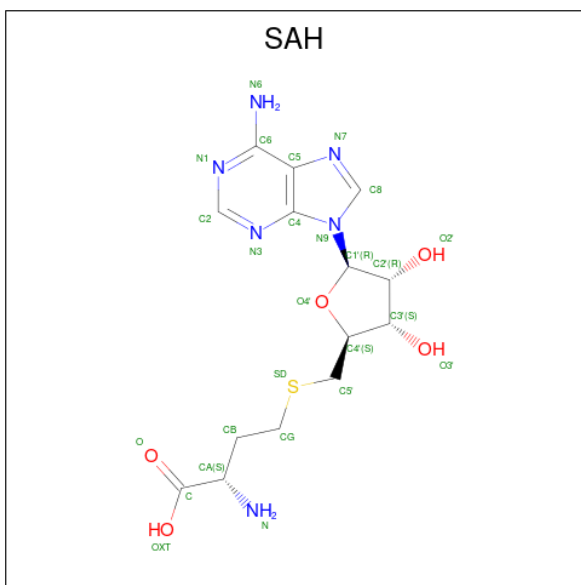
- Molecule 1 is a protein called hypothetical protein PH1948.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	196	Total	C	N	O	S	Se	0	0	0
			1571	1025	269	275	1	1			
1	B	198	Total	C	N	O	S	Se	0	0	0
			1582	1033	267	279	1	2			
1	C	201	Total	C	N	O	S	Se	0	0	0
			1610	1046	276	284	1	3			
1	D	204	Total	C	N	O	S	Se	0	0	0
			1644	1077	278	285	1	3			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP O59611
A	2	MSE	MET	modified residue	UNP O59611
A	118	MSE	MET	modified residue	UNP O59611
B	1	MSE	MET	modified residue	UNP O59611
B	2	MSE	MET	modified residue	UNP O59611
B	118	MSE	MET	modified residue	UNP O59611
C	1	MSE	MET	modified residue	UNP O59611
C	2	MSE	MET	modified residue	UNP O59611
C	118	MSE	MET	modified residue	UNP O59611
D	1	MSE	MET	modified residue	UNP O59611
D	2	MSE	MET	modified residue	UNP O59611
D	118	MSE	MET	modified residue	UNP O59611

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



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		
2	D	1	Total	C	N	O	S	0	0
			26	14	6	5	1		

- Molecule 3 is water.

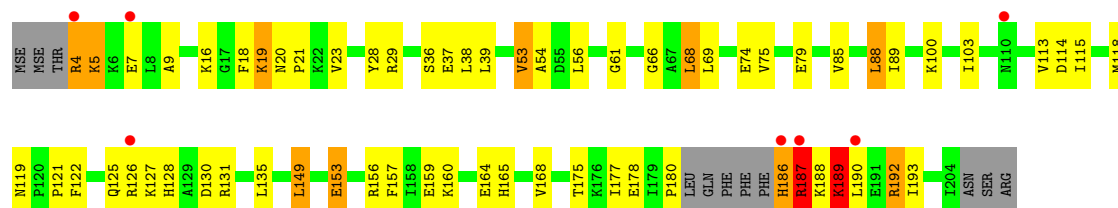
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	100	Total	O	0	0
			100	100		
3	B	85	Total	O	0	0
			85	85		
3	C	64	Total	O	0	0
			64	64		
3	D	63	Total	O	0	0
			63	63		

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.

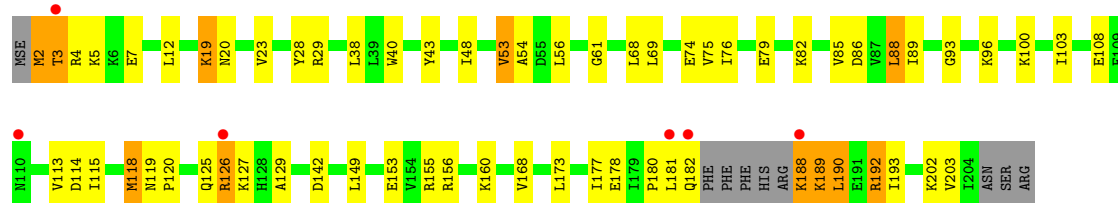
- Molecule 1: hypothetical protein PH1948

Chain A: 



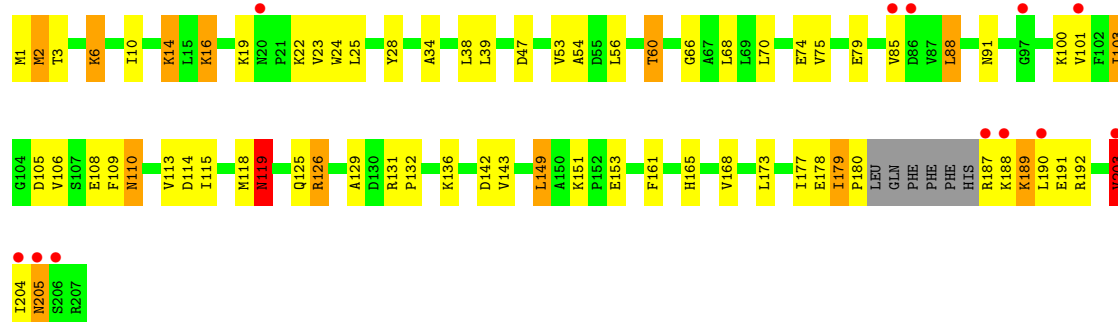
- Molecule 1: hypothetical protein PH1948

Chain B: 

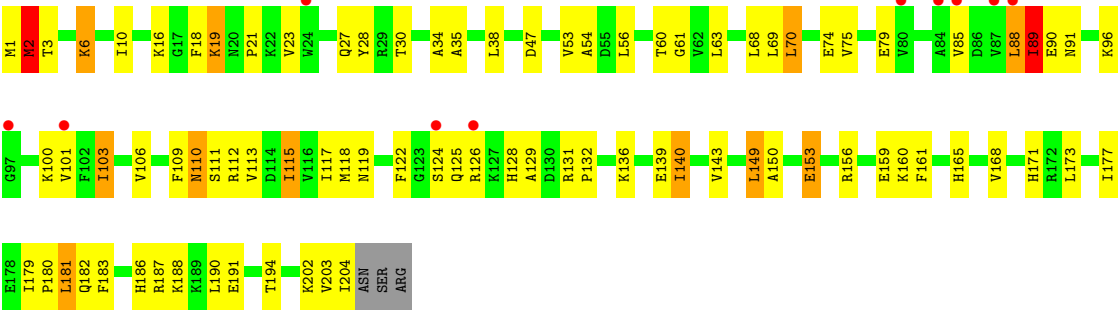


- Molecule 1: hypothetical protein PH1948

Chain C: 



- Molecule 1: hypothetical protein PH1948



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	206.97Å 43.14Å 118.24Å 90.00° 92.08° 90.00°	Depositor
Resolution (Å)	10.00 – 2.20 10.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.3 (10.00-2.20) 98.2 (10.00-2.20)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.59 (at 2.19Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.242 , 0.268 0.244 , 0.270	Depositor DCC
R_{free} test set	5285 reflections (9.99%)	wwPDB-VP
Wilson B-factor (Å ²)	38.2	Xtriage
Anisotropy	0.437	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.44 , 67.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.055 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6823	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.62% 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

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	0.59	2/1603 (0.1%)	0.99	6/2157 (0.3%)
1	B	0.53	2/1613 (0.1%)	0.89	2/2171 (0.1%)
1	C	0.51	4/1640 (0.2%)	0.88	3/2202 (0.1%)
1	D	0.52	4/1679 (0.2%)	0.93	6/2258 (0.3%)
All	All	0.54	12/6535 (0.2%)	0.93	17/8788 (0.2%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	186	HIS	CB-CG	8.91	1.62	1.50
1	B	118	MSE	SE-CE	-8.69	1.69	1.95
1	A	118	MSE	SE-CE	-6.59	1.75	1.95
1	D	118	MSE	SE-CE	-6.29	1.76	1.95
1	C	1	MSE	SE-CE	-6.01	1.77	1.95
1	D	2	MSE	SE-CE	-5.50	1.78	1.95
1	D	1	MSE	SE-CE	-5.47	1.79	1.95
1	B	2	MSE	SE-CE	-5.46	1.79	1.95
1	C	2	MSE	SE-CE	-5.43	1.79	1.95
1	C	118	MSE	CG-SE	-5.36	1.79	1.95
1	C	118	MSE	SE-CE	-5.26	1.79	1.95
1	D	2	MSE	CG-SE	-5.22	1.79	1.95

