



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 03:39 PM UTC

PDB ID : 9TB4 / pdb_00009tb4
Title : SAM-dependent C6-FPP methyltransferase from *Streptomyces varsoviensis* in complex with SAH and FPP
Authors : Li, H.; Schnakenburg, G.; Groll, M.; Dickschat, J.S.
Deposited on : 2025-11-19
Resolution : 2.20 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

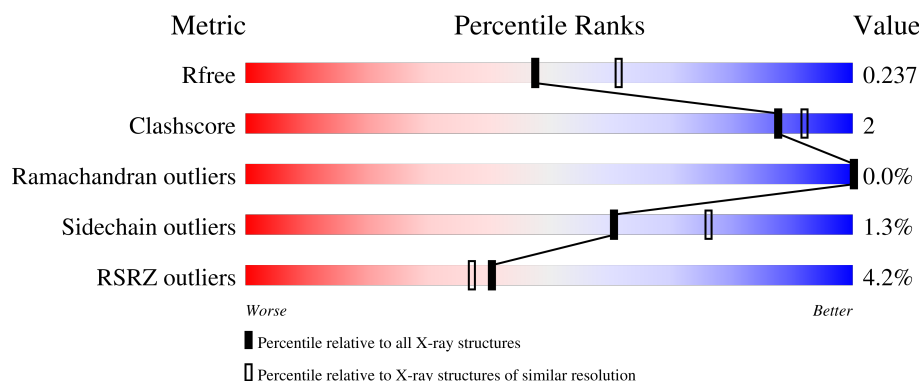
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	303	3% 86% 5% 9%
1	B	303	6% 87% • 9%
1	C	303	• 85% 5% 9%
1	D	303	3% 85% 5% 9%
1	E	303	6% 86% • 10%

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Mol	Chain	Length	Quality of chain
1	F	303	
1	G	303	
1	H	303	
1	I	303	
1	J	303	
1	K	303	
1	L	303	

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
6	PE8	E	902	-	-	X	-

2 Entry composition

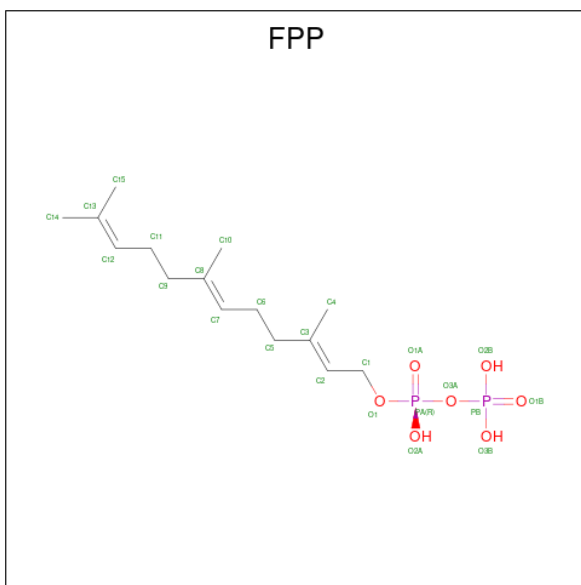
There are 9 unique types of molecules in this entry. The entry contains 27128 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 C6-FPP methyltransferase.

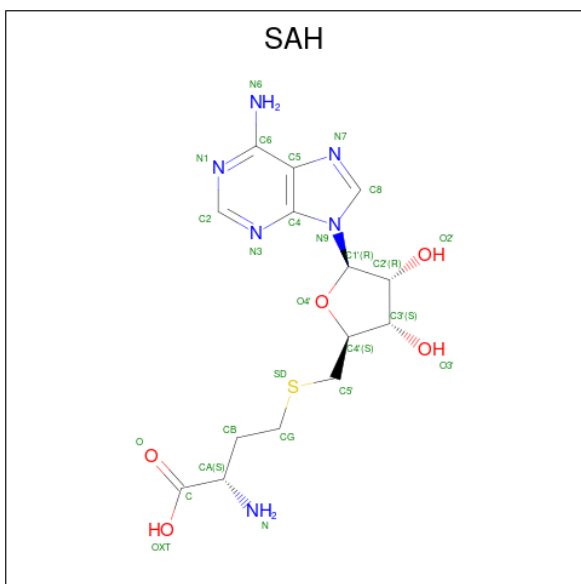
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	277	Total	C	N	O	S	0	0	0
			2170	1378	387	397	8			
1	B	276	Total	C	N	O	S	0	0	0
			2154	1364	386	396	8			
1	C	275	Total	C	N	O	S	0	0	0
			2146	1360	385	393	8			
1	D	275	Total	C	N	O	S	0	0	0
			2146	1360	385	393	8			
1	E	274	Total	C	N	O	S	0	0	0
			2140	1357	384	391	8			
1	F	278	Total	C	N	O	S	0	0	0
			2178	1382	388	400	8			
1	G	274	Total	C	N	O	S	0	0	0
			2137	1355	381	393	8			
1	H	274	Total	C	N	O	S	0	0	0
			2137	1355	381	393	8			
1	I	271	Total	C	N	O	S	0	0	0
			2115	1342	375	390	8			
1	J	276	Total	C	N	O	S	0	0	0
			2154	1364	386	396	8			
1	K	274	Total	C	N	O	S	0	0	0
			2137	1355	381	393	8			
1	L	268	Total	C	N	O	S	0	0	0
			2090	1327	374	381	8			

- Molecule 2 is FARNESYL DIPHOSPHATE (CCD ID: FPP) (formula: $C_{15}H_{28}O_7P_2$) (labeled as "Ligand of Interest" by depositor).



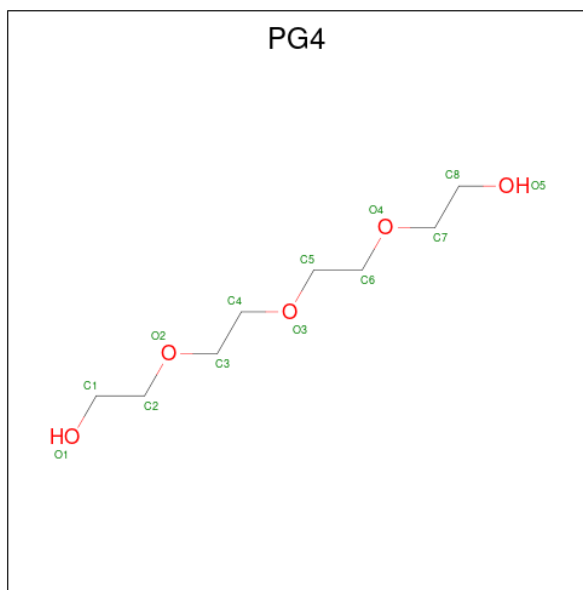
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			24	15	7	2		
2	F	1	Total	C	O	P	0	0
			24	15	7	2		
2	G	1	Total	C	O	P	0	0
			24	15	7	2		

- Molecule 3 is S-ADENOSYL-L-HOMOCYSTEINE (CCD ID: SAH) (formula: $C_{14}H_{20}N_6O_5S$) (labeled as "Ligand of Interest" by depositor).



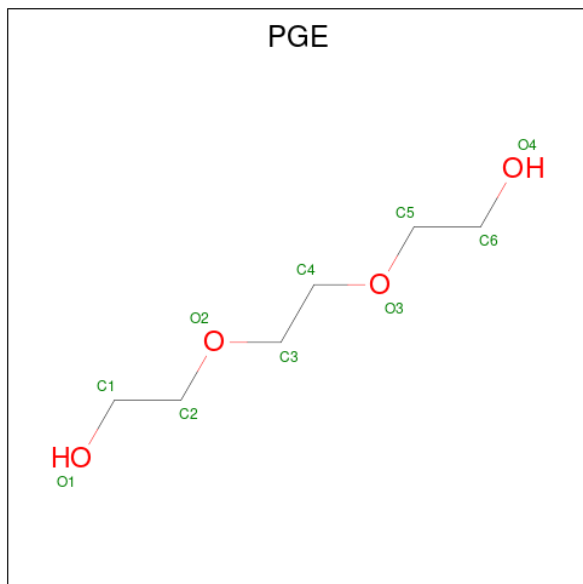
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
3	B	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
3	C	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
3	D	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
3	E	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
3	F	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
3	G	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
3	H	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
3	I	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
3	J	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
3	K	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
3	L	1	Total	C	N	O	S	0	0
			26	14	6	5	1		

