



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:49 PM UTC

PDB ID : 7UCI / pdb_00007uci
Title : SxtA Methyltransferase and decarboxylase didomain in complex with Mn²⁺ and SAH
Authors : Lao, Y.; Skiba, M.A.; Smith, J.L.
Deposited on : 2022-03-16
Resolution : 2.60 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

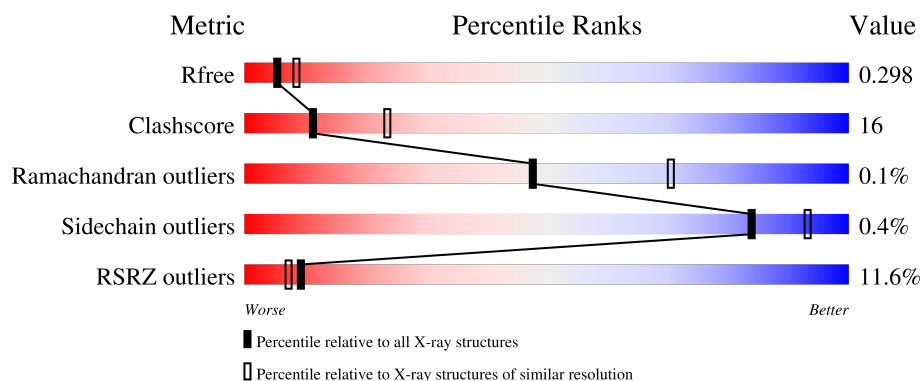
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	713	8% 66% 30% .
1	B	713	14% 65% 28% 6%

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	GOL	B	803	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 21540 atoms, of which 10658 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.

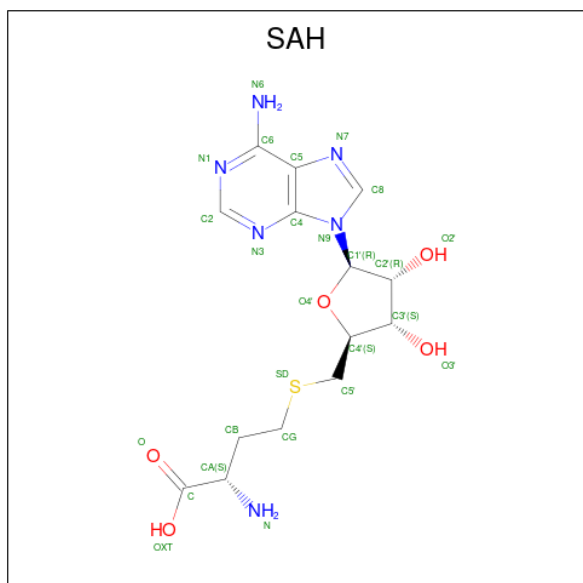
- Molecule 1 is a protein called Polyketide synthase-related protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	686	Total	C	H	N	O	S	0	2	0
			10869	3485	5415	942	1002	25			
1	B	670	Total	C	H	N	O	S	0	2	0
			10479	3380	5198	912	965	24			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP B3EYF9
A	-1	ASN	-	expression tag	UNP B3EYF9
A	0	ALA	-	expression tag	UNP B3EYF9
B	-2	SER	-	expression tag	UNP B3EYF9
B	-1	ASN	-	expression tag	UNP B3EYF9
B	0	ALA	-	expression tag	UNP B3EYF9

- Molecule 2 is S-ADENOSYL-L-HOMOCYSTEINE (CCD ID: SAH) (formula: $C_{14}H_{20}N_6O_5S$) (labeled as "Ligand of Interest" by depositor).

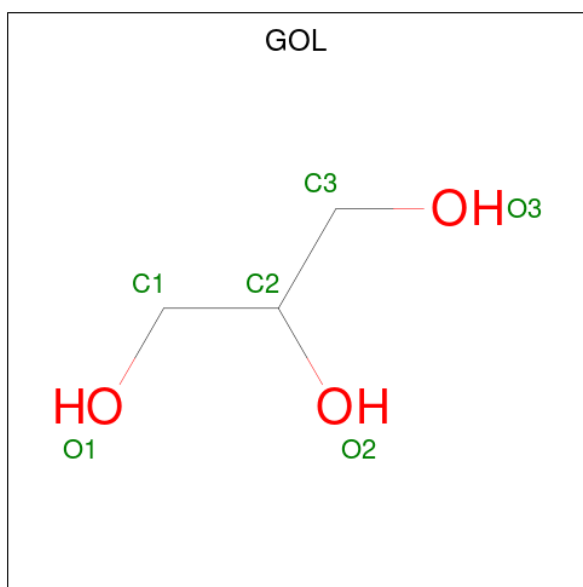


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	S	
			45	14	19	6	5	1	
2	B	1	Total	C	H	N	O	S	
			44	14	18	6	5	1	

- Molecule 3 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mn		
			1	1	0	0
3	B	1	Total	Mn		
			1	1	0	0

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	H	O		
			14	3	8	3	0	0

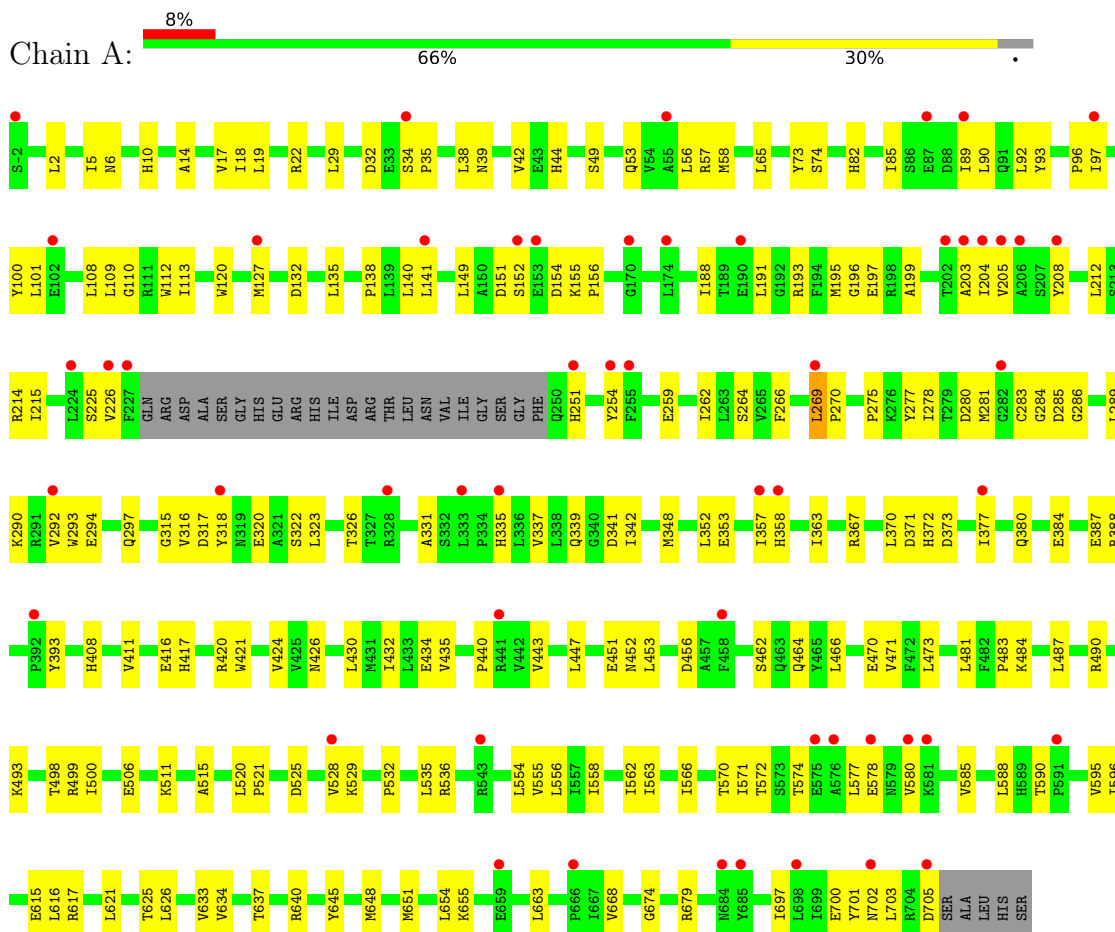
- Molecule 5 is water.