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	186	HIS	CA-CB-CG	10.11	123.91	113.80
1	A	187	ARG	CA-C-N	-9.00	108.04	122.53
1	A	187	ARG	C-N-CA	-9.00	108.04	122.53
1	D	187	ARG	N-CA-C	7.58	120.62	111.82
1	C	205	ASN	N-CA-C	6.39	124.42	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	140	ILE	N-CA-C	6.12	118.89	112.83
1	A	189	LYS	N-CA-C	5.78	123.11	110.80
1	D	89	ILE	CB-CA-C	-5.76	104.36	112.14
1	D	179	ILE	N-CA-C	5.52	115.07	108.45
1	D	182	GLN	N-CA-C	-5.45	107.00	113.97
1	A	119	ASN	N-CA-C	-5.40	99.14	108.97
1	B	119	ASN	N-CA-C	-5.35	99.24	108.97
1	D	119	ASN	N-CA-C	-5.29	99.33	108.97
1	A	53	VAL	N-CA-C	5.29	116.19	108.53
1	C	203	VAL	N-CA-C	5.20	120.15	109.34
1	C	119	ASN	N-CA-C	-5.15	99.60	108.97
1	B	53	VAL	N-CA-C	5.01	115.80	108.53

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	1571	0	1629	96	4
1	B	1582	0	1644	72	1
1	C	1610	0	1674	73	0
1	D	1644	0	1704	73	0
2	A	26	0	19	3	0
2	B	26	0	19	1	0
2	C	26	0	19	0	0
2	D	26	0	19	2	0
3	A	100	0	0	1	0
3	B	85	0	0	7	3
3	C	64	0	0	3	0
3	D	63	0	0	4	0
All	All	6823	0	6727	310	4

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 23.

