



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 07:10 AM UTC

PDB ID : 2C2B / pdb_00002c2b
Title : Crystallographic structure of Arabidopsis thaliana Threonine synthase complexed with pyridoxal phosphate and S-adenosylmethionine
Authors : Mas-Droux, C.; Biou, V.; Dumas, R.
Deposited on : 2005-09-27
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)
Xtrriage (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

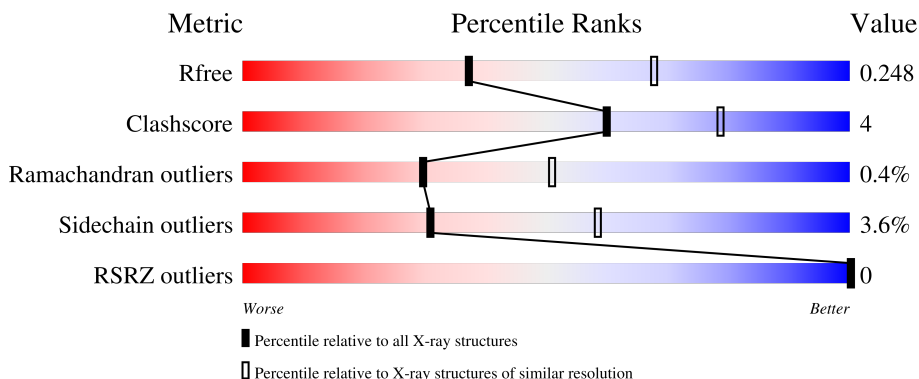
1 Overall quality at a glance

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	486	
1	B	486	
1	C	486	
1	D	486	
1	E	486	

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Mol	Chain	Length	Quality of chain
1	F	486	 76% 14% • 8%

2 Entry composition [i](#)

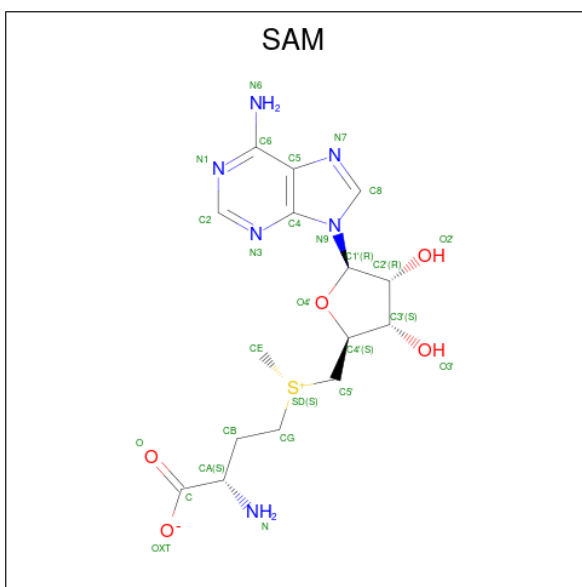
There are 5 unique types of molecules in this entry. The entry contains 20958 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 THREONINE SYNTHASE 1, CHLOROPLASTIC.

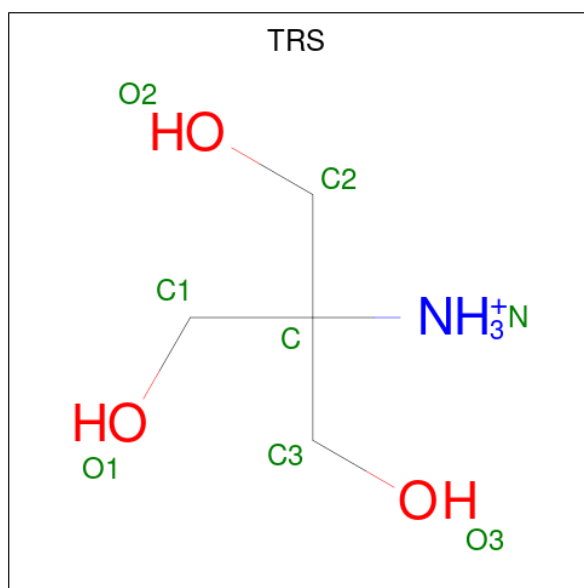
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	445	Total 3428	C 2189	N 576	O 640	S 23	0	0	1
1	B	429	Total 3305	C 2108	N 558	O 617	S 22	0	0	1
1	C	444	Total 3421	C 2184	N 575	O 639	S 23	0	0	1
1	D	429	Total 3305	C 2108	N 558	O 617	S 22	0	0	1
1	E	431	Total 3320	C 2116	N 561	O 621	S 22	0	0	1
1	F	445	Total 3428	C 2189	N 576	O 640	S 23	0	0	1

- Molecule 2 is S-ADENOSYLMETHIONINE (CCD ID: SAM) (formula: C₁₅H₂₂N₆O₅S).



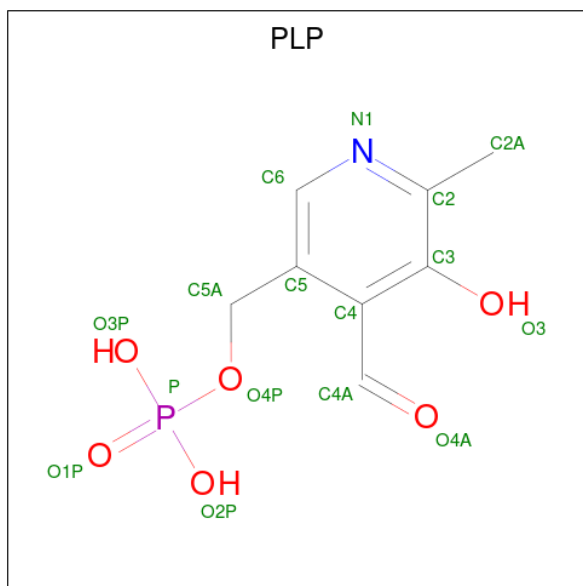
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
2	A	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
2	B	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
2	B	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
2	C	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
2	C	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
2	D	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
2	D	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
2	E	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
2	E	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
2	F	1	Total	C	N	O	S	0	0
			27	15	6	5	1		
2	F	1	Total	C	N	O	S	0	0
			27	15	6	5	1		

- Molecule 3 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			8	4	1	3		
3	C	1	Total	C	N	O	0	0
			8	4	1	3		
3	F	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 4 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: C₈H₁₀NO₆P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
4	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
4	C	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
4	D	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
4	E	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
4	F	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	51	Total	O	0	0
			51	51		

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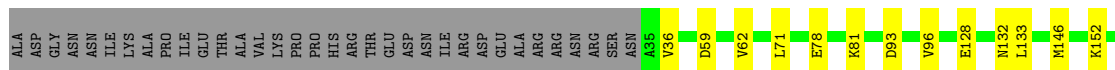
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	55	Total 55	O 55	0	0
5	C	52	Total 52	O 52	0	0
5	D	55	Total 55	O 55	0	0
5	E	44	Total 44	O 44	0	0
5	F	56	Total 56	O 56	0	0



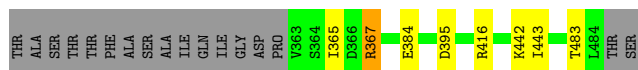
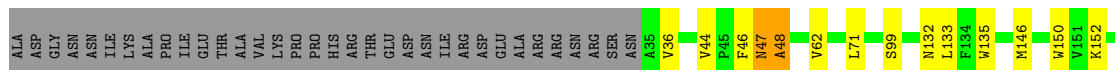
- Molecule 1: THREONINE SYNTHASE 1, CHLOROPLASTIC

Chain D: 78% 10% 12%



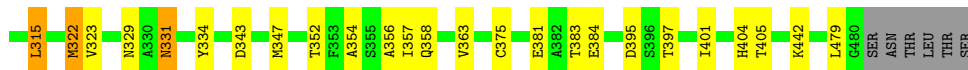
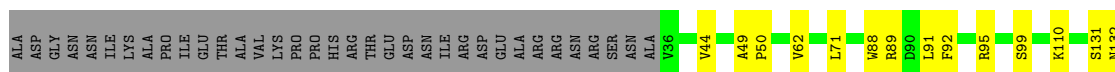
- Molecule 1: THREONINE SYNTHASE 1, CHLOROPLASTIC

Chain E: 78% 10% 11%



- Molecule 1: THREONINE SYNTHASE 1, CHLOROPLASTIC

Chain F: 76% 14% 8%



4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	191.92Å 110.85Å 152.78Å 90.00° 89.98° 90.00°	Depositor
Resolution (Å)	29.89 – 2.60 29.89 – 2.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.89-2.60) 99.6 (29.89-2.60)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 2.61Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.206 , 0.248 0.206 , 0.248	Depositor DCC
R_{free} test set	4904 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	53.8	Xtrriage
Anisotropy	0.086	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 38.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.027 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.027 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.459 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.467 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.027 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	20958	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SAM, PLP, TRS

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	0.44	1/3509 (0.0%)	0.77	0/4763
1	B	0.44	1/3381 (0.0%)	0.76	0/4586
1	C	0.45	1/3502 (0.0%)	0.76	0/4753
1	D	0.43	1/3381 (0.0%)	0.76	0/4586
1	E	0.45	1/3396 (0.0%)	0.76	0/4607
1	F	0.44	1/3509 (0.0%)	0.77	0/4763
All	All	0.44	6/20678 (0.0%)	0.76	0/28058

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	479	LEU	C-N	-6.98	1.23	1.33
1	A	479	LEU	C-N	-6.95	1.23	1.33
1	E	483	THR	C-N	-6.85	1.23	1.33
1	C	479	LEU	C-N	-6.84	1.23	1.33
1	D	481	SER	C-N	-6.68	1.24	1.33
1	B	481	SER	C-N	-6.68	1.24	1.33

There are no bond angle outliers.

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	3428	0	3382	25	0
1	B	3305	0	3257	22	0
1	C	3421	0	3373	27	0
1	D	3305	0	3257	28	0
1	E	3320	0	3270	33	0
1	F	3428	0	3382	41	0
2	A	54	0	44	1	0
2	B	54	0	44	0	0
2	C	54	0	44	2	0
2	D	54	0	44	1	0
2	E	54	0	44	6	0
2	F	54	0	44	8	0
3	A	8	0	12	0	0
3	C	8	0	12	0	0
3	F	8	0	12	0	0
4	A	15	0	6	1	0
4	B	15	0	6	0	0
4	C	15	0	6	1	0
4	D	15	0	6	0	0
4	E	15	0	6	1	0
4	F	15	0	6	1	0
5	A	51	0	0	0	0
5	B	55	0	0	1	0
5	C	52	0	0	0	0
5	D	55	0	0	1	0
5	E	44	0	0	0	0
5	F	56	0	0	0	0
All	All	20958	0	20257	175	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 4.