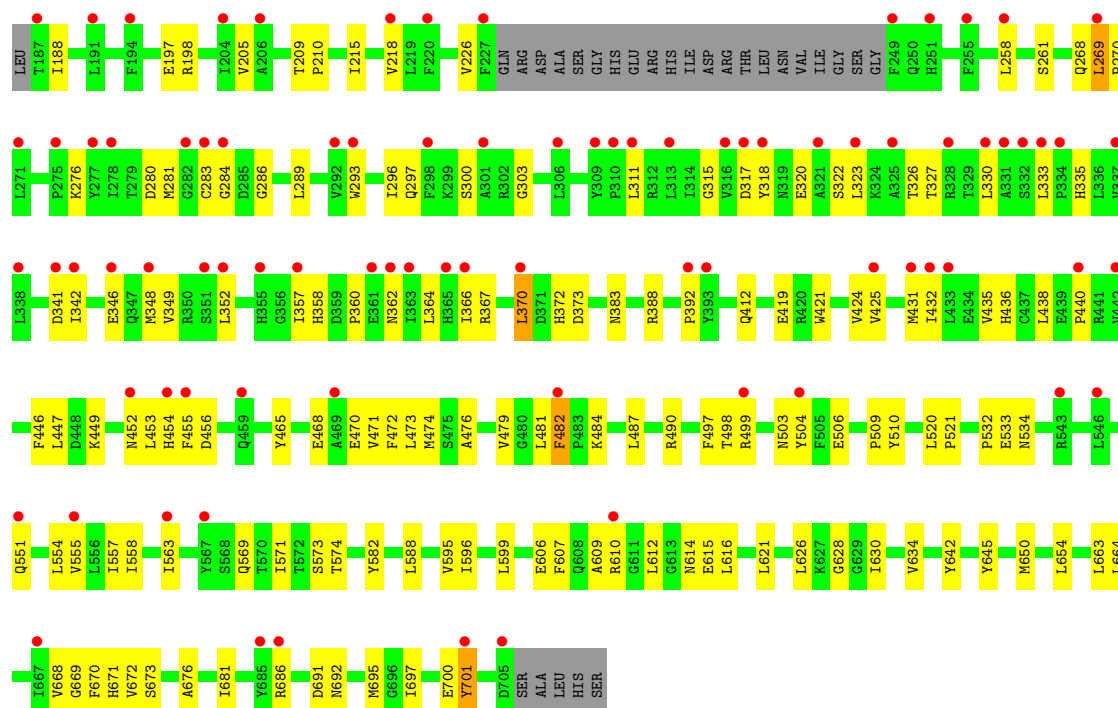
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	47	Total	O		
			47	47	0	0
5	B	40	Total	O		
			40	40	0	0

3 Residue-property plots [i](#)

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

- Molecule 1: Polyketide synthase-related protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	43.80Å 69.67Å 127.61Å 86.44° 83.26° 83.95°	Depositor
Resolution (Å)	42.43 – 2.60 42.43 – 2.60	Depositor EDS
% Data completeness (in resolution range)	97.8 (42.43-2.60) 83.4 (42.43-2.60)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.68 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.237 , 0.298 0.238 , 0.298	Depositor DCC
R_{free} test set	2006 reflections (4.36%)	wwPDB-VP
Wilson B-factor (Å ²)	41.4	Xtriage
Anisotropy	0.400	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 45.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	21540	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.53% 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: GOL, MN, 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	0.12	0/5582	0.32	0/7578
1	B	0.13	0/5404	0.32	0/7338
All	All	0.13	0/10986	0.32	0/14916

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	5454	5415	5406	169	0
1	B	5281	5198	5189	181	0
2	A	26	19	19	6	0
2	B	26	18	19	6	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	B	6	8	8	4	0
5	A	47	0	0	9	0
5	B	40	0	0	12	0
All	All	10882	10658	10641	347	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 16.