- Molecule 4 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	8	5		
4	B	1	Total	C	O	0	0
			13	8	5		
4	E	1	Total	C	O	0	0
			13	8	5		
4	E	1	Total	C	O	0	0
			13	8	5		
4	F	1	Total	C	O	0	0
			13	8	5		
4	H	1	Total	C	O	0	0
			13	8	5		

- Molecule 5 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



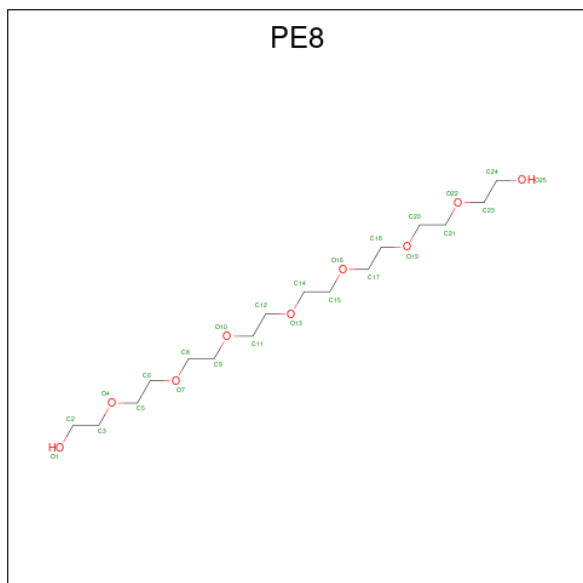
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			10	6	4		
5	B	1	Total	C	O	0	0
			10	6	4		
5	C	1	Total	C	O	0	0
			10	6	4		
5	D	1	Total	C	O	0	0
			10	6	4		
5	F	1	Total	C	O	0	0
			10	6	4		
5	I	1	Total	C	O	0	0
			10	6	4		

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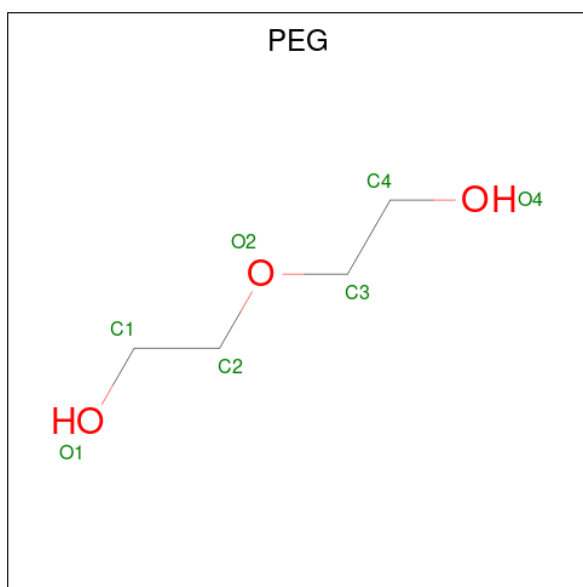
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	K	1	Total	C	O	0	0
			10	6	4		
5	K	1	Total	C	O	0	0
			10	6	4		

- Molecule 6 is 3,6,9,12,15,18,21-HEPTAOXATRICOSANE-1,23-DIOL (CCD ID: PE8) (formula: $C_{16}H_{34}O_9$).



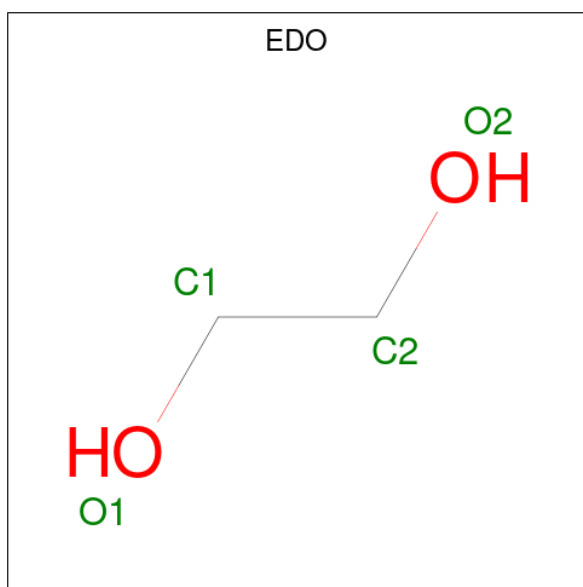
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	E	1	Total	C	O	0	0
			25	16	9		

- Molecule 7 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	E	1	Total	C	O	0	0
			7	4	3		
7	K	1	Total	C	O	0	0
			7	4	3		
7	K	1	Total	C	O	0	0
			7	4	3		

- Molecule 8 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

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

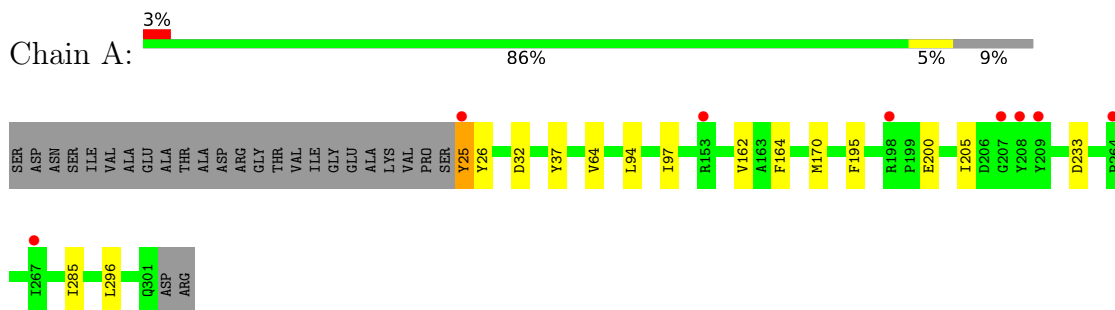
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	61	Total	O	0	0
			61	61		
9	B	93	Total	O	0	0
			93	93		
9	C	79	Total	O	0	0
			79	79		
9	D	70	Total	O	0	0
			70	70		
9	E	85	Total	O	0	0
			85	85		
9	F	93	Total	O	0	0
			93	93		
9	G	67	Total	O	0	0
			67	67		
9	H	78	Total	O	0	0
			78	78		
9	I	45	Total	O	0	0
			45	45		
9	J	44	Total	O	0	0
			44	44		
9	K	49	Total	O	0	0
			49	49		
9	L	56	Total	O	0	0
			56	56		

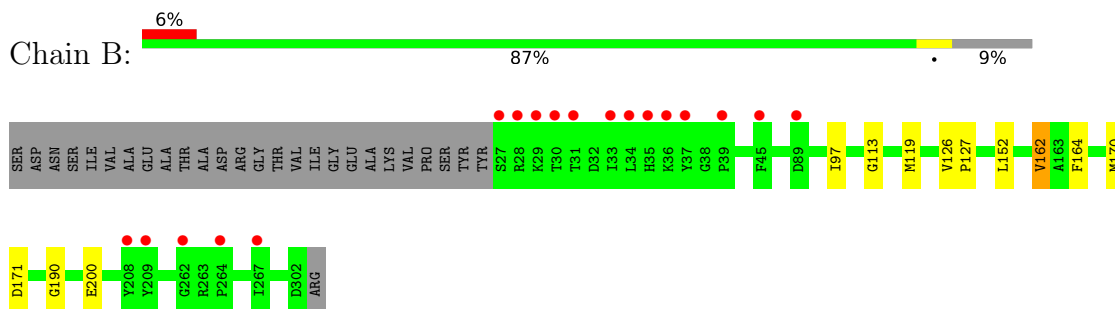
3 Residue-property plots [i](#)

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.