All (310) 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:187:ARG:N	1:A:187:ARG:CD	1.70	1.46
1:A:186:HIS:CB	1:A:187:ARG:HD2	1.52	1.39
1:A:187:ARG:CD	1:A:187:ARG:H	1.32	1.29
1:A:186:HIS:HB3	1:A:187:ARG:CD	1.65	1.24
1:A:187:ARG:CZ	1:A:188:LYS:HD2	1.72	1.19
1:A:187:ARG:N	1:A:187:ARG:HD3	0.88	1.03
1:A:187:ARG:NH2	1:A:188:LYS:HD2	1.75	1.00
1:A:186:HIS:C	1:A:187:ARG:HD3	1.88	0.98
1:A:186:HIS:C	1:A:187:ARG:CD	2.42	0.93
1:C:47:ASP:O	1:C:115:ILE:HD13	1.68	0.93
1:C:178:GLU:HG2	1:C:192:ARG:HG2	1.54	0.88
1:A:180:PRO:HA	1:A:190:LEU:HD23	1.56	0.88
1:D:28:TYR:HB3	1:D:181:LEU:HD11	1.57	0.87
1:A:187:ARG:NH1	1:A:188:LYS:HD2	1.89	0.86
1:A:128:HIS:HB3	1:A:131:ARG:HD3	1.56	0.86
1:C:125:GLN:HE21	1:C:125:GLN:HA	1.41	0.85
1:C:126:ARG:HG3	1:C:129:ALA:HB2	1.56	0.85
1:A:187:ARG:CZ	1:A:188:LYS:CD	2.52	0.84
1:A:178:GLU:HB3	1:A:190:LEU:HD13	1.59	0.84
1:B:19:LYS:HD3	1:B:20:ASN:N	1.93	0.84
1:A:189:LYS:HB2	1:A:189:LYS:NZ	1.92	0.84
1:A:186:HIS:CA	1:A:187:ARG:HD2	2.08	0.84
1:D:203:VAL:HG12	1:D:204:ILE:HG13	1.58	0.83
1:A:178:GLU:HG2	1:A:192:ARG:HG3	1.60	0.83
1:C:74:GLU:OE1	1:C:100:LYS:HE3	1.79	0.82
1:B:180:PRO:HB3	1:B:188:LYS:NZ	1.94	0.81
1:D:85:VAL:O	1:D:89:ILE:HG13	1.81	0.80
1:C:187:ARG:HG3	1:C:188:LYS:H	1.47	0.79
1:A:189:LYS:HB2	1:A:189:LYS:HZ3	1.45	0.79
1:A:180:PRO:HG3	1:A:190:LEU:CD2	2.12	0.79
1:C:125:GLN:HA	1:C:125:GLN:NE2	1.97	0.79
1:B:189:LYS:HZ2	1:B:189:LYS:HA	1.47	0.78
1:D:126:ARG:HG3	1:D:129:ALA:HB2	1.66	0.78
1:C:178:GLU:HB3	1:C:190:LEU:HD21	1.65	0.78
1:C:187:ARG:HG3	1:C:188:LYS:N	1.99	0.78
1:A:187:ARG:HB3	1:B:7:GLU:OE2	1.84	0.77
1:A:16:LYS:HD2	1:A:16:LYS:N	1.99	0.77
1:B:3:THR:HG23	1:B:4:ARG:H	1.49	0.77
1:D:101:VAL:HG12	1:D:103:ILE:CD1	2.16	0.75
1:B:189:LYS:HA	1:B:189:LYS:NZ	2.00	0.75
1:D:74:GLU:OE1	1:D:100:LYS:HE3	1.88	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:GLU:CB	1:A:190:LEU:HD13	2.18	0.73
1:C:189:LYS:CD	1:C:189:LYS:H	2.02	0.72
1:B:190:LEU:HD22	1:B:190:LEU:O	1.90	0.72
1:C:23:VAL:HG22	3:C:1059:HOH:O	1.90	0.72
1:A:189:LYS:NZ	1:A:189:LYS:CB	2.50	0.72
1:C:101:VAL:HG12	1:C:103:ILE:HD11	1.70	0.71
1:C:28:TYR:CD1	1:C:179:ILE:HG21	2.25	0.70
1:C:103:ILE:N	1:C:103:ILE:HD12	2.07	0.70
1:A:19:LYS:HD3	1:A:19:LYS:N	2.06	0.69
1:A:125:GLN:C	1:A:126:ARG:HD2	2.17	0.69
1:B:126:ARG:HG3	1:B:129:ALA:HB2	1.73	0.69
1:A:180:PRO:HG3	1:A:190:LEU:HD21	1.74	0.69
1:A:128:HIS:HB3	1:A:131:ARG:CD	2.22	0.68
1:A:186:HIS:CA	1:A:187:ARG:CD	2.70	0.68
1:D:3:THR:HG21	3:D:1051:HOH:O	1.93	0.68
1:B:190:LEU:HD13	1:B:190:LEU:N	2.09	0.67
1:C:3:THR:HG23	1:C:6:LYS:H	1.59	0.67
1:C:14:LYS:HZ3	1:C:14:LYS:HA	1.60	0.67
1:C:203:VAL:HG12	1:C:204:ILE:HG13	1.76	0.67
1:A:186:HIS:N	1:A:187:ARG:NH1	2.43	0.67
1:B:2:MSE:HE2	1:B:5:LYS:NZ	2.10	0.67
1:C:189:LYS:H	1:C:189:LYS:CE	2.09	0.66
1:B:180:PRO:HB3	1:B:188:LYS:HZ2	1.61	0.66
1:D:16:LYS:HG3	1:D:90:GLU:OE2	1.95	0.66
1:B:74:GLU:OE1	1:B:100:LYS:HE2	1.95	0.66
1:A:187:ARG:C	1:A:188:LYS:CG	2.69	0.66
1:A:187:ARG:NH2	1:A:188:LYS:CD	2.57	0.65
1:B:2:MSE:HE3	1:B:2:MSE:N	2.11	0.65
1:C:101:VAL:HG12	1:C:103:ILE:CD1	2.25	0.65
1:A:186:HIS:HB3	1:A:187:ARG:HD2	0.73	0.65
1:D:180:PRO:O	1:D:181:LEU:HD13	1.97	0.64
1:A:125:GLN:O	1:A:126:ARG:HD2	1.96	0.64
1:D:3:THR:HG23	1:D:6:LYS:H	1.62	0.64
1:A:19:LYS:HD3	1:A:19:LYS:H	1.60	0.64
1:B:188:LYS:HD3	1:B:189:LYS:N	2.13	0.64
1:A:85:VAL:O	1:A:89:ILE:HG12	1.97	0.64
1:D:101:VAL:HG12	1:D:103:ILE:HD13	1.80	0.62
1:A:186:HIS:C	1:A:187:ARG:HD2	2.17	0.62
1:A:19:LYS:CD	1:A:20:ASN:H	2.12	0.61
1:D:16:LYS:O	1:D:60:THR:HG21	2.00	0.61
1:D:150:ALA:HB3	1:D:194:THR:HG22	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:181:LEU:O	1:D:181:LEU:HD22	2.00	0.61
1:D:28:TYR:HB3	1:D:181:LEU:CD1	2.29	0.61
1:B:190:LEU:HD13	1:B:190:LEU:H	1.64	0.60
1:B:190:LEU:HD21	1:B:192:ARG:HH11	1.66	0.60
1:B:4:ARG:HG2	1:B:4:ARG:HH11	1.66	0.60
1:C:189:LYS:H	1:C:189:LYS:HE3	1.66	0.60
1:A:187:ARG:NH1	1:A:188:LYS:HG3	2.17	0.60
1:D:161:PHE:O	1:D:165:HIS:HD2	1.85	0.60
1:C:14:LYS:HA	1:C:14:LYS:NZ	2.16	0.60
1:A:38:LEU:HD23	1:A:38:LEU:C	2.27	0.60
1:B:190:LEU:O	1:B:192:ARG:HD2	2.02	0.59
1:A:156:ARG:HG2	1:A:160:LYS:HE2	1.85	0.59
1:C:2:MSE:SE	1:C:10:ILE:HD12	2.53	0.59
1:D:168:VAL:HG13	1:D:203:VAL:HG22	1.85	0.58
1:A:29:ARG:HA	2:A:1001:SAH:N	2.18	0.58
1:A:114:ASP:C	1:A:115:ILE:HD12	2.29	0.58
1:B:2:MSE:HE2	1:B:5:LYS:HZ1	1.69	0.58
1:B:180:PRO:HB3	1:B:188:LYS:HZ1	1.68	0.58
1:A:187:ARG:C	1:A:188:LYS:HG3	2.29	0.57
1:A:189:LYS:HZ3	1:A:189:LYS:CB	2.13	0.57
1:D:180:PRO:C	1:D:181:LEU:HD13	2.30	0.57
1:D:150:ALA:CB	1:D:194:THR:HG22	2.34	0.57
1:D:183:PHE:HB2	1:D:186:HIS:O	2.04	0.57
1:B:190:LEU:CD2	1:B:192:ARG:HH11	2.18	0.57
1:C:54:ALA:HB2	1:C:113:VAL:HG11	1.85	0.57
1:C:34:ALA:HB2	1:C:177:ILE:HG21	1.87	0.57
1:D:69:LEU:O	1:D:69:LEU:HD13	2.04	0.57
1:C:189:LYS:H	1:C:189:LYS:HD3	1.70	0.57
1:B:188:LYS:HD3	1:B:188:LYS:C	2.30	0.56
1:A:180:PRO:CA	1:A:190:LEU:HD23	2.33	0.56
1:A:16:LYS:HD2	1:A:16:LYS:H	1.70	0.56
1:A:5:LYS:HE2	3:A:1020:HOH:O	2.06	0.56
1:C:101:VAL:CG1	1:C:103:ILE:HD11	2.35	0.55
1:C:110:ASN:O	1:C:110:ASN:ND2	2.32	0.55
1:A:19:LYS:HG2	1:A:20:ASN:N	2.21	0.55
1:B:142:ASP:OD2	1:B:202:LYS:HD2	2.06	0.55
1:B:114:ASP:C	1:B:115:ILE:HD12	2.32	0.55
1:A:28:TYR:O	2:A:1001:SAH:HG1	2.07	0.55
1:B:85:VAL:O	1:B:89:ILE:HG12	2.07	0.55
1:C:47:ASP:C	1:C:115:ILE:HD13	2.31	0.54
1:A:187:ARG:NH1	1:A:188:LYS:CD	2.65	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:GLU:OE1	1:A:100:LYS:NZ	2.41	0.54
1:A:135:LEU:HD21	1:A:165:HIS:CE1	2.43	0.54
1:A:19:LYS:CG	1:A:20:ASN:N	2.71	0.54
1:B:40:TRP:NE1	3:B:1037:HOH:O	2.30	0.54
1:B:156:ARG:HD3	1:B:160:LYS:NZ	2.24	0.53
1:D:61:GLY:HA3	1:D:88:LEU:HD22	1.90	0.53
1:A:177:ILE:HD11	1:A:193:ILE:HD11	1.90	0.53
1:B:38:LEU:C	1:B:38:LEU:HD23	2.33	0.53
1:B:54:ALA:HB2	1:B:113:VAL:HG11	1.91	0.53
1:A:54:ALA:HB2	1:A:113:VAL:HG11	1.90	0.53
1:A:159:GLU:HG2	3:C:1023:HOH:O	2.09	0.53
1:B:173:LEU:HD22	1:D:171:HIS:CD2	2.44	0.53
1:B:126:ARG:CG	1:B:129:ALA:HB2	2.38	0.53
1:D:136:LYS:O	1:D:140:ILE:HG12	2.08	0.53
1:C:114:ASP:C	1:C:115:ILE:HG13	2.34	0.53
1:B:188:LYS:HD2	1:B:190:LEU:HD12	1.90	0.52
1:B:19:LYS:HD3	1:B:20:ASN:H	1.74	0.52
1:D:79:GLU:O	1:D:103:ILE:HA	2.09	0.52
1:B:178:GLU:HB3	1:B:190:LEU:HD23	1.91	0.52
1:C:16:LYS:NZ	1:C:16:LYS:HB3	2.24	0.52
1:A:160:LYS:O	1:A:164:GLU:HG3	2.09	0.52
1:A:186:HIS:N	1:A:187:ARG:HH11	2.08	0.52
1:D:27:GLN:HG2	2:D:1004:SAH:H3'	1.92	0.52
1:D:18:PHE:HB2	1:D:21:PRO:HG3	1.92	0.52
1:D:103:ILE:HD13	1:D:103:ILE:N	2.25	0.52
1:D:38:LEU:C	1:D:38:LEU:HD23	2.36	0.52
1:D:153:GLU:CD	1:D:153:GLU:H	2.17	0.51
1:D:183:PHE:CE1	1:D:191:GLU:HB2	2.44	0.51
1:A:4:ARG:O	1:A:5:LYS:C	2.53	0.51
1:A:180:PRO:HA	1:A:190:LEU:CD2	2.35	0.51
1:C:60:THR:HG23	1:C:91:ASN:ND2	2.25	0.51
1:C:125:GLN:HE21	1:C:125:GLN:CA	2.13	0.51
1:D:203:VAL:C	1:D:204:ILE:HG13	2.35	0.51
1:C:178:GLU:HB3	1:C:190:LEU:CD2	2.39	0.51