All (175) 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:367:ARG:HH11	1:B:367:ARG:HG2	1.05	1.16
1:E:367:ARG:HH11	1:E:367:ARG:HG2	1.08	1.09
1:C:291:ILE:HG13	1:C:323:VAL:HB	1.57	0.85
1:B:367:ARG:HG2	1:B:367:ARG:NH1	1.85	0.84
1:E:367:ARG:HG2	1:E:367:ARG:NH1	1.88	0.83
2:F:500:SAM:H5'2	2:F:501:SAM:H2	1.61	0.82
1:F:99:SER:HB2	2:F:500:SAM:HE1	1.62	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:150:TRP:CH2	2:F:500:SAM:HG2	2.14	0.81
1:B:367:ARG:HH11	1:B:367:ARG:CG	1.94	0.78
2:E:500:SAM:HG2	1:F:150:TRP:CH2	2.21	0.75
1:E:99:SER:HB2	2:E:500:SAM:HE1	1.69	0.74
1:F:251:ILE:HG23	1:F:255:LEU:HD12	1.68	0.74
1:E:367:ARG:HH11	1:E:367:ARG:CG	1.96	0.73
1:F:219:ILE:HD11	1:F:223:GLN:HB2	1.70	0.72
1:B:46:PHE:O	1:B:47:ASN:HB2	1.90	0.72
1:F:331:ASN:HD22	1:F:331:ASN:H	1.38	0.72
1:C:331:ASN:HD22	1:C:331:ASN:H	1.37	0.72
1:C:195:THR:HG23	1:C:267:ARG:HH21	1.54	0.71
1:E:331:ASN:HB2	1:E:332:PRO:CD	2.20	0.71
1:D:331:ASN:HB2	1:D:332:PRO:CD	2.22	0.70
1:F:291:ILE:HG13	1:F:323:VAL:HB	1.73	0.70
1:A:331:ASN:H	1:A:331:ASN:HD22	1.38	0.70
1:C:251:ILE:HG23	1:C:255:LEU:HD12	1.75	0.68
1:D:446:HIS:HD2	1:D:459:ASN:H	1.38	0.68
1:D:367:ARG:HG2	1:D:367:ARG:HH11	1.57	0.68
1:E:46:PHE:O	1:E:47:ASN:HB2	1.92	0.68
1:A:251:ILE:HG23	1:A:255:LEU:HD12	1.75	0.68
1:A:131:SER:HB3	1:A:152:LYS:HE2	1.78	0.66
1:A:291:ILE:HG13	1:A:323:VAL:HB	1.78	0.66
1:B:331:ASN:HB2	1:B:332:PRO:CD	2.25	0.66
1:E:331:ASN:HB2	1:E:332:PRO:HD2	1.78	0.66
1:D:331:ASN:HB2	1:D:332:PRO:HD2	1.79	0.63
1:A:170:LEU:HD11	1:A:258:TYR:CE2	2.34	0.62
1:C:170:LEU:HD11	1:C:258:TYR:CE2	2.35	0.61
1:D:367:ARG:HG2	1:D:367:ARG:NH1	2.13	0.61
1:B:331:ASN:HB2	1:B:332:PRO:HD2	1.86	0.58
1:E:443:ILE:HD11	1:F:222:ALA:HA	1.84	0.58
2:C:500:SAM:H8	1:D:132:ASN:HD21	1.71	0.56
1:C:386:GLU:HB3	1:C:410:THR:HG21	1.89	0.55
1:E:187:VAL:HG11	1:E:199:LEU:HD11	1.88	0.55
1:E:133:LEU:HD13	1:E:152:LYS:HD3	1.88	0.55
1:C:296:ASN:O	1:C:363:VAL:O	2.24	0.55
2:E:500:SAM:HG2	1:F:150:TRP:HH2	1.71	0.55
1:B:177:LEU:O	1:B:182:ARG:HB2	2.07	0.54
1:D:187:VAL:HG11	1:D:199:LEU:HD11	1.89	0.54
1:F:226:GLN:HB2	1:F:227:PRO:HD3	1.91	0.53
1:F:290:VAL:HB	1:F:322:MET:HE2	1.91	0.53
1:F:110:LYS:NZ	1:F:280:GLN:OE1	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:367:ARG:HH11	1:D:367:ARG:CG	2.21	0.53
2:E:500:SAM:H5'2	2:E:501:SAM:H2	1.92	0.52
1:F:356:ALA:H	1:F:404:HIS:HE1	1.58	0.52
1:B:187:VAL:HG11	1:B:199:LEU:HD11	1.92	0.52
1:E:135:TRP:CH2	2:F:501:SAM:HG2	2.45	0.51
1:E:163:LYS:NZ	4:E:1163:PLP:O3	2.44	0.51
1:F:296:ASN:O	1:F:363:VAL:O	2.29	0.50
1:D:446:HIS:CD2	1:D:459:ASN:H	2.25	0.50
1:B:133:LEU:HD13	1:B:152:LYS:HD3	1.93	0.50
1:E:150:TRP:HH2	2:F:500:SAM:HG2	1.73	0.50
1:A:163:LYS:NZ	4:A:1163:PLP:O3	2.44	0.49
1:A:214:LEU:HD22	1:A:219:ILE:HD11	1.94	0.49
1:F:354:ALA:O	1:F:357:ILE:HG22	2.11	0.49
1:C:356:ALA:H	1:C:404:HIS:HE1	1.60	0.49
1:A:356:ALA:H	1:A:404:HIS:HE1	1.60	0.49
1:E:132:ASN:ND2	2:F:500:SAM:H1'	2.27	0.49
1:D:244:CYS:O	1:D:248:ILE:HG12	2.13	0.49
1:F:49:ALA:HB1	1:F:50:PRO:HD2	1.96	0.48
1:C:163:LYS:NZ	4:C:1163:PLP:O3	2.45	0.48
1:D:177:LEU:O	1:D:182:ARG:HB2	2.14	0.48
1:C:336:HIS:CD2	1:C:344:PHE:HB3	2.49	0.48
1:C:354:ALA:O	1:C:357:ILE:HG22	2.14	0.48
1:F:91:LEU:O	1:F:95:ARG:HG3	2.14	0.48
1:C:85:GLY:HA2	1:C:315:LEU:HD23	1.96	0.48
1:F:167:MET:HE2	1:F:267:ARG:HE	1.78	0.48
1:A:397:THR:O	1:A:397:THR:HG22	2.13	0.48
1:F:88:TRP:CB	1:F:315:LEU:HD11	2.44	0.48
1:C:88:TRP:HB2	1:C:315:LEU:CD2	2.44	0.48
1:F:331:ASN:HB2	1:F:334:TYR:HB3	1.95	0.47
1:A:128:GLU:H	1:A:128:GLU:CD	2.23	0.47
1:B:159:THR:HG21	1:B:198:ALA:HA	1.97	0.47
1:C:88:TRP:HB2	1:C:315:LEU:HD21	1.97	0.47
1:E:395:ASP:OD2	1:E:442:LYS:NZ	2.46	0.47
1:A:222:ALA:HA	1:B:443:ILE:HD11	1.97	0.47
1:C:331:ASN:HB2	1:C:334:TYR:HB3	1.96	0.47
1:E:205:SER:HA	1:F:397:THR:O	2.15	0.46
2:A:500:SAM:H8	1:B:132:ASN:HD21	1.80	0.46
1:F:163:LYS:NZ	4:F:1163:PLP:O3	2.48	0.46
1:B:46:PHE:O	1:B:47:ASN:CB	2.62	0.46
1:F:170:LEU:HD11	1:F:258:TYR:CE2	2.51	0.46
1:C:91:LEU:O	1:C:95:ARG:HG3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:ASN:HB2	1:A:334:TYR:HB3	1.97	0.46
1:B:128:GLU:H	1:B:128:GLU:CD	2.24	0.45
1:E:247:LEU:O	1:E:251:ILE:HD12	2.15	0.45
1:E:132:ASN:HD21	2:F:500:SAM:H8	1.82	0.45
1:D:220:SER:HB3	1:D:223:GLN:HB2	1.98	0.45
1:F:187:VAL:HG11	1:F:199:LEU:HD11	1.99	0.45
1:F:138:ARG:O	1:F:142:GLN:HB2	2.17	0.45
1:F:401:ILE:HD13	1:F:405:THR:HB	1.99	0.45
1:C:132:ASN:HD21	2:D:500:SAM:H8	1.82	0.45
1:D:247:LEU:O	1:D:251:ILE:HG12	2.17	0.45
1:E:230:ASN:HA	1:F:395:ASP:O	2.16	0.45
1:A:226:GLN:HB2	1:A:227:PRO:HD3	1.98	0.45
1:A:352:THR:HG22	1:A:354:ALA:H	1.81	0.45
1:D:128:GLU:H	1:D:128:GLU:CD	2.25	0.44
1:A:91:LEU:O	1:A:95:ARG:HG3	2.17	0.44
1:A:358:GLN:O	1:A:358:GLN:HG3	2.17	0.44
1:A:397:THR:O	1:A:397:THR:CG2	2.66	0.44
1:F:88:TRP:HB2	1:F:315:LEU:HD11	2.00	0.44
1:B:221:MET:HA	1:B:221:MET:HE2	1.99	0.44
1:C:50:PRO:HG3	1:E:319:ILE:O	2.18	0.44
1:C:214:LEU:HD22	1:C:219:ILE:HD11	2.00	0.44
1:E:188:GLY:HA2	1:E:211:ILE:O	2.18	0.44
1:F:322:MET:HB2	1:F:375:CYS:SG	2.58	0.44
1:F:352:THR:HG22	1:F:354:ALA:H	1.82	0.44
1:E:177:LEU:O	1:E:182:ARG:HB2	2.17	0.44
2:F:500:SAM:HN1	2:F:501:SAM:HE1	1.83	0.44
1:D:395:ASP:OD2	1:D:442:LYS:NZ	2.49	0.43
1:E:297:LEU:HD21	1:E:326:GLN:NE2	2.33	0.43
1:A:137:GLU:HG3	1:A:141:LYS:HE2	1.99	0.43
1:B:391:MET:SD	1:B:442:LYS:NZ	2.86	0.43
1:D:203:CYS:HB3	1:D:208:ILE:O	2.18	0.43
1:D:291:ILE:HG12	1:D:323:VAL:HB	2.01	0.43
1:E:269:GLU:O	1:E:272:LYS:HG2	2.18	0.43
1:A:191:SER:HB3	1:A:195:THR:HB	2.00	0.43
1:E:159:THR:HG21	1:E:198:ALA:HA	2.00	0.43
1:F:261:ASN:HD22	1:F:261:ASN:HA	1.65	0.43
1:B:93:ASP:O	1:B:96:VAL:HG12	2.18	0.43
1:D:391:MET:HE2	1:D:403:PRO:HA	2.01	0.43
1:D:93:ASP:O	1:D:96:VAL:HG12	2.19	0.43
1:C:401:ILE:HD13	1:C:405:THR:HB	2.01	0.43
2:E:500:SAM:HN1	2:E:501:SAM:H5'1	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:500:SAM:HB8	1:F:132:ASN:HD21	1.84	0.42
1:B:296:ASN:ND2	5:B:2035:HOH:O	2.51	0.42
1:D:296:ASN:ND2	5:D:2038:HOH:O	2.52	0.42
1:C:85:GLY:HA3	1:C:314:GLY:O	2.19	0.42
1:A:77:MET:HE1	1:A:309:MET:HE3	2.01	0.42
1:D:395:ASP:HA	1:D:399:MET:O	2.19	0.42
1:F:191:SER:HA	1:F:241:PHE:HE1	1.85	0.42
1:F:138:ARG:HB3	1:F:397:THR:HG23	2.01	0.42
1:F:292:VAL:HB	1:F:300:ILE:HG12	2.01	0.42
1:F:358:GLN:HG3	1:F:358:GLN:O	2.19	0.42
1:D:390:ALA:HA	1:D:393:GLN:HE21	1.85	0.42
1:E:255:LEU:HA	1:E:256:PRO:HD3	1.93	0.42
1:B:146:MET:SD	1:B:416:ARG:HD2	2.60	0.41
1:E:176:ARG:O	1:E:180:MET:HG2	2.19	0.41
1:F:331:ASN:HD22	1:F:331:ASN:N	2.08	0.41
1:A:89:ARG:HD2	1:A:284:TRP:CD1	2.55	0.41
1:C:245:MET:HE1	1:C:261:ASN:HD21	1.85	0.41
1:C:279:LEU:HD13	1:C:307:PHE:HE1	1.85	0.41
1:E:146:MET:SD	1:E:416:ARG:HD2	2.60	0.41
1:C:191:SER:HB3	1:C:195:THR:HB	2.02	0.41
1:D:59:ASP:HA	1:D:177:LEU:HD21	2.01	0.41
2:C:500:SAM:HB8	1:D:132:ASN:ND2	2.35	0.41
1:D:81:LYS:HG3	1:D:313:LEU:HD23	2.01	0.41
1:A:167:MET:HE2	1:A:267:ARG:HE	1.84	0.41
1:B:135:TRP:HB2	1:B:150:TRP:CZ3	2.56	0.41
1:E:47:ASN:HB3	1:E:48:ALA:H	1.58	0.41
1:A:354:ALA:O	1:A:357:ILE:HG22	2.20	0.41
1:C:226:GLN:HB2	1:C:227:PRO:HD3	2.03	0.41
1:D:133:LEU:HD13	1:D:152:LYS:HD3	2.03	0.41
1:D:446:HIS:HD2	1:D:459:ASN:N	2.12	0.41
1:A:401:ILE:HD13	1:A:405:THR:HB	2.03	0.41
1:E:203:CYS:HB3	1:E:208:ILE:O	2.20	0.41
1:F:89:ARG:HD2	1:F:284:TRP:CD1	2.56	0.41
1:C:203:CYS:HB3	1:C:208:ILE:O	2.21	0.41
1:F:329:ASN:HD21	1:F:384:GLU:HG3	1.86	0.41
1:E:135:TRP:HB2	1:E:150:TRP:CZ3	2.55	0.40
1:F:92:PHE:C	1:F:280:GLN:HE22	2.28	0.40
1:F:395:ASP:OD2	1:F:442:LYS:NZ	2.53	0.40
1:B:336:HIS:HD2	1:B:343:ASP:C	2.29	0.40
1:C:89:ARG:HD2	1:C:284:TRP:CD1	2.55	0.40
1:D:146:MET:SD	1:D:416:ARG:HD2	2.61	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:ARG:HH11	1:A:208:ILE:HG12	1.86	0.40
1:B:188:GLY:HA2	1:B:211:ILE:O	2.22	0.40
1:E:46:PHE:O	1:E:47:ASN:CB	2.67	0.40
1:F:224:LEU:O	1:F:228:ILE:HG12	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	443/486 (91%)	424 (96%)	18 (4%)	1 (0%)	43	66
1	B	425/486 (87%)	406 (96%)	17 (4%)	2 (0%)	24	46
1	C	442/486 (91%)	422 (96%)	18 (4%)	2 (0%)	24	46
1	D	425/486 (87%)	399 (94%)	25 (6%)	1 (0%)	43	66
1	E	427/486 (88%)	400 (94%)	24 (6%)	3 (1%)	18	38
1	F	443/486 (91%)	420 (95%)	21 (5%)	2 (0%)	24	46
All	All	2605/2916 (89%)	2471 (95%)	123 (5%)	11 (0%)	30	51