All (347) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:317:ASP:OD1	2:B:801:SAH:O2'	1.71	1.08
1:B:317:ASP:CG	2:B:801:SAH:O2'	2.03	1.01
1:A:317:ASP:OD1	2:A:801:SAH:O2'	1.89	0.90
1:B:373:ASP:OD1	5:B:901:HOH:O	1.91	0.87
1:A:331:ALA:O	1:B:610:ARG:NH2	2.13	0.82
1:B:293:TRP:NE1	1:B:297:GLN:OE1	2.14	0.81
1:A:556:LEU:HD22	1:A:563:ILE:HD11	1.61	0.81
1:A:490:ARG:NH2	1:A:493:LYS:O	2.14	0.81
1:A:585:VAL:HA	1:A:588:LEU:HD13	1.64	0.80
1:B:447:LEU:HD11	1:B:453:LEU:HD22	1.61	0.80
1:B:362:ASN:O	5:B:903:HOH:O	2.01	0.79
1:A:112:TRP:N	5:A:902:HOH:O	2.17	0.78
1:A:283:CYS:HB3	1:A:326:THR:HG21	1.64	0.78
1:B:383:ASN:ND2	5:B:907:HOH:O	2.19	0.76
1:B:19:LEU:HD11	1:B:89:ILE:HD11	1.67	0.75
1:B:468:GLU:OE2	5:B:904:HOH:O	2.05	0.74
1:A:615:GLU:OE2	5:A:901:HOH:O	2.05	0.73
1:B:97:ILE:HD12	1:B:138:PRO:HB2	1.70	0.72
1:A:556:LEU:HB2	1:A:616:LEU:HD12	1.72	0.71
1:A:571:ILE:HD12	1:A:588:LEU:HB3	1.70	0.71
1:B:456:ASP:OD1	5:B:905:HOH:O	2.09	0.71
1:B:366:ILE:HG12	1:B:431:MET:HG2	1.73	0.71
1:A:528:VAL:HG12	1:A:536:ARG:HD3	1.72	0.71
1:A:97:ILE:HD12	1:A:138:PRO:HB2	1.73	0.70
1:B:317:ASP:OD2	2:B:801:SAH:O2'	2.09	0.70
1:A:572:THR:HG23	1:A:590:THR:HG21	1.75	0.69
1:A:331:ALA:HB1	1:B:509:PRO:HG3	1.74	0.69
1:B:645:TYR:HE2	1:B:664:LEU:HD21	1.58	0.69
1:B:276:LYS:O	1:B:311:LEU:HD12	1.92	0.68
1:B:34:SER:HB3	1:B:35:PRO:HD3	1.76	0.68
1:A:262:ILE:HD13	1:A:292:VAL:HG22	1.76	0.68
1:B:634:VAL:HG12	1:B:700:GLU:HG3	1.77	0.67
1:B:197:GLU:OE1	1:B:198:ARG:NH1	2.29	0.66
1:B:33:GLU:HB3	1:B:36:LEU:HD11	1.77	0.66
1:B:676:ALA:HB2	1:B:701:TYR:CE1	2.31	0.65
1:B:490:ARG:NH2	1:B:497:PHE:O	2.29	0.65
1:A:284:GLY:HA2	2:A:801:SAH:O3'	1.97	0.65
1:B:474:MET:HE1	1:B:554:LEU:HD22	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:ASN:O	1:A:10:HIS:ND1	2.23	0.65
1:B:342:ILE:HD13	1:B:370:LEU:HD13	1.79	0.65
1:B:691:ASP:OD2	5:B:906:HOH:O	2.15	0.65
1:A:214:ARG:NH2	1:A:225:SER:OG	2.30	0.64
1:A:470:GLU:OE2	5:A:903:HOH:O	2.15	0.64
1:B:654:LEU:HD11	1:B:697:ILE:HD12	1.79	0.64
1:A:637:THR:HG23	1:A:697:ILE:HG23	1.80	0.63
4:B:803:GOL:O1	5:B:902:HOH:O	1.96	0.63
1:B:681:ILE:HG21	1:B:695:MET:HE2	1.80	0.63
1:A:34:SER:HB2	1:A:35:PRO:HD3	1.80	0.63
1:A:555:VAL:HG13	1:A:562:ILE:HG23	1.81	0.63
1:B:473:LEU:HD22	1:B:626:LEU:HD12	1.80	0.63
1:A:451:GLU:OE1	1:A:451:GLU:N	2.31	0.63
1:A:342:ILE:HD11	1:A:370:LEU:CD2	2.28	0.63
1:B:555:VAL:HG12	1:B:557:ILE:HG13	1.81	0.62
1:A:204:ILE:HG23	1:A:205:VAL:HG13	1.79	0.62
1:A:283:CYS:SG	1:A:316:VAL:N	2.72	0.62
1:B:609:ALA:N	5:B:902:HOH:O	2.19	0.62
1:B:671:HIS:HD1	1:B:701:TYR:HH	1.45	0.62
1:A:572:THR:CG2	1:A:590:THR:HG21	2.30	0.62
1:B:81:ILE:HD11	1:B:122:LEU:HD22	1.81	0.61
1:B:571:ILE:HD12	1:B:588:LEU:HB3	1.82	0.61
1:B:35:PRO:O	1:B:36:LEU:HD12	2.01	0.61
1:B:22:ARG:NH1	1:B:82:HIS:O	2.34	0.60
1:A:44:HIS:NE2	5:A:907:HOH:O	2.31	0.60
4:B:803:GOL:O2	5:B:902:HOH:O	2.17	0.60
1:B:621:LEU:HD11	1:B:701:TYR:CD2	2.37	0.60
1:B:100:TYR:OH	1:B:168:GLU:OE1	2.13	0.59
1:B:330:LEU:HD23	1:B:333:LEU:HD12	1.83	0.59
1:A:39:ASN:HA	1:A:42:VAL:HG22	1.84	0.59
1:A:570:THR:O	1:A:571:ILE:HD13	2.03	0.59
1:A:280:ASP:HB2	1:A:289:LEU:HD21	1.83	0.58
1:B:296:ILE:O	1:B:300:SER:OG	2.14	0.58
1:B:574:THR:HG22	1:B:634:VAL:HG11	1.83	0.58
1:A:352:LEU:HD22	1:A:357:ILE:HD12	1.83	0.58
1:B:188:ILE:HD12	1:B:188:ILE:H	1.68	0.58
1:B:498:THR:HG21	1:B:626:LEU:HD22	1.85	0.58
1:A:151:ASP:HB3	1:A:193:ARG:HH21	1.69	0.58
1:A:331:ALA:O	1:B:610:ARG:CZ	2.51	0.58
1:A:498:THR:HG21	1:A:626:LEU:HD22	1.84	0.57
1:B:341:ASP:OD1	1:B:342:ILE:N	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:621:LEU:HD21	1:B:701:TYR:CD2	2.40	0.57
1:A:286:GLY:N	1:A:326:THR:HG22	2.19	0.57
1:A:317:ASP:CG	2:A:801:SAH:O2'	2.47	0.57
1:B:281:MET:HG2	1:B:342:ILE:CG2	2.35	0.57
1:A:281:MET:HE2	1:A:348:MET:CE	2.35	0.57
1:A:317:ASP:OD2	2:A:801:SAH:O3'	2.13	0.57
1:A:408:HIS:HA	1:A:411:VAL:HG22	1.86	0.56
1:A:484:LYS:HG3	1:A:506:GLU:HG2	1.87	0.56
1:B:330:LEU:O	1:B:335:HIS:ND1	2.38	0.56
1:A:637:THR:CG2	1:A:697:ILE:HG23	2.35	0.56
1:A:154:ASP:OD1	1:A:155:LYS:N	2.36	0.56
1:B:283:CYS:HB2	1:B:326:THR:HG21	1.87	0.56
1:B:30:LEU:HD11	1:B:41:MET:SD	2.45	0.56
1:A:317:ASP:O	1:A:339:GLN:HA	2.06	0.56
1:A:269:LEU:CB	1:A:270:PRO:HD3	2.36	0.56
1:B:209:THR:HB	1:B:210:PRO:HD3	1.88	0.56
1:A:140:LEU:HB3	1:A:199:ALA:HB1	1.86	0.56
1:A:262:ILE:CD1	1:A:292:VAL:HG22	2.36	0.56
1:B:352:LEU:HD22	1:B:357:ILE:HD12	1.87	0.55
1:A:19:LEU:HD11	1:A:89:ILE:HD11	1.89	0.55
1:A:85:ILE:HD11	1:A:127:MET:HE2	1.88	0.55
1:B:317:ASP:OD1	2:B:801:SAH:N3	2.40	0.55
1:B:35:PRO:C	1:B:36:LEU:HD12	2.31	0.55
1:A:17:VAL:HG12	1:A:215:ILE:HG12	1.89	0.55
1:A:141:LEU:HD12	1:A:203:ALA:HB1	1.89	0.54
1:B:36:LEU:O	1:B:73:TYR:N	2.30	0.54
1:B:474:MET:HE1	1:B:554:LEU:CD2	2.37	0.54
1:A:38:LEU:HD22	1:A:53:GLN:OE1	2.06	0.54
1:B:327:THR:HG23	1:B:335:HIS:NE2	2.23	0.54
1:B:286:GLY:N	1:B:326:THR:HG22	2.23	0.54
1:B:498:THR:HG21	1:B:626:LEU:CD2	2.38	0.54
1:A:285:ASP:C	1:A:326:THR:HG22	2.33	0.54
1:B:342:ILE:HD13	1:B:370:LEU:CD1	2.37	0.53
1:A:499:ARG:HD3	1:A:500:ILE:HG13	1.90	0.53
1:B:388:ARG:NE	1:B:412:GLN:OE1	2.36	0.53
1:A:318:TYR:HA	1:A:339:GLN:HG3	1.90	0.53
1:A:113:ILE:N	5:A:902:HOH:O	2.11	0.53
