



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 21, 2020 – 05:13 AM BST

PDB ID : 4NDN
Title : Structural insights of MAT enzymes: MATa2b complexed with SAM and PPNP
Authors : Murray, B.; Antonyuk, S.V.; Marina, A.; Lu, S.C.; Mato, J.M.; Hasnain, S.S.; Rojas, A.L.
Deposited on : 2013-10-27
Resolution : 2.34 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

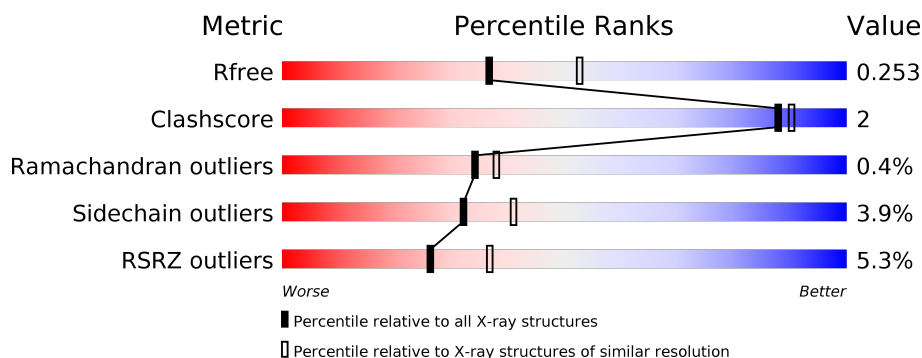
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.34 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}	130704	2096 (2.36-2.32)
Clashscore	141614	2193 (2.36-2.32)
Ramachandran outliers	138981	2159 (2.36-2.32)
Sidechain outliers	138945	2160 (2.36-2.32)
RSRZ outliers	127900	2067 (2.36-2.32)

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 on 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	395	<div> <div>2%</div> <div> <div></div> <div>87%</div> <div>5%</div> <div>8%</div> </div> </div>
1	B	395	<div> <div>3%</div> <div> <div></div> <div>88%</div> <div>8%</div> <div>• •</div> </div> </div>
1	C	395	<div> <div>7%</div> <div> <div></div> <div>83%</div> <div>8%</div> <div>9%</div> </div> </div>
1	D	395	<div> <div>4%</div> <div> <div></div> <div>88%</div> <div>7%</div> <div>• •</div> </div> </div>
2	E	323	<div> <div>2%</div> <div> <div></div> <div>92%</div> <div>• • •</div> </div> </div>
2	F	323	<div> <div>13%</div> <div> <div></div> <div>83%</div> <div>10%</div> <div>• 6%</div> </div> </div>

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 17113 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 S-adenosylmethionine synthase isoform type-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	362	Total	C	N	O	S	0	0	0
			2822	1790	492	529	11			
1	B	383	Total	C	N	O	S	0	0	0
			2980	1883	520	566	11			
1	C	360	Total	C	N	O	S	0	0	0
			2809	1782	490	526	11			
1	D	382	Total	C	N	O	S	0	1	0
			2980	1882	522	565	11			

- Molecule 2 is a protein called Methionine adenosyltransferase 2 subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	309	Total	C	N	O	S	0	1	0
			2456	1551	449	446	10			
2	F	304	Total	C	N	O	S	0	1	0
			2416	1527	442	437	10			

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Mg	0	0
			2	2		
3	D	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



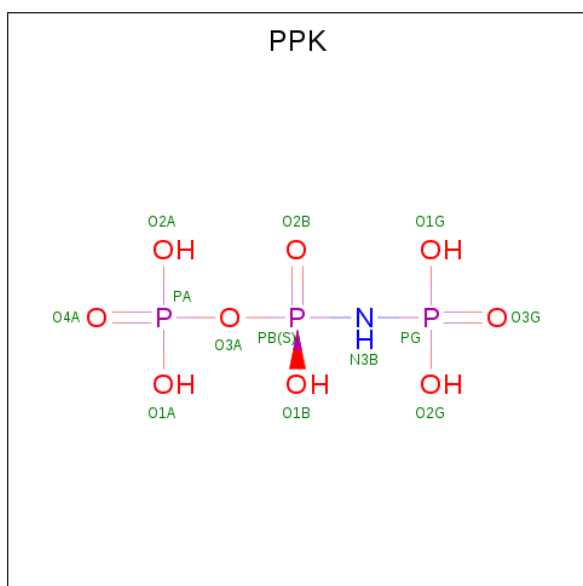
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		

