



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:37 AM UTC

PDB ID : 1P91 / pdb_00001p91
Title : Crystal Structure of RlmA(I) enzyme: 23S rRNA n1-G745 methyltransferase (NORTHEAST STRUCTURAL GENOMICS CONSORTIUM TARGET ER19)
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Deposited on : 2003-05-08
Resolution : 2.80 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](#) ①) 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

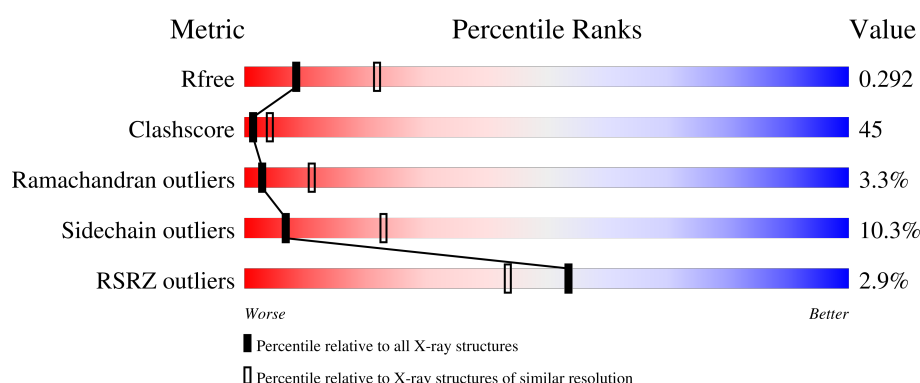
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	269	
1	B	269	

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
4	SAM	B	2401	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4342 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 Ribosomal RNA large subunit methyltransferase A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	Se	0	0	0
			2130	1351	379	385	8	7			
1	B	258	Total	C	N	O	S	Se	0	0	0
			2046	1303	359	369	8	7			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	29	MSE	MET	modified residue	UNP P36999
A	54	MSE	MET	modified residue	UNP P36999
A	55	MSE	MET	modified residue	UNP P36999
A	149	MSE	MET	modified residue	UNP P36999
A	185	MSE	MET	modified residue	UNP P36999
A	220	MSE	MET	modified residue	UNP P36999
A	233	MSE	MET	modified residue	UNP P36999
B	29	MSE	MET	modified residue	UNP P36999
B	54	MSE	MET	modified residue	UNP P36999
B	55	MSE	MET	modified residue	UNP P36999
B	149	MSE	MET	modified residue	UNP P36999
B	185	MSE	MET	modified residue	UNP P36999
B	220	MSE	MET	modified residue	UNP P36999
B	233	MSE	MET	modified residue	UNP P36999

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

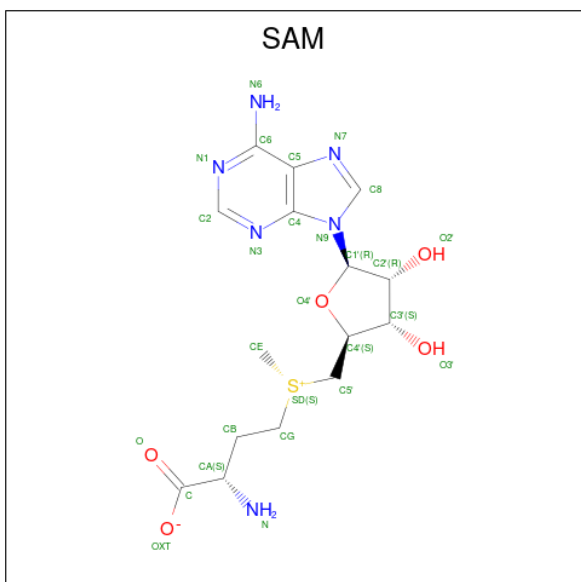
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 4 is S-ADENOSYLMETHIONINE (CCD ID: SAM) (formula: $C_{15}H_{22}N_6O_5S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
4	B	1	Total	C	N	O	S	0	0
			27	15	6	5	1		

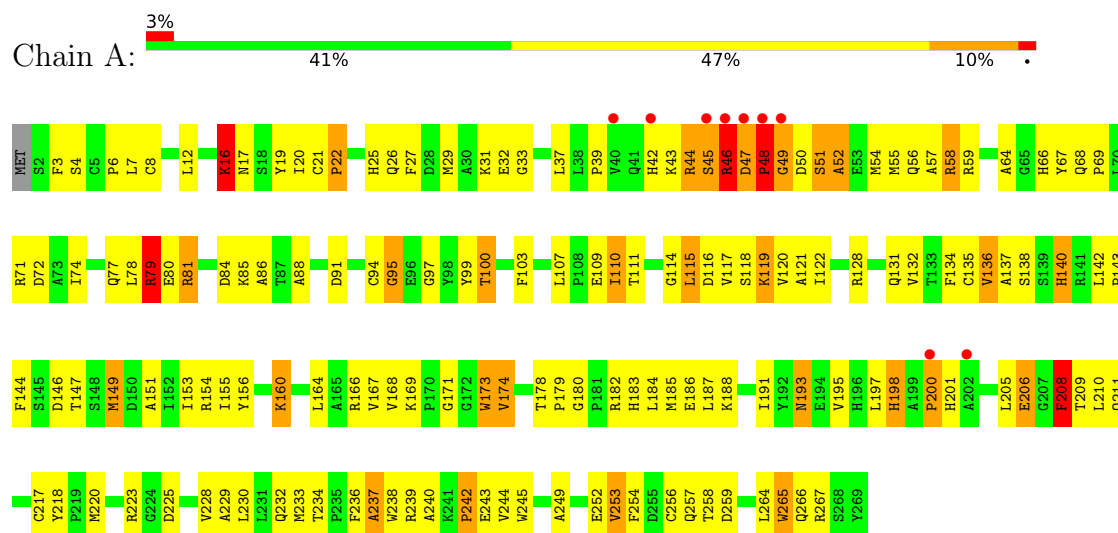
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	54	Total	O	0	0
			54	54		
5	B	31	Total	O	0	0
			31	31		

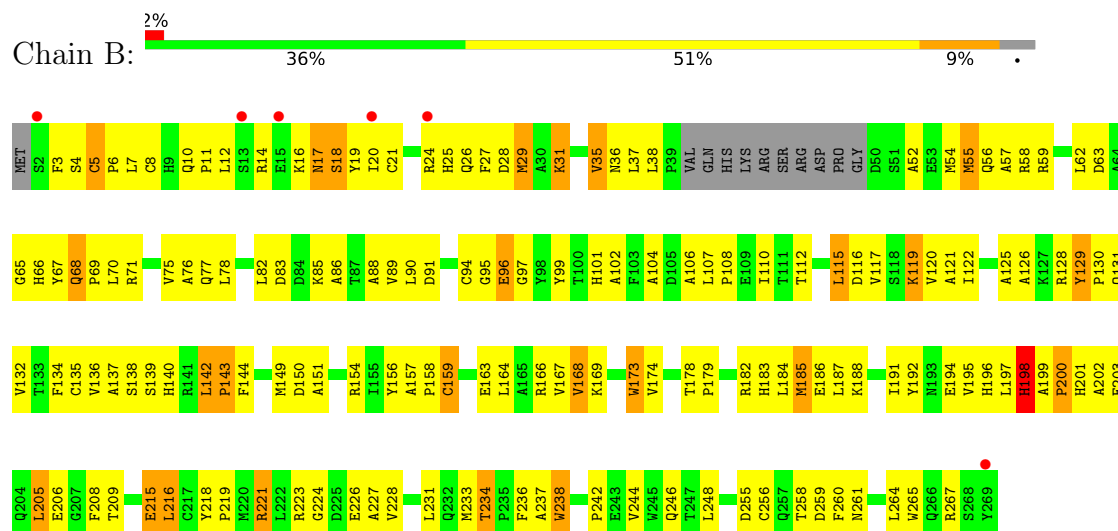
3 Residue-property plots

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: Ribosomal RNA large subunit methyltransferase A