1:A:484:LYS:HE3	1:A:506:GLU:CD	2.34	0.53
1:B:358:HIS:C	1:B:360:PRO:HD3	2.34	0.53
1:A:596:ILE:HB	1:A:633:VAL:HG22	1.91	0.53
1:A:96:PRO:HD2	1:A:108:LEU:HD21	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:446:PHE:HA	1:B:449:LYS:HE2	1.91	0.53
1:A:17:VAL:CG2	1:A:58:MET:HE1	2.39	0.53
1:B:60:GLU:HG2	1:B:440:PRO:HB2	1.91	0.53
1:A:254:TYR:CZ	1:A:500:ILE:HD12	2.45	0.52
1:B:669:GLY:HA2	1:B:672:VAL:HG22	1.90	0.52
1:B:421:TRP:O	1:B:425:VAL:HG23	2.09	0.52
1:A:335:HIS:CE1	1:B:610:ARG:HD2	2.45	0.52
1:A:342:ILE:HD11	1:A:370:LEU:HD21	1.89	0.52
1:B:269:LEU:CB	1:B:270:PRO:HD3	2.40	0.52
1:B:606:GLU:HB2	1:B:607:PHE:HD1	1.74	0.52
1:A:323:LEU:HD22	1:A:337:VAL:HG13	1.92	0.52
1:A:384:GLU:O	1:A:387:GLU:HG2	2.10	0.51
1:A:367:ARG:HD2	1:A:371:ASP:OD2	2.10	0.51
1:A:251:HIS:HB2	1:A:254:TYR:CZ	2.45	0.51
1:A:372:HIS:ND1	1:A:434:GLU:OE2	2.44	0.51
1:B:258:LEU:HG	1:B:431:MET:HE2	1.92	0.51
1:B:89:ILE:HD12	1:B:90:LEU:N	2.26	0.51
1:A:563:ILE:C	1:A:563:ILE:HD12	2.36	0.51
1:B:30:LEU:O	1:B:34:SER:HA	2.11	0.51
1:B:281:MET:HG2	1:B:342:ILE:HG23	1.92	0.50
1:A:269:LEU:CB	1:A:270:PRO:CD	2.89	0.50
1:B:320:GLU:HA	1:B:323:LEU:HD12	1.94	0.50
1:B:481:LEU:C	1:B:482:PHE:HD1	2.19	0.50
1:B:280:ASP:HB2	1:B:289:LEU:HD21	1.92	0.50
1:A:651:MET:O	1:A:655:LYS:HG3	2.11	0.50
1:B:6:ASN:HA	1:B:9:THR:HG22	1.93	0.50
1:A:22:ARG:NH1	1:A:82:HIS:O	2.44	0.50
1:B:432:ILE:HD12	1:B:432:ILE:O	2.12	0.50
1:B:558:ILE:HD12	1:B:563:ILE:HG13	1.94	0.50
1:B:431:MET:HA	1:B:503:ASN:O	2.11	0.49
1:A:2:LEU:HD23	1:A:5:ILE:HD12	1.92	0.49
1:B:33:GLU:O	1:B:36:LEU:HD13	2.12	0.49
1:B:86:SER:O	1:B:89:ILE:HG13	2.12	0.49
1:B:571:ILE:HG22	1:B:573:SER:H	1.78	0.49
1:B:672:VAL:HG23	1:B:673:SER:N	2.28	0.49
1:A:29:LEU:O	1:A:32:ASP:N	2.45	0.49
1:A:269:LEU:HB3	1:A:270:PRO:HD3	1.94	0.49
1:A:637:THR:HG23	1:A:697:ILE:CG2	2.43	0.49
1:A:341:ASP:OD1	1:A:342:ILE:N	2.46	0.49
1:A:435:VAL:HG13	1:A:466:LEU:HD21	1.95	0.48
1:A:640:ARG:NH2	5:A:908:HOH:O	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:ALA:CB	1:B:509:PRO:HG3	2.42	0.48
1:B:616:LEU:C	1:B:616:LEU:HD23	2.38	0.48
1:A:266:PHE:CZ	1:A:278:ILE:HD11	2.47	0.48
1:A:571:ILE:HB	1:A:595:VAL:HG13	1.95	0.48
1:B:645:TYR:HE2	1:B:664:LEU:CD2	2.23	0.48
1:A:208:TYR:O	1:A:212:LEU:HG	2.13	0.48
1:B:269:LEU:HB2	1:B:270:PRO:HD3	1.95	0.48
1:B:571:ILE:HB	1:B:595:VAL:HG23	1.95	0.48
1:A:264:SER:HB3	1:A:487:LEU:HD21	1.95	0.48
1:A:616:LEU:C	1:A:616:LEU:HD23	2.39	0.48
1:A:152:SER:HA	1:A:156:PRO:HA	1.95	0.48
1:B:12:PHE:CD1	1:B:205:VAL:HG22	2.49	0.48
1:A:318:TYR:HE1	2:A:801:SAH:N7	2.11	0.48
1:A:353:GLU:OE2	1:A:358:HIS:NE2	2.47	0.48
1:A:353:GLU:OE2	1:A:358:HIS:CD2	2.67	0.48
1:B:364:LEU:HD21	1:B:366:ILE:HD11	1.96	0.48
1:B:424:VAL:HG12	1:B:424:VAL:O	2.13	0.47
1:B:681:ILE:HG21	1:B:695:MET:CE	2.43	0.47
1:B:67:ARG:NH2	5:B:909:HOH:O	2.41	0.47
1:A:515:ALA:HB2	1:A:555:VAL:HG23	1.95	0.47
1:B:297:GLN:O	1:B:303:GLY:HA3	2.14	0.47
1:B:330:LEU:CD2	1:B:333:LEU:HD12	2.43	0.47
1:B:435:VAL:HG23	1:B:499:ARG:HD2	1.96	0.47
1:B:621:LEU:HD11	1:B:701:TYR:CE2	2.49	0.47
1:B:663:LEU:HD12	1:B:672:VAL:HG21	1.96	0.47
1:A:65:LEU:HA	1:A:74:SER:O	2.15	0.47
1:A:109:LEU:HD22	1:A:135:LEU:HD11	1.97	0.47
1:A:191:LEU:HD11	1:A:195:MET:HE2	1.97	0.47
1:A:293:TRP:O	1:A:297:GLN:N	2.40	0.47
1:A:56:LEU:HD13	1:A:73:TYR:CE2	2.48	0.47
1:B:23:GLU:OE2	1:B:87:GLU:OE2	2.31	0.47
1:B:17:VAL:HG21	1:B:58:MET:SD	2.55	0.47
1:B:38:LEU:O	1:B:42:VAL:HG13	2.14	0.47
1:B:269:LEU:CB	1:B:270:PRO:CD	2.93	0.47
1:A:323:LEU:HD22	1:A:337:VAL:CG1	2.45	0.47
1:A:280:ASP:CB	1:A:289:LEU:HD21	2.44	0.47
1:A:621:LEU:O	1:A:625:THR:HG23	2.15	0.46
1:B:14:ALA:O	1:B:18:ILE:HG13	2.14	0.46
1:B:621:LEU:HD11	1:B:701:TYR:HD2	1.81	0.46
1:A:668:VAL:HG21	1:A:697:ILE:HD13	1.97	0.46
1:A:266:PHE:CZ	1:A:275:PRO:HG2	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:LEU:HD23	1:A:196:GLY:HA3	1.97	0.46
1:A:421:TRP:O	1:A:424:VAL:HG22	2.15	0.46
1:B:14:ALA:O	1:B:17:VAL:HG22	2.16	0.46
1:B:89:ILE:O	1:B:92:LEU:HD23	2.14	0.46
1:B:510:TYR:OH	1:B:615:GLU:OE1	2.25	0.46
1:A:367:ARG:NH1	1:A:371:ASP:OD1	2.48	0.46
1:A:580:VAL:HG11	1:A:588:LEU:HD11	1.98	0.46
1:B:284:GLY:N	2:B:801:SAH:O3'	2.49	0.46
1:A:387:GLU:HG3	1:A:388:ARG:N	2.30	0.46
1:B:297:GLN:HA	1:B:303:GLY:HA3	1.95	0.46
1:A:645:TYR:O	1:A:648:MET:HG2	2.16	0.46
1:B:7:ARG:HH21	1:B:126:LEU:HA	1.80	0.46
1:A:14:ALA:O	1:A:18:ILE:HG13	2.16	0.46
1:A:17:VAL:HG23	1:A:58:MET:HE1	1.98	0.46
1:A:320:GLU:HA	1:A:323:LEU:HD12	1.98	0.46
1:A:705:ASP:N	1:A:705:ASP:OD1	2.46	0.46
1:B:482:PHE:HD1	1:B:482:PHE:N	2.14	0.46
1:A:484:LYS:HE3	1:A:506:GLU:CG	2.46	0.46
1:B:654:LEU:O	1:B:663:LEU:HD22	2.16	0.45
1:A:470:GLU:HG2	1:A:471:VAL:N	2.29	0.45
1:B:283:CYS:SG	1:B:315:GLY:HA2	2.56	0.45
1:B:621:LEU:CD1	1:B:701:TYR:HD2	2.29	0.45
1:B:70:GLU:O	1:B:71:LEU:HB2	2.16	0.45
1:B:484:LYS:HE3	1:B:506:GLU:OE2	2.16	0.45
1:A:283:CYS:SG	1:A:315:GLY:CA	3.04	0.45
1:B:15:VAL:HB	1:B:16:PRO:HD3	1.97	0.45
1:A:89:ILE:O	1:A:92:LEU:HD23	2.16	0.45
1:A:38:LEU:O	1:A:42:VAL:HG13	2.16	0.45
1:A:100:TYR:HD2	1:A:101:LEU:HD22	1.80	0.45
1:B:333:LEU:O	1:B:335:HIS:N	2.48	0.45
1:B:596:ILE:HG12	1:B:630:ILE:HG12	1.99	0.45
1:A:490:ARG:HD2	1:A:498:THR:OG1	2.17	0.45
1:B:37:SER:OG	5:B:908:HOH:O	2.21	0.45