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	47	ASN
1	E	47	ASN
1	A	220	SER
1	C	220	SER
1	F	220	SER
1	E	48	ALA
1	B	36	VAL
1	C	346	PRO

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Mol	Chain	Res	Type
1	E	36	VAL
1	F	343	ASP
1	D	36	VAL

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	368/403 (91%)	354 (96%)	14 (4%)	29	56
1	B	354/403 (88%)	342 (97%)	12 (3%)	32	60
1	C	367/403 (91%)	353 (96%)	14 (4%)	29	56
1	D	354/403 (88%)	344 (97%)	10 (3%)	38	66
1	E	356/403 (88%)	345 (97%)	11 (3%)	35	63
1	F	368/403 (91%)	352 (96%)	16 (4%)	26	51
All	All	2167/2418 (90%)	2090 (96%)	77 (4%)	31	58

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	VAL
1	A	62	VAL
1	A	71	LEU
1	A	131	SER
1	A	144	LEU
1	A	218	LYS
1	A	224	LEU
1	A	239	THR
1	A	266	LEU
1	A	279	LEU
1	A	300	ILE
1	A	315	LEU
1	A	331	ASN
1	A	479	LEU
1	B	44	VAL

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Mol	Chain	Res	Type
1	B	62	VAL
1	B	71	LEU
1	B	239	THR
1	B	247	LEU
1	B	279	LEU
1	B	315	LEU
1	B	365	ILE
1	B	367	ARG
1	B	384	GLU
1	B	464	VAL
1	B	465	LYS
1	C	44	VAL
1	C	62	VAL
1	C	71	LEU
1	C	237	ILE
1	C	239	THR
1	C	245	MET
1	C	266	LEU
1	C	267	ARG
1	C	279	LEU
1	C	300	ILE
1	C	312	GLU
1	C	322	MET
1	C	331	ASN
1	C	342	LYS
1	D	62	VAL
1	D	71	LEU
1	D	78	GLU
1	D	181	LYS
1	D	239	THR
1	D	279	LEU
1	D	315	LEU
1	D	365	ILE
1	D	367	ARG
1	D	464	VAL
1	E	44	VAL
1	E	62	VAL
1	E	71	LEU
1	E	239	THR
1	E	247	LEU
1	E	279	LEU
1	E	300	ILE

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Mol	Chain	Res	Type
1	E	315	LEU
1	E	365	ILE
1	E	367	ARG
1	E	384	GLU
1	F	44	VAL
1	F	62	VAL
1	F	71	LEU
1	F	131	SER
1	F	144	LEU
1	F	245	MET
1	F	252	THR
1	F	266	LEU
1	F	279	LEU
1	F	300	ILE
1	F	315	LEU
1	F	322	MET
1	F	331	ASN
1	F	347	MET
1	F	381	GLU
1	F	383	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (60) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	132	ASN
1	A	226	GLN
1	A	230	ASN
1	A	261	ASN
1	A	280	GLN
1	A	299	ASN
1	A	326	GLN
1	A	331	ASN
1	A	376	ASN
1	A	404	HIS
1	A	417	ASN
1	B	132	ASN
1	B	230	ASN
1	B	264	ASN
1	B	326	GLN
1	B	329	ASN
1	B	336	HIS
1	B	376	ASN

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Mol	Chain	Res	Type
1	B	393	GLN
1	B	448	ASN
1	C	37	ASN
1	C	132	ASN
1	C	226	GLN
1	C	230	ASN
1	C	261	ASN
1	C	271	GLN
1	C	280	GLN
1	C	299	ASN
1	C	326	GLN
1	C	331	ASN
1	C	376	ASN
1	C	417	ASN
1	C	448	ASN
1	C	459	ASN
1	D	132	ASN
1	D	230	ASN
1	D	264	ASN
1	D	311	GLN
1	D	326	GLN
1	D	329	ASN
1	D	393	GLN
1	D	404	HIS
1	D	434	HIS
1	D	446	HIS
1	E	132	ASN
1	E	226	GLN
1	E	230	ASN
1	E	329	ASN
1	E	331	ASN
1	E	404	HIS
1	F	132	ASN
1	F	230	ASN
1	F	261	ASN
1	F	299	ASN
1	F	326	GLN
1	F	329	ASN
1	F	331	ASN
1	F	404	HIS
1	F	417	ASN
1	F	448	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](#)

21 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
4	PLP	A	1163	1	15,15,16	1.87	3 (20%)	21,22,23	1.62	3 (14%)
4	PLP	C	1163	1	15,15,16	1.87	3 (20%)	21,22,23	1.69	3 (14%)
2	SAM	B	500	-	27,29,29	1.13	4 (14%)	34,42,42	2.09	8 (23%)
2	SAM	C	501	-	27,29,29	1.12	3 (11%)	34,42,42	1.96	8 (23%)
4	PLP	B	1163	1	15,15,16	1.90	3 (20%)	21,22,23	1.49	2 (9%)
3	TRS	C	800	-	7,7,7	0.31	0	9,9,9	0.36	0
4	PLP	F	1163	1	15,15,16	1.85	3 (20%)	21,22,23	1.72	4 (19%)
2	SAM	C	500	-	27,29,29	1.15	4 (14%)	34,42,42	2.02	9 (26%)
2	SAM	F	501	-	27,29,29	1.13	5 (18%)	34,42,42	1.96	8 (23%)
2	SAM	B	501	-	27,29,29	1.16	5 (18%)	34,42,42	2.01	9 (26%)
2	SAM	A	500	-	27,29,29	1.14	4 (14%)	34,42,42	2.02	10 (29%)
4	PLP	E	1163	1	15,15,16	1.90	3 (20%)	21,22,23	1.70	4 (19%)
2	SAM	E	500	-	27,29,29	1.11	4 (14%)	34,42,42	2.12	9 (26%)
2	SAM	A	501	-	27,29,29	1.12	3 (11%)	34,42,42	1.98	8 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SAM	D	500	-	27,29,29	1.13	4 (14%)	34,42,42	2.11	8 (23%)
2	SAM	F	500	-	27,29,29	1.13	3 (11%)	34,42,42	2.09	8 (23%)
4	PLP	D	1163	1	15,15,16	1.90	3 (20%)	21,22,23	1.58	2 (9%)
3	TRS	F	800	-	7,7,7	0.29	0	9,9,9	0.33	0
3	TRS	A	800	-	7,7,7	0.31	0	9,9,9	0.19	0
2	SAM	E	501	-	27,29,29	1.17	5 (18%)	34,42,42	1.97	9 (26%)
2	SAM	D	501	-	27,29,29	1.18	5 (18%)	34,42,42	1.98	9 (26%)

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	PLP	A	1163	1	-	0/6/6/8	0/1/1/1
4	PLP	C	1163	1	-	1/6/6/8	0/1/1/1
2	SAM	B	500	-	-	1/17/33/33	0/3/3/3
2	SAM	C	501	-	-	3/17/33/33	0/3/3/3
4	PLP	B	1163	1	-	3/6/6/8	0/1/1/1
3	TRS	C	800	-	-	1/9/9/9	-
4	PLP	F	1163	1	-	0/6/6/8	0/1/1/1
2	SAM	C	500	-	-	1/17/33/33	0/3/3/3
2	SAM	F	501	-	-	8/17/33/33	0/3/3/3
2	SAM	B	501	-	-	6/17/33/33	0/3/3/3
2	SAM	A	500	-	-	1/17/33/33	0/3/3/3
4	PLP	E	1163	1	-	0/6/6/8	0/1/1/1
2	SAM	E	500	-	-	8/17/33/33	0/3/3/3
2	SAM	A	501	-	-	8/17/33/33	0/3/3/3
2	SAM	D	500	-	-	5/17/33/33	0/3/3/3
2	SAM	F	500	-	-	6/17/33/33	0/3/3/3
4	PLP	D	1163	1	-	0/6/6/8	0/1/1/1
3	TRS	F	800	-	-	3/9/9/9	-
3	TRS	A	800	-	-	0/9/9/9	-
2	SAM	E	501	-	-	7/17/33/33	0/3/3/3
2	SAM	D	501	-	-	5/17/33/33	0/3/3/3

All (67) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1163	PLP	O3-C3	-5.84	1.23	1.36
4	E	1163	PLP	O3-C3	-5.81	1.23	1.36
4	C	1163	PLP	O3-C3	-5.79	1.23	1.36
4	B	1163	PLP	O3-C3	-5.77	1.23	1.36
4	D	1163	PLP	O3-C3	-5.75	1.23	1.36
4	F	1163	PLP	O3-C3	-5.70	1.23	1.36
2	C	501	SAM	C2-N1	2.87	1.39	1.33
2	A	500	SAM	C2-N3	2.86	1.39	1.33
2	C	500	SAM	C2-N3	2.86	1.39	1.33
2	E	501	SAM	C2-N1	2.81	1.38	1.33
2	A	501	SAM	C2-N1	2.79	1.38	1.33
2	D	501	SAM	C2-N1	2.78	1.38	1.33
2	B	501	SAM	C2-N3	2.77	1.38	1.33
2	D	501	SAM	C2-N3	2.77	1.38	1.33
2	D	500	SAM	C2-N3	2.76	1.38	1.33
2	F	500	SAM	C2-N3	2.76	1.38	1.33
2	B	501	SAM	C2-N1	2.74	1.38	1.33
4	D	1163	PLP	C2-N1	2.74	1.38	1.33
2	E	501	SAM	C2-N3	2.72	1.38	1.33
2	B	500	SAM	C2-N3	2.71	1.38	1.33
4	E	1163	PLP	C2-N1	2.70	1.38	1.33
4	B	1163	PLP	C2-N1	2.68	1.38	1.33
2	F	501	SAM	C2-N1	2.65	1.38	1.33
2	A	501	SAM	C2-N3	2.65	1.38	1.33
2	C	501	SAM	C2-N3	2.64	1.38	1.33
2	B	500	SAM	C2-N1	2.63	1.38	1.33
2	A	500	SAM	C2-N1	2.60	1.38	1.33
2	C	500	SAM	C2-N1	2.60	1.38	1.33
2	E	500	SAM	C2-N3	2.60	1.38	1.33
2	E	500	SAM	C2-N1	2.60	1.38	1.33
2	F	501	SAM	C2-N3	2.60	1.38	1.33
2	F	500	SAM	C2-N1	2.58	1.38	1.33
4	C	1163	PLP	C2-N1	2.53	1.38	1.33
2	D	500	SAM	C2-N1	2.52	1.38	1.33
4	F	1163	PLP	C2-N1	2.46	1.38	1.33
4	A	1163	PLP	C2-N1	2.45	1.38	1.33
2	F	500	SAM	OXT-C	-2.29	1.23	1.30
2	A	500	SAM	OXT-C	-2.28	1.23	1.30
2	C	500	SAM	OXT-C	-2.26	1.23	1.30
2	A	501	SAM	OXT-C	-2.22	1.23	1.30
2	C	500	SAM	C5-N7	-2.21	1.35	1.39
2	D	501	SAM	C8-N7	2.21	1.35	1.31
2	E	501	SAM	OXT-C	-2.21	1.23	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	SAM	OXT-C	-2.21	1.23	1.30
2	C	501	SAM	OXT-C	-2.20	1.23	1.30
2	D	501	SAM	OXT-C	-2.19	1.23	1.30
2	F	501	SAM	OXT-C	-2.19	1.23	1.30
4	D	1163	PLP	C6-N1	2.19	1.38	1.34
2	D	500	SAM	C8-N7	2.18	1.35	1.31
2	D	500	SAM	OXT-C	-2.17	1.23	1.30
2	B	500	SAM	OXT-C	-2.17	1.23	1.30
2	E	501	SAM	C5-N7	-2.16	1.35	1.39
4	B	1163	PLP	C6-N1	2.14	1.38	1.34
2	E	501	SAM	C8-N7	2.14	1.35	1.31
2	E	500	SAM	OXT-C	-2.14	1.23	1.30
4	E	1163	PLP	C6-N1	2.14	1.38	1.34
2	A	500	SAM	C5-N7	-2.13	1.35	1.39
2	F	501	SAM	C8-N7	2.12	1.35	1.31
4	F	1163	PLP	C6-N1	2.11	1.38	1.34
2	B	501	SAM	C8-N7	2.09	1.35	1.31
2	B	501	SAM	C5-N7	-2.08	1.35	1.39
4	C	1163	PLP	C6-N1	2.07	1.38	1.34
2	E	500	SAM	C8-N7	2.06	1.35	1.31
2	D	501	SAM	C5-N7	-2.05	1.35	1.39
2	B	500	SAM	C8-N7	2.04	1.35	1.31
4	A	1163	PLP	C6-N1	2.04	1.38	1.34
2	F	501	SAM	C5-N7	-2.00	1.35	1.39