- Molecule 1: Ribosomal RNA large subunit methyltransferase A



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	107.19Å 122.28Å 143.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.80 20.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	92.9 (20.00-2.80) 92.6 (20.00-2.80)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 2.81Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.248 , 0.296 0.247 , 0.292	Depositor DCC
R_{free} test set	1143 reflections (5.22%)	wwPDB-VP
Wilson B-factor (Å ²)	50.6	Xtriage
Anisotropy	0.645	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 44.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4342	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% 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: SAM, ZN, SO4

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	0.71	0/2179	1.28	28/2944 (1.0%)
1	B	0.69	0/2092	1.19	18/2828 (0.6%)
All	All	0.70	0/4271	1.24	46/5772 (0.8%)

There are no bond length outliers.

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	46	ARG	N-CA-C	13.11	125.51	108.34
1	A	49	GLY	N-CA-C	-12.74	98.88	113.79
1	A	44	ARG	N-CA-C	-9.14	101.31	111.28
1	B	203	GLU	N-CA-C	-8.70	95.94	109.25
1	B	201	HIS	N-CA-C	7.93	122.41	109.72
1	B	142	LEU	CA-C-N	7.52	129.24	119.84
1	B	142	LEU	C-N-CA	7.52	129.24	119.84
1	B	52	ALA	N-CA-C	-7.50	104.66	113.88
1	B	234	THR	CA-C-N	7.34	127.20	119.05
1	B	234	THR	C-N-CA	7.34	127.20	119.05
1	B	26	GLN	N-CA-C	7.34	120.22	108.41
1	B	55	MSE	N-CA-C	-7.30	102.50	111.33
1	B	129	TYR	CA-C-N	7.07	126.70	119.56
1	B	129	TYR	C-N-CA	7.07	126.70	119.56
1	A	64	ALA	N-CA-C	-6.75	104.79	113.16
1	A	198	HIS	N-CA-C	6.69	119.31	108.41
1	A	167	VAL	N-CA-C	6.65	119.40	111.09
1	B	198	HIS	N-CA-C	6.32	118.03	110.19
1	A	136	VAL	CB-CA-C	-6.30	103.33	111.53
1	A	200	PRO	CA-C-N	6.10	129.37	120.71
1	A	200	PRO	C-N-CA	6.10	129.37	120.71
1	A	47	ASP	N-CA-C	-6.05	98.30	108.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	173	TRP	N-CA-C	5.84	118.73	109.50
1	A	48	PRO	N-CA-C	5.83	124.47	112.47
1	B	200	PRO	CA-C-N	-5.78	113.73	122.81
1	B	200	PRO	C-N-CA	-5.78	113.73	122.81
1	A	208	PHE	N-CA-C	5.53	117.03	109.18
1	B	168	VAL	N-CA-C	5.50	116.64	108.23
1	A	206	GLU	N-CA-C	-5.49	101.92	110.10
1	A	81	ARG	N-CA-C	5.42	117.19	111.28
1	A	218	TYR	CA-C-N	5.41	127.55	120.25
1	A	218	TYR	C-N-CA	5.41	127.55	120.25
1	A	45	SER	N-CA-C	-5.39	102.06	110.10
1	A	197	LEU	CA-C-N	-5.38	115.42	123.00
1	A	197	LEU	C-N-CA	-5.38	115.42	123.00
1	B	215	GLU	N-CA-C	-5.34	101.17	109.72
1	A	174	VAL	N-CA-C	-5.23	100.53	108.17
1	A	55	MSE	N-CA-C	-5.17	105.54	111.07
1	A	200	PRO	N-CA-C	5.15	118.94	110.55
1	A	173	TRP	N-CA-C	5.11	117.44	109.52
1	A	140	HIS	N-CA-C	-5.09	106.57	112.89
1	A	149	MSE	N-CA-C	5.09	117.79	109.59
1	B	139	SER	N-CA-C	-5.09	107.07	113.28
1	A	7	LEU	N-CA-C	5.06	119.41	112.68
1	A	195	VAL	N-CA-C	5.05	115.51	108.84
1	A	223	ARG	N-CA-C	-5.00	104.08	110.53

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2130	0	2089	182	0
1	B	2046	0	2002	199	0
2	A	1	0	0	1	0
2	B	1	0	0	0	0
3	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	20	0	0	0	0
4	A	27	0	22	5	0
4	B	27	0	22	15	0
5	A	54	0	0	3	0
5	B	31	0	0	4	0
All	All	4342	0	4135	373	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 45.