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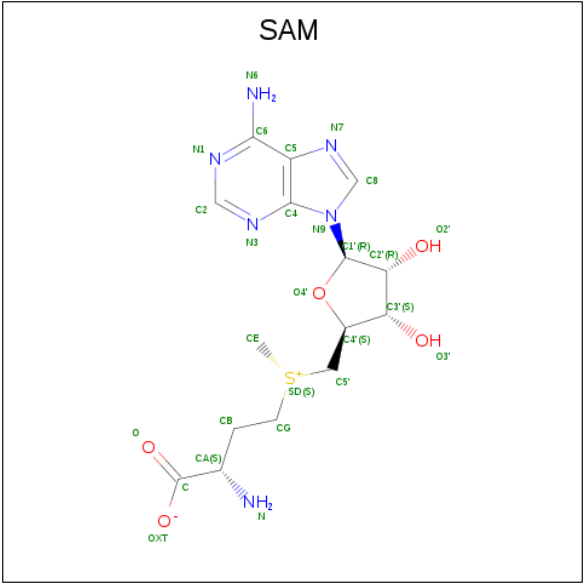
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	E	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is (DIPHOSPHONO)AMINOPHOSPHONIC ACID (three-letter code: PPK) (formula: $\text{H}_6\text{NO}_9\text{P}_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	N	O	P	0	0
			13	1	9	3		
5	C	1	Total	N	O	P	0	0
			13	1	9	3		

- Molecule 6 is S-ADENOSYLMETHIONINE (three-letter code: SAM) (formula: $\text{C}_{15}\text{H}_{22}\text{N}_6\text{O}_5\text{S}$).



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

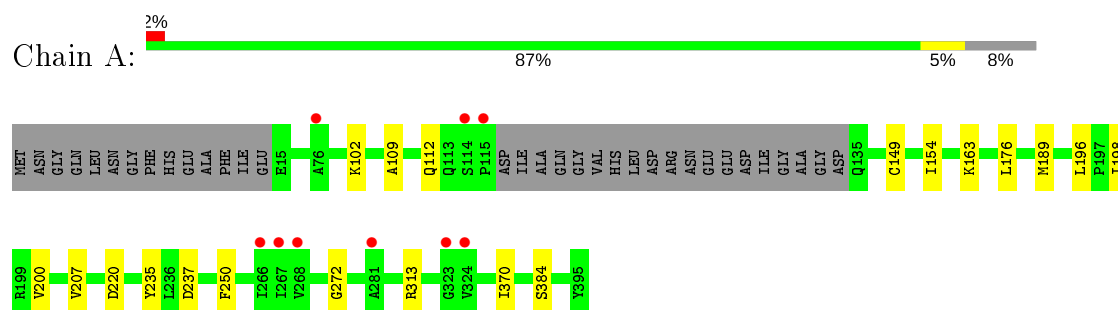
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	109	Total	O	0	0
			109	109		
7	B	97	Total	O	0	0
			97	97		
7	C	91	Total	O	0	0
			91	91		
7	D	116	Total	O	0	0
			116	116		
7	E	67	Total	O	0	0
			67	67		
7	F	18	Total	O	0	0
			18	18		

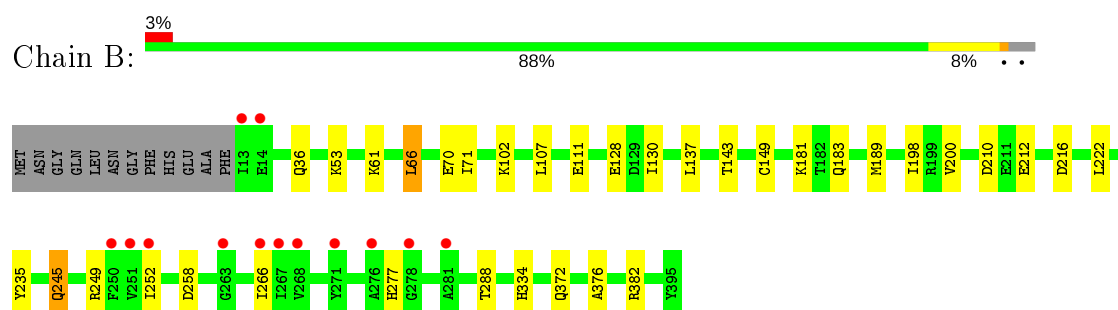
3 Residue-property plots [i](#)

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.

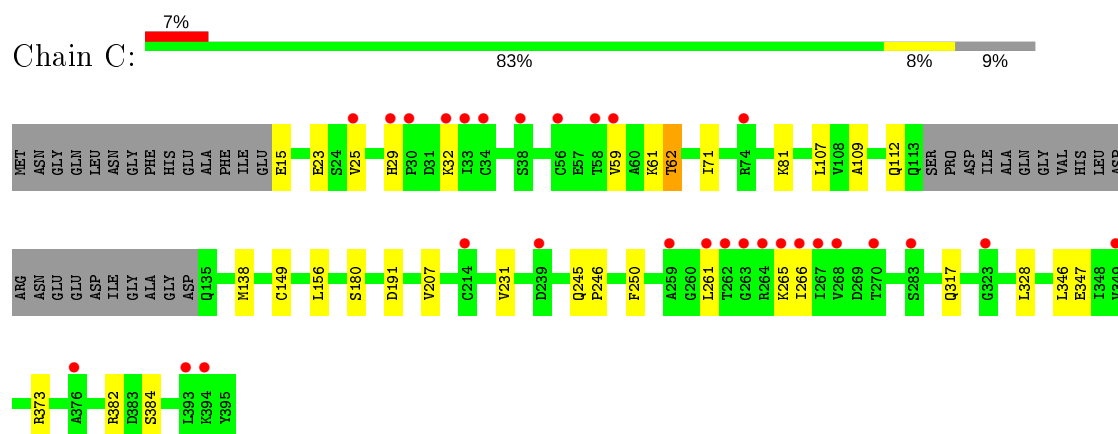
- Molecule 1: S-adenosylmethionine synthase isoform type-2