1:B:76:ILE:HG12	1:B:100:LYS:HE3	1.91	0.50
1:C:151:LYS:HE2	1:C:153:GLU:OE1	2.10	0.50
1:D:2:MSE:SE	1:D:10:ILE:HD12	2.61	0.50
1:D:18:PHE:CE2	1:D:27:GLN:HB3	2.47	0.50
1:D:30:THR:HG22	1:D:35:ALA:HB2	1.92	0.50
1:C:34:ALA:HB2	1:C:177:ILE:CG2	2.42	0.50
1:C:189:LYS:HE3	1:C:190:LEU:H	1.77	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:PHE:CG	1:A:149:LEU:HD13	2.46	0.50
1:A:187:ARG:O	1:A:188:LYS:HG2	2.11	0.50
1:C:60:THR:HG23	1:C:91:ASN:HD22	1.75	0.50
1:C:189:LYS:HD3	1:C:189:LYS:N	2.27	0.50
1:A:79:GLU:O	1:A:103:ILE:HA	2.12	0.49
1:C:2:MSE:HG3	1:C:6:LYS:HZ3	1.77	0.49
1:D:131:ARG:HB2	1:D:132:PRO:HD3	1.95	0.49
1:D:181:LEU:HD12	3:D:1022:HOH:O	2.12	0.49
1:A:68:LEU:HD22	1:A:75:VAL:HG21	1.95	0.49
1:B:190:LEU:N	1:B:190:LEU:CD1	2.74	0.49
1:C:38:LEU:C	1:C:38:LEU:HD23	2.38	0.49
1:C:189:LYS:CD	1:C:189:LYS:N	2.73	0.49
1:A:180:PRO:HG3	1:A:190:LEU:HD22	1.91	0.49
1:B:93:GLY:O	1:B:96:LYS:HD3	2.12	0.49
1:C:114:ASP:O	1:C:115:ILE:HG13	2.12	0.49
1:C:79:GLU:O	1:C:103:ILE:HA	2.13	0.49
1:D:168:VAL:HG13	1:D:203:VAL:CG2	2.43	0.49
1:B:43:TYR:HB2	1:B:48:ILE:HD12	1.95	0.49
1:A:178:GLU:HB3	1:A:190:LEU:HB3	1.96	0.48
1:D:16:LYS:O	1:D:60:THR:CG2	2.60	0.48
1:B:82:LYS:HE2	1:B:86:ASP:OD1	2.13	0.48
1:D:109:PHE:CZ	1:D:111:SER:HB3	2.49	0.48
1:B:3:THR:HG23	1:B:4:ARG:N	2.23	0.48
1:A:4:ARG:HG3	1:A:37:GLU:OE1	2.13	0.48
1:C:6:LYS:NZ	1:C:6:LYS:HB3	2.28	0.48
1:D:63:LEU:HB3	1:D:117:ILE:HD12	1.96	0.48
1:A:88:LEU:HD12	1:A:89:ILE:HD13	1.95	0.48
1:C:88:LEU:HD13	1:C:88:LEU:C	2.39	0.48
1:B:4:ARG:NH1	3:B:1016:HOH:O	2.47	0.47
1:B:61:GLY:HA3	1:B:88:LEU:HD22	1.95	0.47
1:C:24:TRP:CD1	1:C:25:LEU:HG	2.49	0.47
1:C:153:GLU:CD	1:C:153:GLU:H	2.22	0.47
1:D:69:LEU:HD13	1:D:69:LEU:C	2.39	0.47
1:C:125:GLN:NE2	1:C:125:GLN:CA	2.69	0.47
1:D:110:ASN:N	1:D:110:ASN:HD22	2.13	0.47
1:D:181:LEU:O	1:D:181:LEU:CD2	2.62	0.47
1:A:122:PHE:CD2	1:A:149:LEU:CD1	2.97	0.47
1:C:6:LYS:HZ3	1:C:6:LYS:HB3	1.80	0.47
1:C:60:THR:HG23	1:C:60:THR:O	2.13	0.47
1:B:7:GLU:HG3	3:B:1051:HOH:O	2.14	0.47
1:B:23:VAL:O	1:B:125:GLN:HG3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:ARG:CD	1:B:129:ALA:HB2	2.46	0.46
1:A:156:ARG:CG	1:A:160:LYS:HE2	2.46	0.46
1:C:105:ASP:HB3	1:C:108:GLU:HG3	1.97	0.46
1:D:79:GLU:OE2	2:D:1004:SAH:O2'	2.30	0.46
1:A:19:LYS:CG	1:A:20:ASN:H	2.27	0.46
1:A:122:PHE:CD2	1:A:149:LEU:HD13	2.50	0.46
1:A:178:GLU:HB3	1:A:190:LEU:CD1	2.38	0.46
1:C:187:ARG:CG	1:C:188:LYS:H	2.13	0.46
1:D:54:ALA:HB2	1:D:113:VAL:HG11	1.97	0.46
1:D:88:LEU:C	1:D:88:LEU:HD13	2.39	0.46
1:A:4:ARG:HB3	1:A:4:ARG:NH1	2.30	0.46
1:B:4:ARG:HG2	1:B:4:ARG:NH1	2.31	0.46
1:B:4:ARG:HD3	1:B:40:TRP:CH2	2.51	0.46
1:D:112:ARG:NH1	1:D:139:GLU:O	2.48	0.46
1:A:187:ARG:H	1:A:187:ARG:HH11	1.63	0.46
1:B:79:GLU:O	1:B:103:ILE:HA	2.15	0.46
1:B:88:LEU:HD13	1:B:88:LEU:C	2.41	0.46
1:B:2:MSE:HA	1:B:5:LYS:HE2	1.98	0.45
1:D:202:LYS:NZ	3:D:1023:HOH:O	2.49	0.45
1:C:2:MSE:HG3	1:C:6:LYS:NZ	2.31	0.45
1:C:131:ARG:HB2	1:C:132:PRO:HD3	1.98	0.45
1:D:53:VAL:HG13	1:D:115:ILE:HG22	1.99	0.45
1:A:180:PRO:CA	1:A:190:LEU:CD2	2.93	0.45
1:C:56:LEU:HG	1:C:106:VAL:HG23	1.98	0.45
1:C:22:LYS:HD3	1:C:24:TRP:CZ2	2.52	0.45
1:D:16:LYS:HE2	1:D:90:GLU:OE1	2.17	0.45
1:A:180:PRO:CG	1:A:190:LEU:CD2	2.88	0.45
1:B:155:ARG:HH22	1:D:159:GLU:CD	2.24	0.45
1:D:60:THR:CG2	1:D:91:ASN:ND2	2.80	0.45
1:C:2:MSE:CG	1:C:6:LYS:HZ3	2.29	0.45
1:D:23:VAL:O	1:D:125:GLN:HG3	2.16	0.45
1:D:60:THR:HG22	1:D:60:THR:O	2.17	0.45
1:A:187:ARG:NH1	1:A:188:LYS:CG	2.80	0.45
1:D:34:ALA:HB2	1:D:177:ILE:CG2	2.47	0.44
1:A:187:ARG:O	1:A:188:LYS:CG	2.64	0.44
1:D:47:ASP:OD1	1:D:143:VAL:HG11	2.18	0.44
1:D:156:ARG:HH11	1:D:160:LYS:HB2	1.82	0.44
1:A:187:ARG:CZ	1:A:188:LYS:CG	2.95	0.44
1:B:2:MSE:HE2	1:B:5:LYS:HZ3	1.80	0.44
1:B:3:THR:C	1:B:5:LYS:H	2.24	0.44
1:A:16:LYS:N	1:A:16:LYS:CD	2.78	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:160:LYS:HB2	1:D:160:LYS:HE3	1.78	0.44
1:A:18:PHE:HB2	1:A:21:PRO:HG3	2.00	0.44
1:C:119:ASN:C	1:C:119:ASN:HD22	2.24	0.44
1:B:190:LEU:HD21	1:B:192:ARG:NH1	2.32	0.44
1:D:89:ILE:HG13	1:D:89:ILE:H	1.63	0.44
1:B:189:LYS:HZ2	1:B:190:LEU:H	1.65	0.43
1:D:122:PHE:CD1	1:D:149:LEU:HD13	2.53	0.43
1:A:53:VAL:O	1:A:75:VAL:HA	2.19	0.43
1:B:177:ILE:HD11	1:B:193:ILE:HD11	2.00	0.43
1:C:85:VAL:HG21	1:C:103:ILE:HG13	2.00	0.43
1:B:4:ARG:HD3	1:B:40:TRP:HH2	1.83	0.43
1:B:156:ARG:HD3	1:B:160:LYS:HZ2	1.83	0.43
1:A:4:ARG:CG	1:A:37:GLU:OE1	2.66	0.43
1:A:153:GLU:H	1:A:153:GLU:CD	2.27	0.43
1:C:39:LEU:HD13	1:C:66:GLY:HA3	2.01	0.43
1:C:109:PHE:O	1:C:136:LYS:HD2	2.19	0.43
1:D:128:HIS:HB3	1:D:131:ARG:HD2	2.00	0.43
1:C:2:MSE:HG3	1:C:6:LYS:HE2	2.00	0.43
1:A:88:LEU:C	1:A:88:LEU:HD13	2.43	0.43
1:A:131:ARG:NH1	1:A:157:PHE:CE2	2.87	0.43
1:D:19:LYS:HA	1:D:19:LYS:CE	2.49	0.43
1:D:190:LEU:N	1:D:190:LEU:HD22	2.34	0.43
1:C:149:LEU:HB2	3:C:1025:HOH:O	2.19	0.42
1:D:53:VAL:O	1:D:75:VAL:HA	2.19	0.42
1:C:103:ILE:N	1:C:103:ILE:CD1	2.77	0.42
1:D:115:ILE:HD11	3:D:1063:HOH:O	2.18	0.42
1:A:61:GLY:HA3	1:A:88:LEU:HD22	2.02	0.42
1:D:109:PHE:O	1:D:136:LYS:NZ	2.47	0.42
1:C:23:VAL:O	1:C:125:GLN:HG3	2.20	0.42
1:C:142:ASP:OD1	1:C:205:ASN:OD1	2.37	0.42
1:B:118:MSE:HG3	1:B:120:PRO:HD3	2.01	0.42
1:A:19:LYS:HE2	1:A:20:ASN:H	1.84	0.42
1:B:126:ARG:HB3	3:B:1063:HOH:O	2.19	0.42
1:B:168:VAL:HG13	1:B:203:VAL:HG22	2.02	0.42
1:C:190:LEU:HD13	1:C:190:LEU:C	2.45	0.42
1:D:19:LYS:HA	1:D:19:LYS:NZ	2.35	0.41
1:D:56:LEU:HG	1:D:106:VAL:HG23	2.02	0.41
1:A:39:LEU:HD13	1:A:66:GLY:HA3	2.02	0.41
1:C:161:PHE:O	1:C:165:HIS:HD2	2.03	0.41
1:A:178:GLU:O	1:A:180:PRO:HD3	2.20	0.41
1:B:53:VAL:O	1:B:75:VAL:HA	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:ARG:HA	2:A:1001:SAH:HN2	1.83	0.41
1:D:128:HIS:HB3	1:D:131:ARG:CD	2.51	0.41
1:A:121:PRO:O	1:A:130:ASP:HB3	2.20	0.41
1:B:189:LYS:HZ2	1:B:189:LYS:CA	2.26	0.41
1:D:70:LEU:HD12	1:D:70:LEU:HA	1.91	0.41
1:D:203:VAL:HG12	1:D:204:ILE:CG1	2.39	0.41
1:A:23:VAL:O	1:A:125:GLN:HG3	2.21	0.41
1:A:9:ALA:HB2	1:A:36:SER:OG	2.21	0.41
1:B:155:ARG:NH2	1:D:159:GLU:CD	2.79	0.41
1:B:188:LYS:HA	3:B:1062:HOH:O	2.21	0.41
1:B:12:LEU:O	1:B:29:ARG:NH1	2.54	0.41
1:B:181:LEU:C	1:B:181:LEU:HD12	2.45	0.41
1:C:85:VAL:CG1	1:C:101:VAL:HG11	2.51	0.41
1:B:3:THR:C	1:B:5:LYS:N	2.80	0.40
1:B:188:LYS:C	1:B:188:LYS:CD	2.92	0.40
1:B:28:TYR:O	2:B:1002:SAH:HG1	2.21	0.40
1:B:108:GLU:OE1	3:B:1060:HOH:O	2.20	0.40
1:C:168:VAL:HG13	1:C:203:VAL:HG22	2.03	0.40
1:A:187:ARG:CZ	1:A:188:LYS:HG3	2.51	0.40
1:B:181:LEU:HB3	3:B:1061:HOH:O	2.20	0.40
1:C:47:ASP:OD1	1:C:143:VAL:HG11	2.21	0.40
1:C:53:VAL:O	1:C:75:VAL:HA	2.20	0.40
1:C:179:ILE:HA	1:C:180:PRO:HD3	1.82	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:LYS:NZ	3:B:1031:HOH:O[1_545]	0.54	1.66
1:A:188:LYS:CE	3:B:1031:HOH:O[1_545]	1.35	0.85
1:A:188:LYS:CD	3:B:1031:HOH:O[1_545]	2.05	0.15
1:A:186:HIS:N	1:B:125:GLN:OE1[1_545]	2.14	0.06