All (373) 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:160:LYS:HE2	1:A:160:LYS:HA	1.28	1.08
1:B:20:ILE:HG12	1:B:25:HIS:O	1.56	1.06
1:A:185:MSE:HE2	1:A:257:GLN:NE2	1.73	1.03
1:B:10:GLN:HB3	1:B:11:PRO:HD2	1.47	0.95
1:A:8:CYS:HG	2:A:1301:ZN:ZN	0.66	0.93
1:A:66:HIS:HB3	1:A:220:MSE:HE2	1.50	0.91
1:B:6:PRO:HG2	1:B:37:LEU:HB3	1.53	0.91
1:B:183:HIS:CD2	1:B:184:LEU:HG	2.05	0.90
1:B:119:LYS:HD3	1:B:119:LYS:N	1.84	0.89
1:A:119:LYS:HD3	1:A:119:LYS:N	1.88	0.89
1:A:21:CYS:SG	1:A:25:HIS:HB2	2.13	0.88
1:A:201:HIS:CD2	1:A:201:HIS:H	1.86	0.88
1:A:122:ILE:HG13	1:A:136:VAL:HG22	1.54	0.88
1:A:232:GLN:HG2	1:B:246:GLN:HE22	1.36	0.88
1:B:35:VAL:HG12	1:B:135:CYS:HB2	1.57	0.87
1:B:38:LEU:HD13	1:B:137:ALA:HA	1.56	0.86
1:A:252:GLU:HG2	1:A:253:VAL:HG12	1.58	0.84
1:A:66:HIS:HB3	1:A:220:MSE:CE	2.09	0.83
1:A:185:MSE:HE2	1:A:257:GLN:HE22	1.39	0.83
1:B:182:ARG:O	1:B:185:MSE:HG3	1.79	0.82
1:A:178:THR:HB	1:A:179:PRO:HD2	1.60	0.81
1:B:191:ILE:HD11	1:B:236:PHE:HD1	1.44	0.81
1:B:18:SER:HB3	1:B:28:ASP:HA	1.64	0.79
4:A:1401:SAM:HB2	4:A:1401:SAM:H4'	1.64	0.79
1:B:20:ILE:HG23	5:B:2402:HOH:O	1.83	0.78
1:B:184:LEU:HD13	1:B:187:LEU:HD12	1.65	0.78
1:A:3:PHE:HB2	1:A:12:LEU:HD12	1.65	0.77
1:B:25:HIS:HB2	5:B:2402:HOH:O	1.84	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:LYS:HD2	1:A:17:ASN:ND2	2.00	0.76
1:B:6:PRO:CG	1:B:37:LEU:HB3	2.16	0.76
1:B:144:PHE:O	1:B:166:ARG:NH2	2.18	0.76
1:B:31:LYS:H	1:B:31:LYS:HD3	1.53	0.74
1:A:117:VAL:HG12	1:A:137:ALA:O	1.88	0.74
1:A:228:VAL:HG13	1:A:245:TRP:CE3	2.22	0.73
1:B:156:TYR:OH	1:B:179:PRO:HG3	1.89	0.73
1:A:107:LEU:HB3	1:A:110:ILE:HD11	1.70	0.72
1:B:198:HIS:CD2	1:B:199:ALA:H	2.08	0.72
1:A:44:ARG:O	1:A:46:ARG:CZ	2.37	0.72
1:A:243:GLU:H	1:A:243:GLU:CD	1.98	0.71
1:B:198:HIS:CD2	1:B:199:ALA:N	2.59	0.71
1:B:78:LEU:HD22	1:B:82:LEU:HD11	1.73	0.71
1:A:185:MSE:CE	1:A:257:GLN:HE22	2.03	0.71
1:B:115:LEU:C	1:B:115:LEU:HD23	2.16	0.70
1:A:119:LYS:N	1:A:119:LYS:CD	2.55	0.70
1:B:35:VAL:HG13	1:B:37:LEU:HD21	1.74	0.69
1:B:191:ILE:HD11	1:B:236:PHE:CD1	2.27	0.69
1:A:119:LYS:HE2	1:A:120:VAL:HG13	1.74	0.69
1:A:122:ILE:CG1	1:A:136:VAL:HG22	2.22	0.69
1:B:156:TYR:CZ	1:B:179:PRO:HG3	2.28	0.69
1:B:128:ARG:HB2	5:B:2425:HOH:O	1.93	0.68
1:A:95:GLY:HA2	1:A:116:ASP:OD1	1.93	0.68
1:A:54:MSE:HB3	1:A:58:ARG:CZ	2.23	0.68
1:B:71:ARG:O	1:B:75:VAL:HG23	1.93	0.67
1:A:115:LEU:HD23	1:A:116:ASP:N	2.10	0.67
1:B:19:TYR:CE1	1:B:35:VAL:HG21	2.29	0.67
1:B:58:ARG:NH1	1:B:233:MSE:HE3	2.08	0.67
1:A:119:LYS:CD	1:A:119:LYS:H	2.06	0.67
1:B:224:GLY:O	1:B:228:VAL:HG23	1.96	0.66
1:A:107:LEU:HD13	1:A:110:ILE:HD11	1.78	0.66
1:A:156:TYR:CD1	1:A:179:PRO:HD3	2.30	0.66
1:A:201:HIS:CD2	1:A:201:HIS:N	2.60	0.66
1:B:96:GLU:C	1:B:125:ALA:HB2	2.21	0.65
1:A:116:ASP:HB3	1:A:122:ILE:HD11	1.77	0.65
1:A:119:LYS:HD3	1:A:119:LYS:H	1.60	0.65
1:B:70:LEU:HD13	1:B:260:PHE:CG	2.31	0.65
1:B:191:ILE:HG13	1:B:191:ILE:O	1.96	0.64
1:A:122:ILE:HG21	1:A:136:VAL:CG2	2.28	0.64
1:B:70:LEU:HD13	1:B:260:PHE:CD1	2.32	0.64
1:B:54:MSE:O	1:B:55:MSE:C	2.41	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:TRP:CZ3	1:A:266:GLN:HG3	2.32	0.64
1:B:156:TYR:CZ	4:B:2401:SAM:HG1	2.32	0.64
1:B:221:ARG:HG2	1:B:255:ASP:OD1	1.97	0.64
1:B:119:LYS:HE2	1:B:120:VAL:HG12	1.80	0.64
1:B:31:LYS:HD3	1:B:31:LYS:N	2.13	0.63
1:B:5:CYS:C	1:B:7:LEU:H	2.06	0.63
1:A:3:PHE:CB	1:A:12:LEU:HD12	2.29	0.63
1:B:117:VAL:HG12	1:B:137:ALA:O	1.98	0.63
1:B:110:ILE:C	1:B:110:ILE:HD12	2.23	0.62
1:A:54:MSE:HB3	1:A:58:ARG:NH1	2.14	0.62
1:A:186:GLU:CD	1:A:256:CYS:HA	2.24	0.62
1:A:166:ARG:HG2	1:A:166:ARG:HH11	1.65	0.62
1:A:31:LYS:HE2	1:A:32:GLU:OE2	2.00	0.62
1:A:228:VAL:HG13	1:A:245:TRP:HE3	1.65	0.62
1:A:107:LEU:HB3	1:A:110:ILE:CD1	2.30	0.61
1:A:94:CYS:O	1:A:95:GLY:C	2.43	0.61
1:A:122:ILE:CB	1:A:136:VAL:HG22	2.29	0.61
1:B:77:GLN:HE21	1:B:77:GLN:HA	1.64	0.61
1:A:156:TYR:CE1	1:A:179:PRO:HD3	2.35	0.61
1:B:126:ALA:HA	1:B:134:PHE:HD1	1.66	0.61
1:B:154:ARG:HD3	1:B:157:ALA:O	2.01	0.61
1:B:7:LEU:HB3	1:B:25:HIS:CE1	2.36	0.61
1:B:17:ASN:O	1:B:29:MSE:HB2	2.00	0.61
1:B:35:VAL:HG13	1:B:37:LEU:CD2	2.31	0.60