1:A:109:LEU:C	5:A:902:HOH:O	2.60	0.45
1:B:281:MET:HE2	1:B:348:MET:CE	2.47	0.45
1:B:261:SER:HA	1:B:487:LEU:HD22	1.99	0.45
1:B:431:MET:HB3	1:B:504:TYR:HA	1.99	0.45
1:A:577:LEU:HG	1:A:578:GLU:H	1.82	0.44
1:A:322:SER:O	1:A:326:THR:HG23	2.18	0.44
1:B:370:LEU:HD13	1:B:370:LEU:O	2.18	0.44
1:B:482:PHE:N	1:B:482:PHE:CD1	2.84	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:701:TYR:CD1	1:B:701:TYR:N	2.85	0.44
1:A:654:LEU:O	1:A:663:LEU:HD22	2.18	0.44
1:B:9:THR:HG23	1:B:454:HIS:CE1	2.52	0.44
1:B:139:LEU:CD2	1:B:143:LEU:HD11	2.46	0.44
1:B:419:GLU:OE1	1:B:479:VAL:HG12	2.18	0.44
1:B:582:TYR:CB	1:B:686:ARG:HG2	2.47	0.44
1:A:380:GLN:H	1:A:416:GLU:CD	2.24	0.44
1:A:373:ASP:OD2	1:A:464:GLN:NE2	2.51	0.44
1:B:62:LEU:HD21	1:B:453:LEU:HD21	2.00	0.44
1:B:628:GLY:O	1:B:630:ILE:HD12	2.18	0.44
1:A:149:LEU:HG	1:A:188:ILE:HD11	2.00	0.44
1:B:109:LEU:HD22	1:B:135:LEU:HD11	2.00	0.44
1:B:470:GLU:HG2	1:B:471:VAL:N	2.33	0.44
1:A:417:HIS:HA	1:A:420:ARG:HH11	1.83	0.43
1:B:34:SER:HB3	1:B:35:PRO:CD	2.47	0.43
1:B:520:LEU:N	1:B:521:PRO:HD2	2.33	0.43
1:A:342:ILE:HG22	1:A:421:TRP:HE1	1.82	0.43
1:A:277:TYR:HB2	1:A:363:ILE:HG12	2.00	0.43
1:A:679:ARG:NH1	1:A:700:GLU:OE1	2.51	0.43
1:A:481:LEU:HA	1:A:506:GLU:O	2.18	0.43
1:A:574:THR:HG22	1:A:634:VAL:HG11	2.01	0.43
1:B:133:GLY:O	1:B:137:ILE:HG12	2.18	0.43
1:B:569:GLN:HE21	1:B:571:ILE:HD11	1.84	0.43
1:B:654:LEU:HD22	1:B:668:VAL:HG11	2.00	0.43
1:A:284:GLY:CA	2:A:801:SAH:O3'	2.64	0.43
1:B:610:ARG:NH2	1:B:612:LEU:HD11	2.34	0.43
1:B:468:GLU:OE1	1:B:470:GLU:OE2	2.37	0.43
1:A:193:ARG:O	1:A:197:GLU:HG3	2.19	0.42
1:A:443:VAL:O	1:A:447:LEU:HA	2.20	0.42
1:B:71:LEU:O	5:B:909:HOH:O	2.21	0.42
1:B:283:CYS:CB	1:B:326:THR:HG21	2.48	0.42
1:B:438:LEU:HD12	1:B:452:ASN:HD21	1.84	0.42
1:B:634:VAL:HG12	1:B:700:GLU:CG	2.47	0.42
1:A:57:ARG:HG3	1:A:440:PRO:HG3	2.01	0.42
1:A:254:TYR:CZ	1:A:500:ILE:CD1	3.03	0.42
1:B:113:ILE:HD12	1:B:168:GLU:HB3	2.00	0.42
1:B:642:TYR:HB3	1:B:692:ASN:OD1	2.19	0.42
1:B:650:MET:HB3	1:B:695:MET:HE1	2.02	0.42
1:A:447:LEU:HD11	1:A:453:LEU:HD22	2.00	0.42
1:A:462[B]:SER:O	1:A:462[B]:SER:OG	2.36	0.42
1:A:520:LEU:HB3	1:A:521:PRO:HD3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:269:LEU:HB2	1:B:270:PRO:CD	2.49	0.42
1:B:372:HIS:HB2	1:B:465:TYR:O	2.19	0.42
1:A:451:GLU:HG2	1:A:499:ARG:NH1	2.35	0.42
1:B:432:ILE:HD12	1:B:432:ILE:C	2.45	0.42
1:B:482:PHE:CE2	1:B:510:TYR:CE2	3.06	0.42
1:A:571:ILE:HG22	1:A:572:THR:N	2.34	0.42
1:B:670:PHE:CE1	4:B:803:GOL:H32	2.55	0.42
1:A:562:ILE:HD12	1:A:562:ILE:H	1.85	0.42
1:A:558:ILE:HD12	1:A:563:ILE:HG12	2.02	0.42
1:B:392:PRO:HD2	1:B:551:GLN:NE2	2.34	0.42
1:B:218:VAL:CG2	1:B:226:VAL:HG11	2.50	0.41
1:A:89:ILE:HD12	1:A:90:LEU:N	2.34	0.41
1:A:393:TYR:CE2	1:A:411:VAL:HG21	2.55	0.41
1:A:570:THR:HG22	1:A:596:ILE:CD1	2.51	0.41
1:A:616:LEU:HD23	1:A:616:LEU:O	2.20	0.41
1:A:120:TRP:NE1	1:A:132:ASP:OD1	2.40	0.41
1:A:269:LEU:HB2	1:A:270:PRO:HD3	2.03	0.41
1:A:377:ILE:HD12	1:A:377:ILE:O	2.21	0.41
1:B:366:ILE:HG22	1:B:367:ARG:N	2.35	0.41
1:A:473:LEU:HD11	1:A:483:PRO:HG3	2.03	0.41
1:B:268:GLN:O	1:B:269:LEU:C	2.63	0.41
1:B:297:GLN:HA	1:B:303:GLY:CA	2.50	0.41
1:A:226:VAL:HG13	1:A:226:VAL:O	2.20	0.41
1:A:259:GLU:HA	1:A:262:ILE:HD12	2.03	0.41
1:A:426:ASN:OD1	5:A:904:HOH:O	2.22	0.41
1:A:525:ASP:OD1	1:A:529:LYS:HE3	2.20	0.41
1:B:17:VAL:HG12	1:B:215:ILE:HG12	2.01	0.41
1:B:342:ILE:CG2	1:B:370:LEU:HD12	2.50	0.41
1:B:346:GLU:O	1:B:349:VAL:HG22	2.21	0.41
1:B:432:ILE:HD11	1:B:472:PHE:HE2	1.85	0.41
1:A:264:SER:CB	1:A:487:LEU:HD21	2.51	0.41
1:A:452:ASN:HB3	1:A:456:ASP:OD2	2.21	0.41
1:A:574:THR:CG2	1:A:634:VAL:HG11	2.51	0.41
1:A:625:THR:HG22	1:A:703:LEU:HD11	2.03	0.41
1:B:318:TYR:CD1	2:B:801:SAH:C4	3.03	0.41
1:B:532:PRO:HG2	1:B:534:ASN:OD1	2.21	0.41
1:A:42:VAL:HG11	1:A:49:SER:HA	2.03	0.41
1:A:430:LEU:HD21	1:A:432:ILE:HD11	2.03	0.41
1:B:258:LEU:C	1:B:258:LEU:HD23	2.45	0.41
1:B:281:MET:HG2	1:B:342:ILE:HG22	2.03	0.41
1:B:436:HIS:N	1:B:499:ARG:O	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:476:ALA:HB1	1:B:481:LEU:HB2	2.02	0.41
1:A:701:TYR:O	1:A:702:ASN:C	2.64	0.41
1:B:366:ILE:HG12	1:B:431:MET:CG	2.47	0.41
1:A:110:GLY:C	5:A:902:HOH:O	2.63	0.40
1:B:34:SER:CB	1:B:35:PRO:HD3	2.49	0.40
1:B:136:VAL:HG13	1:B:178:TRP:HH2	1.85	0.40
1:B:317:ASP:OD2	1:B:322:SER:OG	2.40	0.40
1:B:452:ASN:HA	1:B:455:PHE:HB3	2.03	0.40
1:A:532:PRO:HD2	1:A:535:LEU:HD12	2.04	0.40
1:B:532:PRO:O	1:B:533:GLU:C	2.65	0.40
1:B:614:ASN:N	4:B:803:GOL:H11	2.36	0.40
1:A:89:ILE:HD12	1:A:89:ILE:C	2.46	0.40
1:A:290:LYS:HE2	1:A:294:GLU:OE2	2.21	0.40
1:A:554:LEU:HB2	1:A:566:ILE:HG22	2.03	0.40
1:A:668:VAL:CG2	1:A:697:ILE:HD13	2.51	0.40
1:A:90:LEU:HD23	1:A:93:TYR:HE2	1.85	0.40
1:B:10:HIS:HA	1:B:454:HIS:HB3	2.04	0.40
1:A:617:ARG:CZ	1:A:674:GLY:HA3	2.51	0.40
1:B:120:TRP:CG	1:B:131:LEU:HD22	2.57	0.40
1:B:435:VAL:HG23	1:B:499:ARG:CD	2.51	0.40
1:B:571:ILE:HD12	1:B:588:LEU:CB	2.50	0.40
1:B:672:VAL:HG23	1:B:673:SER:H	1.87	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	684/713 (96%)	652 (95%)	31 (4%)	1 (0%)	48	70
1	B	664/713 (93%)	634 (96%)	29 (4%)	1 (0%)	43	66
All	All	1348/1426 (94%)	1286 (95%)	60 (4%)	2 (0%)	48	70