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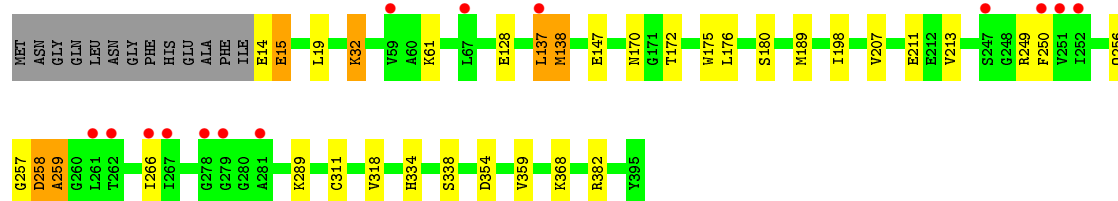


- Molecule 1: S-adenosylmethionine synthase isoform type-2




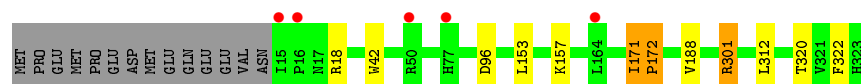
- Molecule 1: S-adenosylmethionine synthase isoform type-2

Chain D: 




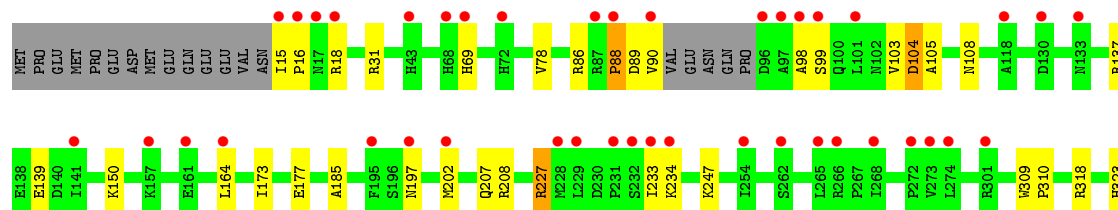
- Molecule 2: Methionine adenosyltransferase 2 subunit beta

Chain E: 



- Molecule 2: Methionine adenosyltransferase 2 subunit beta

Chain F: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.14Å 122.18Å 298.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.78 – 2.34 47.74 – 2.34	Depositor EDS
% Data completeness (in resolution range)	96.9 (47.78-2.34) 96.9 (47.74-2.34)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 2.34Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.211 , 0.251 0.212 , 0.253	Depositor DCC
R_{free} test set	5409 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	47.5	Xtriage
Anisotropy	0.220	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 32.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17113	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.37% 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: PPK, MG, EDO, SAM

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	0.32	0/2879	0.54	0/3893
1	B	0.31	0/3039	0.53	0/4111
1	C	0.29	0/2865	0.53	0/3873
1	D	0.32	0/3042	0.54	0/4114
2	E	0.30	0/2518	0.50	0/3415
2	F	0.30	0/2476	0.50	0/3355
All	All	0.31	0/16819	0.52	0/22761

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	F	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	F	88	PRO	Peptide

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	2822	0	2832	11	0
1	B	2980	0	2973	16	0
1	C	2809	0	2820	13	0
1	D	2980	0	2975	22	0
2	E	2456	0	2431	4	0
2	F	2416	0	2394	10	0
3	A	2	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	12	0	18	0	0
4	B	20	0	30	0	0
4	C	8	0	12	0	0
4	D	24	0	36	0	0
4	E	4	0	6	0	0
5	A	13	0	1	0	0
5	C	13	0	1	0	0
6	A	27	0	22	0	0
6	C	27	0	22	1	0
7	A	109	0	0	0	0
7	B	97	0	0	0	0
7	C	91	0	0	0	0
7	D	116	0	0	0	0
7	E	67	0	0	0	0
7	F	18	0	0	0	0
All	All	17113	0	16573	67	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 2.

The worst 5 of 67 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:D:189:MET:HB2	1:D:198:ILE:HD11	1.77	0.66
1:B:36:GLN:HE21	1:B:372:GLN:HE21	1.52	0.56
1:A:313:ARG:NH2	2:F:323:HIS:O	2.35	0.56
1:C:59:VAL:HG13	1:C:261:LEU:HD23	1.88	0.55
1:A:109:ALA:HB1	1:B:61:LYS:CE	2.37	0.54

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	358/395 (91%)	345 (96%)	13 (4%)	0	100	100
1	B	381/395 (96%)	365 (96%)	16 (4%)	0	100	100
1	C	356/395 (90%)	343 (96%)	12 (3%)	1 (0%)	41	47
1	D	381/395 (96%)	370 (97%)	10 (3%)	1 (0%)	41	47
2	E	308/323 (95%)	294 (96%)	13 (4%)	1 (0%)	41	47
2	F	301/323 (93%)	284 (94%)	12 (4%)	5 (2%)	9	6
All	All	2085/2226 (94%)	2001 (96%)	76 (4%)	8 (0%)	34	38