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	192/207 (93%)	186 (97%)	4 (2%)	2 (1%)	12	11
1	B	194/207 (94%)	188 (97%)	5 (3%)	1 (0%)	24	27
1	C	197/207 (95%)	192 (98%)	4 (2%)	1 (0%)	24	27
1	D	202/207 (98%)	194 (96%)	8 (4%)	0	100	100
All	All	785/828 (95%)	760 (97%)	21 (3%)	4 (0%)	24	27

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	189	LYS
1	C	203	VAL
1	B	3	THR
1	A	5	LYS

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	169/177 (96%)	154 (91%)	15 (9%)	9	10
1	B	171/177 (97%)	157 (92%)	14 (8%)	10	12
1	C	174/177 (98%)	157 (90%)	17 (10%)	7	8
1	D	177/177 (100%)	160 (90%)	17 (10%)	8	8
All	All	691/708 (98%)	628 (91%)	63 (9%)	9	9

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ARG
1	A	7	GLU
1	A	19	LYS
1	A	56	LEU

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Mol	Chain	Res	Type
1	A	68	LEU
1	A	69	LEU
1	A	88	LEU
1	A	127	LYS
1	A	149	LEU
1	A	153	GLU
1	A	168	VAL
1	A	175	THR
1	A	187	ARG
1	A	189	LYS
1	A	192	ARG
1	B	19	LYS
1	B	56	LEU
1	B	68	LEU
1	B	69	LEU
1	B	88	LEU
1	B	126	ARG
1	B	127	LYS
1	B	149	LEU
1	B	153	GLU
1	B	182	GLN
1	B	188	LYS
1	B	189	LYS
1	B	190	LEU
1	B	192	ARG
1	C	6	LYS
1	C	14	LYS
1	C	16	LYS
1	C	19	LYS
1	C	60	THR
1	C	68	LEU
1	C	70	LEU
1	C	88	LEU
1	C	103	ILE
1	C	110	ASN
1	C	119	ASN
1	C	126	ARG
1	C	149	LEU
1	C	173	LEU
1	C	179	ILE
1	C	189	LYS
1	C	191	GLU