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	269	LEU
1	B	269	LEU

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	593/629 (94%)	592 (100%)	1 (0%)	87	95
1	B	564/629 (90%)	560 (99%)	4 (1%)	76	89
All	All	1157/1258 (92%)	1152 (100%)	5 (0%)	84	93

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	511	LYS
1	B	370	LEU
1	B	482	PHE
1	B	599	LEU
1	B	701	TYR

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

Mol	Chain	Res	Type
1	A	146	HIS
1	A	250	GLN
1	A	297	GLN
1	A	408	HIS
1	A	569	GLN
1	A	589	HIS
1	A	597	GLN
1	A	657	HIS
1	B	48	ASN
1	B	175	HIS

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Mol	Chain	Res	Type
1	B	390	HIS
1	B	569	GLN
1	B	597	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 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 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
2	SAH	B	801	-	27,28,28	0.90	1 (3%)	36,40,40	2.20	14 (38%)
4	GOL	B	803	-	5,5,5	0.46	0	5,5,5	0.98	0
2	SAH	A	801	-	27,28,28	0.73	1 (3%)	36,40,40	1.92	11 (30%)

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	801	-	-	3/15/31/31	0/3/3/3
4	GOL	B	803	-	-	2/4/4/4	-
2	SAH	A	801	-	-	5/15/31/31	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	SAH	OXT-C	-2.56	1.22	1.30
2	A	801	SAH	OXT-C	-2.23	1.23	1.30

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	SAH	N3-C2-N1	-4.47	121.82	128.58
2	B	801	SAH	N3-C2-N1	-4.45	121.84	128.58
2	B	801	SAH	C2'-C3'-C4'	-4.24	94.42	102.61
2	A	801	SAH	C5-N7-C8	3.94	109.64	103.45
2	B	801	SAH	C5-N7-C8	3.93	109.63	103.45
2	A	801	SAH	C5-C4-N3	-3.81	121.47	126.72
2	B	801	SAH	O4'-C1'-C2'	-3.77	98.55	106.62
2	B	801	SAH	N9-C8-N7	-3.74	108.63	113.94
2	A	801	SAH	N9-C8-N7	-3.70	108.69	113.94
2	B	801	SAH	C5-C4-N3	-3.55	121.82	126.72
2	B	801	SAH	O2'-C2'-C3'	-3.46	100.73	111.82
2	B	801	SAH	O4'-C1'-N9	3.05	113.95	108.09
2	A	801	SAH	N3-C4-N9	3.05	132.35	127.17
2	B	801	SAH	N3-C4-N9	3.04	132.34	127.17
2	A	801	SAH	O3'-C3'-C2'	-2.97	102.31	111.82
2	A	801	SAH	C2-N3-C4	2.91	118.94	111.83
2	B	801	SAH	C2-N3-C4	2.83	118.75	111.83
2	B	801	SAH	C4-C5-N7	-2.63	107.57	110.58
2	A	801	SAH	C4-C5-N7	-2.56	107.66	110.58
2	B	801	SAH	C4-N9-C8	2.47	108.33	105.74
2	A	801	SAH	OXT-C-O	-2.39	118.67	124.08
2	B	801	SAH	OXT-C-O	-2.38	118.68	124.08
2	B	801	SAH	C3'-C2'-C1'	2.31	105.84	101.46
2	A	801	SAH	O4'-C4'-C5'	2.06	114.15	108.83
2	A	801	SAH	C4-N9-C8	2.00	107.84	105.74

There are no chirality outliers.

All (10) torsion outliers are listed below:

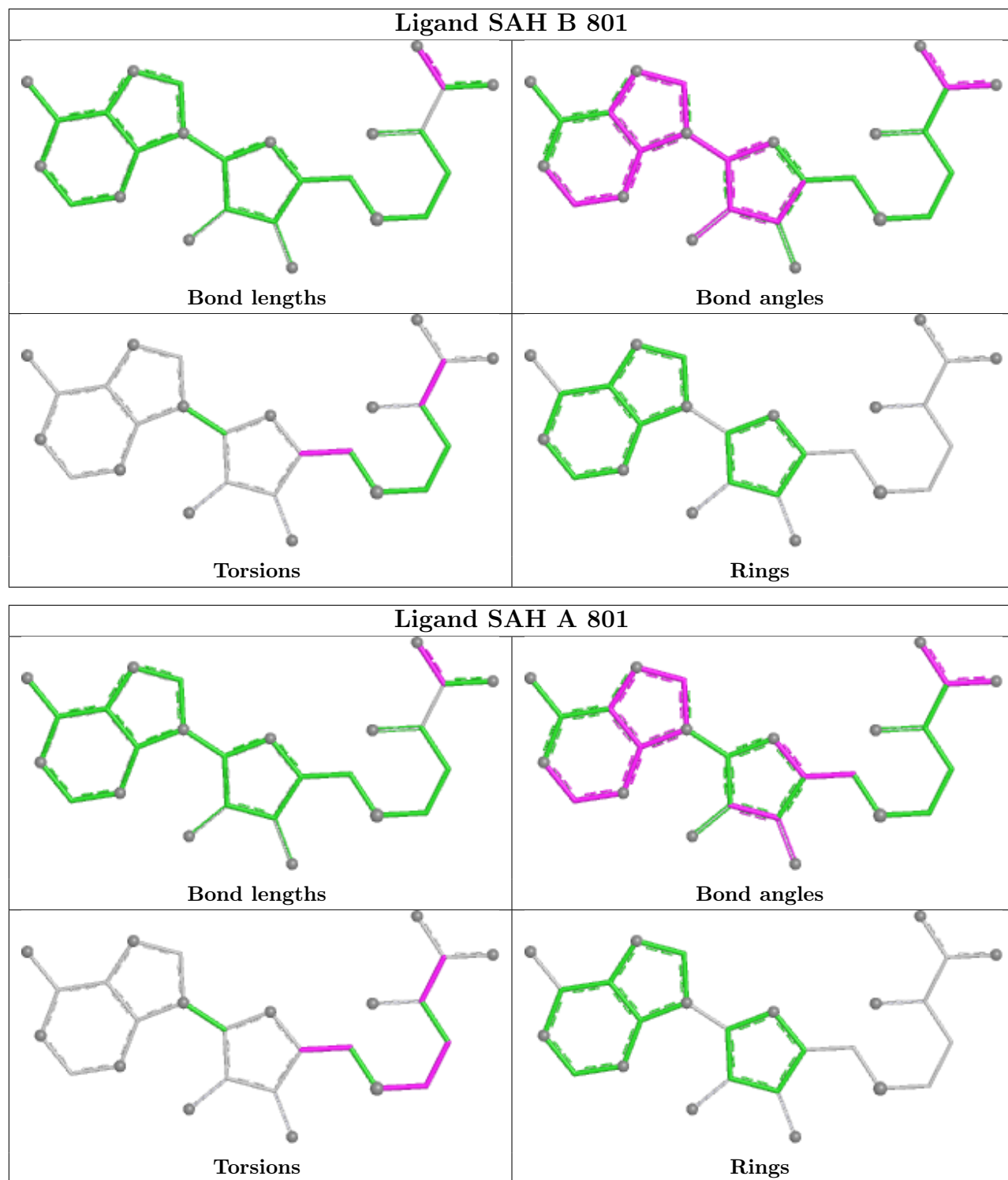
Mol	Chain	Res	Type	Atoms
2	A	801	SAH	O4'-C4'-C5'-SD
2	A	801	SAH	C3'-C4'-C5'-SD
2	A	801	SAH	CA-CB-CG-SD
2	B	801	SAH	OXT-C-CA-N
4	B	803	GOL	O1-C1-C2-C3
2	B	801	SAH	O-C-CA-N
4	B	803	GOL	O1-C1-C2-O2
2	B	801	SAH	C3'-C4'-C5'-SD
2	A	801	SAH	CB-CG-SD-C5'
2	A	801	SAH	OXT-C-CA-N