All (121) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1163	PLP	O4P-C5A-C5	5.98	120.57	109.36
2	D	500	SAM	N3-C2-N1	-5.97	119.54	128.58
2	E	500	SAM	N3-C2-N1	-5.97	119.54	128.58
2	F	500	SAM	N3-C2-N1	-5.89	119.67	128.58
2	C	500	SAM	N3-C2-N1	-5.84	119.74	128.58
2	B	500	SAM	N3-C2-N1	-5.81	119.79	128.58
2	A	500	SAM	N3-C2-N1	-5.79	119.81	128.58
2	A	501	SAM	N3-C2-N1	-5.67	120.00	128.58
2	E	501	SAM	N3-C2-N1	-5.65	120.03	128.58
2	B	501	SAM	N3-C2-N1	-5.62	120.08	128.58
2	D	501	SAM	N3-C2-N1	-5.61	120.09	128.58
2	F	501	SAM	N3-C2-N1	-5.54	120.19	128.58
2	C	501	SAM	N3-C2-N1	-5.51	120.25	128.58
4	F	1163	PLP	O4P-C5A-C5	5.46	119.59	109.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1163	PLP	O4P-C5A-C5	5.27	119.23	109.36
4	E	1163	PLP	O4P-C5A-C5	5.20	119.10	109.36
2	E	500	SAM	C5-C4-N3	-5.11	119.67	126.72
2	D	500	SAM	C5-C4-N3	-5.06	119.75	126.72
2	B	500	SAM	C5-C4-N3	-5.05	119.77	126.72
2	F	500	SAM	C5-C4-N3	-5.01	119.81	126.72
2	B	501	SAM	C5-C4-N3	-4.86	120.02	126.72
2	D	501	SAM	C5-C4-N3	-4.83	120.07	126.72
2	A	500	SAM	C5-C4-N3	-4.82	120.08	126.72
2	C	500	SAM	C5-C4-N3	-4.73	120.20	126.72
4	D	1163	PLP	O4P-C5A-C5	4.73	118.22	109.36
2	C	501	SAM	C5-C4-N3	-4.73	120.21	126.72
2	E	501	SAM	C5-C4-N3	-4.72	120.21	126.72
2	A	501	SAM	C5-C4-N3	-4.69	120.26	126.72
2	F	501	SAM	C5-C4-N3	-4.67	120.28	126.72
4	B	1163	PLP	O4P-C5A-C5	4.45	117.70	109.36
2	E	500	SAM	C2-N3-C4	3.91	121.38	111.83
2	D	500	SAM	C2-N3-C4	3.89	121.32	111.83
2	F	500	SAM	C2-N3-C4	3.77	121.03	111.83
2	D	500	SAM	N9-C8-N7	-3.76	108.61	113.94
2	B	500	SAM	C2-N3-C4	3.74	120.96	111.83
2	B	500	SAM	N9-C8-N7	-3.68	108.72	113.94
2	A	501	SAM	N9-C8-N7	-3.63	108.78	113.94
2	A	500	SAM	C2-N3-C4	3.62	120.67	111.83
2	F	501	SAM	N9-C8-N7	-3.62	108.80	113.94
2	C	500	SAM	C2-N3-C4	3.61	120.64	111.83
2	D	501	SAM	C2-N3-C4	3.55	120.51	111.83
2	A	501	SAM	C2-N3-C4	3.55	120.51	111.83
2	B	501	SAM	C2-N3-C4	3.54	120.47	111.83
2	B	500	SAM	C5-N7-C8	3.53	109.00	103.45
2	C	501	SAM	N9-C8-N7	-3.50	108.98	113.94
2	E	501	SAM	C2-N3-C4	3.49	120.36	111.83
2	F	500	SAM	N9-C8-N7	-3.48	108.99	113.94
2	C	501	SAM	C2-N3-C4	3.48	120.33	111.83
2	D	500	SAM	C5-N7-C8	3.48	108.92	103.45
2	F	501	SAM	C2-N3-C4	3.47	120.32	111.83
2	E	500	SAM	N9-C8-N7	-3.47	109.01	113.94
2	F	500	SAM	C5-N7-C8	3.45	108.86	103.45
2	B	500	SAM	N3-C4-N9	3.37	132.91	127.17
2	E	500	SAM	C5-N7-C8	3.37	108.75	103.45
2	D	500	SAM	N3-C4-N9	3.36	132.89	127.17
2	F	501	SAM	C5-N7-C8	3.36	108.73	103.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	500	SAM	N3-C4-N9	3.35	132.87	127.17
2	C	500	SAM	N9-C8-N7	-3.34	109.19	113.94
2	F	500	SAM	N3-C4-N9	3.34	132.84	127.17
2	E	501	SAM	N9-C8-N7	-3.33	109.21	113.94
2	A	501	SAM	C5-N7-C8	3.33	108.68	103.45
2	B	501	SAM	N9-C8-N7	-3.30	109.25	113.94
2	C	501	SAM	C5-N7-C8	3.30	108.64	103.45
2	A	500	SAM	N9-C8-N7	-3.28	109.28	113.94
2	B	501	SAM	C5-N7-C8	3.27	108.58	103.45
2	E	501	SAM	C5-N7-C8	3.25	108.55	103.45
2	D	501	SAM	N9-C8-N7	-3.23	109.35	113.94
2	C	500	SAM	C5-N7-C8	3.22	108.52	103.45
2	A	500	SAM	N3-C4-N9	3.22	132.64	127.17
2	D	501	SAM	C5-N7-C8	3.20	108.48	103.45
2	A	501	SAM	N3-C4-N9	3.18	132.58	127.17
2	C	500	SAM	N3-C4-N9	3.17	132.56	127.17
2	A	500	SAM	C5-N7-C8	3.17	108.43	103.45
2	C	501	SAM	N3-C4-N9	3.14	132.51	127.17
2	B	501	SAM	N3-C4-N9	3.07	132.39	127.17
2	F	501	SAM	N3-C4-N9	3.02	132.30	127.17
2	D	501	SAM	N3-C4-N9	3.02	132.30	127.17
2	E	501	SAM	N3-C4-N9	2.99	132.26	127.17
2	E	500	SAM	OXT-C-O	-2.97	117.33	124.08
4	E	1163	PLP	C4A-C4-C5	-2.82	118.03	120.94
2	D	500	SAM	OXT-C-O	-2.80	117.72	124.08
2	A	500	SAM	OXT-C-O	-2.79	117.74	124.08
2	B	500	SAM	OXT-C-O	-2.79	117.75	124.08
2	B	501	SAM	OXT-C-O	-2.76	117.81	124.08
2	C	500	SAM	OXT-C-O	-2.76	117.83	124.08
2	D	501	SAM	OXT-C-O	-2.73	117.88	124.08
4	F	1163	PLP	C6-C5-C4	2.69	120.31	118.10
2	F	501	SAM	OXT-C-O	-2.68	118.01	124.08
2	A	501	SAM	OXT-C-O	-2.65	118.06	124.08
2	C	501	SAM	OXT-C-O	-2.60	118.18	124.08
2	E	501	SAM	OXT-C-O	-2.60	118.19	124.08
2	F	500	SAM	C4-C5-N7	-2.59	107.62	110.58
2	F	500	SAM	OXT-C-O	-2.55	118.29	124.08
4	A	1163	PLP	C6-C5-C4	2.48	120.13	118.10
2	B	500	SAM	C4-C5-N7	-2.47	107.76	110.58
2	C	500	SAM	O4 ² -C1 ¹ -N9	2.43	112.75	108.09
4	B	1163	PLP	C5-C6-N1	-2.41	119.90	123.83
2	A	500	SAM	O4 ² -C1 ¹ -N9	2.41	112.72	108.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	SAM	C4-C5-N7	-2.39	107.84	110.58
2	E	500	SAM	C4-C5-N7	-2.39	107.85	110.58
2	D	500	SAM	C4-C5-N7	-2.39	107.85	110.58
2	F	501	SAM	C4-C5-N7	-2.37	107.87	110.58
2	E	500	SAM	O4'-C1'-N9	2.37	112.64	108.09
2	D	501	SAM	C4-C5-N7	-2.36	107.88	110.58
4	D	1163	PLP	C5-C6-N1	-2.36	119.99	123.83
2	C	500	SAM	C4-C5-N7	-2.35	107.89	110.58
2	C	501	SAM	C4-C5-N7	-2.32	107.93	110.58
4	F	1163	PLP	C5-C6-N1	-2.31	120.08	123.83
2	E	501	SAM	C4-C5-N7	-2.30	107.95	110.58
2	A	500	SAM	C4-C5-N7	-2.27	107.98	110.58
4	F	1163	PLP	C4A-C4-C5	-2.26	118.61	120.94
2	E	501	SAM	C3'-C2'-C1'	2.24	105.71	101.46
2	A	501	SAM	C4-C5-N7	-2.24	108.02	110.58
4	E	1163	PLP	C5-C6-N1	-2.23	120.21	123.83
4	A	1163	PLP	C5-C6-N1	-2.21	120.23	123.83
4	C	1163	PLP	C6-C5-C4	2.20	119.90	118.10
2	B	501	SAM	C3'-C2'-C1'	2.11	105.46	101.46
4	E	1163	PLP	C6-C5-C4	2.11	119.83	118.10
2	A	500	SAM	C2'-C1'-N9	-2.10	108.08	113.30
4	C	1163	PLP	C5-C6-N1	-2.08	120.45	123.83
2	D	501	SAM	C3'-C2'-C1'	2.02	105.28	101.46

There are no chirality outliers.

All (67) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	SAM	N-CA-CB-CG
2	A	501	SAM	C-CA-CB-CG
2	A	501	SAM	C4'-C5'-SD-CE
2	B	500	SAM	CA-CB-CG-SD
2	B	501	SAM	CB-CG-SD-C5'
2	C	500	SAM	C4'-C5'-SD-CE
2	D	500	SAM	O-C-CA-N
2	D	500	SAM	CA-CB-CG-SD
2	D	501	SAM	CB-CG-SD-C5'
2	E	500	SAM	C4'-C5'-SD-CE
2	E	500	SAM	O4'-C4'-C5'-SD
2	E	500	SAM	C3'-C4'-C5'-SD
2	E	501	SAM	O-C-CA-N
2	E	501	SAM	C4'-C5'-SD-CE

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Mol	Chain	Res	Type	Atoms
2	F	500	SAM	C4'-C5'-SD-CE
2	F	500	SAM	O4'-C4'-C5'-SD
2	F	500	SAM	C3'-C4'-C5'-SD
2	F	501	SAM	N-CA-CB-CG
2	F	501	SAM	C-CA-CB-CG
2	F	501	SAM	C4'-C5'-SD-CE
4	B	1163	PLP	C5A-O4P-P-O2P
4	B	1163	PLP	C5A-O4P-P-O3P
2	D	500	SAM	OXT-C-CA-N
2	E	500	SAM	OXT-C-CA-N
2	F	501	SAM	OXT-C-CA-N
2	E	501	SAM	OXT-C-CA-N
2	D	501	SAM	CB-CG-SD-CE
2	D	501	SAM	C2'-C1'-N9-C8
2	F	500	SAM	OXT-C-CA-N
2	C	501	SAM	C2'-C1'-N9-C8
2	E	501	SAM	C2'-C1'-N9-C8
4	B	1163	PLP	C5A-O4P-P-O1P
2	A	501	SAM	C2'-C1'-N9-C8
2	B	501	SAM	C2'-C1'-N9-C8
2	F	501	SAM	C2'-C1'-N9-C8
2	B	501	SAM	CB-CG-SD-CE
3	F	800	TRS	C1-C-C3-O3
3	F	800	TRS	C2-C-C3-O3
2	D	501	SAM	C2'-C1'-N9-C4
2	E	500	SAM	O-C-CA-N
2	E	500	SAM	O-C-CA-CB
2	E	500	SAM	OXT-C-CA-CB
2	E	500	SAM	CA-CB-CG-SD
2	E	501	SAM	CA-CB-CG-SD
2	F	500	SAM	O-C-CA-CB
2	C	501	SAM	C2'-C1'-N9-C4
2	E	501	SAM	C2'-C1'-N9-C4
4	C	1163	PLP	C5A-O4P-P-O1P
2	A	501	SAM	C2'-C1'-N9-C4
2	B	501	SAM	C2'-C1'-N9-C4
2	F	501	SAM	C2'-C1'-N9-C4
2	F	500	SAM	OXT-C-CA-CB
3	C	800	TRS	C1-C-C3-O3
3	F	800	TRS	N-C-C3-O3
2	A	501	SAM	OXT-C-CA-N
2	A	501	SAM	O4'-C1'-N9-C8