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Mol	Chain	Res	Type
1	D	2	MSE
1	D	6	LYS
1	D	19	LYS
1	D	68	LEU
1	D	70	LEU
1	D	88	LEU
1	D	89	ILE
1	D	96	LYS
1	D	103	ILE
1	D	110	ASN
1	D	115	ILE
1	D	124	SER
1	D	149	LEU
1	D	153	GLU
1	D	173	LEU
1	D	181	LEU
1	D	188	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	ASN
1	A	125	GLN
1	A	165	HIS
1	B	110	ASN
1	B	119	ASN
1	B	125	GLN
1	B	128	HIS
1	B	165	HIS
1	C	33	ASN
1	C	119	ASN
1	C	125	GLN
1	C	165	HIS
1	D	110	ASN
1	D	119	ASN
1	D	125	GLN
1	D	165	HIS
1	D	186	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 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	A	1001	-	27,28,28	0.76	1 (3%)	36,40,40	1.39	3 (8%)
2	SAH	B	1002	-	27,28,28	0.81	1 (3%)	36,40,40	1.27	3 (8%)
2	SAH	C	1003	-	27,28,28	0.68	0	36,40,40	0.97	2 (5%)
2	SAH	D	1004	-	27,28,28	0.82	1 (3%)	36,40,40	1.06	3 (8%)

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	A	1001	-	-	3/15/31/31	0/3/3/3
2	SAH	B	1002	-	-	4/15/31/31	0/3/3/3
2	SAH	C	1003	-	-	3/15/31/31	0/3/3/3
2	SAH	D	1004	-	-	6/15/31/31	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	SAH	C4-N3	2.22	1.38	1.34
2	D	1004	SAH	C4-N3	2.15	1.38	1.34
2	B	1002	SAH	C4-N3	2.13	1.38	1.34

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	SAH	CB-CG-SD	4.67	123.88	113.45
2	B	1002	SAH	CB-CG-SD	3.66	121.61	113.45
2	B	1002	SAH	O2'-C2'-C3'	3.18	122.00	111.82
2	A	1001	SAH	O2'-C2'-C3'	3.04	121.58	111.82
2	D	1004	SAH	O2'-C2'-C3'	2.88	121.05	111.82
2	B	1002	SAH	O3'-C3'-C4'	-2.82	102.98	111.08
2	C	1003	SAH	O2'-C2'-C3'	2.68	120.40	111.82
2	A	1001	SAH	O3'-C3'-C4'	-2.59	103.65	111.08
2	D	1004	SAH	O3'-C3'-C4'	-2.47	103.99	111.08
2	D	1004	SAH	C2'-C3'-C4'	2.13	106.73	102.61
2	C	1003	SAH	O3'-C3'-C4'	-2.09	105.09	111.08

There are no chirality outliers.

All (16) torsion outliers are listed below:

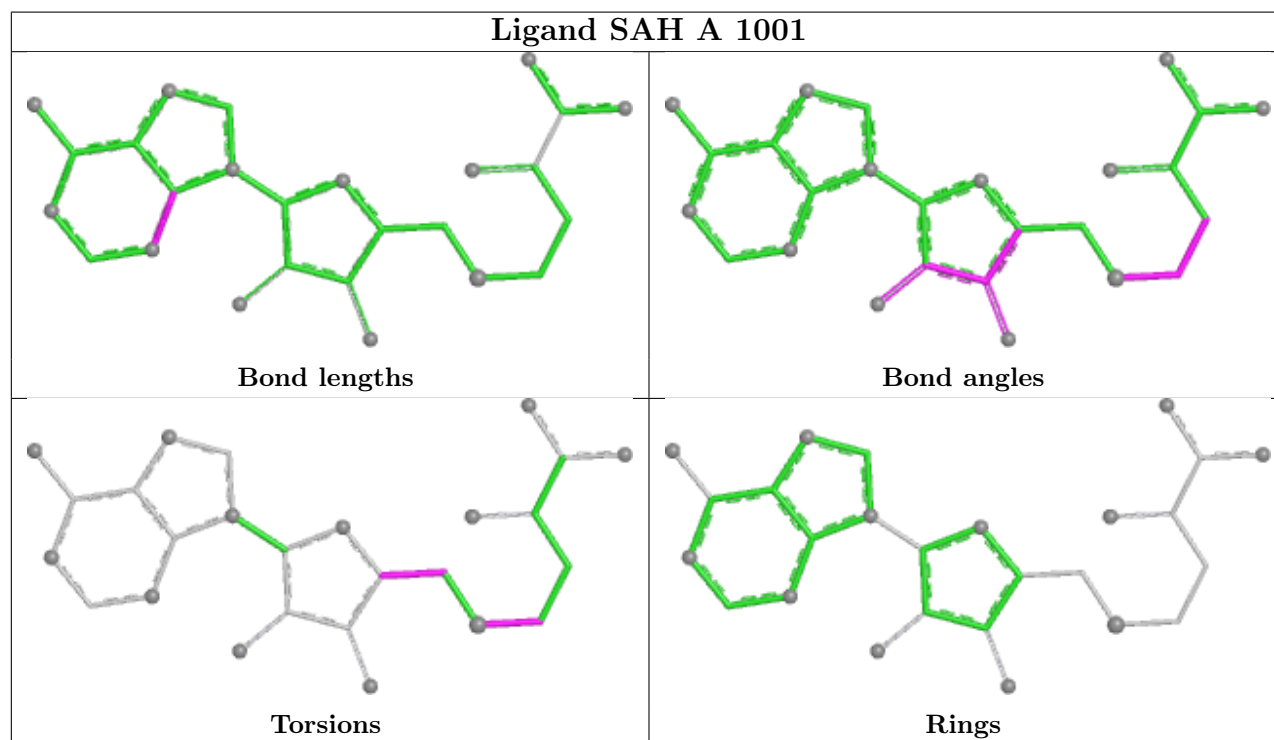
Mol	Chain	Res	Type	Atoms
2	A	1001	SAH	CB-CG-SD-C5'
2	A	1001	SAH	O4'-C4'-C5'-SD
2	A	1001	SAH	C3'-C4'-C5'-SD
2	B	1002	SAH	N-CA-CB-CG
2	B	1002	SAH	C-CA-CB-CG
2	B	1002	SAH	O4'-C4'-C5'-SD
2	B	1002	SAH	C3'-C4'-C5'-SD
2	D	1004	SAH	O4'-C4'-C5'-SD
2	D	1004	SAH	C3'-C4'-C5'-SD
2	C	1003	SAH	C-CA-CB-CG
2	D	1004	SAH	O-C-CA-CB
2	C	1003	SAH	N-CA-CB-CG
2	D	1004	SAH	OXT-C-CA-CB
2	C	1003	SAH	O4'-C4'-C5'-SD
2	D	1004	SAH	CB-CG-SD-C5'
2	D	1004	SAH	OXT-C-CA-N

There are no ring outliers.

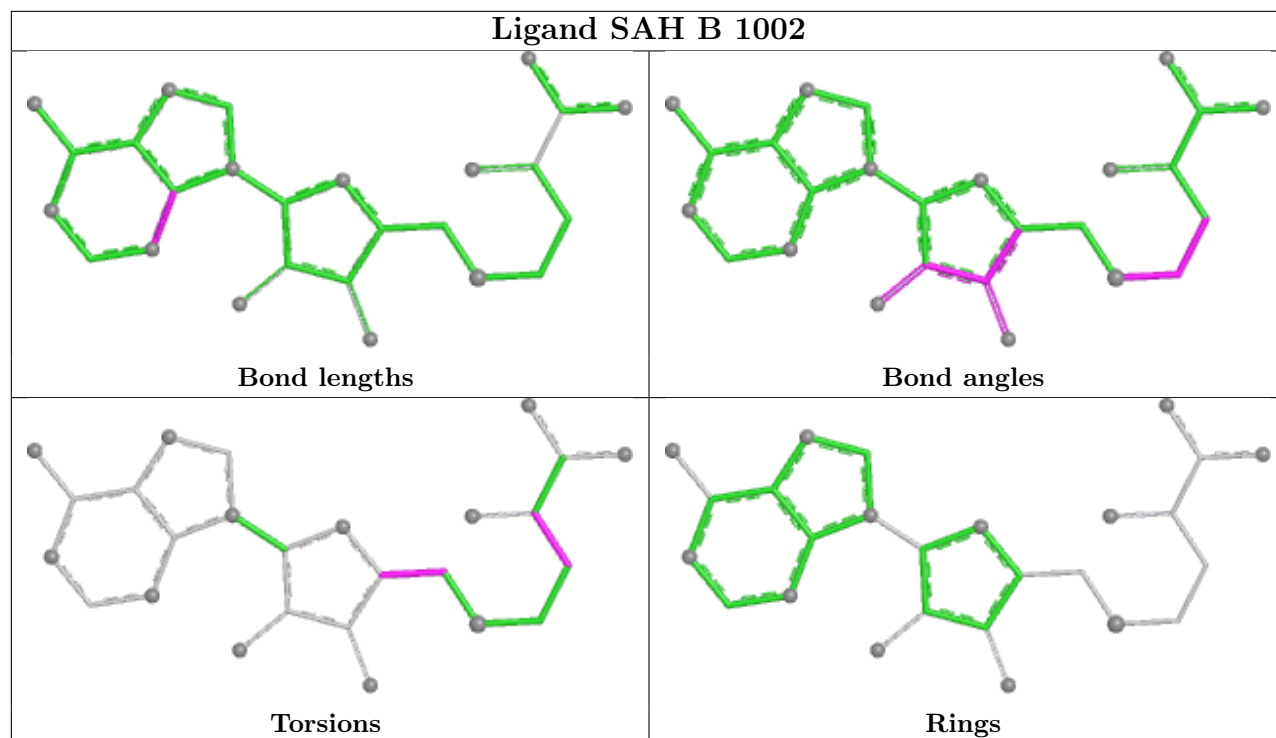
3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1001	SAH	3	0
2	B	1002	SAH	1	0
2	D	1004	SAH	2	0