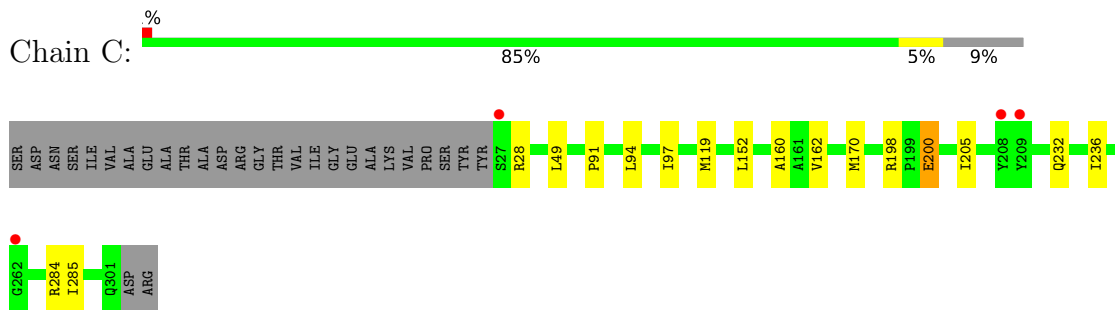
- Molecule 1: C6-FPP methytransferase



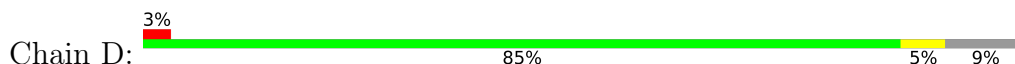
- Molecule 1: C6-FPP methytransferase

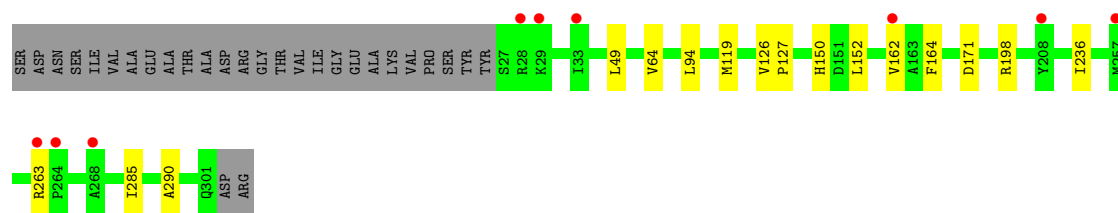


- Molecule 1: C6-FPP methytransferase

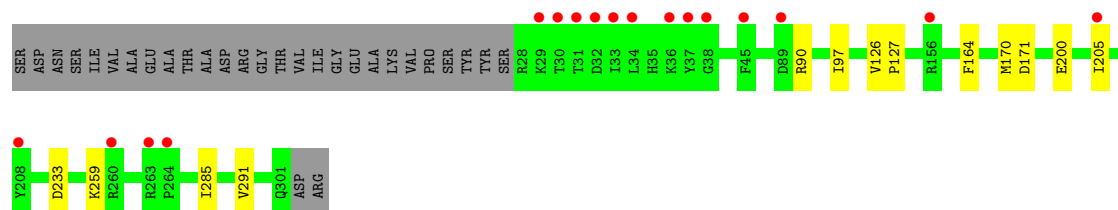
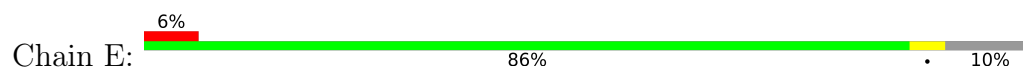


- Molecule 1: C6-FPP methytransferase

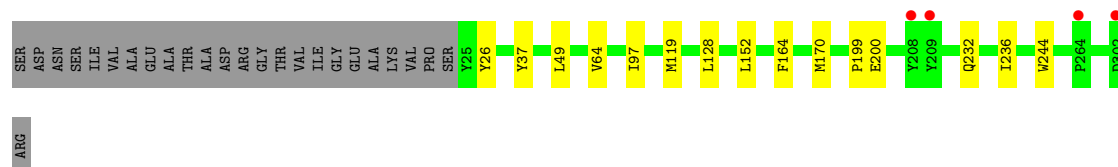
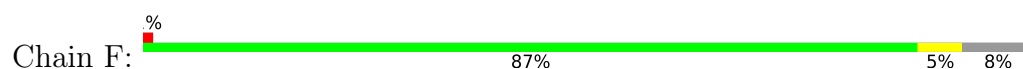




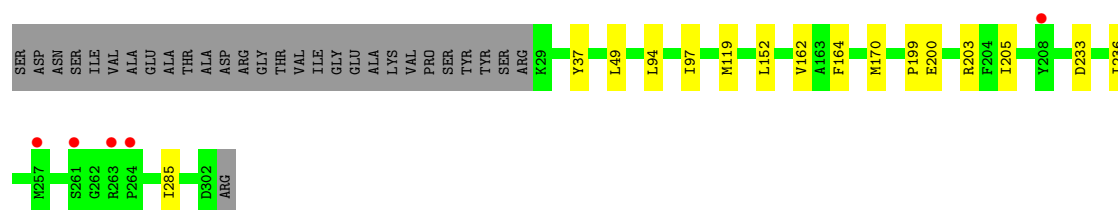
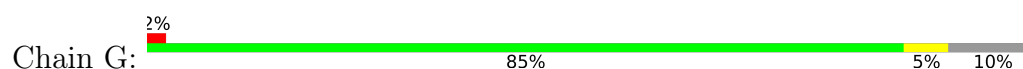
● Molecule 1: C6-FPP methytransferase



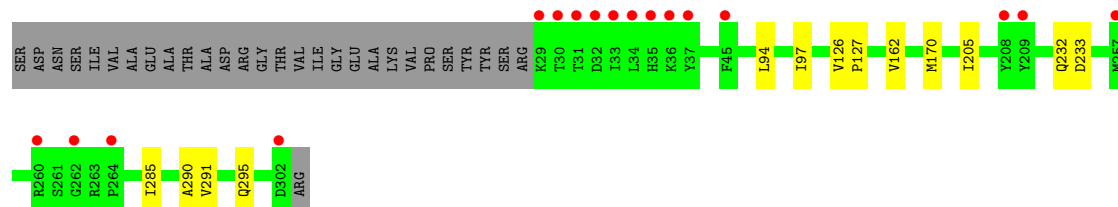
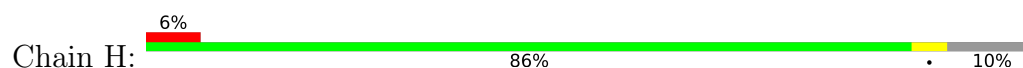
● Molecule 1: C6-FPP methytransferase



● Molecule 1: C6-FPP methytransferase

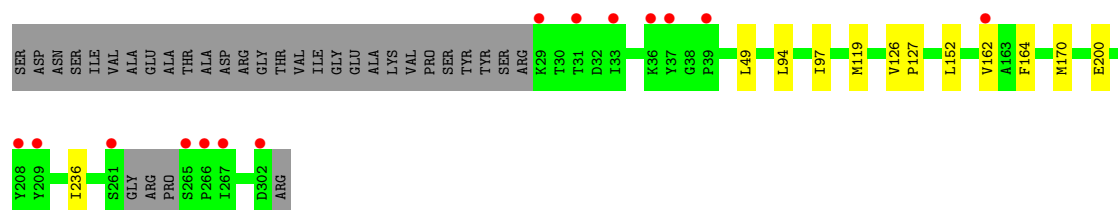


● Molecule 1: C6-FPP methytransferase



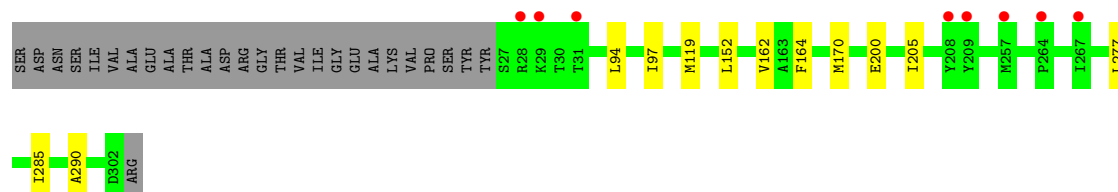
- Molecule 1: C6-FPP methytransferase

Chain I: 



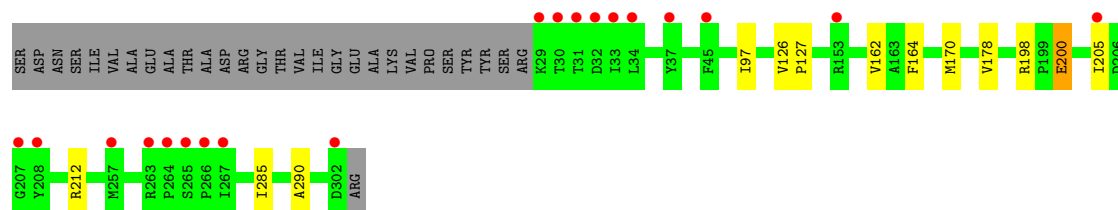
- Molecule 1: C6-FPP methytransferase

Chain J: 