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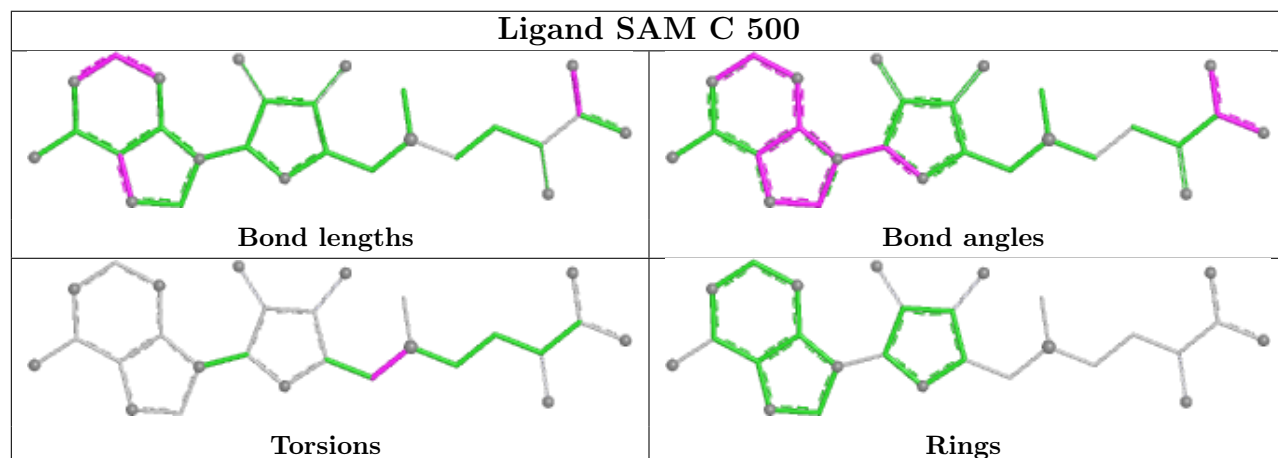
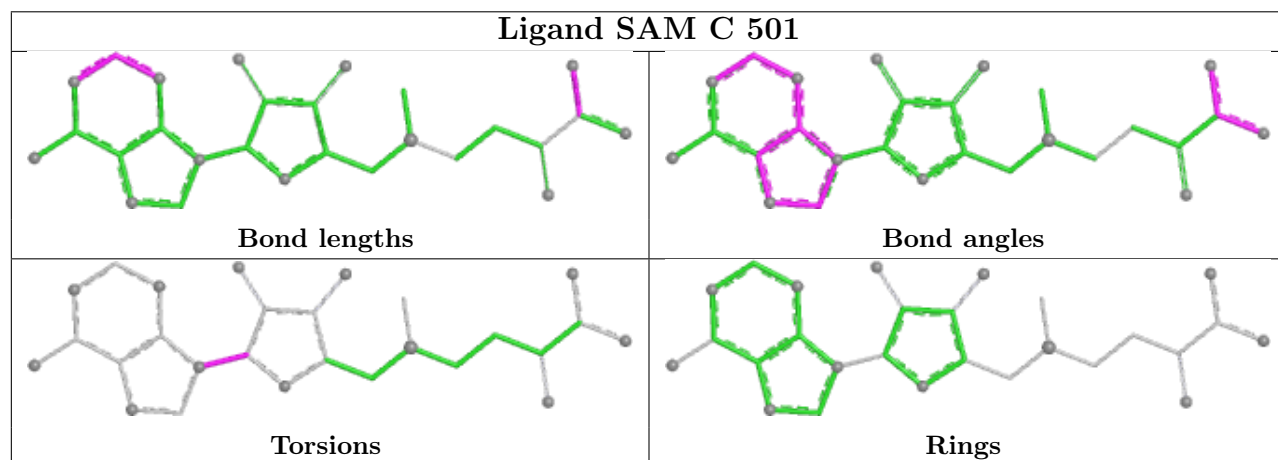
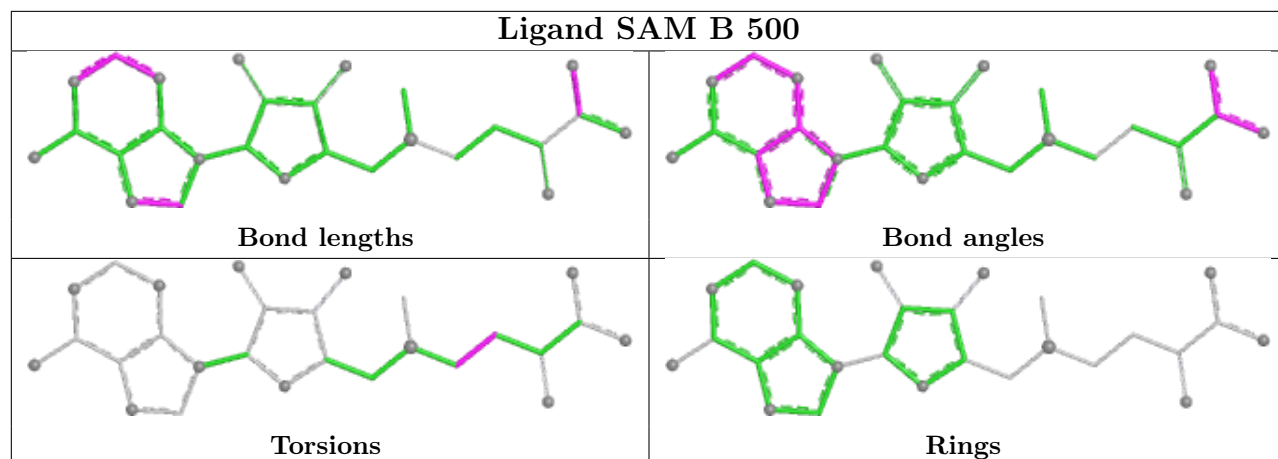
Mol	Chain	Res	Type	Atoms
2	B	501	SAM	O4'-C1'-N9-C8
2	C	501	SAM	O4'-C1'-N9-C8
2	E	501	SAM	O4'-C1'-N9-C8
2	F	501	SAM	O4'-C1'-N9-C8
2	A	500	SAM	CA-CB-CG-SD
2	B	501	SAM	OXT-C-CA-N
2	D	500	SAM	CB-CG-SD-C5'
2	D	501	SAM	O4'-C1'-N9-C8
2	A	501	SAM	O-C-CA-N
2	F	501	SAM	O-C-CA-N
2	D	500	SAM	C-CA-CB-CG

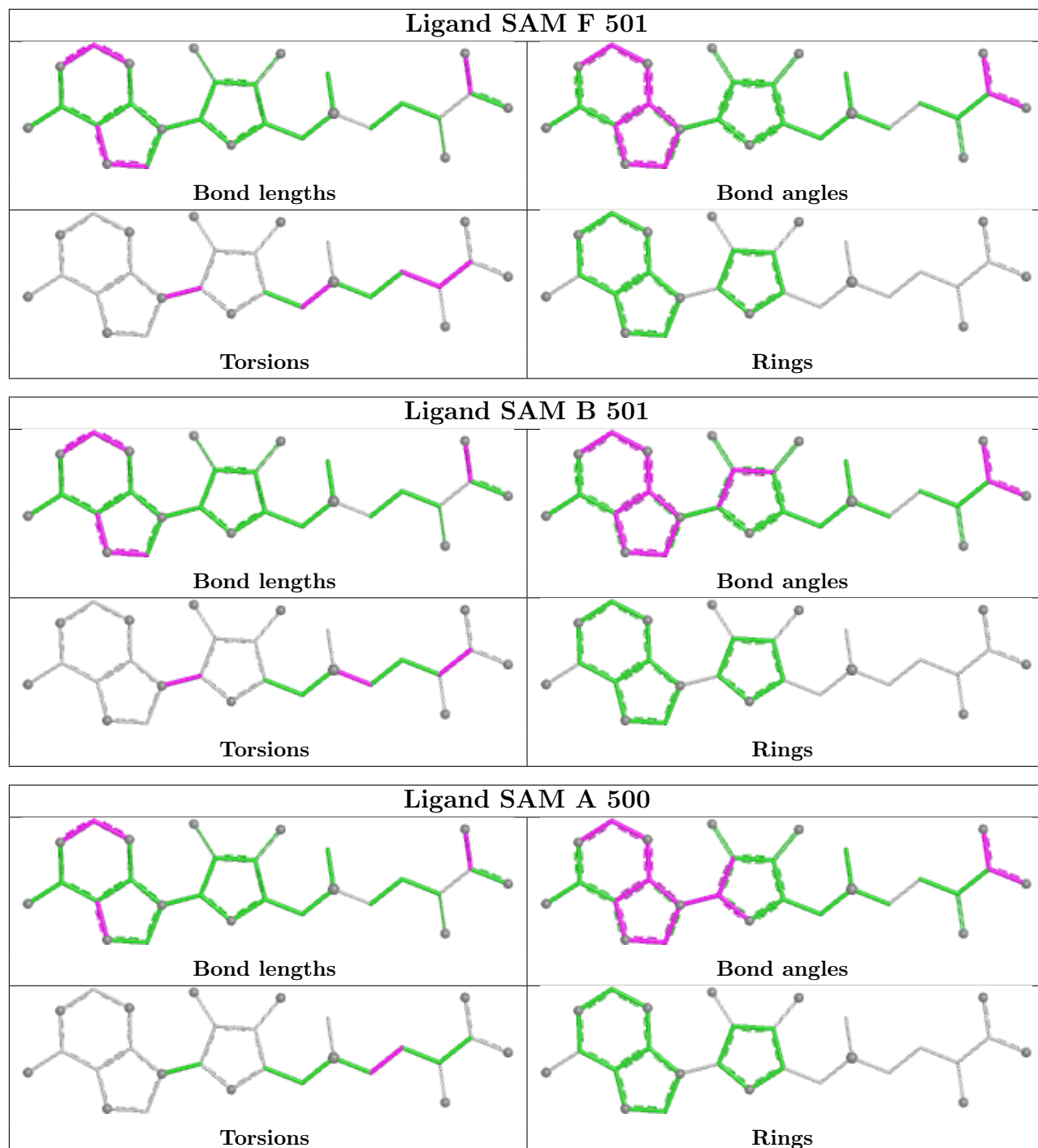
There are no ring outliers.