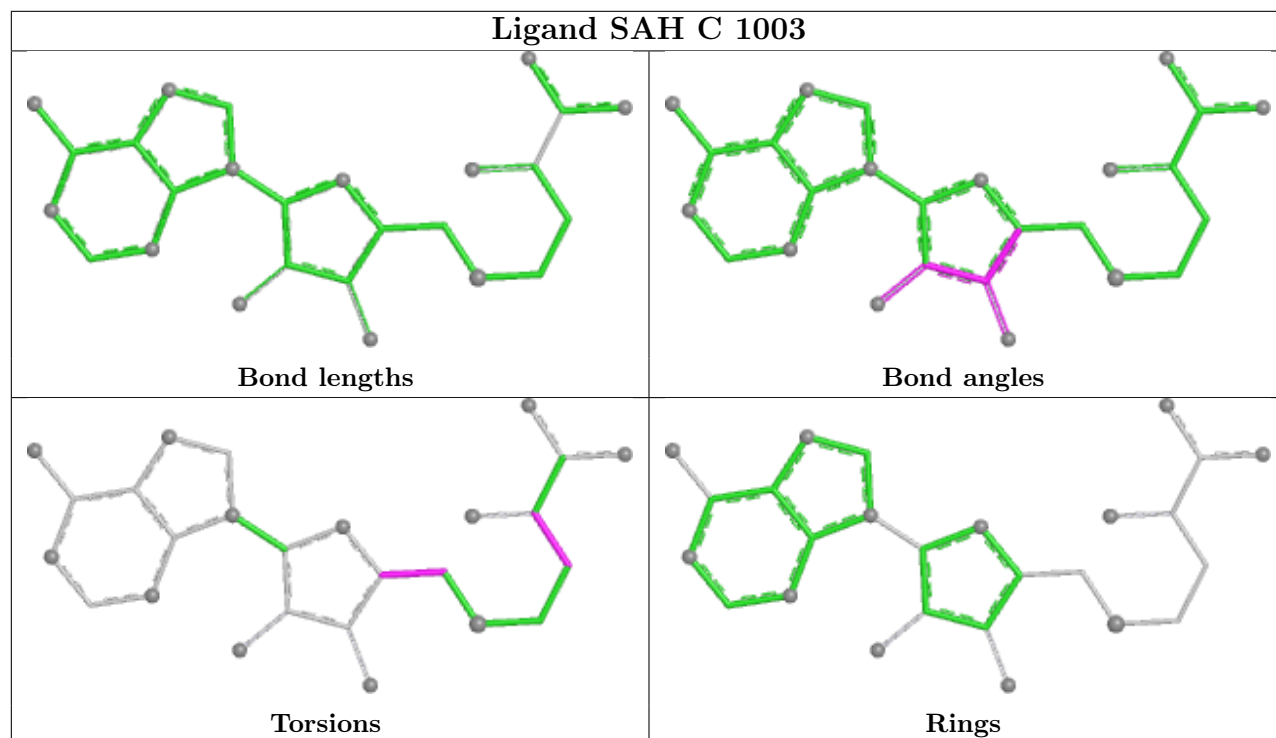
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