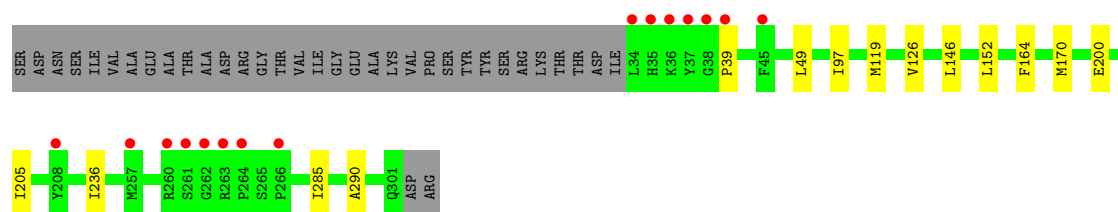
- Molecule 1: C6-FPP methytransferase

Chain K: 



- Molecule 1: C6-FPP methytransferase

Chain L: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	130.56Å 133.33Å 231.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.20 30.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.6 (30.00-2.20) 98.6 (30.00-2.20)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.209 , 0.237 0.208 , 0.237	Depositor DCC
R_{free} test set	10080 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	34.6	Xtriage
Anisotropy	0.131	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 43.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	27128	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.25% 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: PE8, PG4, SAH, EDO, FPP, PEG, PGE

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	0.43	0/2224	0.90	0/3021
1	B	0.43	0/2206	0.92	0/2996
1	C	0.43	0/2198	0.92	0/2985
1	D	0.43	0/2198	0.93	0/2985
1	E	0.43	0/2192	0.91	0/2977
1	F	0.43	0/2232	0.91	0/3032
1	G	0.43	0/2189	0.92	0/2974
1	H	0.43	0/2189	0.92	0/2974
1	I	0.43	0/2165	0.91	0/2940
1	J	0.43	0/2206	0.92	0/2996
1	K	0.43	0/2189	0.93	0/2974
1	L	0.43	0/2142	0.90	0/2910
All	All	0.43	0/26330	0.92	0/35764

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 [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	2170	0	2108	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2154	0	2094	6	0
1	C	2146	0	2090	12	0
1	D	2146	0	2090	8	0
1	E	2140	0	2085	7	0
1	F	2178	0	2112	11	0
1	G	2137	0	2076	10	0
1	H	2137	0	2076	9	0
1	I	2115	0	2052	5	0
1	J	2154	0	2094	4	0
1	K	2137	0	2076	4	0
1	L	2090	0	2030	5	0
2	A	24	0	25	6	0
2	F	24	0	25	4	0
2	G	24	0	25	3	0
3	A	26	0	19	0	0
3	B	26	0	19	0	0
3	C	26	0	19	0	0
3	D	26	0	19	0	0
3	E	26	0	19	0	0
3	F	26	0	19	0	0
3	G	26	0	19	0	0
3	H	26	0	19	0	0
3	I	26	0	19	0	0
3	J	26	0	19	0	0
3	K	26	0	19	0	0
3	L	26	0	19	0	0
4	A	13	0	18	0	0
4	B	13	0	18	1	0
4	E	26	0	36	0	0
4	F	13	0	18	0	0
4	H	13	0	18	0	0
5	B	20	0	28	1	0
5	C	10	0	14	0	0
5	D	10	0	14	0	0
5	F	10	0	14	0	0
5	I	10	0	14	0	0
5	K	20	0	28	0	0
6	E	25	0	34	10	0
7	E	7	0	10	1	0
7	K	14	0	20	0	0
8	E	12	0	18	3	0
8	H	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	A	61	0	0	0	0
9	B	93	0	0	0	0
9	C	79	0	0	0	0
9	D	70	0	0	0	0
9	E	85	0	0	0	0
9	F	93	0	0	0	0
9	G	67	0	0	0	0
9	H	78	0	0	0	0
9	I	45	0	0	0	0
9	J	44	0	0	0	0
9	K	49	0	0	0	0
9	L	56	0	0	0	0
All	All	27128	0	25594	101	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 101 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)
6:E:902:PE8:H22	6:E:902:PE8:H61	1.31	1.11
1:A:32:ASP:HB3	2:A:401:FPP:O3B	1.66	0.96
1:A:37:TYR:OH	2:A:401:FPP:H52	1.73	0.88
1:G:37:TYR:OH	2:G:401:FPP:H52	1.79	0.81
2:A:401:FPP:H101	2:A:401:FPP:C12	2.10	0.79

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	275/303 (91%)	268 (98%)	7 (2%)	0	100	100
1	B	274/303 (90%)	268 (98%)	6 (2%)	0	100	100
1	C	273/303 (90%)	266 (97%)	7 (3%)	0	100	100
1	D	273/303 (90%)	265 (97%)	8 (3%)	0	100	100
1	E	272/303 (90%)	267 (98%)	5 (2%)	0	100	100
1	F	276/303 (91%)	271 (98%)	5 (2%)	0	100	100
1	G	272/303 (90%)	266 (98%)	6 (2%)	0	100	100
1	H	272/303 (90%)	267 (98%)	5 (2%)	0	100	100
1	I	267/303 (88%)	258 (97%)	9 (3%)	0	100	100
1	J	274/303 (90%)	269 (98%)	5 (2%)	0	100	100
1	K	272/303 (90%)	267 (98%)	5 (2%)	0	100	100
1	L	266/303 (88%)	255 (96%)	10 (4%)	1 (0%)	30	34
All	All	3266/3636 (90%)	3187 (98%)	78 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	39	PRO

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	214/234 (92%)	211 (99%)	3 (1%)	59	75
1	B	213/234 (91%)	210 (99%)	3 (1%)	59	75
1	C	212/234 (91%)	210 (99%)	2 (1%)	70	84
1	D	212/234 (91%)	210 (99%)	2 (1%)	70	84
1	E	211/234 (90%)	207 (98%)	4 (2%)	50	66
1	F	215/234 (92%)	214 (100%)	1 (0%)	81	90
1	G	211/234 (90%)	209 (99%)	2 (1%)	70	84

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	211/234 (90%)	211 (100%)	0	100	100
1	I	209/234 (89%)	207 (99%)	2 (1%)	68	81
1	J	213/234 (91%)	209 (98%)	4 (2%)	50	66
1	K	211/234 (90%)	205 (97%)	6 (3%)	38	52
1	L	205/234 (88%)	202 (98%)	3 (2%)	57	73
All	All	2537/2808 (90%)	2505 (99%)	32 (1%)	61	76

5 of 32 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	K	212	ARG
1	L	164	PHE
1	E	200	GLU
1	E	164	PHE
1	L	200	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 27 such sidechains are listed below:

Mol	Chain	Res	Type
1	H	35	HIS
1	I	110	GLN
1	L	192	GLN
1	I	35	HIS
1	I	232	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 oligosaccharides in this entry.