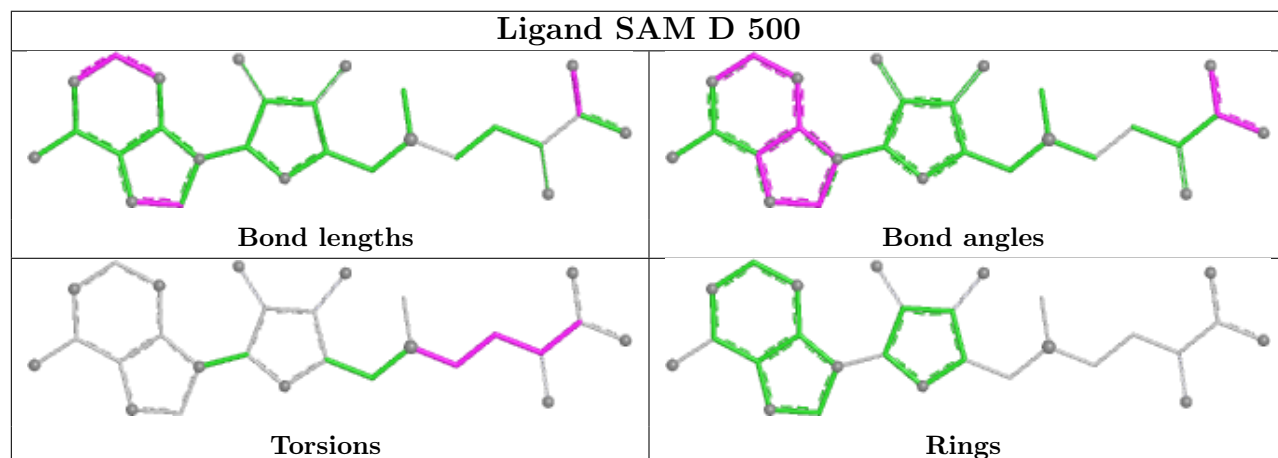
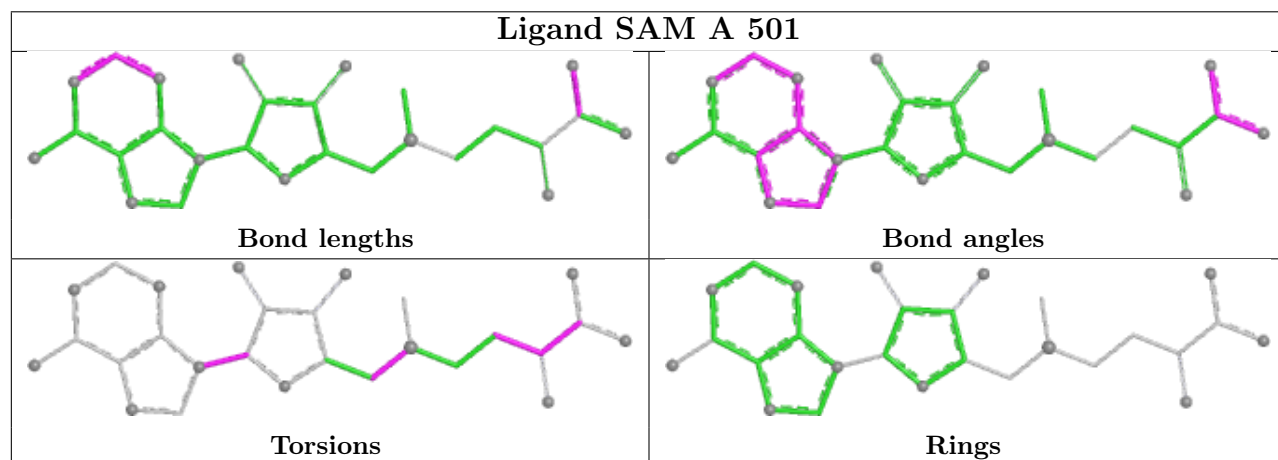
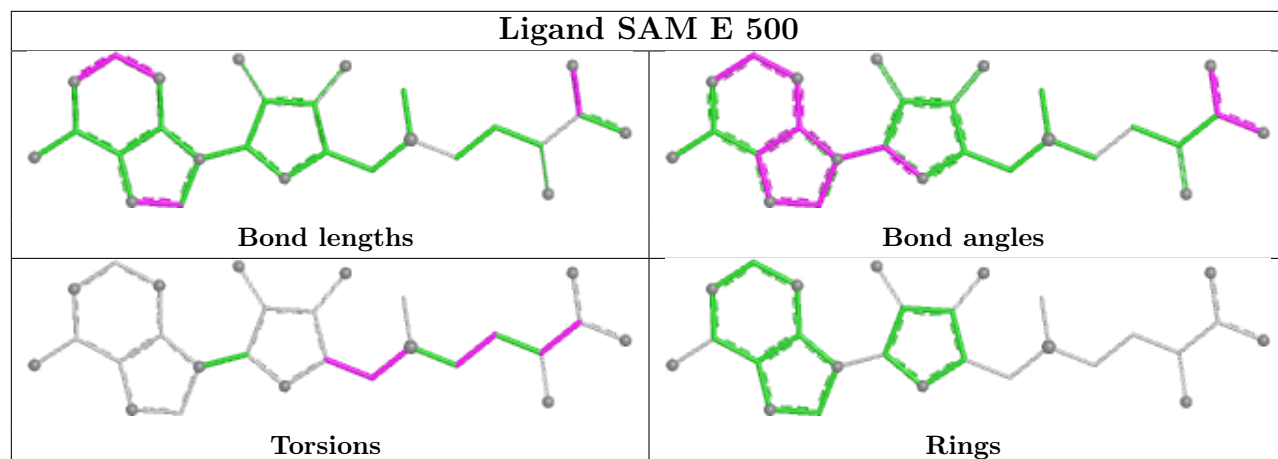
11 monomers are involved in 22 short contacts:

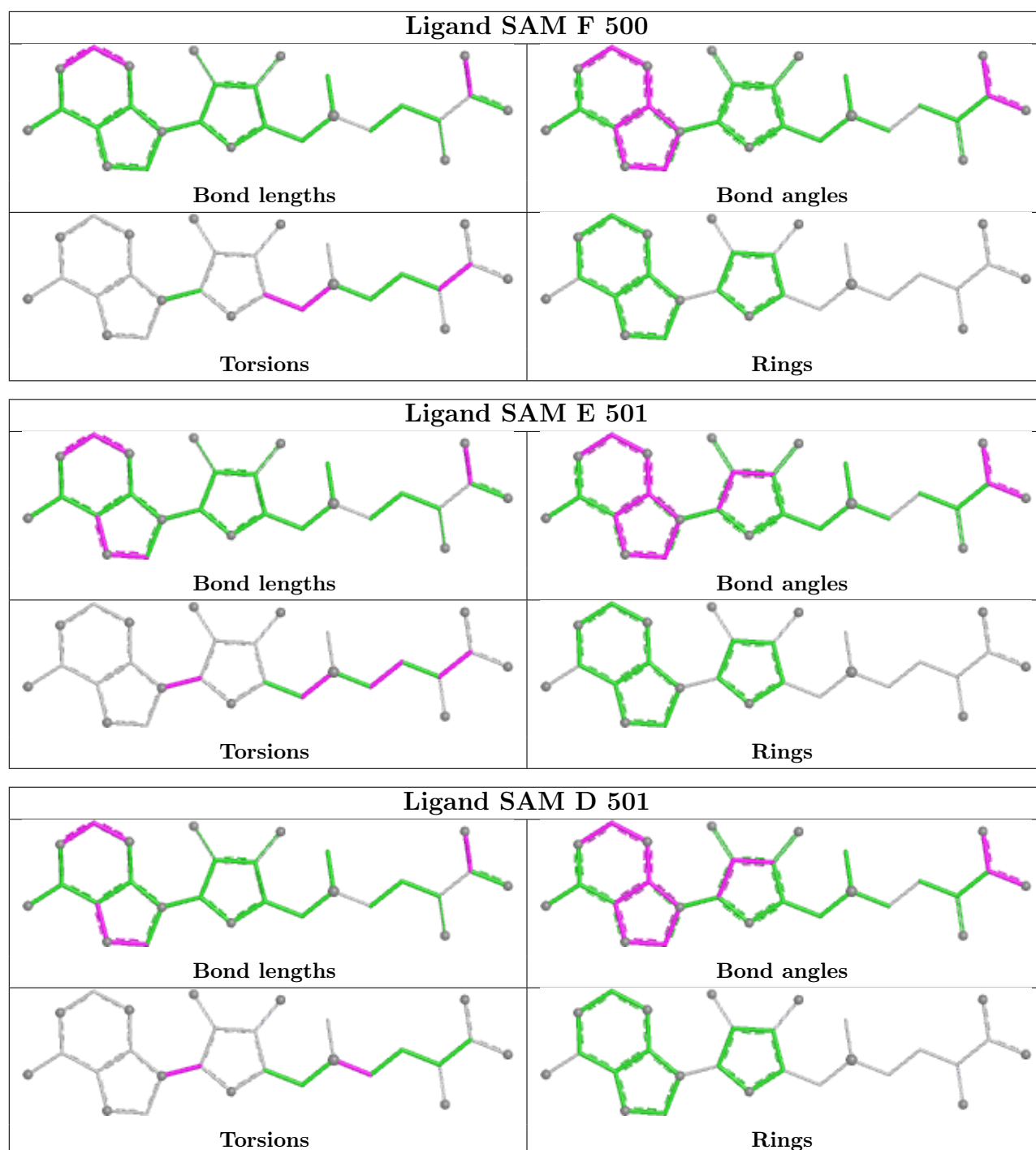
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1163	PLP	1	0
4	C	1163	PLP	1	0
4	F	1163	PLP	1	0
2	C	500	SAM	2	0
2	F	501	SAM	3	0
2	A	500	SAM	1	0
4	E	1163	PLP	1	0
2	E	500	SAM	6	0
2	D	500	SAM	1	0
2	F	500	SAM	7	0
2	E	501	SAM	2	0

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.









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	445/486 (91%)	-1.03	0 100 100	48, 58, 80, 94	0
1	B	429/486 (88%)	-1.15	0 100 100	51, 59, 77, 89	0
1	C	444/486 (91%)	-1.05	0 100 100	48, 58, 79, 95	0
1	D	429/486 (88%)	-1.15	0 100 100	50, 59, 77, 89	0
1	E	431/486 (88%)	-1.16	0 100 100	51, 59, 77, 91	0
1	F	445/486 (91%)	-1.04	0 100 100	48, 58, 81, 94	0
All	All	2623/2916 (89%)	-1.10	0 100 100	48, 59, 79, 95	0

There are no RSRZ outliers to report.

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	TRS	A	800	8/8	0.96	0.07	96,96,96,96	0

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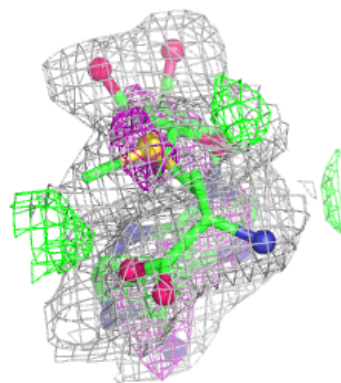
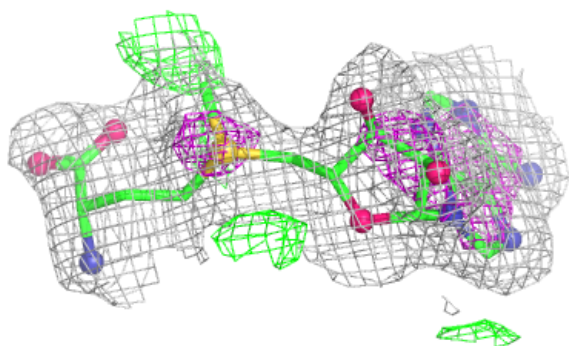
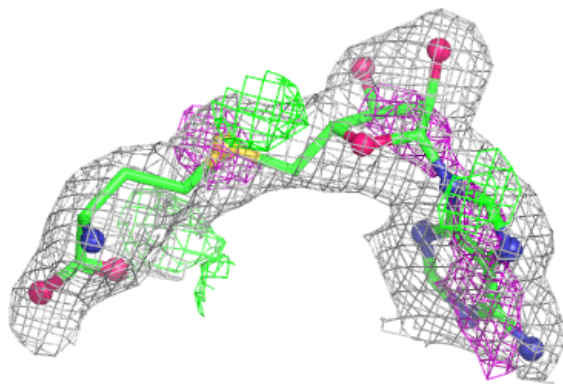
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	TRS	F	800	8/8	0.96	0.06	96,96,96,97	0
2	SAM	E	500	27/27	0.97	0.07	46,52,77,78	0
2	SAM	B	500	27/27	0.97	0.07	47,55,75,76	0
3	TRS	C	800	8/8	0.97	0.05	96,96,96,97	0
2	SAM	D	500	27/27	0.97	0.07	42,50,83,83	0
4	PLP	B	1163	15/16	0.97	0.07	55,57,59,59	15
4	PLP	E	1163	15/16	0.97	0.06	56,58,60,60	15
4	PLP	F	1163	15/16	0.97	0.06	48,54,56,57	0
2	SAM	E	501	27/27	0.98	0.06	52,63,81,81	0
2	SAM	F	500	27/27	0.98	0.06	36,43,64,65	0
2	SAM	F	501	27/27	0.98	0.05	40,55,68,69	0
2	SAM	B	501	27/27	0.98	0.06	62,70,81,82	0
2	SAM	C	500	27/27	0.98	0.06	31,41,57,58	0
2	SAM	C	501	27/27	0.98	0.05	49,58,65,65	0
4	PLP	A	1163	15/16	0.98	0.05	51,53,57,58	0
2	SAM	A	501	27/27	0.98	0.05	45,56,68,68	0
4	PLP	C	1163	15/16	0.98	0.06	49,52,53,55	0
4	PLP	D	1163	15/16	0.98	0.05	57,58,60,60	15
2	SAM	D	501	27/27	0.98	0.06	53,65,82,82	0
2	SAM	A	500	27/27	0.98	0.06	32,43,60,62	0