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	89	ASP
2	F	104	ASP
1	D	259	ALA
2	F	88	PRO
1	C	62	THR

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	302/327 (92%)	293 (97%)	9 (3%)	41	50
1	B	318/327 (97%)	306 (96%)	12 (4%)	33	41
1	C	300/327 (92%)	288 (96%)	12 (4%)	31	40

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	318/327 (97%)	306 (96%)	12 (4%)	33	41
2	E	263/276 (95%)	255 (97%)	8 (3%)	41	50
2	F	258/276 (94%)	242 (94%)	16 (6%)	18	20
All	All	1759/1860 (95%)	1690 (96%)	69 (4%)	32	41

5 of 69 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	382	ARG
1	D	147	GLU
2	F	202	MET
1	C	384	SER
1	D	128	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	372	GLN
2	F	207	GLN
1	C	352	ASN
1	B	245	GLN
1	C	48	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 25 ligands modelled in this entry, 4 are monoatomic - leaving 21 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	EDO	D	407	-	3,3,3	0.50	0	2,2,2	0.14	0
5	PPK	C	404	3	11,12,12	2.63	4 (36%)	15,20,20	1.04	1 (6%)
4	EDO	B	403	-	3,3,3	0.46	0	2,2,2	0.29	0
4	EDO	B	402	-	3,3,3	0.67	0	2,2,2	0.10	0
4	EDO	D	402	-	3,3,3	0.48	0	2,2,2	0.25	0
4	EDO	C	402	-	3,3,3	0.42	0	2,2,2	0.33	0
4	EDO	D	406	-	3,3,3	0.46	0	2,2,2	0.34	0
6	SAM	A	407	-	21,29,29	0.99	1 (4%)	18,42,42	1.68	4 (22%)
4	EDO	C	403	-	3,3,3	0.56	0	2,2,2	0.12	0
4	EDO	B	401	-	3,3,3	0.45	0	2,2,2	0.32	0
5	PPK	A	406	3	11,12,12	2.66	4 (36%)	15,20,20	1.07	0
4	EDO	D	405	-	3,3,3	0.59	0	2,2,2	0.21	0
4	EDO	B	405	-	3,3,3	0.45	0	2,2,2	0.30	0
6	SAM	C	405	-	21,29,29	1.03	1 (4%)	18,42,42	1.53	3 (16%)
4	EDO	D	404	-	3,3,3	0.47	0	2,2,2	0.38	0
4	EDO	A	404	-	3,3,3	0.44	0	2,2,2	0.31	0
4	EDO	B	404	-	3,3,3	0.50	0	2,2,2	0.24	0
4	EDO	E	401	-	3,3,3	0.48	0	2,2,2	0.33	0
4	EDO	A	405	-	3,3,3	0.46	0	2,2,2	0.31	0
4	EDO	D	403	-	3,3,3	0.49	0	2,2,2	0.31	0
4	EDO	A	403	-	3,3,3	0.47	0	2,2,2	0.22	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	EDO	D	407	-	-	1/1/1/1	-
5	PPK	C	404	3	-	3/8/12/12	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	403	-	-	0/1/1/1	-
4	EDO	B	402	-	-	0/1/1/1	-
4	EDO	D	402	-	-	1/1/1/1	-
4	EDO	C	402	-	-	1/1/1/1	-
4	EDO	D	406	-	-	1/1/1/1	-
6	SAM	A	407	-	-	0/8/33/33	0/3/3/3
4	EDO	C	403	-	-	1/1/1/1	-
4	EDO	B	401	-	-	1/1/1/1	-
5	PPK	A	406	3	-	2/8/12/12	-
4	EDO	D	405	-	-	1/1/1/1	-
4	EDO	B	405	-	-	1/1/1/1	-
6	SAM	C	405	-	-	4/8/33/33	0/3/3/3
4	EDO	D	404	-	-	1/1/1/1	-
4	EDO	A	404	-	-	1/1/1/1	-
4	EDO	B	404	-	-	0/1/1/1	-
4	EDO	E	401	-	-	0/1/1/1	-
4	EDO	A	405	-	-	1/1/1/1	-
4	EDO	D	403	-	-	1/1/1/1	-
4	EDO	A	403	-	-	1/1/1/1	-

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	406	PPK	PB-N3B	5.08	1.76	1.63
5	A	406	PPK	PG-N3B	4.94	1.76	1.63
5	C	404	PPK	PB-N3B	4.87	1.76	1.63
5	C	404	PPK	PG-N3B	4.67	1.75	1.63
5	C	404	PPK	PB-O3A	3.19	1.63	1.59

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	407	SAM	N3-C2-N1	-3.81	122.72	128.68
6	C	405	SAM	N3-C2-N1	-3.61	123.03	128.68
6	C	405	SAM	C4-C5-N7	-2.79	106.50	109.40
6	A	407	SAM	C2-N1-C6	2.67	123.31	118.75
6	C	405	SAM	C3'-C2'-C1'	2.42	104.62	100.98

There are no chirality outliers.