5.6 Ligand geometry

37 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$
8	EDO	H	903	-	3,3,3	0.06	0	2,2,2	0.11	0
4	PG4	H	902	-	12,12,12	0.16	0	11,11,11	0.11	0
7	PEG	K	904	-	6,6,6	0.12	0	5,5,5	0.07	0
3	SAH	A	402	-	27,28,28	0.49	1 (3%)	36,40,40	0.58	1 (2%)
4	PG4	E	903	-	12,12,12	0.17	0	11,11,11	0.10	0
3	SAH	D	901	-	27,28,28	0.49	1 (3%)	36,40,40	0.55	1 (2%)
3	SAH	E	901	-	27,28,28	0.49	1 (3%)	36,40,40	0.58	1 (2%)
2	FPP	A	401	-	22,23,23	0.62	0	27,31,31	0.64	0
5	PGE	K	903	-	9,9,9	0.19	0	8,8,8	0.10	0
3	SAH	B	901	-	27,28,28	0.49	1 (3%)	36,40,40	0.59	1 (2%)
7	PEG	K	905	-	6,6,6	0.13	0	5,5,5	0.08	0
3	SAH	L	901	-	27,28,28	0.49	1 (3%)	36,40,40	0.55	1 (2%)
3	SAH	J	901	-	27,28,28	0.49	1 (3%)	36,40,40	0.56	1 (2%)
4	PG4	E	904	-	12,12,12	0.16	0	11,11,11	0.11	0
6	PE8	E	902	-	24,24,24	0.16	0	23,23,23	0.20	0
5	PGE	C	902	-	9,9,9	0.14	0	8,8,8	0.11	0
2	FPP	G	401	-	22,23,23	0.66	1 (4%)	27,31,31	0.66	0
5	PGE	K	902	-	9,9,9	0.14	0	8,8,8	0.11	0
8	EDO	E	906	-	3,3,3	0.06	0	2,2,2	0.12	0
5	PGE	F	404	-	9,9,9	0.14	0	8,8,8	0.10	0
5	PGE	B	904	-	9,9,9	0.14	0	8,8,8	0.11	0
7	PEG	E	905	-	6,6,6	0.12	0	5,5,5	0.06	0
3	SAH	H	901	-	27,28,28	0.49	1 (3%)	36,40,40	0.58	1 (2%)
3	SAH	F	402	-	27,28,28	0.49	1 (3%)	36,40,40	0.55	1 (2%)
4	PG4	F	403	-	12,12,12	0.15	0	11,11,11	0.13	0
3	SAH	K	901	-	27,28,28	0.49	1 (3%)	36,40,40	0.57	1 (2%)
2	FPP	F	401	-	22,23,23	0.60	0	27,31,31	0.67	0
3	SAH	I	901	-	27,28,28	0.48	1 (3%)	36,40,40	0.54	1 (2%)
5	PGE	D	902	-	9,9,9	0.14	0	8,8,8	0.11	0
8	EDO	E	908	-	3,3,3	0.04	0	2,2,2	0.16	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	PGE	I	902	-	9,9,9	0.15	0	8,8,8	0.10	0
3	SAH	G	402	-	27,28,28	0.49	1 (3%)	36,40,40	0.56	1 (2%)
5	PGE	B	903	-	9,9,9	0.15	0	8,8,8	0.11	0
8	EDO	E	907	-	3,3,3	0.06	0	2,2,2	0.15	0
4	PG4	B	902	-	12,12,12	0.15	0	11,11,11	0.11	0
4	PG4	A	403	-	12,12,12	0.21	0	11,11,11	0.18	0
3	SAH	C	901	-	27,28,28	0.49	1 (3%)	36,40,40	0.58	1 (2%)

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
8	EDO	H	903	-	-	0/1/1/1	-
4	PG4	H	902	-	-	10/10/10/10	-
7	PEG	K	904	-	-	4/4/4/4	-
3	SAH	A	402	-	-	0/15/31/31	0/3/3/3
4	PG4	E	903	-	-	6/10/10/10	-
3	SAH	D	901	-	-	0/15/31/31	0/3/3/3
3	SAH	E	901	-	-	2/15/31/31	0/3/3/3
2	FPP	A	401	-	-	9/25/25/25	-
5	PGE	K	903	-	-	4/7/7/7	-
3	SAH	B	901	-	-	2/15/31/31	0/3/3/3
7	PEG	K	905	-	-	3/4/4/4	-
3	SAH	L	901	-	-	0/15/31/31	0/3/3/3
3	SAH	J	901	-	-	1/15/31/31	0/3/3/3
4	PG4	E	904	-	-	8/10/10/10	-
6	PE8	E	902	-	-	12/22/22/22	-
5	PGE	C	902	-	-	6/7/7/7	-
2	FPP	G	401	-	-	12/25/25/25	-
5	PGE	K	902	-	-	4/7/7/7	-
8	EDO	E	906	-	-	1/1/1/1	-
5	PGE	F	404	-	-	3/7/7/7	-
5	PGE	B	904	-	-	4/7/7/7	-
7	PEG	E	905	-	-	3/4/4/4	-
3	SAH	H	901	-	-	0/15/31/31	0/3/3/3
3	SAH	F	402	-	-	0/15/31/31	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	PG4	F	403	-	-	8/10/10/10	-
3	SAH	K	901	-	-	0/15/31/31	0/3/3/3
2	FPP	F	401	-	-	13/25/25/25	-
3	SAH	I	901	-	-	1/15/31/31	0/3/3/3
5	PGE	D	902	-	-	5/7/7/7	-
8	EDO	E	908	-	-	1/1/1/1	-
5	PGE	I	902	-	-	5/7/7/7	-
3	SAH	G	402	-	-	0/15/31/31	0/3/3/3
5	PGE	B	903	-	-	4/7/7/7	-
8	EDO	E	907	-	-	0/1/1/1	-
4	PG4	B	902	-	-	3/10/10/10	-
4	PG4	A	403	-	-	7/10/10/10	-
3	SAH	C	901	-	-	1/15/31/31	0/3/3/3

The worst 5 of 13 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	401	FPP	PA-O3A	2.29	1.62	1.59
3	K	901	SAH	OXT-C	-2.26	1.23	1.30
3	J	901	SAH	OXT-C	-2.26	1.23	1.30
3	G	402	SAH	OXT-C	-2.25	1.23	1.30
3	C	901	SAH	OXT-C	-2.25	1.23	1.30

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	901	SAH	OXT-C-O	-2.67	118.02	124.08
3	B	901	SAH	OXT-C-O	-2.66	118.05	124.08
3	K	901	SAH	OXT-C-O	-2.65	118.06	124.08
3	A	402	SAH	OXT-C-O	-2.65	118.07	124.08
3	C	901	SAH	OXT-C-O	-2.65	118.07	124.08

There are no chirality outliers.

5 of 142 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	FPP	C1-O1-PA-O2A
2	A	401	FPP	PA-O3A-PB-O2B
2	A	401	FPP	PA-O3A-PB-O3B
2	F	401	FPP	C1-O1-PA-O1A

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Mol	Chain	Res	Type	Atoms
2	F	401	FPP	C1-O1-PA-O2A

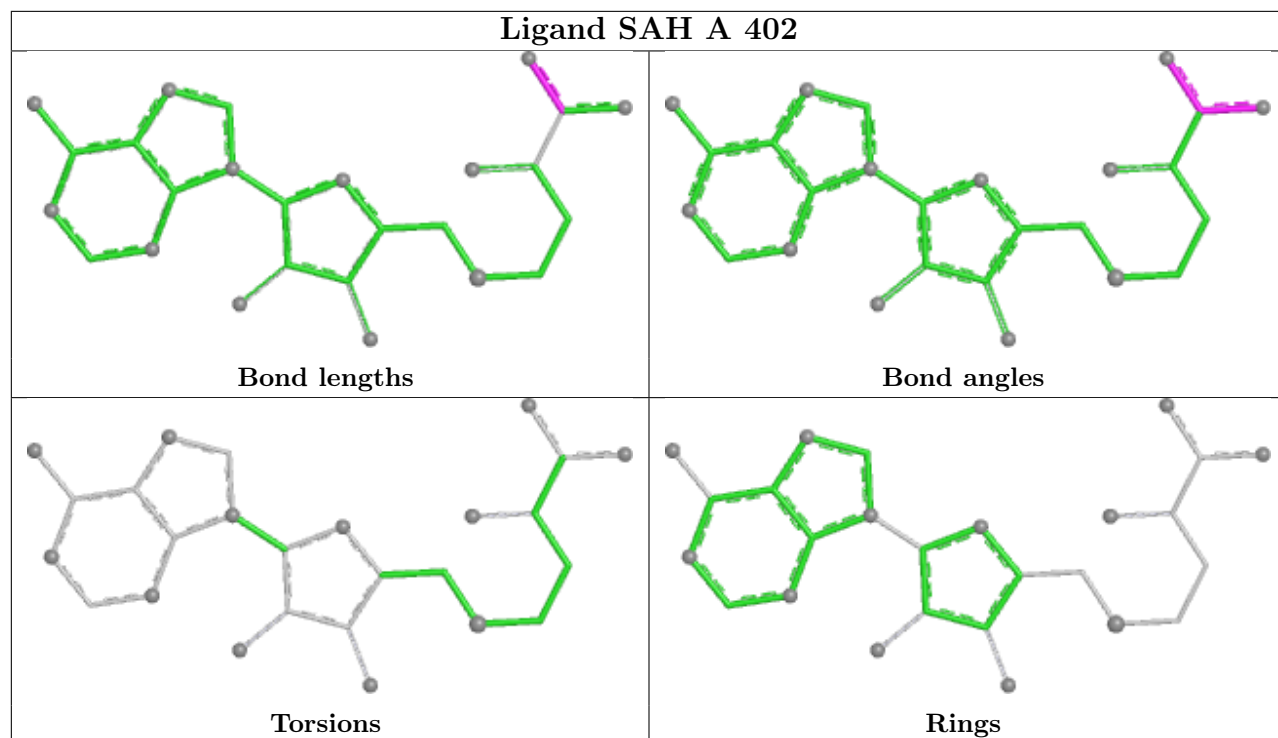
There are no ring outliers.