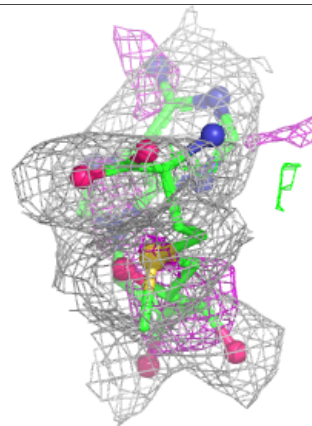
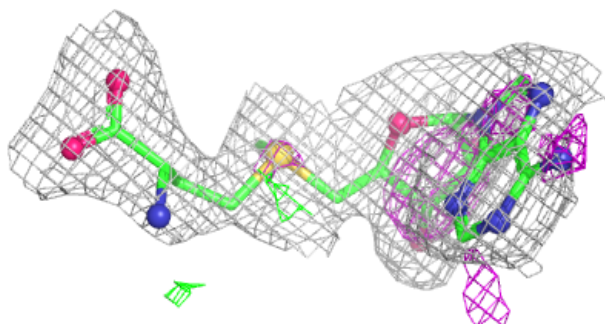
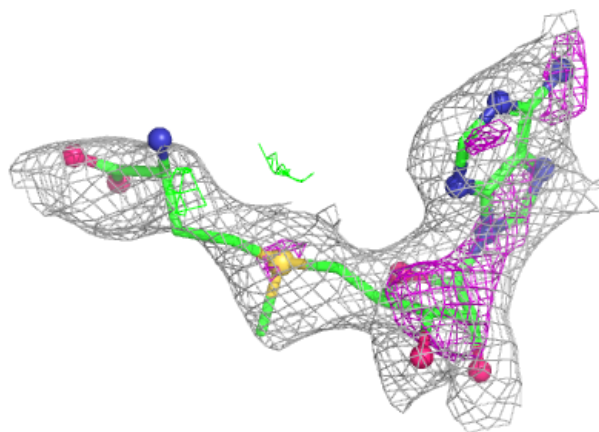
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around SAM E 500:

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

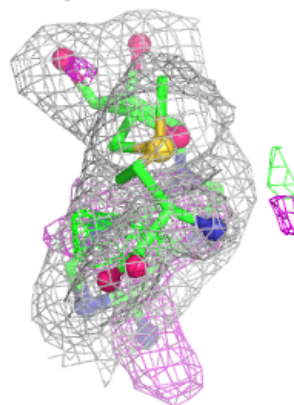
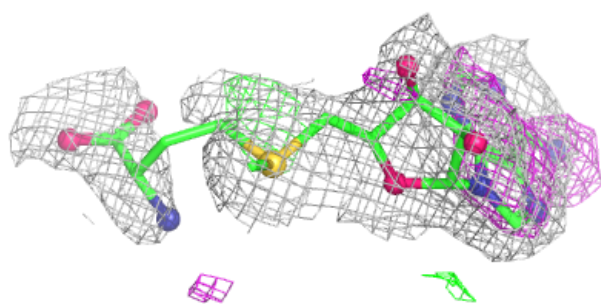
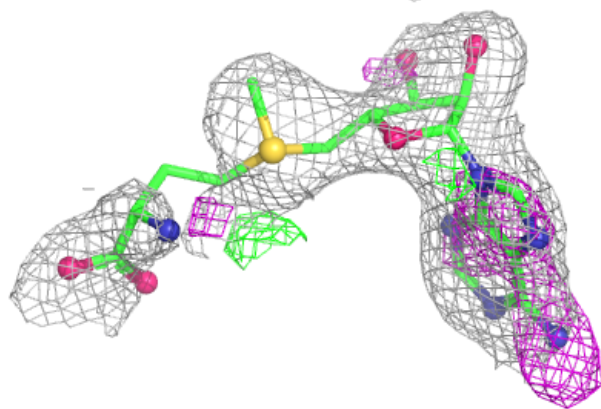
**Electron density around SAM B 500:**

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



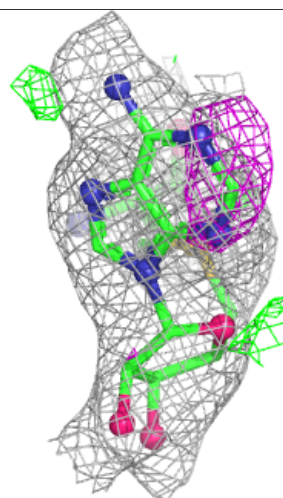
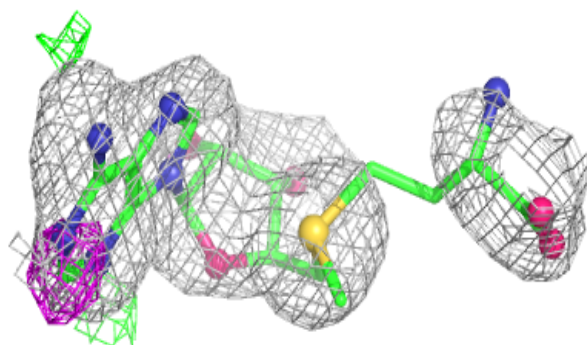
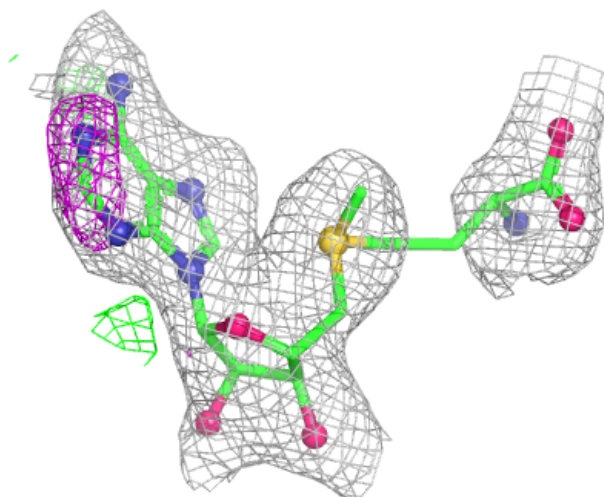
Electron density around SAM D 500:

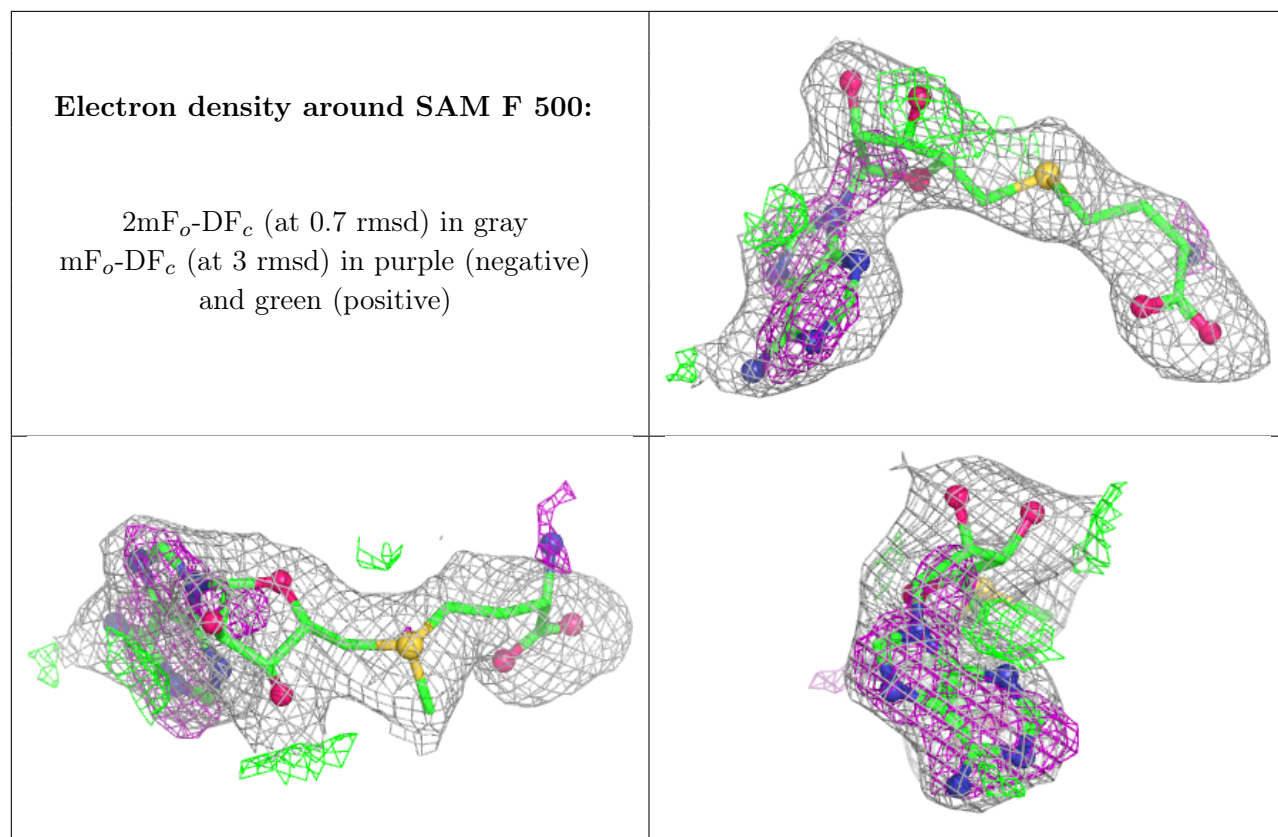
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around SAM E 501:

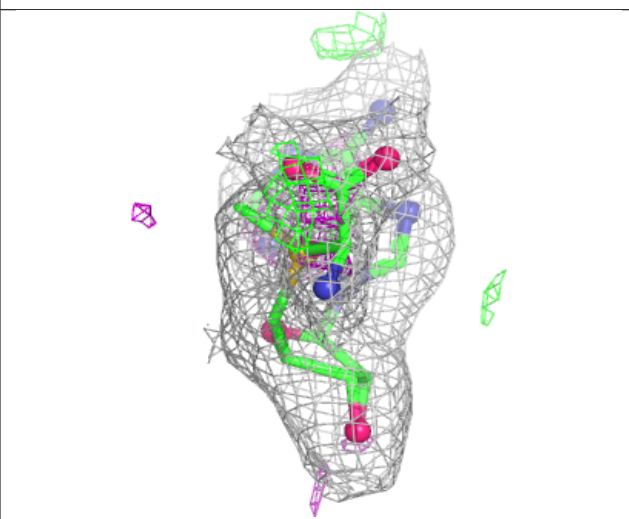
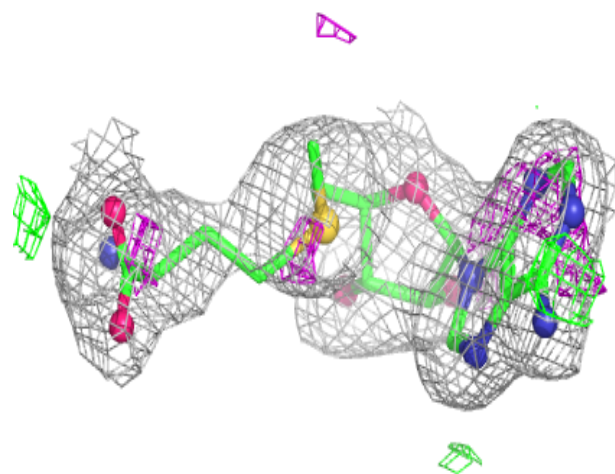
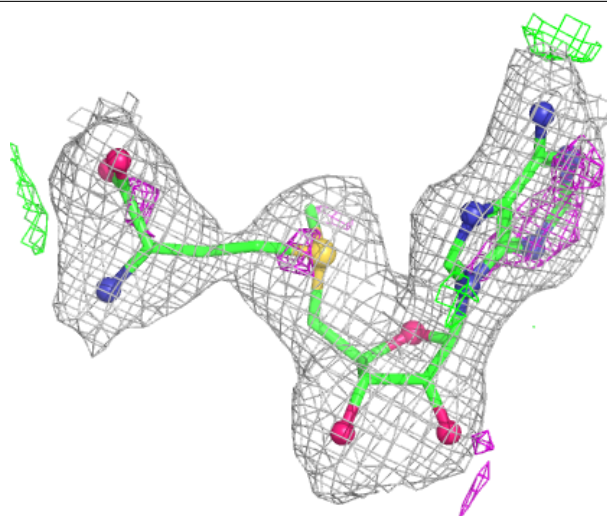
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