5 of 22 torsion outliers are listed below:

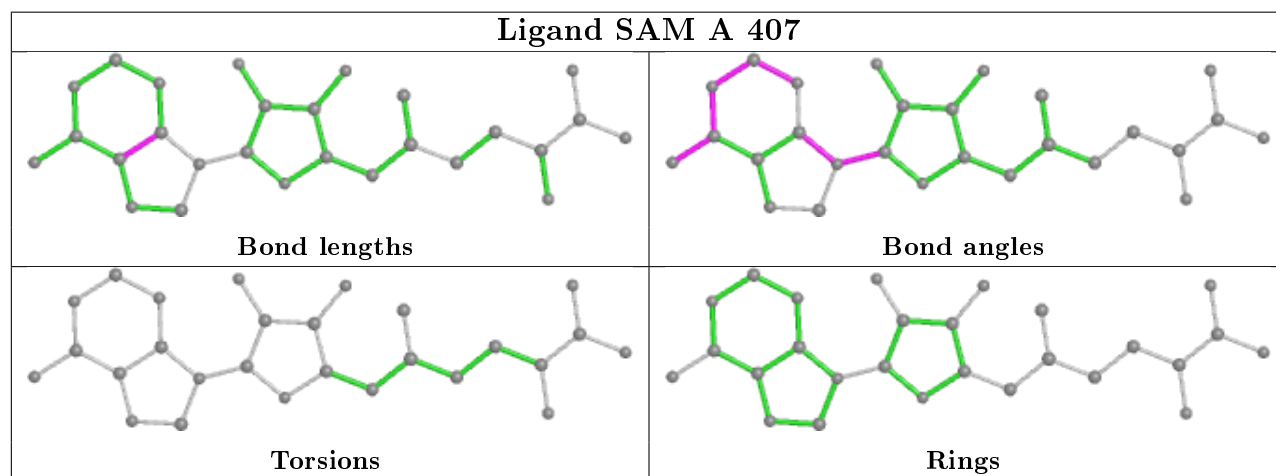
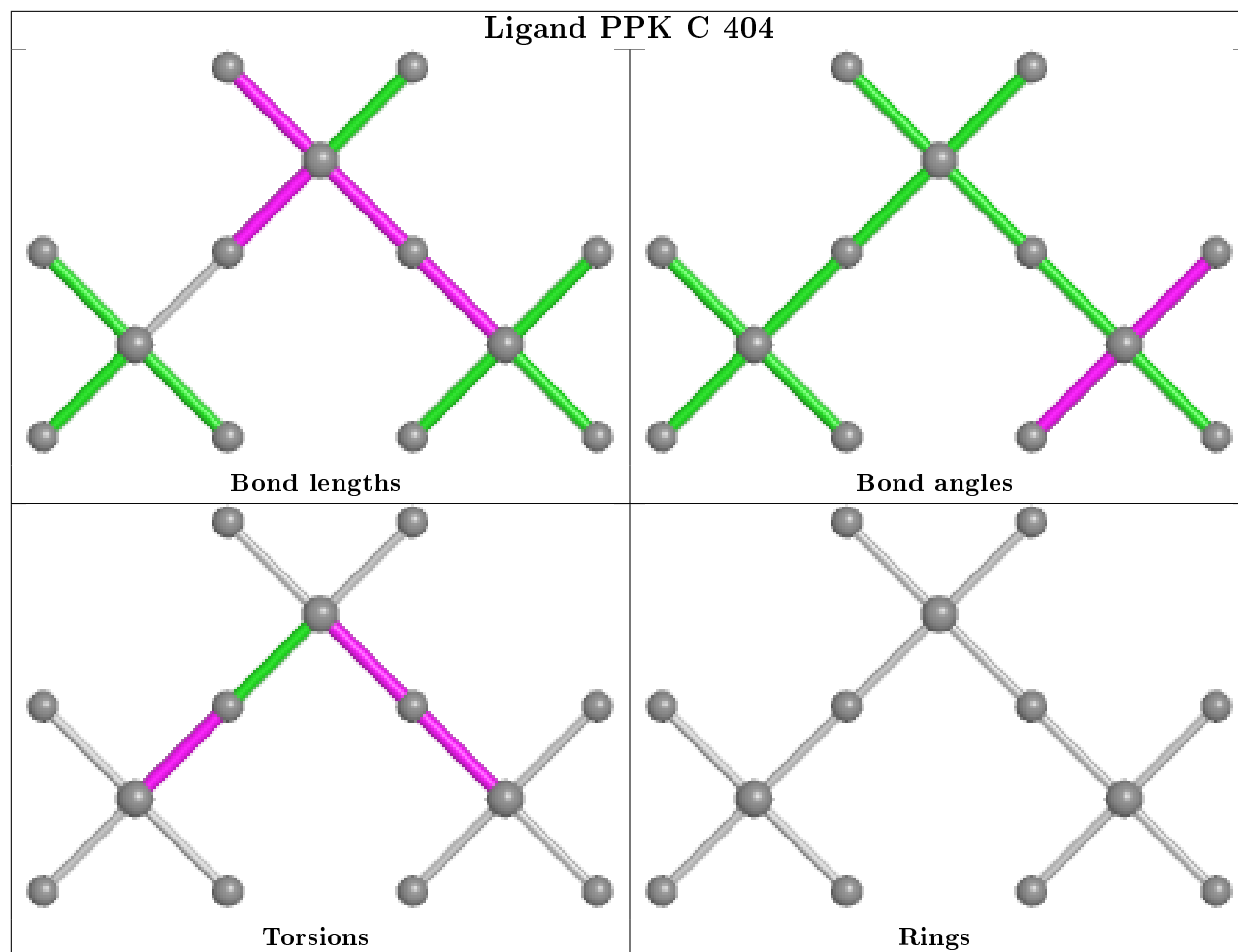
Mol	Chain	Res	Type	Atoms
5	C	404	PPK	PB-N3B-PG-O3G
5	C	404	PPK	PG-N3B-PB-O2B
5	C	404	PPK	PB-O3A-PA-O2A
5	A	406	PPK	PB-N3B-PG-O3G
5	A	406	PPK	PG-N3B-PB-O2B