1:B:119:LYS:HD3	1:B:119:LYS:H	1.64	0.60
1:A:57:ALA:O	1:A:58:ARG:C	2.44	0.60
1:B:31:LYS:H	1:B:31:LYS:CD	2.10	0.60
1:A:107:LEU:HB3	1:A:110:ILE:CG1	2.31	0.60
1:B:10:GLN:CB	1:B:11:PRO:HD2	2.29	0.60
1:A:153:ILE:N	1:A:153:ILE:HD12	2.17	0.60
1:B:198:HIS:CG	1:B:199:ALA:N	2.69	0.59
1:A:52:ALA:O	1:A:56:GLN:HG3	2.01	0.59
1:A:185:MSE:HE2	1:A:257:GLN:HE21	1.64	0.59
1:A:54:MSE:CE	1:B:242:PRO:HG3	2.32	0.59
1:B:107:LEU:HB3	1:B:110:ILE:HG13	1.84	0.59
1:A:48:PRO:HG2	1:B:238:TRP:HB3	1.84	0.59
1:A:209:THR:O	1:A:211:GLN:HG2	2.01	0.59
1:A:59:ARG:NH2	1:A:128:ARG:HD3	2.17	0.59
1:A:94:CYS:SG	1:A:114:GLY:C	2.86	0.58
1:A:147:THR:HG22	1:A:147:THR:O	2.02	0.58
1:B:107:LEU:HB2	1:B:110:ILE:HD11	1.83	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:35:VAL:CG1	1:B:37:LEU:HD21	2.33	0.58
1:B:66:HIS:O	1:B:67:TYR:HD2	1.87	0.58
1:A:78:LEU:O	1:A:80:GLU:N	2.37	0.58
1:A:198:HIS:O	1:A:200:PRO:HD3	2.04	0.58
1:B:19:TYR:O	1:B:20:ILE:HG13	2.04	0.58
1:A:228:VAL:HG13	1:A:245:TRP:CZ3	2.39	0.57
1:B:106:ALA:O	1:B:108:PRO:HD3	2.03	0.57
1:B:119:LYS:HG2	1:B:120:VAL:N	2.19	0.57
1:A:85:LYS:O	1:A:85:LYS:HG2	2.04	0.57
1:B:198:HIS:O	1:B:200:PRO:HD3	2.04	0.57
1:A:79:ARG:HH11	1:A:79:ARG:HG2	1.68	0.57
1:B:29:MSE:HG3	1:B:35:VAL:HG23	1.87	0.57
1:A:122:ILE:HG21	1:A:136:VAL:HG23	1.86	0.57
1:B:71:ARG:HD2	1:B:99:TYR:HA	1.87	0.57
1:B:184:LEU:O	1:B:185:MSE:C	2.48	0.57
1:A:31:LYS:HG3	1:A:32:GLU:OE2	2.05	0.56
1:A:37:LEU:HB2	1:A:137:ALA:HB2	1.87	0.56
1:A:74:ILE:HD12	1:A:155:ILE:HD11	1.88	0.56
1:A:230:LEU:O	1:A:234:THR:HG23	2.06	0.56
1:B:63:ASP:C	1:B:65:GLY:H	2.12	0.56
1:B:122:ILE:HG13	1:B:136:VAL:HG22	1.88	0.56
1:B:164:LEU:HD13	1:B:205:LEU:HD11	1.86	0.56
1:B:36:ASN:C	1:B:37:LEU:HD23	2.31	0.56
1:A:115:LEU:HD23	1:A:115:LEU:C	2.31	0.56
1:B:71:ARG:HG3	1:B:99:TYR:HD2	1.69	0.56
1:A:191:ILE:CD1	1:A:236:PHE:HB2	2.36	0.56
1:B:156:TYR:CE2	1:B:179:PRO:HG3	2.41	0.56
1:B:183:HIS:NE2	1:B:184:LEU:HG	2.21	0.55
1:B:96:GLU:O	1:B:96:GLU:HG2	2.07	0.55
1:A:156:TYR:O	1:A:156:TYR:CD2	2.60	0.55
1:B:6:PRO:HD2	1:B:27:PHE:CE1	2.42	0.55
1:A:107:LEU:HD22	1:A:110:ILE:HD11	1.87	0.55
1:A:160:LYS:HA	1:A:160:LYS:CE	2.12	0.55
1:A:103:PHE:CE1	1:A:153:ILE:HG21	2.41	0.55
1:B:242:PRO:O	1:B:246:GLN:HG2	2.06	0.54
1:B:107:LEU:HB3	1:B:110:ILE:CG1	2.36	0.54
1:B:218:TYR:H	1:B:258:THR:HG22	1.73	0.54
1:A:217:CYS:HA	1:A:258:THR:O	2.07	0.54
1:A:147:THR:HG23	1:A:169:LYS:HA	1.90	0.54
1:A:81:ARG:HB3	1:A:173:TRP:CE2	2.42	0.54
1:B:88:ALA:HB1	1:B:149:MSE:SE	2.58	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:TYR:CD1	4:A:1401:SAM:HE2	2.43	0.54
1:B:104:ALA:HA	1:B:112:THR:OG1	2.08	0.54
1:A:166:ARG:HG2	1:A:166:ARG:NH1	2.23	0.54
1:A:46:ARG:NH1	1:A:119:LYS:HE3	2.23	0.53
1:A:54:MSE:CB	1:A:58:ARG:NH1	2.71	0.53
1:A:171:GLY:N	1:A:267:ARG:O	2.34	0.53
1:A:174:VAL:HG21	1:A:208:PHE:CE2	2.43	0.53
1:A:116:ASP:CB	1:A:122:ILE:HD11	2.37	0.53
1:B:10:GLN:HB3	1:B:11:PRO:CD	2.30	0.53
1:B:90:LEU:HD12	1:B:91:ASP:N	2.24	0.53
1:B:5:CYS:C	1:B:7:LEU:N	2.66	0.53
1:A:187:LEU:HD13	1:A:254:PHE:CE2	2.44	0.53
1:B:19:TYR:HE1	1:B:35:VAL:HG21	1.74	0.52
1:B:95:GLY:HA2	1:B:116:ASP:OD1	2.09	0.52
1:B:119:LYS:CE	1:B:120:VAL:HG12	2.38	0.52
1:A:46:ARG:CG	1:A:48:PRO:HG3	2.39	0.52
1:A:156:TYR:HD1	4:A:1401:SAM:HE2	1.74	0.52
1:A:71:ARG:O	1:A:72:ASP:C	2.51	0.52
1:A:138:SER:HG	1:A:140:HIS:CE1	2.28	0.52
1:B:85:LYS:O	1:B:85:LYS:HG3	2.09	0.52
1:A:78:LEU:C	1:A:80:GLU:N	2.68	0.52
1:A:208:PHE:N	1:A:208:PHE:CD1	2.76	0.52
1:A:178:THR:HB	1:A:179:PRO:CD	2.35	0.52
1:B:119:LYS:CG	1:B:120:VAL:N	2.72	0.52
1:B:97:GLY:N	4:B:2401:SAM:HN62	2.06	0.52
1:A:132:VAL:HB	1:A:134:PHE:CE1	2.45	0.52
1:B:96:GLU:N	4:B:2401:SAM:N1	2.51	0.52
1:A:91:ASP:OD1	1:A:100:THR:HG23	2.09	0.51
1:B:117:VAL:HG12	1:B:138:SER:HA	1.91	0.51
1:B:119:LYS:CG	1:B:120:VAL:H	2.23	0.51
1:B:231:LEU:O	1:B:237:ALA:HB2	2.11	0.51
1:B:191:ILE:CD1	1:B:236:PHE:HD1	2.20	0.51
1:B:227:ALA:HB1	1:B:248:LEU:HD11	1.93	0.51
1:A:185:MSE:CE	1:A:257:GLN:NE2	2.57	0.51
1:B:188:LYS:HD3	1:B:197:LEU:HD12	1.93	0.51
1:B:188:LYS:HB2	1:B:192:TYR:CD1	2.46	0.50
1:B:12:LEU:HD22	1:B:19:TYR:HB3	1.92	0.50