There are no ring outliers.

3 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	801	SAH	6	0
4	B	803	GOL	4	0
2	A	801	SAH	6	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 ⓘ

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	686/713 (96%)	0.88	54 (7%) 18 14	28, 63, 95, 119	1 (0%)
1	B	670/713 (93%)	1.13	103 (15%) 5 4	29, 69, 132, 241	1 (0%)
All	All	1356/1426 (95%)	1.00	157 (11%) 9 7	28, 66, 121, 241	2 (0%)

All (157) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	332	SER	6.5
1	A	203	ALA	5.6
1	B	321	ALA	5.0
1	A	255	PHE	4.7
1	A	392	PRO	4.6
1	A	152	SER	4.3
1	A	227	PHE	4.3
1	B	366	ILE	4.3
1	B	325	ALA	4.2
1	B	271	LEU	4.2
1	A	528	VAL	4.1
1	B	313	LEU	4.1
1	B	227	PHE	3.9
1	B	204	ILE	3.8
1	A	205	VAL	3.8
1	A	357	ILE	3.7
1	A	458	PHE	3.7
1	B	251	HIS	3.7
1	B	393	TYR	3.7
1	B	147	ASN	3.6
1	B	425	VAL	3.5
1	A	702	ASN	3.4
1	B	258	LEU	3.4
1	A	204	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	81	ILE	3.4
1	B	348	MET	3.4
1	B	311	LEU	3.4
1	B	206	ALA	3.4
1	B	293	TRP	3.3
1	B	365	HIS	3.3
1	A	251	HIS	3.3
1	B	74	SER	3.3
1	A	89	ILE	3.2
1	B	361	GLU	3.2
1	B	218	VAL	3.1
1	B	11	GLY	3.1
1	B	316	VAL	3.1
1	B	284	GLY	3.1
1	A	659	GLU	3.1
1	A	34	SER	3.1
1	A	335	HIS	3.0
1	B	341	ASP	3.0
1	A	170	GLY	3.0
1	A	269	LEU	3.0
1	B	338	LEU	3.0
1	A	581	LYS	3.0
1	B	187	THR	3.0
1	B	283	CYS	2.9
1	B	334	PRO	2.9
1	B	370	LEU	2.9
1	B	-1	ASN	2.9
1	A	206	ALA	2.8
1	B	333	LEU	2.8
1	B	432	ILE	2.8
1	A	705	ASP	2.8
1	B	442	VAL	2.7
1	B	504	TYR	2.7
1	A	543	ARG	2.7
1	B	452	ASN	2.7
1	B	169	LEU	2.7
1	B	292	VAL	2.7
1	B	275	PRO	2.7
1	B	45	LEU	2.6
1	A	377	ILE	2.6
1	B	126	LEU	2.6
1	B	330	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	292	VAL	2.6
1	B	34	SER	2.6
1	B	194	PHE	2.6
1	B	298	PHE	2.6
1	B	440	PRO	2.6
1	B	148	LEU	2.6
1	B	277	TYR	2.6
1	B	323	LEU	2.6
1	B	355	HIS	2.6
1	A	87	GLU	2.6
1	B	123	ASP	2.6
1	B	433	LEU	2.6
1	B	499	ARG	2.6
1	B	454	HIS	2.6
1	B	282	GLY	2.6
1	B	310	PRO	2.6
1	B	431	MET	2.5
1	B	342	ILE	2.5
1	B	363	ILE	2.5
1	A	141	LEU	2.5
1	B	686	ARG	2.5
1	B	567	TYR	2.5
1	A	666	PRO	2.5
1	B	249	PHE	2.5
1	B	301	ALA	2.5
1	B	705	ASP	2.5
1	B	346	GLU	2.5
1	B	309	TYR	2.5
1	B	331	ALA	2.5
1	A	202	THR	2.4
1	A	254	TYR	2.4
1	A	358	HIS	2.4
1	B	269	LEU	2.4
1	B	551	GLN	2.4
1	B	328	ARG	2.4
1	B	685	TYR	2.4
1	B	667	ILE	2.4
1	B	362	ASN	2.4
1	A	580	VAL	2.4
1	A	174	LEU	2.3
1	A	685	TYR	2.3
1	B	108	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	576	ALA	2.3
1	B	317	ASP	2.3
1	A	190	GLU	2.3
1	A	441	ARG	2.3
1	B	255	PHE	2.3
1	B	29	LEU	2.2
1	B	546	LEU	2.2
1	B	278	ILE	2.2
1	A	698	LEU	2.2
1	B	220	PHE	2.2
1	A	153	GLU	2.2
1	A	328	ARG	2.2
1	B	610	ARG	2.2
1	B	306	LEU	2.2
1	B	111	ARG	2.2
1	B	318	TYR	2.2
1	B	337	VAL	2.2
1	B	392	PRO	2.1
1	B	482	PHE	2.1
1	A	575	GLU	2.1
1	A	226	VAL	2.1
1	A	55	ALA	2.1
1	B	469	ALA	2.1
1	B	33	GLU	2.1
1	A	591	PRO	2.1
1	A	127	MET	2.1
1	B	701	TYR	2.1
1	B	357	ILE	2.1
1	B	455	PHE	2.1
1	B	351	SER	2.1
1	A	578	GLU	2.1
1	B	459	GLN	2.1
1	B	191	LEU	2.1
1	B	352	LEU	2.1
1	B	144	HIS	2.1
1	A	208	TYR	2.1
1	A	102	GLU	2.0
1	A	224	LEU	2.0
1	A	97	ILE	2.0
1	B	555	VAL	2.0
1	B	543	ARG	2.0
1	B	79	ALA	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	-2	SER	2.0
1	A	333	LEU	2.0
1	A	282	GLY	2.0
1	A	684	ASN	2.0
1	B	89	ILE	2.0
1	B	563	ILE	2.0
1	A	318	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

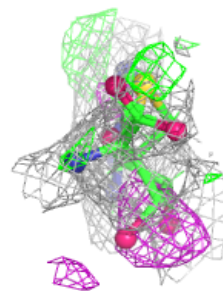
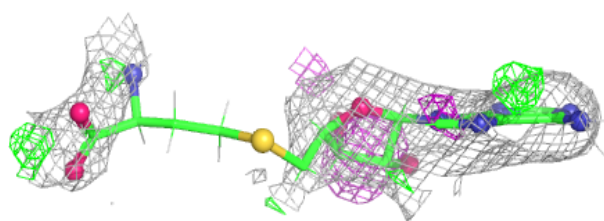
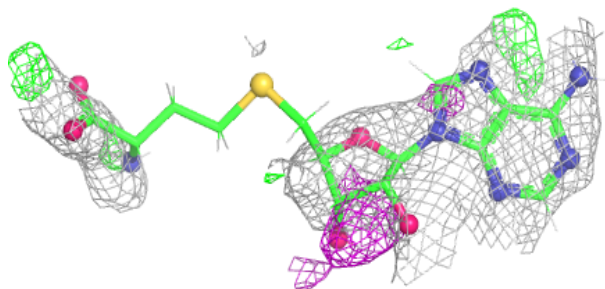
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SAH	B	801	26/26	0.51	0.17	37,52,63,80	7
2	SAH	A	801	26/26	0.68	0.15	28,46,57,87	4
4	GOL	B	803	6/6	0.90	0.15	34,58,75,75	0
3	MN	B	802	1/1	0.98	0.07	56,56,56,56	0
3	MN	A	802	1/1	0.99	0.04	48,48,48,48	0