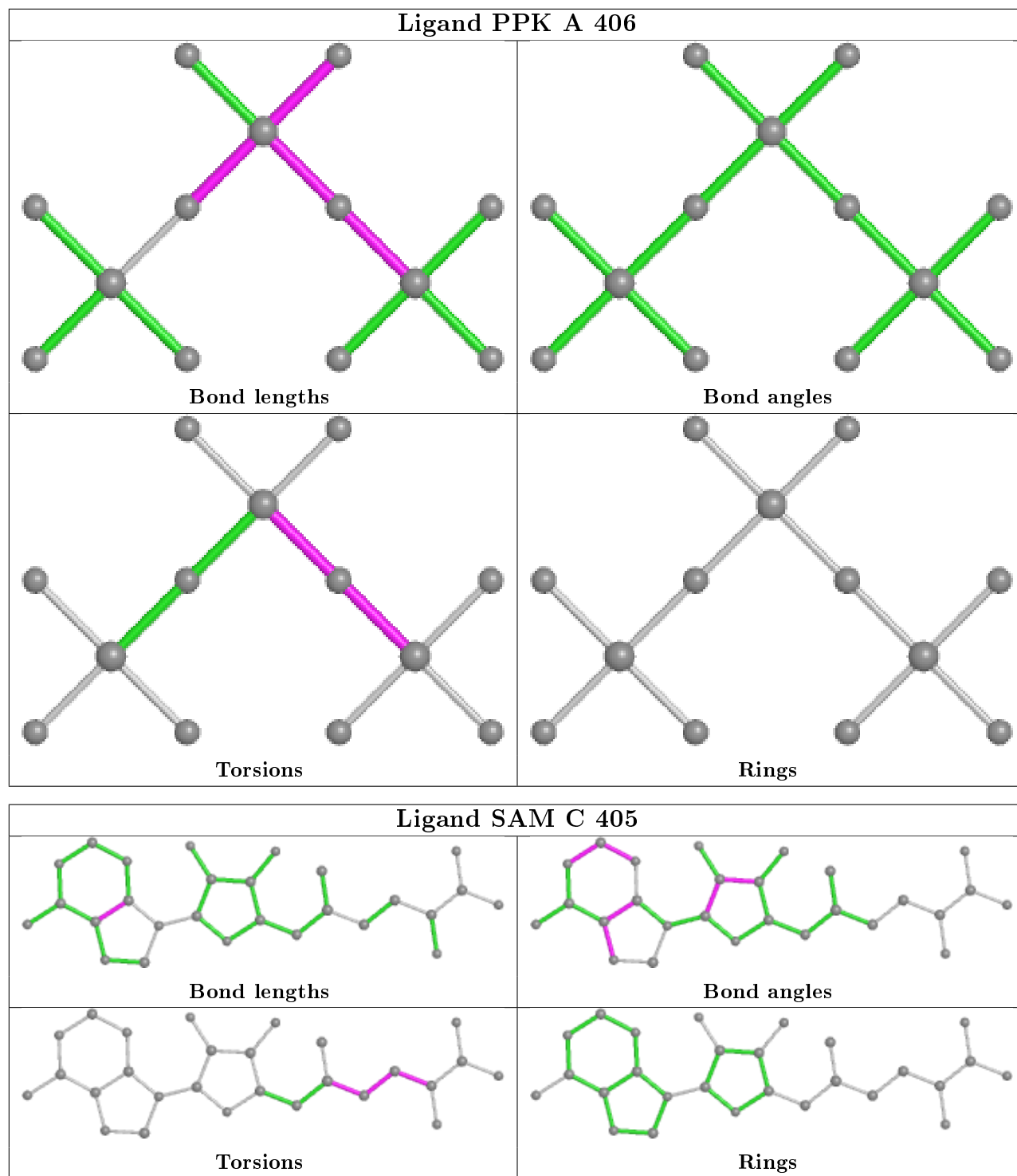
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	405	SAM	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers ⓘ

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 [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	362/395 (91%)	0.07	9 (2%) 57 66	32, 47, 71, 109	0
1	B	383/395 (96%)	0.21	13 (3%) 45 55	34, 48, 74, 120	0
1	C	360/395 (91%)	0.43	29 (8%) 12 18	36, 56, 81, 93	0
1	D	382/395 (96%)	0.09	14 (3%) 41 52	32, 44, 69, 104	0
2	E	309/323 (95%)	-0.01	5 (1%) 72 80	41, 57, 83, 106	0
2	F	304/323 (94%)	0.76	41 (13%) 3 5	46, 83, 111, 130	0
All	All	2100/2226 (94%)	0.25	111 (5%) 26 37	32, 52, 92, 130	0

The worst 5 of 111 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	15	ILE	6.1
2	E	15	ILE	6.0
2	F	99	SER	6.0
2	F	96	ASP	5.5
1	B	13	ILE	5.3

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 monosaccharides in this entry.