1:A:46:ARG:HG2	1:A:48:PRO:HG3	1.93	0.50
1:B:8:CYS:SG	1:B:10:GLN:HB2	2.51	0.50
1:B:20:ILE:CG2	1:B:21:CYS:N	2.74	0.50
1:B:63:ASP:C	1:B:65:GLY:N	2.69	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:95:GLY:HA3	4:B:2401:SAM:N1	2.26	0.50
1:A:182:ARG:O	1:A:185:MSE:HG3	2.11	0.50
1:A:156:TYR:HD1	1:A:179:PRO:HD3	1.77	0.50
1:A:79:ARG:HG2	1:A:79:ARG:NH1	2.26	0.50
1:A:42:HIS:C	1:A:44:ARG:H	2.20	0.50
1:A:180:GLY:N	1:A:259:ASP:O	2.35	0.49
1:B:58:ARG:HH11	1:B:233:MSE:HE3	1.73	0.49
1:B:77:GLN:HA	1:B:77:GLN:NE2	2.27	0.49
1:B:110:ILE:HD12	1:B:110:ILE:O	2.12	0.49
1:B:59:ARG:HD2	1:B:128:ARG:NH1	2.27	0.49
1:B:234:THR:HG22	4:B:2401:SAM:OXT	2.12	0.49
1:B:66:HIS:C	1:B:67:TYR:HD2	2.19	0.49
1:B:150:ASP:CG	1:B:169:LYS:HD2	2.38	0.49
1:A:228:VAL:HG21	1:A:249:ALA:HB2	1.95	0.49
1:B:156:TYR:CD2	4:B:2401:SAM:H5'1	2.47	0.49
1:A:54:MSE:HE2	1:B:242:PRO:HG3	1.94	0.49
1:B:86:ALA:C	1:B:110:ILE:HG22	2.38	0.49
1:B:119:LYS:CD	1:B:120:VAL:H	2.25	0.49
1:B:182:ARG:CB	1:B:182:ARG:HH11	2.25	0.49
1:A:103:PHE:HE1	1:A:153:ILE:HG21	1.77	0.49
1:A:6:PRO:HD2	1:A:27:PHE:CZ	2.48	0.49
1:B:16:LYS:HD2	1:B:17:ASN:ND2	2.28	0.49
1:A:119:LYS:HD3	1:A:120:VAL:H	1.77	0.49
1:A:209:THR:O	1:A:265:TRP:HA	2.13	0.49
1:A:48:PRO:C	1:A:50:ASP:H	2.20	0.48
1:B:188:LYS:CD	1:B:197:LEU:HD12	2.43	0.48
1:A:118:SER:HB3	1:A:121:ALA:HB3	1.95	0.48
1:B:14:ARG:HG3	1:B:14:ARG:HH11	1.78	0.48
1:B:158:PRO:O	1:B:159:CYS:HB2	2.12	0.48
1:B:168:VAL:HG12	1:B:169:LYS:O	2.13	0.48
1:B:54:MSE:C	1:B:56:GLN:N	2.70	0.48
1:B:67:TYR:HE1	4:B:2401:SAM:HG2	1.78	0.48
1:B:119:LYS:HG2	1:B:120:VAL:H	1.78	0.48
1:B:156:TYR:CG	4:B:2401:SAM:H5'1	2.48	0.48
1:A:259:ASP:C	1:A:259:ASP:OD2	2.54	0.48
1:B:131:GLN:OE1	1:B:131:GLN:N	2.44	0.48
1:B:151:ALA:HA	1:B:173:TRP:O	2.13	0.48
1:B:38:LEU:HD13	1:B:137:ALA:CA	2.36	0.48
1:B:58:ARG:HD3	4:B:2401:SAM:N3	2.28	0.48
1:A:19:TYR:O	1:A:26:GLN:HA	2.13	0.48
1:A:78:LEU:C	1:A:80:GLU:H	2.22	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:ALA:O	1:A:232:GLN:HB3	2.12	0.48
1:B:71:ARG:HG3	1:B:99:TYR:CD2	2.48	0.48
1:A:42:HIS:CE1	1:A:43:LYS:HG3	2.49	0.48
1:B:156:TYR:CE2	1:B:179:PRO:CG	2.97	0.48
1:A:107:LEU:HB3	1:A:110:ILE:HG12	1.94	0.48
1:A:156:TYR:HB3	4:A:1401:SAM:H3'	1.96	0.48
1:B:208:PHE:CZ	1:B:267:ARG:HD3	2.49	0.48
1:A:29:MSE:HE3	1:A:33:GLY:HA2	1.95	0.47
1:A:54:MSE:HE1	1:B:242:PRO:HG3	1.96	0.47
1:A:86:ALA:O	1:A:110:ILE:HG22	2.14	0.47
1:A:151:ALA:HA	1:A:173:TRP:O	2.15	0.47
1:B:77:GLN:HE21	1:B:77:GLN:CA	2.25	0.47
1:B:37:LEU:HD12	1:B:143:PRO:HB3	1.96	0.47
1:A:144:PHE:CD2	1:A:149:MSE:CE	2.98	0.47
1:B:95:GLY:HA3	4:B:2401:SAM:C2	2.45	0.47
1:B:132:VAL:HG11	1:B:134:PHE:CE2	2.50	0.47
1:A:68:GLN:O	1:A:69:PRO:C	2.57	0.47
1:B:68:GLN:O	1:B:69:PRO:C	2.55	0.47
1:B:120:VAL:HG13	1:B:121:ALA:N	2.30	0.47
1:B:208:PHE:CE2	1:B:267:ARG:HD3	2.49	0.47
1:B:231:LEU:HD12	1:B:234:THR:OG1	2.14	0.47
1:A:174:VAL:CG2	1:A:208:PHE:CE2	2.98	0.47
1:A:240:ALA:HB3	1:A:245:TRP:NE1	2.30	0.47
1:B:7:LEU:HD13	1:B:25:HIS:CD2	2.50	0.47
1:B:216:LEU:C	1:B:216:LEU:HD12	2.40	0.47
1:A:173:TRP:CH2	1:A:266:GLN:HG3	2.50	0.47
1:B:186:GLU:CD	1:B:256:CYS:HA	2.40	0.46
1:B:194:GLU:OE2	1:B:196:HIS:HB3	2.16	0.46
1:B:117:VAL:CG1	1:B:138:SER:HA	2.46	0.46
1:A:97:GLY:H	4:A:1401:SAM:HN62	1.62	0.46
1:B:115:LEU:HD23	1:B:116:ASP:N	2.30	0.46
1:B:188:LYS:HB2	1:B:192:TYR:CE1	2.50	0.46
1:A:59:ARG:HH21	1:A:128:ARG:HD3	1.80	0.46
1:A:67:TYR:CD1	1:A:220:MSE:HE1	2.51	0.46
1:A:74:ILE:O	1:A:77:GLN:HB3	2.16	0.46
1:B:71:ARG:CD	1:B:99:TYR:HA	2.46	0.46
1:A:191:ILE:HD11	1:A:236:PHE:HB2	1.98	0.45
1:B:14:ARG:HD2	1:B:19:TYR:CE2	2.50	0.45
1:B:198:HIS:CE1	5:B:2430:HOH:O	2.68	0.45
1:B:17:ASN:ND2	1:B:17:ASN:C	2.74	0.45
1:A:20:ILE:HA	1:A:25:HIS:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:HIS:HB3	1:A:220:MSE:HE1	1.97	0.45
1:A:122:ILE:HG21	1:A:136:VAL:HG22	1.98	0.45
1:A:173:TRP:CE3	1:A:264:LEU:HD11	2.51	0.45
1:B:69:PRO:HB2	1:B:218:TYR:CE2	2.51	0.45
1:B:167:VAL:HG12	1:B:167:VAL:O	2.15	0.45
1:B:174:VAL:HB	1:B:265:TRP:HB2	1.99	0.45
1:A:265:TRP:N	1:A:265:TRP:CD1	2.84	0.45
1:B:85:LYS:O	1:B:85:LYS:CG	2.64	0.45