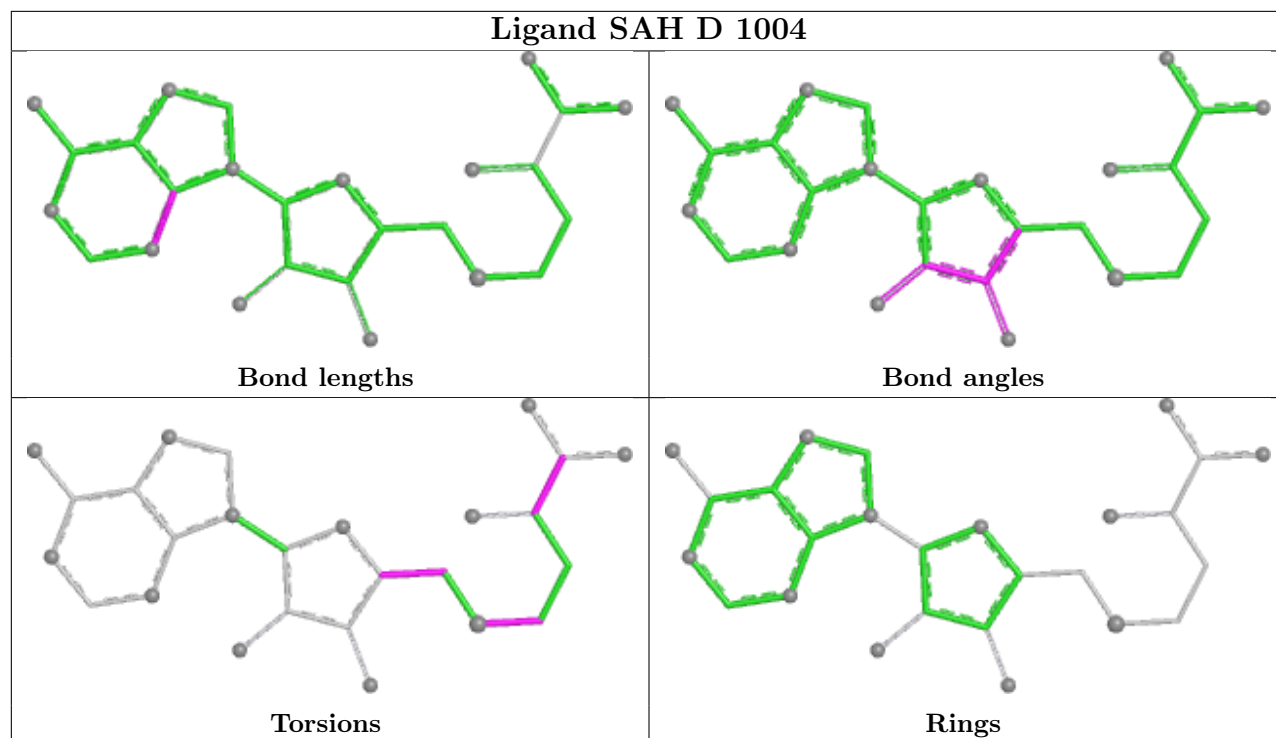


Ligand SAH B 1002



Ligand SAH C 1003





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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	A	195/207 (94%)	-0.06	7 (3%)	46	43	24, 37, 63, 98	0
1	B	196/207 (94%)	-0.02	6 (3%)	51	48	22, 38, 67, 88	0
1	C	198/207 (95%)	0.29	12 (6%)	27	24	29, 46, 70, 96	0
1	D	201/207 (97%)	0.31	10 (4%)	34	31	28, 47, 67, 80	0
All	All	790/828 (95%)	0.13	35 (4%)	39	36	22, 42, 68, 98	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	124	SER	5.6
1	D	97	GLY	4.6
1	B	181	LEU	4.2
1	A	187	ARG	4.1
1	C	86	ASP	3.6
1	A	186	HIS	3.6
1	A	190	LEU	3.5
1	C	190	LEU	3.3
1	D	80	VAL	2.8
1	C	97	GLY	2.8
1	C	187	ARG	2.8
1	B	3	THR	2.8
1	D	87	VAL	2.7
1	B	182	GLN	2.6
1	C	203	VAL	2.6
1	D	101	VAL	2.6
1	D	126	ARG	2.6
1	C	85	VAL	2.6
1	C	101	VAL	2.5
1	A	7	GLU	2.5
1	B	110	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	188	LYS	2.5
1	B	188	LYS	2.5
1	C	206	SER	2.4
1	C	205	ASN	2.4
1	D	85	VAL	2.3
1	A	126	ARG	2.3
1	C	20	ASN	2.3
1	D	84	ALA	2.2
1	C	204	ILE	2.2
1	A	4	ARG	2.1
1	B	126	ARG	2.1
1	A	110	ASN	2.1
1	D	24	TRP	2.1
1	D	88	LEU	2.1

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](#)

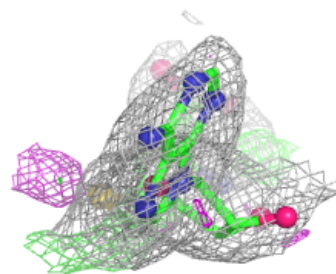
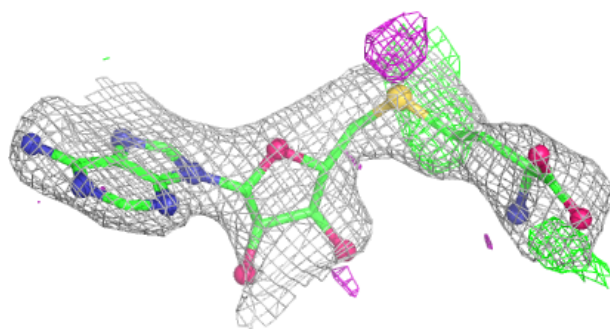
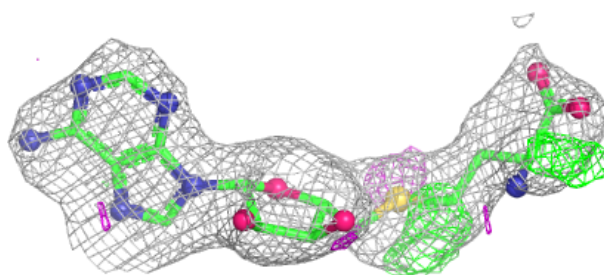
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SAH	D	1004	26/26	0.77	0.13	49,62,71,72	0
2	SAH	C	1003	26/26	0.84	0.11	40,52,68,70	0
2	SAH	B	1002	26/26	0.87	0.10	32,37,54,57	0
2	SAH	A	1001	26/26	0.88	0.10	33,38,50,52	0

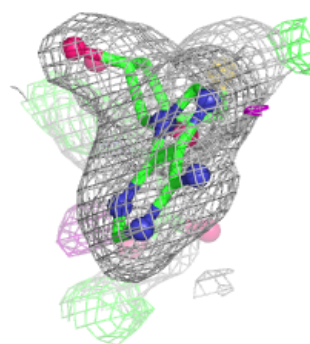
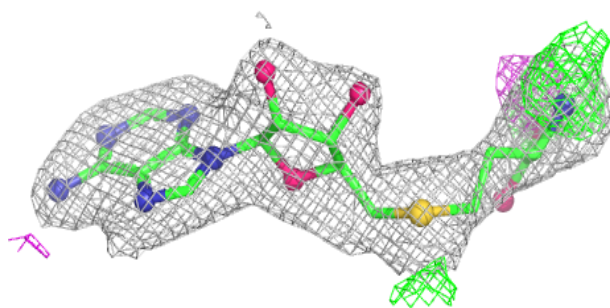
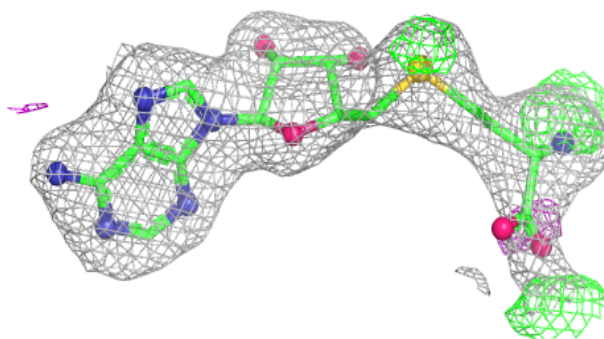
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around SAH D 1004:

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

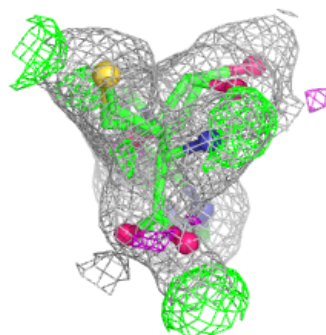
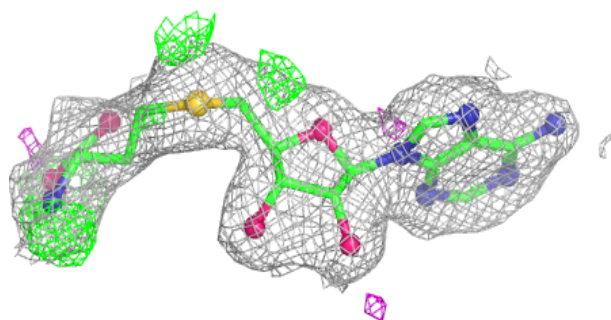
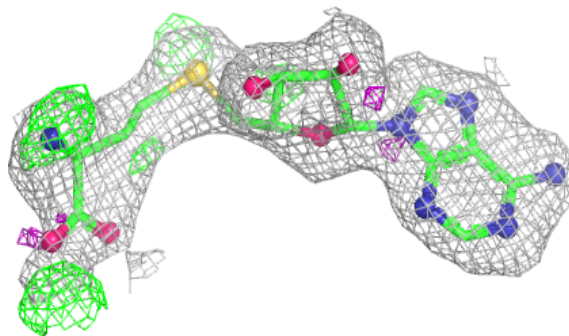
**Electron density around SAH C 1003:**

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

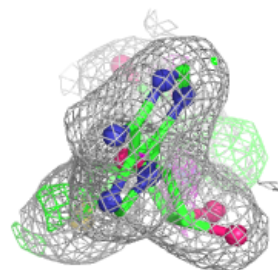
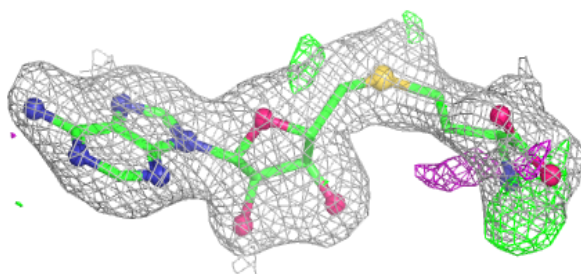
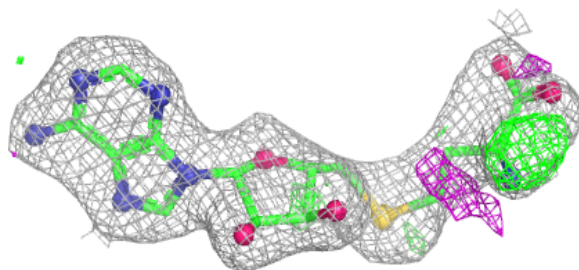


Electron density around SAH B 1002:

$2mF_o-DF_c$ (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.