9 monomers are involved in 28 short contacts:

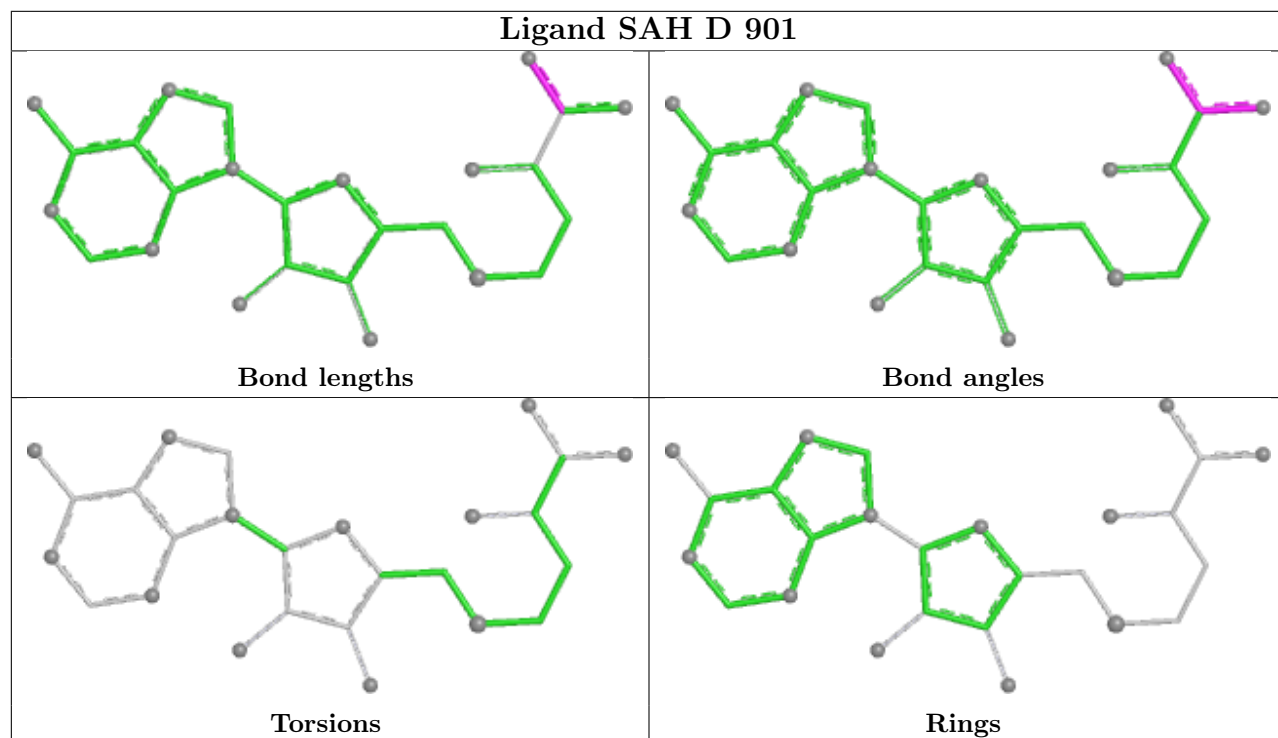
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	FPP	6	0
6	E	902	PE8	10	0
2	G	401	FPP	3	0
7	E	905	PEG	1	0
2	F	401	FPP	4	0
8	E	908	EDO	2	0
5	B	903	PGE	1	0
8	E	907	EDO	1	0
4	B	902	PG4	1	0

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.