1:B:183:HIS:N	1:B:259:ASP:HB3	2.32	0.45
1:A:146:ASP:N	1:A:146:ASP:OD1	2.49	0.45
1:B:4:SER:N	1:B:143:PRO:O	2.48	0.45
1:A:154:ARG:NH2	5:A:1427:HOH:O	2.48	0.45
1:B:12:LEU:HD22	1:B:19:TYR:CB	2.46	0.45
1:A:142:LEU:HA	1:A:143:PRO:HD3	1.77	0.45
1:B:77:GLN:NE2	1:B:264:LEU:HD22	2.32	0.45
1:A:99:TYR:HB2	5:A:1434:HOH:O	2.16	0.44
1:B:83:ASP:OD1	1:B:83:ASP:N	2.50	0.44
1:B:163:GLU:O	1:B:167:VAL:HG23	2.17	0.44
1:A:16:LYS:HD2	1:A:17:ASN:CG	2.42	0.44
1:A:51:SER:O	1:A:52:ALA:C	2.60	0.44
1:A:115:LEU:C	1:A:115:LEU:CD2	2.90	0.44
1:B:20:ILE:HG22	1:B:21:CYS:N	2.32	0.44
1:B:223:ARG:N	1:B:226:GLU:OE1	2.49	0.44
1:A:188:LYS:HG2	1:A:236:PHE:CZ	2.52	0.44
1:B:119:LYS:N	1:B:119:LYS:CD	2.67	0.44
1:B:198:HIS:CD2	1:B:199:ALA:HB3	2.51	0.44
1:A:91:ASP:CB	1:A:100:THR:HG23	2.47	0.44
1:A:187:LEU:HD13	1:A:254:PHE:HE2	1.83	0.44
1:A:187:LEU:HD12	1:A:187:LEU:HA	1.79	0.44
1:A:252:GLU:CD	1:A:252:GLU:H	2.25	0.44
1:A:228:VAL:O	1:A:229:ALA:C	2.61	0.44
1:B:58:ARG:NH1	1:B:233:MSE:CE	2.80	0.44
1:B:68:GLN:HB3	1:B:69:PRO:HD3	2.00	0.44
1:B:142:LEU:HD12	1:B:167:VAL:HG21	2.00	0.44
1:A:16:LYS:HB3	1:A:17:ASN:H	1.73	0.43
1:A:119:LYS:CD	1:A:120:VAL:H	2.31	0.43
1:B:3:PHE:HB3	1:B:12:LEU:HD12	1.98	0.43
1:B:188:LYS:NZ	1:B:197:LEU:HD13	2.33	0.43
1:A:54:MSE:HE1	1:B:242:PRO:HA	2.00	0.43
1:A:67:TYR:CE1	1:A:220:MSE:HE1	2.54	0.43
1:B:119:LYS:HE2	1:B:120:VAL:CG1	2.47	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:PRO:HB2	1:A:49:GLY:H	1.59	0.43
1:A:54:MSE:HE1	1:B:242:PRO:CA	2.49	0.43
1:A:78:LEU:O	1:A:79:ARG:C	2.60	0.43
1:A:239:ARG:HH22	1:B:56:GLN:NE2	2.17	0.43
1:A:107:LEU:CB	1:A:110:ILE:HD11	2.43	0.43
1:A:109:GLU:OE1	1:A:109:GLU:HA	2.18	0.43
1:A:186:GLU:OE1	1:A:256:CYS:HA	2.17	0.43
1:B:129:TYR:HA	1:B:130:PRO:HD2	1.90	0.43
1:A:164:LEU:O	1:A:168:VAL:HG23	2.19	0.43
1:B:197:LEU:HD23	1:B:197:LEU:HA	1.80	0.43
1:A:8:CYS:SG	1:A:21:CYS:SG	3.16	0.42
1:A:12:LEU:HD23	1:A:21:CYS:HB3	2.00	0.42
1:A:131:GLN:OE1	1:A:131:GLN:N	2.49	0.42
1:B:20:ILE:HG22	1:B:21:CYS:O	2.19	0.42
1:B:76:ALA:O	1:B:77:GLN:C	2.62	0.42
1:A:54:MSE:O	1:A:57:ALA:HB3	2.19	0.42
1:A:183:HIS:O	1:A:184:LEU:HB2	2.19	0.42
1:B:138:SER:HB3	1:B:140:HIS:HD2	1.84	0.42
1:B:215:GLU:OE1	1:B:259:ASP:OD1	2.37	0.42
1:B:188:LYS:HZ3	1:B:197:LEU:HD13	1.85	0.42
1:B:3:PHE:CB	1:B:12:LEU:HD12	2.50	0.42
1:B:17:ASN:N	1:B:17:ASN:HD22	2.17	0.42
1:B:156:TYR:CB	4:B:2401:SAM:H3'	2.50	0.42
1:A:201:HIS:H	1:A:201:HIS:HD2	1.58	0.42
1:B:17:ASN:HD22	1:B:17:ASN:H	1.66	0.42
1:B:164:LEU:HD23	1:B:164:LEU:HA	1.88	0.42
1:A:205:LEU:O	1:A:206:GLU:C	2.58	0.42
1:B:138:SER:HB3	1:B:140:HIS:CD2	2.54	0.42
1:B:227:ALA:CB	1:B:248:LEU:HD11	2.50	0.42
1:A:107:LEU:HD13	1:A:110:ILE:CD1	2.49	0.41
1:B:4:SER:HB2	1:B:144:PHE:C	2.45	0.41
1:A:117:VAL:HG12	1:A:138:SER:HA	2.03	0.41
1:A:238:TRP:CD2	1:B:57:ALA:HB2	2.55	0.41
1:B:58:ARG:HD3	4:B:2401:SAM:C2	2.50	0.41
1:B:110:ILE:C	1:B:110:ILE:CD1	2.90	0.41
1:A:88:ALA:HB2	1:A:111:THR:HB	2.03	0.41
1:A:16:LYS:C	1:A:16:LYS:HD3	2.42	0.41
1:A:16:LYS:HD3	1:A:16:LYS:HA	1.88	0.41
1:B:67:TYR:HE1	4:B:2401:SAM:CG	2.33	0.41
1:B:69:PRO:HG2	1:B:218:TYR:CZ	2.56	0.41
1:B:199:ALA:HA	1:B:200:PRO:HD2	1.95	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:PRO:HD3	5:A:1411:HOH:O	2.20	0.41
1:A:46:ARG:HG2	1:A:48:PRO:CD	2.51	0.41
1:A:107:LEU:CD1	1:A:110:ILE:HD11	2.48	0.41
1:A:156:TYR:CE1	1:A:179:PRO:CD	3.03	0.41
1:A:184:LEU:HA	1:A:256:CYS:SG	2.61	0.41
1:A:240:ALA:HB3	1:A:245:TRP:HE1	1.86	0.41
1:A:44:ARG:O	1:A:46:ARG:NE	2.54	0.40
1:B:101:HIS:CG	1:B:102:ALA:N	2.89	0.40
1:B:150:ASP:OD1	1:B:169:LYS:HD2	2.20	0.40
1:A:94:CYS:O	1:A:97:GLY:N	2.52	0.40
1:B:62:LEU:HD23	1:B:62:LEU:HA	1.88	0.40
1:A:236:PHE:O	1:A:237:ALA:C	2.64	0.40
1:B:184:LEU:HD11	4:B:2401:SAM:O	2.21	0.40
1:A:42:HIS:C	1:A:44:ARG:N	2.79	0.40
1:A:242:PRO:O	1:A:243:GLU:C	2.65	0.40
1:B:3:PHE:HD2	1:B:37:LEU:HD11	1.87	0.40
1:B:67:TYR:CE1	4:B:2401:SAM:HG2	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	266/269 (99%)	226 (85%)	31 (12%)	9 (3%)	3	11
1	B	254/269 (94%)	211 (83%)	35 (14%)	8 (3%)	3	12
All	All	520/538 (97%)	437 (84%)	66 (13%)	17 (3%)	3	11