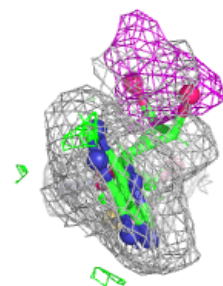
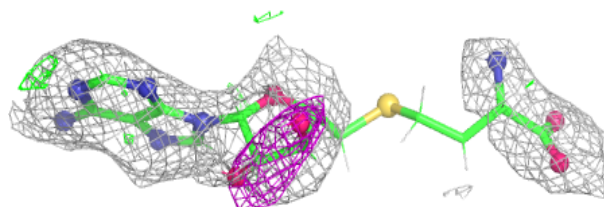
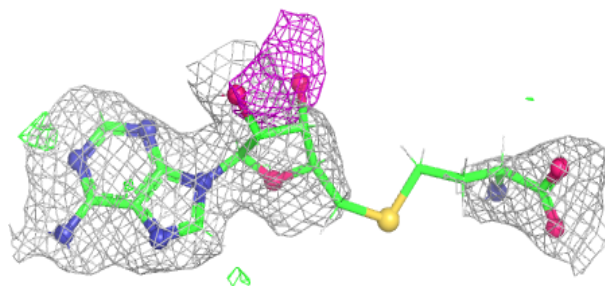
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around SAH B 801:

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

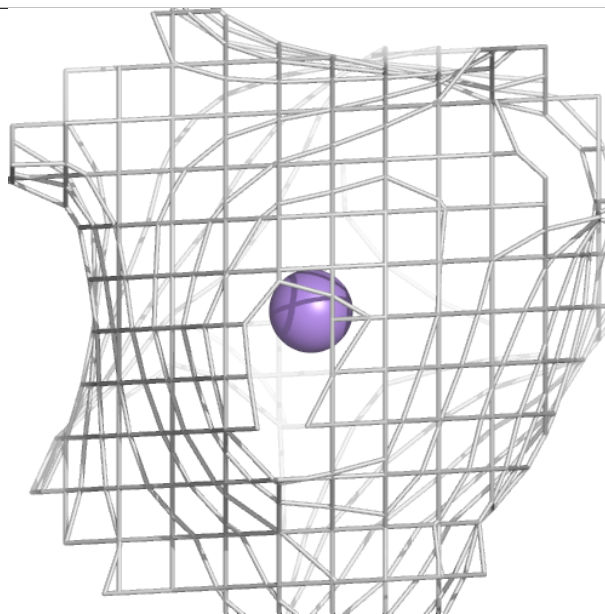
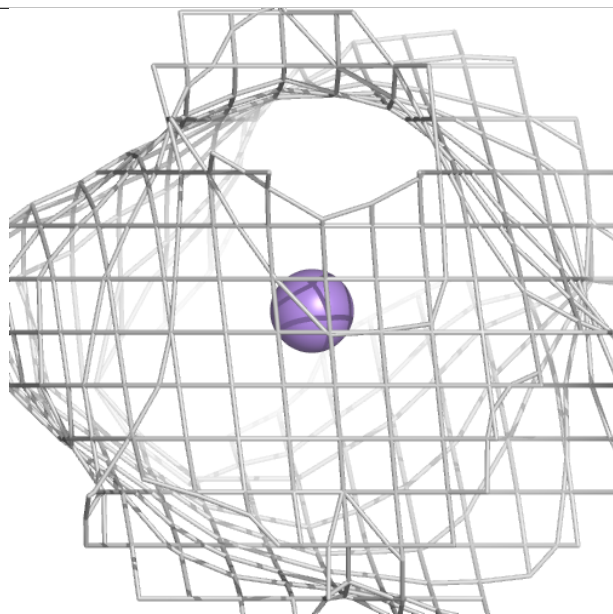
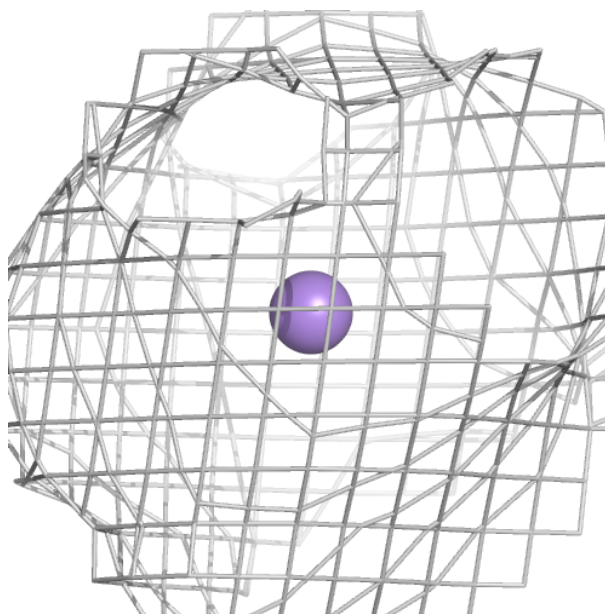
**Electron density around SAH A 801:**

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



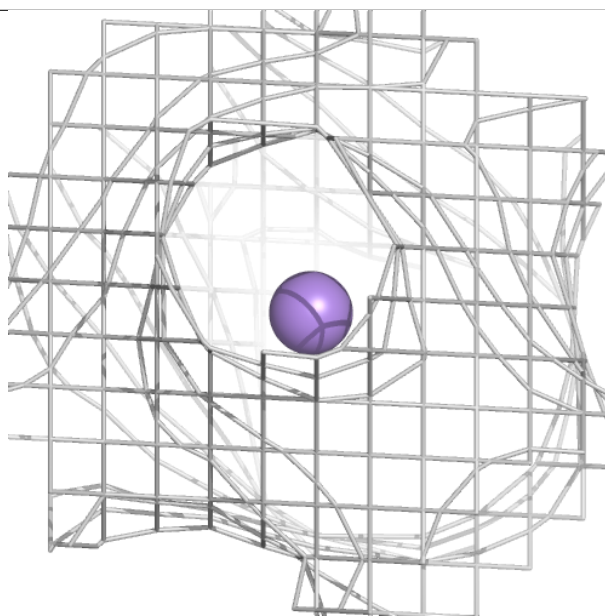
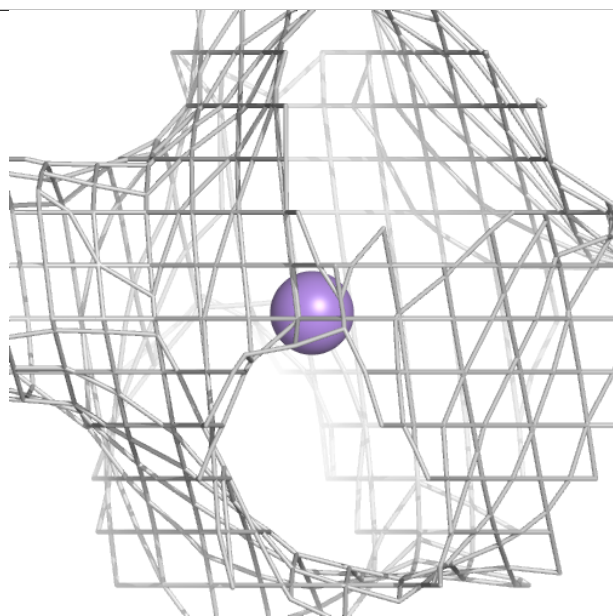
Electron density around MN B 802:

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



Electron density around MN A 802:

$2mF_o-DF_c$ (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.