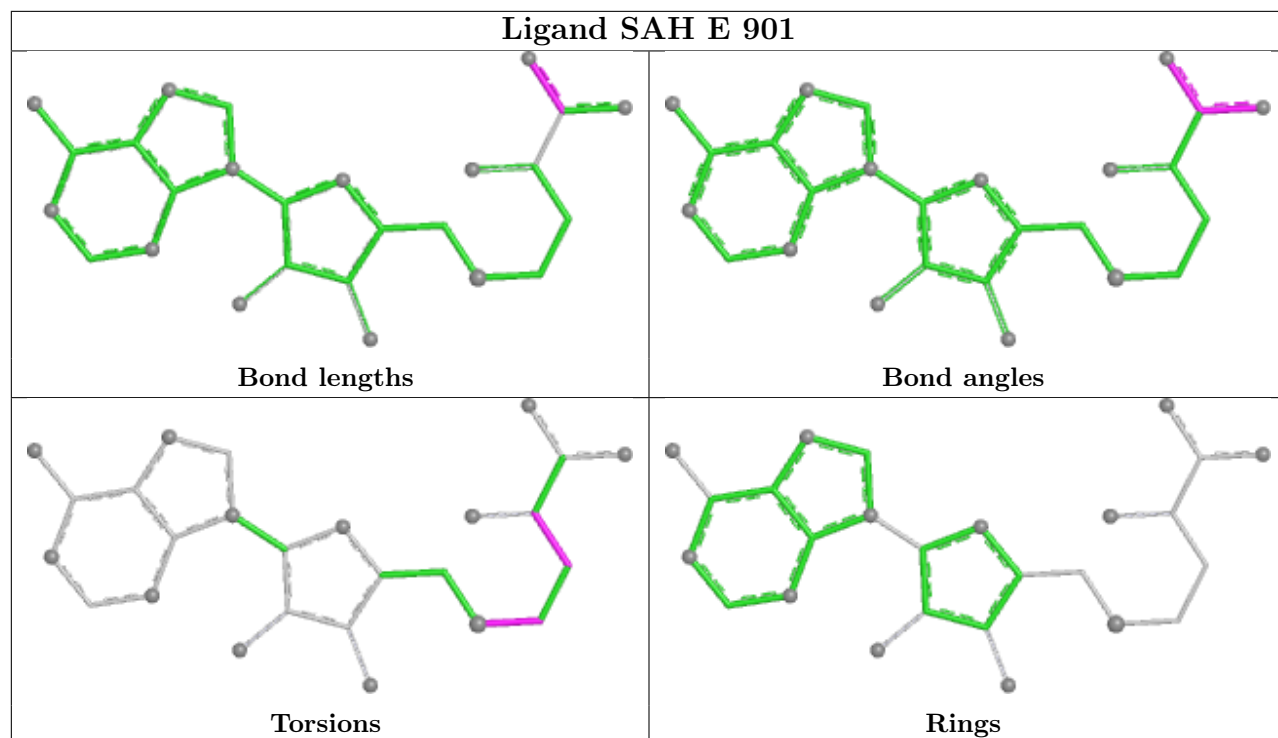
Ligand SAH A 402



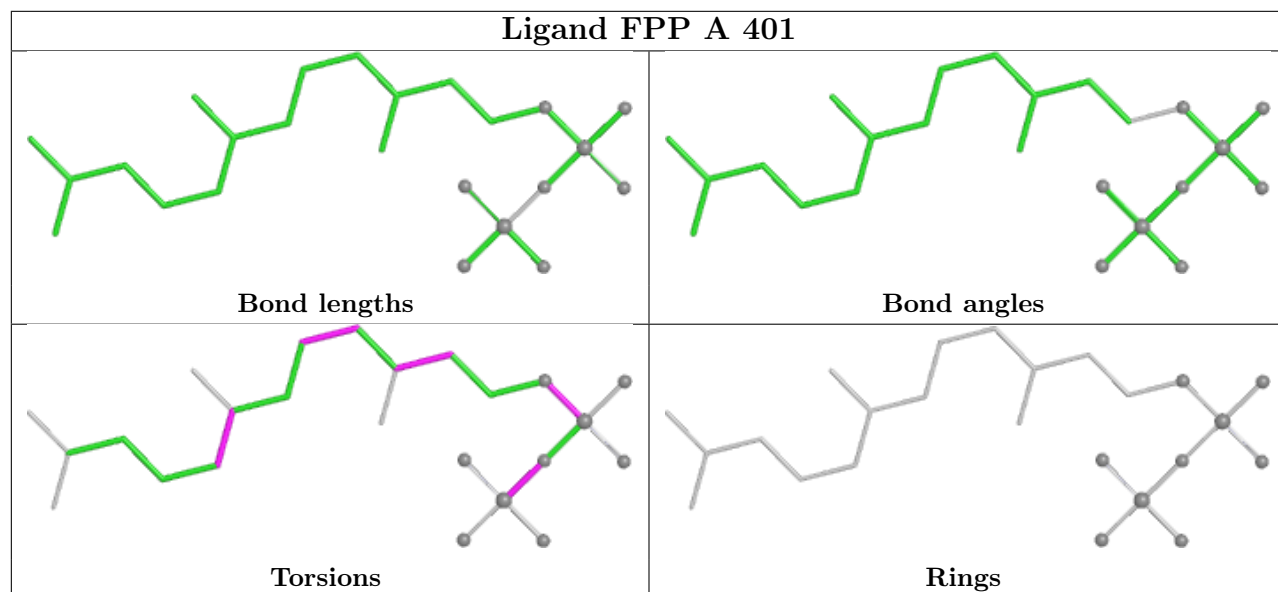
Ligand SAH D 901



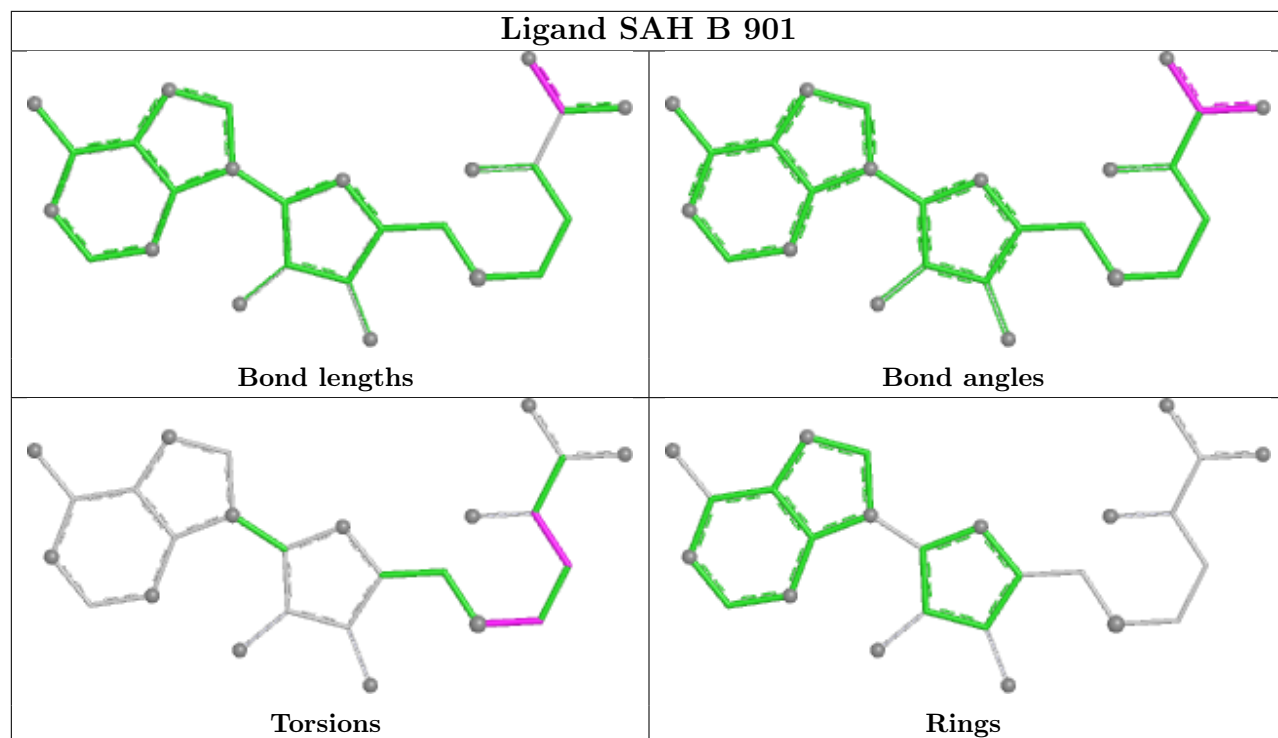
Ligand SAH E 901



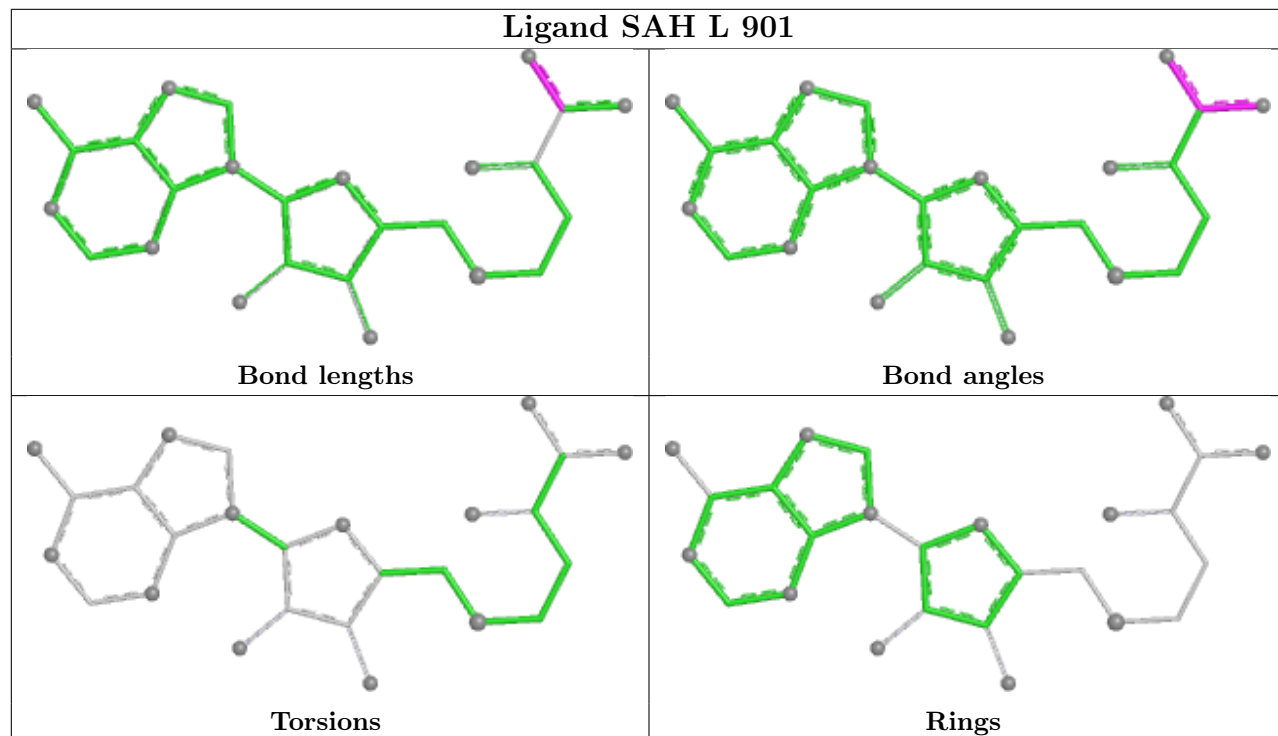
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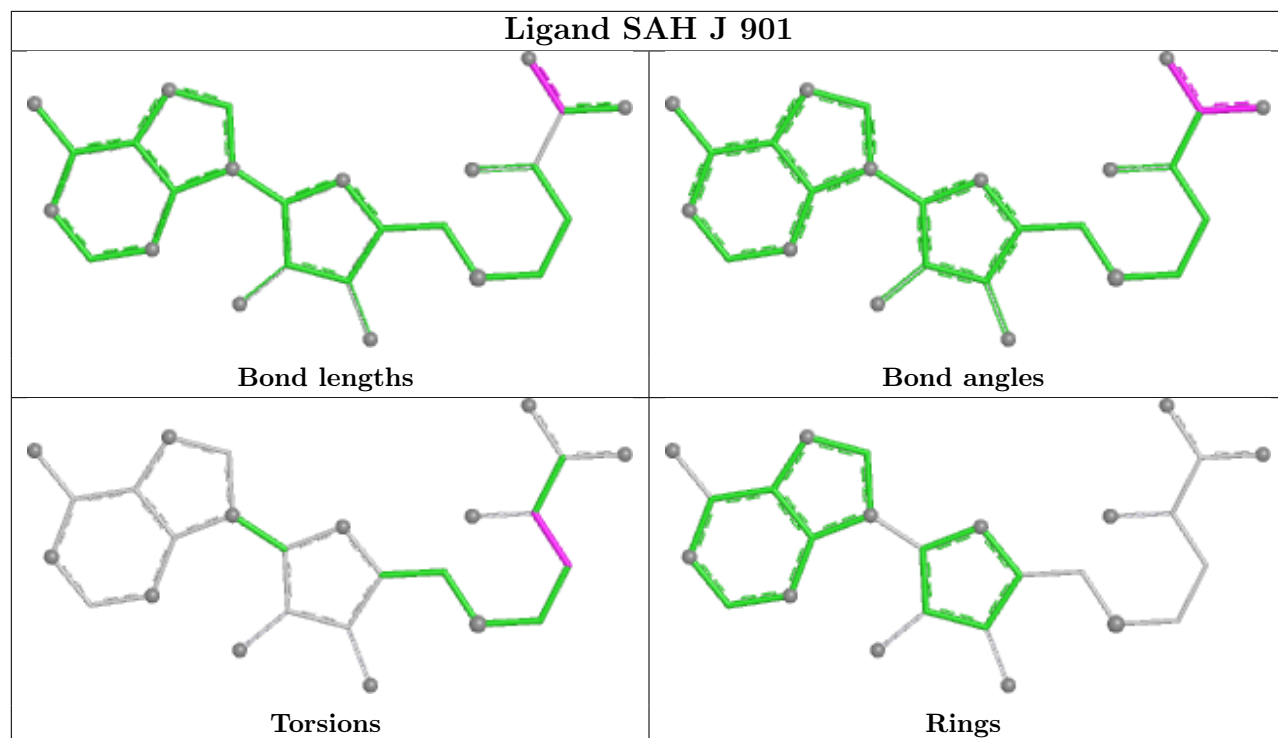
Ligand SAH B 901