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	48	PRO

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Mol	Chain	Res	Type
1	A	52	ALA
1	A	237	ALA
1	A	79	ARG
1	A	95	GLY
1	A	193	ASN
1	B	202	ALA
1	A	16	LYS
1	B	94	CYS
1	B	159	CYS
1	B	185	MSE
1	B	238	TRP
1	A	39	PRO
1	B	89	VAL
1	B	143	PRO
1	A	244	VAL
1	B	195	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	224/218 (103%)	200 (89%)	24 (11%)	6	21
1	B	214/218 (98%)	193 (90%)	21 (10%)	7	25
All	All	438/436 (100%)	393 (90%)	45 (10%)	7	23

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

Mol	Chain	Res	Type
1	A	4	SER
1	A	16	LYS
1	A	22	PRO
1	A	45	SER
1	A	46	ARG
1	A	47	ASP
1	A	51	SER

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Mol	Chain	Res	Type
1	A	58	ARG
1	A	79	ARG
1	A	84	ASP
1	A	100	THR
1	A	110	ILE
1	A	115	LEU
1	A	119	LYS
1	A	135	CYS
1	A	160	LYS
1	A	193	ASN
1	A	208	PHE
1	A	210	LEU
1	A	225	ASP
1	A	233	MSE
1	A	242	PRO
1	A	253	VAL
1	A	265	TRP
1	B	5	CYS
1	B	17	ASN
1	B	18	SER
1	B	24	ARG
1	B	29	MSE
1	B	31	LYS
1	B	35	VAL
1	B	68	GLN
1	B	96	GLU
1	B	115	LEU
1	B	119	LYS
1	B	178	THR
1	B	198	HIS
1	B	205	LEU
1	B	206	GLU
1	B	209	THR
1	B	216	LEU
1	B	219	PRO
1	B	221	ARG
1	B	244	VAL
1	B	261	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	17	ASN
1	A	23	GLN
1	A	42	HIS
1	A	66	HIS
1	A	68	GLN
1	A	198	HIS
1	A	201	HIS
1	A	232	GLN
1	A	257	GLN
1	A	263	HIS
1	B	17	ASN
1	B	56	GLN
1	B	68	GLN
1	B	77	GLN
1	B	140	HIS
1	B	198	HIS
1	B	212	GLN
1	B	261	ASN

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 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 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	SAM	A	1401	-	27,29,29	2.63	9 (33%)	34,42,42	2.88	13 (38%)
4	SAM	B	2401	-	27,29,29	2.47	8 (29%)	34,42,42	2.85	9 (26%)
3	SO4	B	502	-	4,4,4	0.19	0	6,6,6	0.57	0
3	SO4	A	501	-	4,4,4	0.23	0	6,6,6	0.55	0
3	SO4	B	503	-	4,4,4	0.26	0	6,6,6	0.59	0
3	SO4	B	504	-	4,4,4	0.43	0	6,6,6	0.62	0
3	SO4	B	505	-	4,4,4	0.37	0	6,6,6	0.63	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
4	SAM	B	2401	-	-	8/17/33/33	0/3/3/3
4	SAM	A	1401	-	-	6/17/33/33	0/3/3/3

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1401	SAM	C2-N1	7.81	1.47	1.33
4	B	2401	SAM	C2-N1	7.24	1.46	1.33
4	A	1401	SAM	C8-N9	5.35	1.46	1.37
4	B	2401	SAM	C6-N6	4.48	1.45	1.34
4	B	2401	SAM	C5-N7	-4.33	1.31	1.39
4	B	2401	SAM	C8-N9	4.32	1.44	1.37
4	A	1401	SAM	C5-N7	-4.18	1.31	1.39
4	A	1401	SAM	C6-N6	4.15	1.44	1.34
4	A	1401	SAM	C4-N9	3.93	1.45	1.37
4	B	2401	SAM	C5-C6	3.22	1.50	1.41
4	B	2401	SAM	C4-N9	3.16	1.44	1.37
4	A	1401	SAM	O4'-C1'	3.04	1.49	1.42
4	A	1401	SAM	C5-C6	2.84	1.48	1.41
4	B	2401	SAM	C5-C4	-2.84	1.34	1.39
4	B	2401	SAM	OXT-C	-2.57	1.22	1.30
4	A	1401	SAM	C5-C4	-2.22	1.35	1.39
4	A	1401	SAM	OXT-C	-2.07	1.24	1.30

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	2401	SAM	N9-C8-N7	-9.62	100.28	113.94
4	A	1401	SAM	N9-C8-N7	-9.33	100.70	113.94
4	B	2401	SAM	C5-N7-C8	8.33	116.54	103.45
4	A	1401	SAM	C5-N7-C8	8.32	116.53	103.45
4	A	1401	SAM	N3-C2-N1	-6.05	119.43	128.58
4	B	2401	SAM	N3-C2-N1	-5.95	119.58	128.58
4	B	2401	SAM	C4-N9-C1'	-4.05	117.15	126.63
4	A	1401	SAM	C4-N9-C1'	-3.54	118.36	126.63
4	B	2401	SAM	C4-N9-C8	3.45	109.36	105.74
4	A	1401	SAM	CG-SD-C5'	2.98	110.73	103.43
4	B	2401	SAM	C1'-N9-C8	2.88	133.48	127.09
4	A	1401	SAM	C4-N9-C8	2.86	108.74	105.74
4	A	1401	SAM	OXT-C-O	-2.75	117.83	124.08
4	A	1401	SAM	C1'-N9-C8	2.61	132.90	127.09
4	B	2401	SAM	C2-N3-C4	2.57	118.11	111.83
4	A	1401	SAM	C2-N3-C4	2.49	117.91	111.83
4	A	1401	SAM	CB-CA-C	2.26	116.43	110.45
4	A	1401	SAM	C2-N1-C6	2.21	122.36	118.73
4	B	2401	SAM	C2-N1-C6	2.19	122.33	118.73
4	B	2401	SAM	O4'-C1'-C2'	-2.14	102.03	106.62
4	A	1401	SAM	N3-C4-N9	2.12	130.78	127.17
4	A	1401	SAM	N6-C6-N1	2.03	122.89	118.38

There are no chirality outliers.

All (14) torsion outliers are listed below:

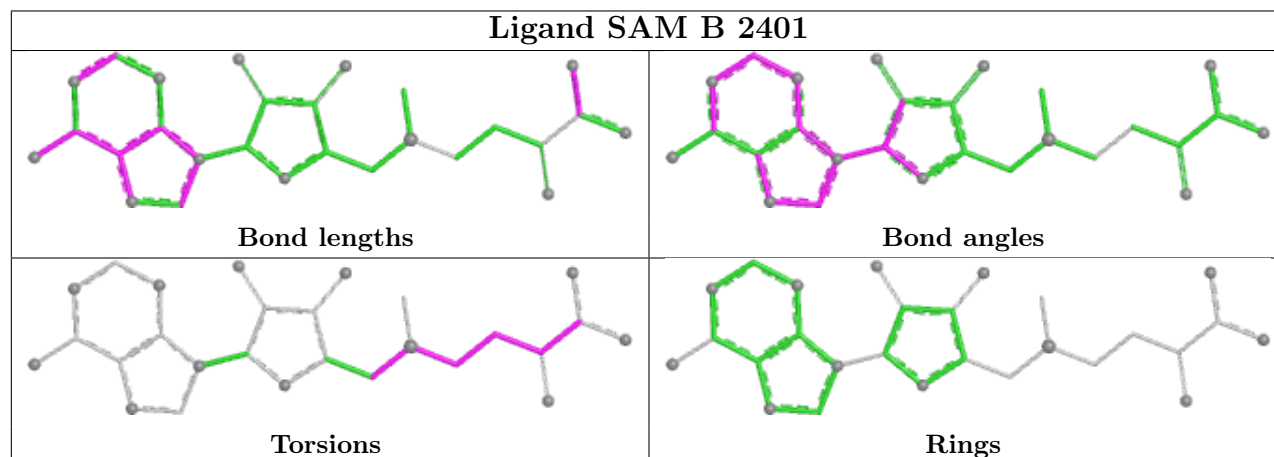
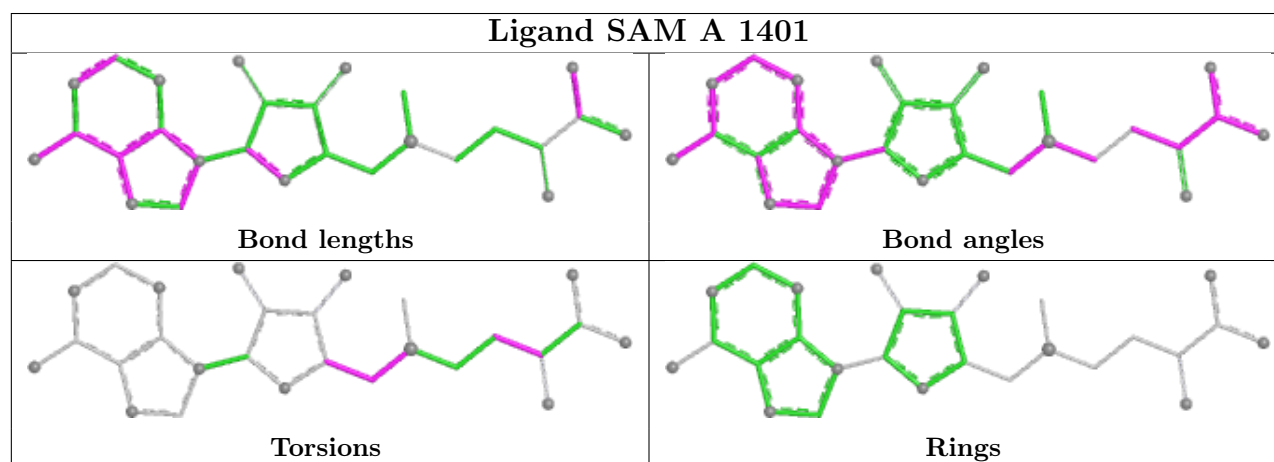
Mol	Chain	Res	Type	Atoms
4	A	1401	SAM	N-CA-CB-CG
4	A	1401	SAM	C-CA-CB-CG
4	A	1401	SAM	C4'-C5'-SD-CG
4	A	1401	SAM	C4'-C5'-SD-CE
4	A	1401	SAM	O4'-C4'-C5'-SD
4	A	1401	SAM	C3'-C4'-C5'-SD
4	B	2401	SAM	N-CA-CB-CG
4	B	2401	SAM	C-CA-CB-CG
4	B	2401	SAM	CA-CB-CG-SD
4	B	2401	SAM	C4'-C5'-SD-CE
4	B	2401	SAM	OXT-C-CA-N
4	B	2401	SAM	O-C-CA-CB
4	B	2401	SAM	OXT-C-CA-CB
4	B	2401	SAM	CB-CG-SD-C5'

There are no ring outliers.

2 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1401	SAM	5	0
4	B	2401	SAM	15	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	261/269 (97%)	-0.17	9 (3%)	48	39	20, 47, 94, 99	0
1	B	251/269 (93%)	0.06	6 (2%)	59	49	17, 56, 93, 99	0
All	All	512/538 (95%)	-0.06	15 (2%)	53	43	17, 51, 94, 99	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	13	SER	6.1
1	B	15	GLU	3.4
1	A	49	GLY	3.3
1	A	47	ASP	3.2
1	A	42	HIS	3.1
1	A	45	SER	2.9
1	A	48	PRO	2.8
1	B	2	SER	2.8
1	A	46	ARG	2.8
1	B	24	ARG	2.7
1	B	20	ILE	2.6
1	A	200	PRO	2.5
1	B	269	TYR	2.5
1	A	202	ALA	2.3
1	A	40	VAL	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](#)

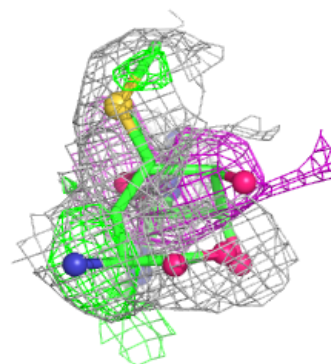
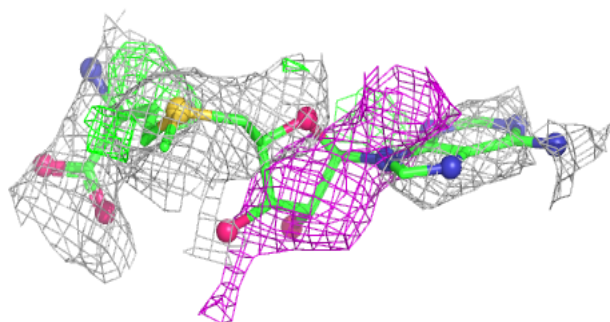
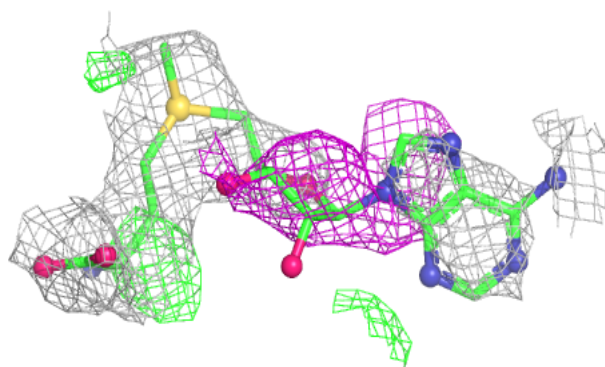
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	SO4	B	504	5/5	0.64	0.31	96,97,97,98	0
3	SO4	B	505	5/5	0.66	0.24	95,95,96,98	0
4	SAM	A	1401	27/27	0.67	0.26	90,96,98,98	0
4	SAM	B	2401	27/27	0.79	0.19	97,98,98,98	0
2	ZN	B	2301	1/1	0.85	0.14	95,95,95,95	0
3	SO4	A	501	5/5	0.87	0.08	73,74,74,76	0
3	SO4	B	502	5/5	0.91	0.09	83,86,87,87	0
3	SO4	B	503	5/5	0.91	0.16	78,80,82,84	0
2	ZN	A	1301	1/1	0.99	0.02	62,62,62,62	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.

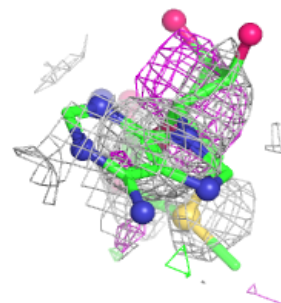
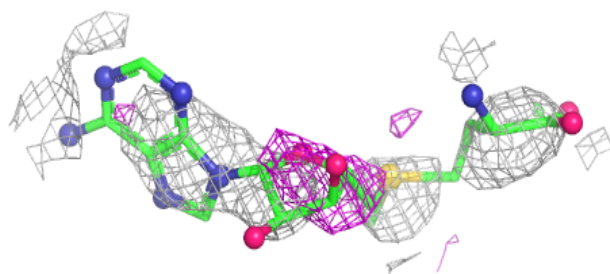
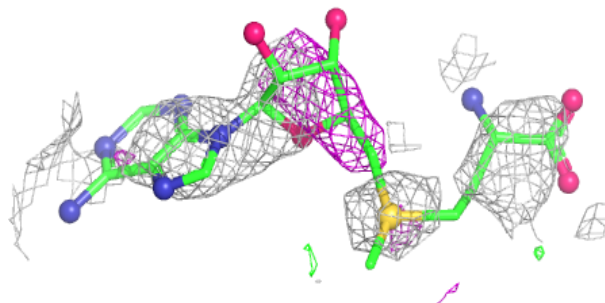
Electron density around SAM A 1401:

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 SAM B 2401:

$2mF_o-DF_c$ (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.