6.4 Ligands ⓘ

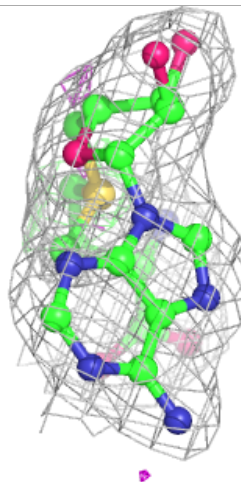
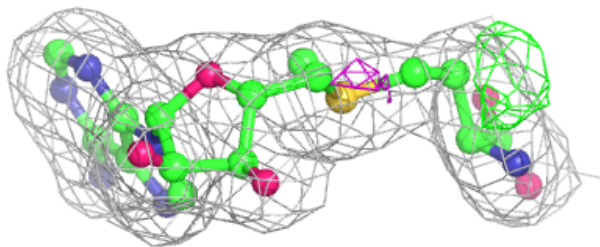
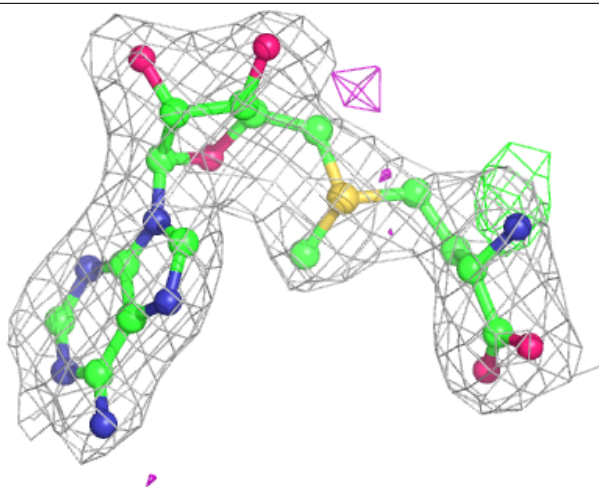
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	EDO	C	403	4/4	0.72	0.25	65,68,68,70	0
4	EDO	D	402	4/4	0.74	0.29	69,70,71,71	0
4	EDO	B	401	4/4	0.74	0.28	75,76,78,81	0
4	EDO	B	402	4/4	0.76	0.26	46,47,47,48	0
4	EDO	B	404	4/4	0.78	0.28	73,75,75,76	0
4	EDO	D	405	4/4	0.79	0.30	57,57,58,60	0
4	EDO	A	405	4/4	0.81	0.22	75,77,77,79	0
4	EDO	D	403	4/4	0.84	0.20	66,67,67,68	0
4	EDO	A	403	4/4	0.85	0.23	63,65,67,71	0
4	EDO	D	406	4/4	0.85	0.29	68,68,70,75	0
4	EDO	A	404	4/4	0.87	0.23	65,67,68,71	0
4	EDO	D	404	4/4	0.88	0.23	69,69,70,70	0
4	EDO	C	402	4/4	0.90	0.58	70,70,71,72	0
6	SAM	A	407	27/27	0.93	0.15	38,43,56,57	0
3	MG	A	401	1/1	0.94	0.04	44,44,44,44	0
5	PPK	C	404	13/13	0.94	0.15	52,61,67,73	0
4	EDO	B	405	4/4	0.94	0.14	56,59,59,61	0
5	PPK	A	406	13/13	0.94	0.13	56,63,72,74	0
4	EDO	B	403	4/4	0.94	0.17	55,56,56,57	0
6	SAM	C	405	27/27	0.94	0.17	41,48,56,57	0
4	EDO	E	401	4/4	0.95	0.12	45,45,45,45	0
4	EDO	D	407	4/4	0.96	0.32	70,71,72,73	0
3	MG	C	401	1/1	0.98	0.15	43,43,43,43	0
3	MG	D	401	1/1	0.98	0.20	32,32,32,32	0
3	MG	A	402	1/1	0.99	0.19	37,37,37,37	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.