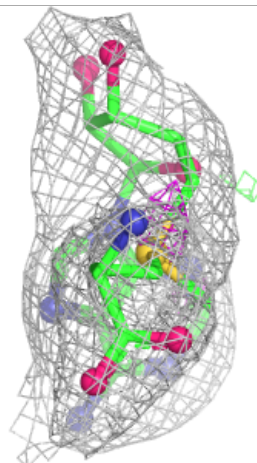
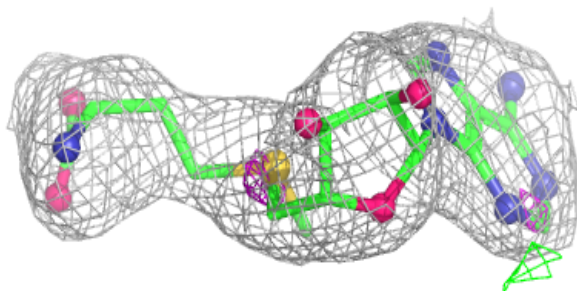
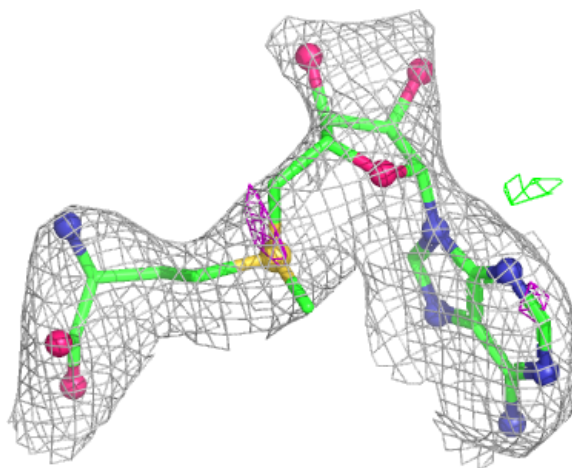
Electron density around SAM F 501:

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



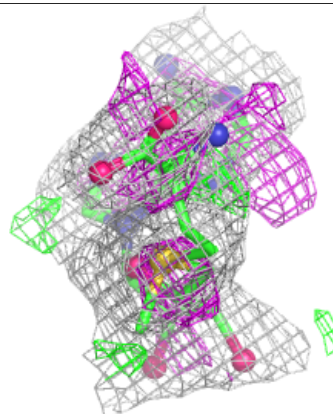
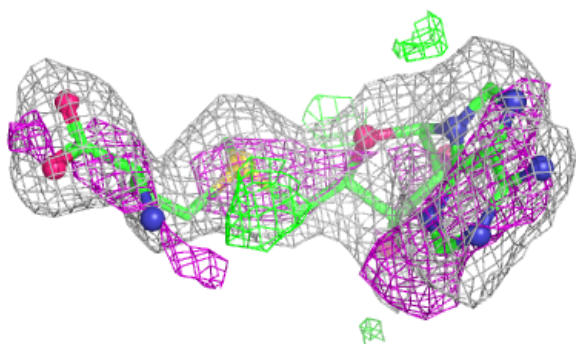
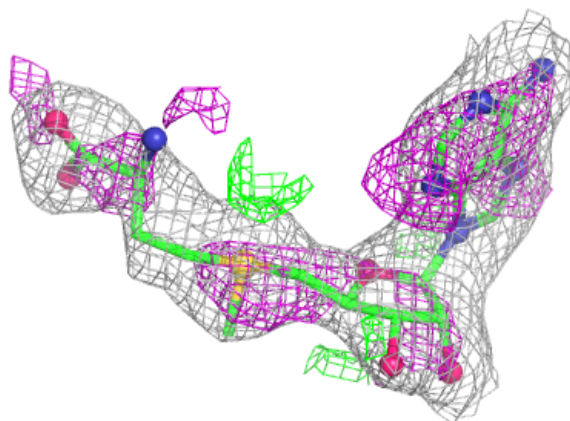
Electron density around SAM B 501:

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

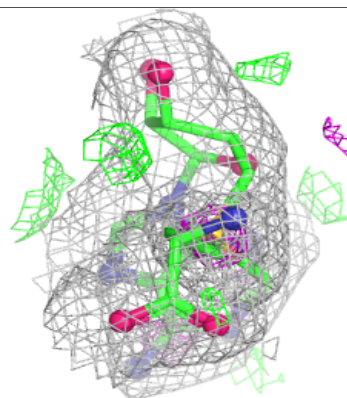
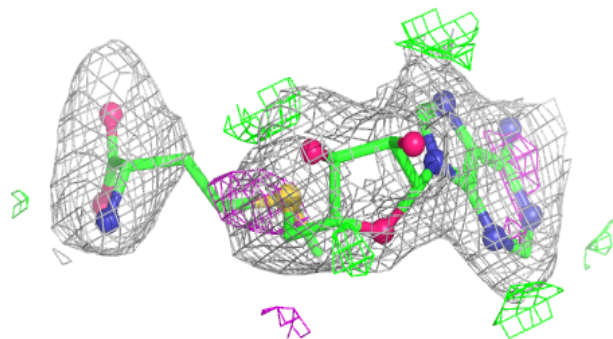
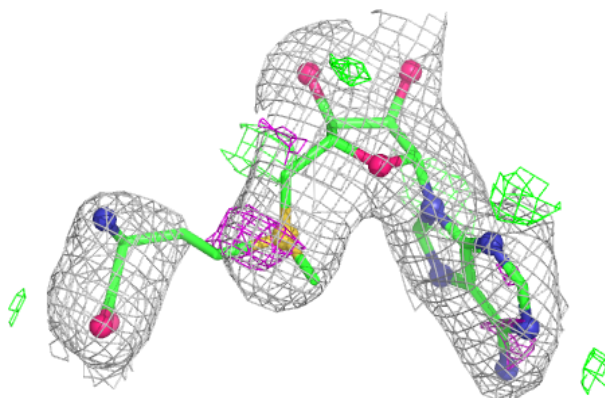


Electron density around SAM C 500:

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

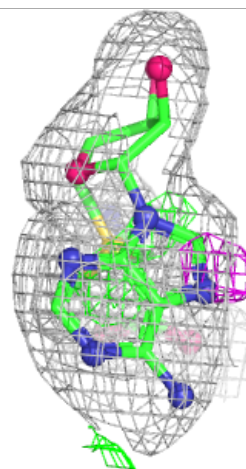
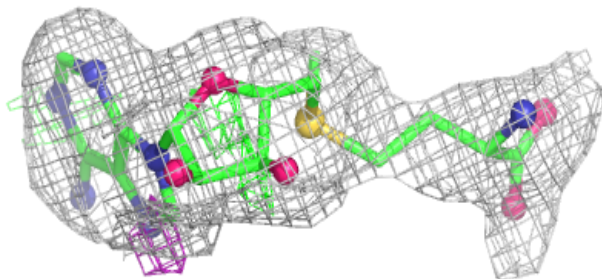
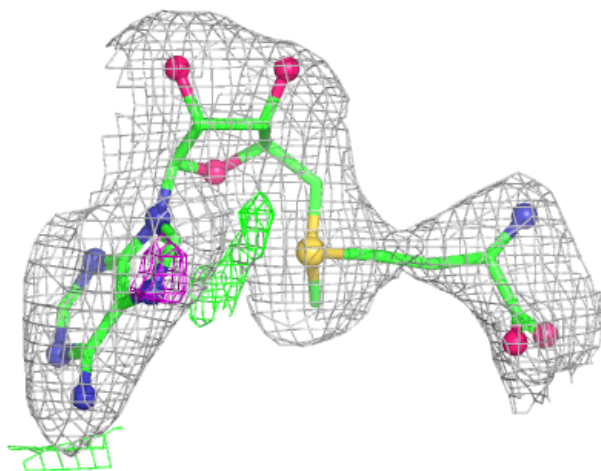
**Electron density around SAM C 501:**

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



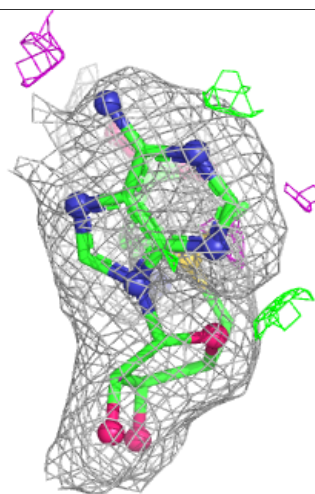
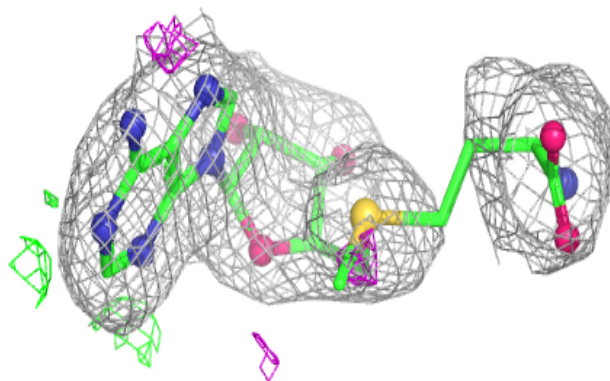
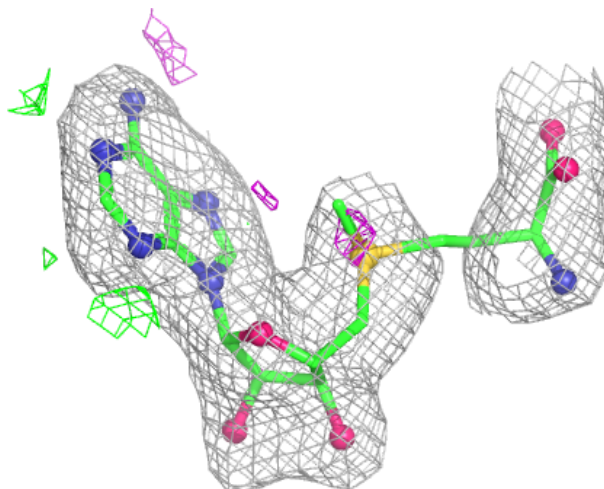
Electron density around SAM A 501:

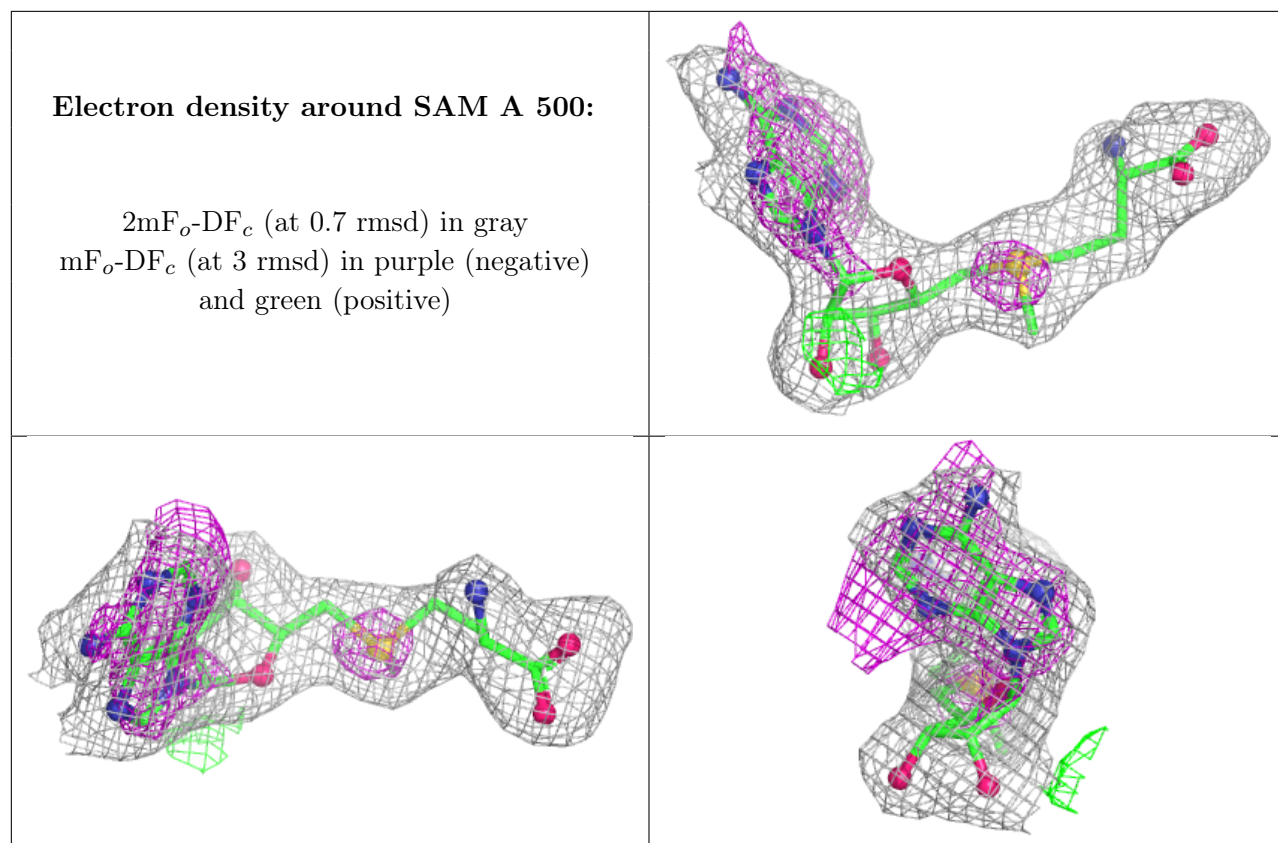
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around SAM D 501:

$2mF_o-DF_c$ (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.