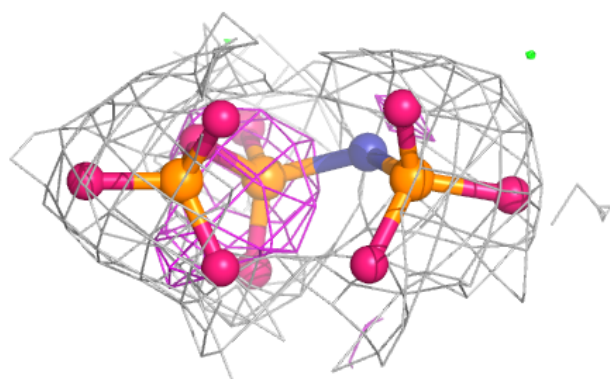
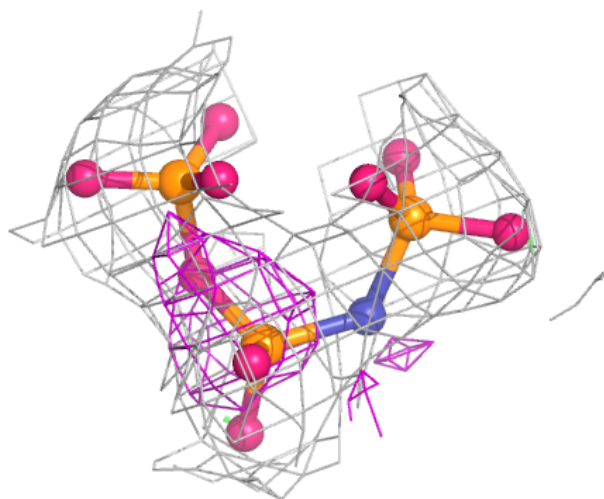
Electron density around SAM A 407:

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



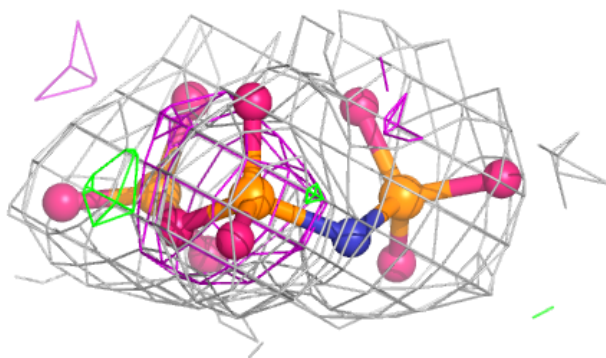
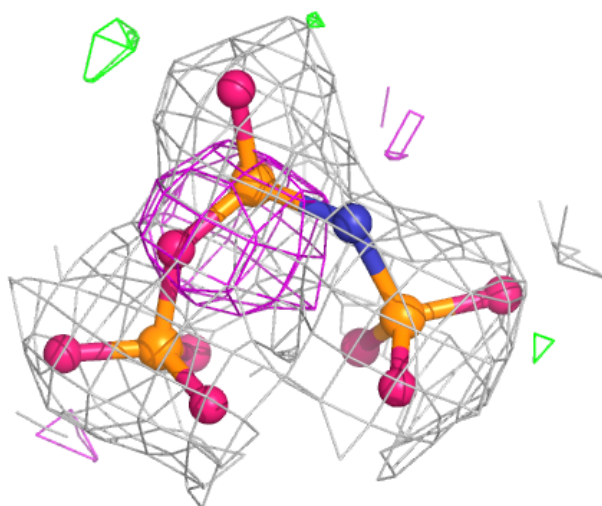
Electron density around PPK C 404:

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



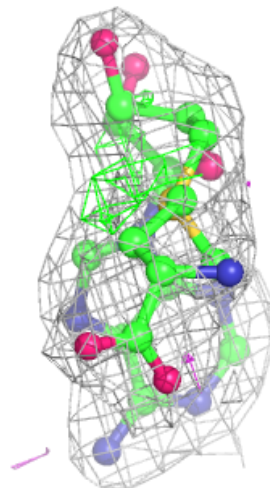
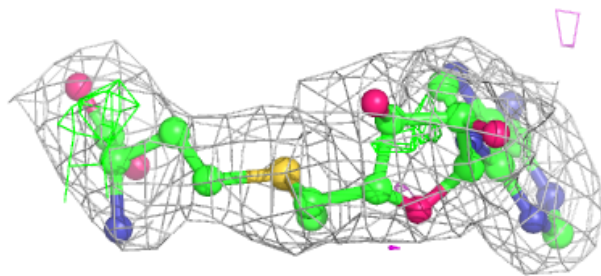
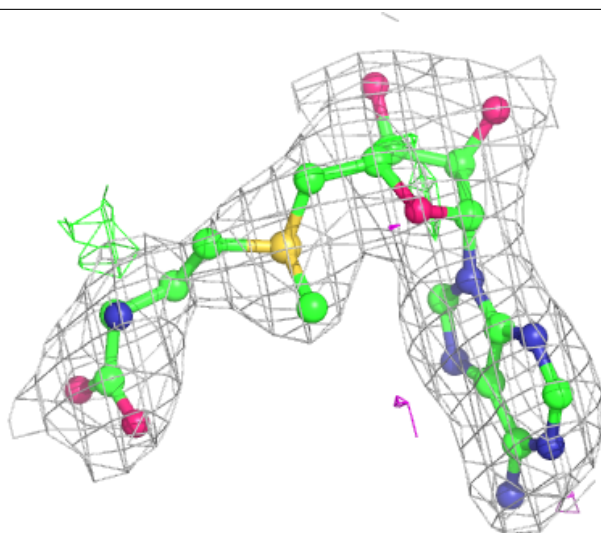
Electron density around PPK A 406:

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



Electron density around SAM C 405:

$2mF_o-DF_c$ (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.