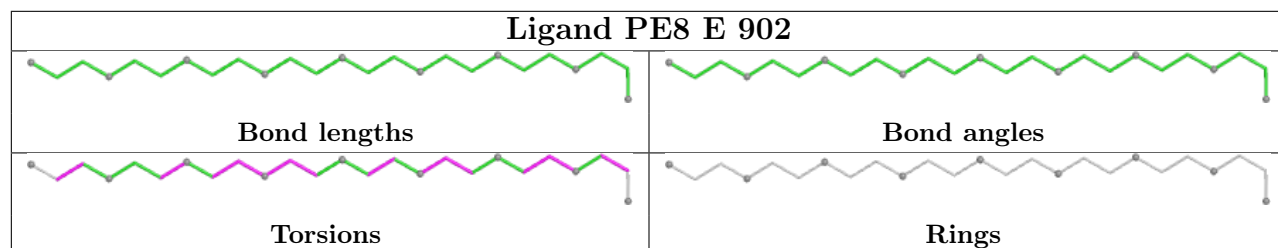
Ligand SAH L 901



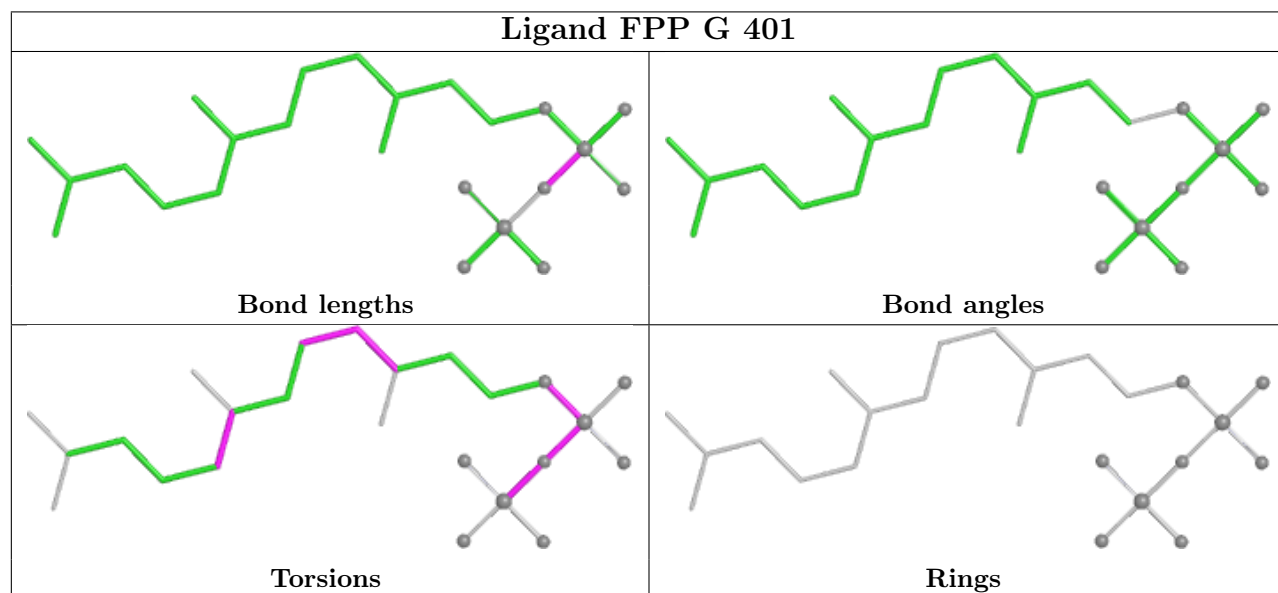
Ligand SAH J 901



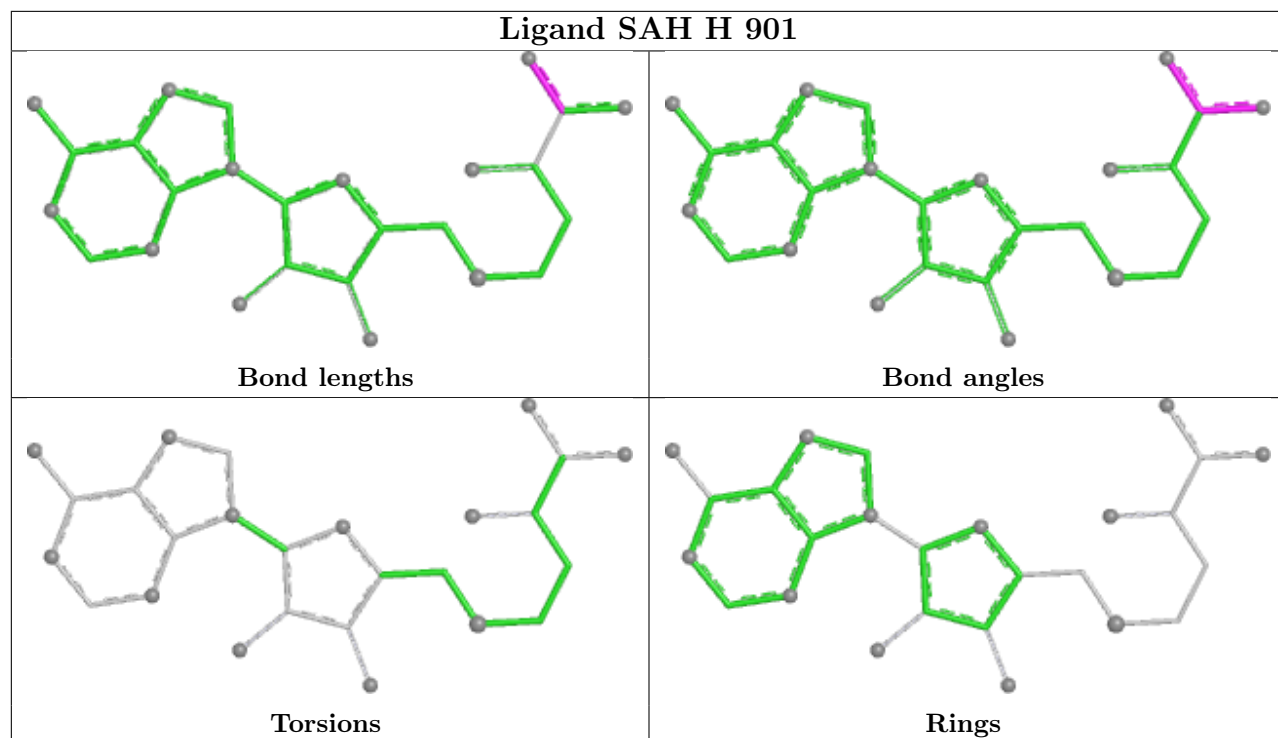
Ligand PE8 E 902



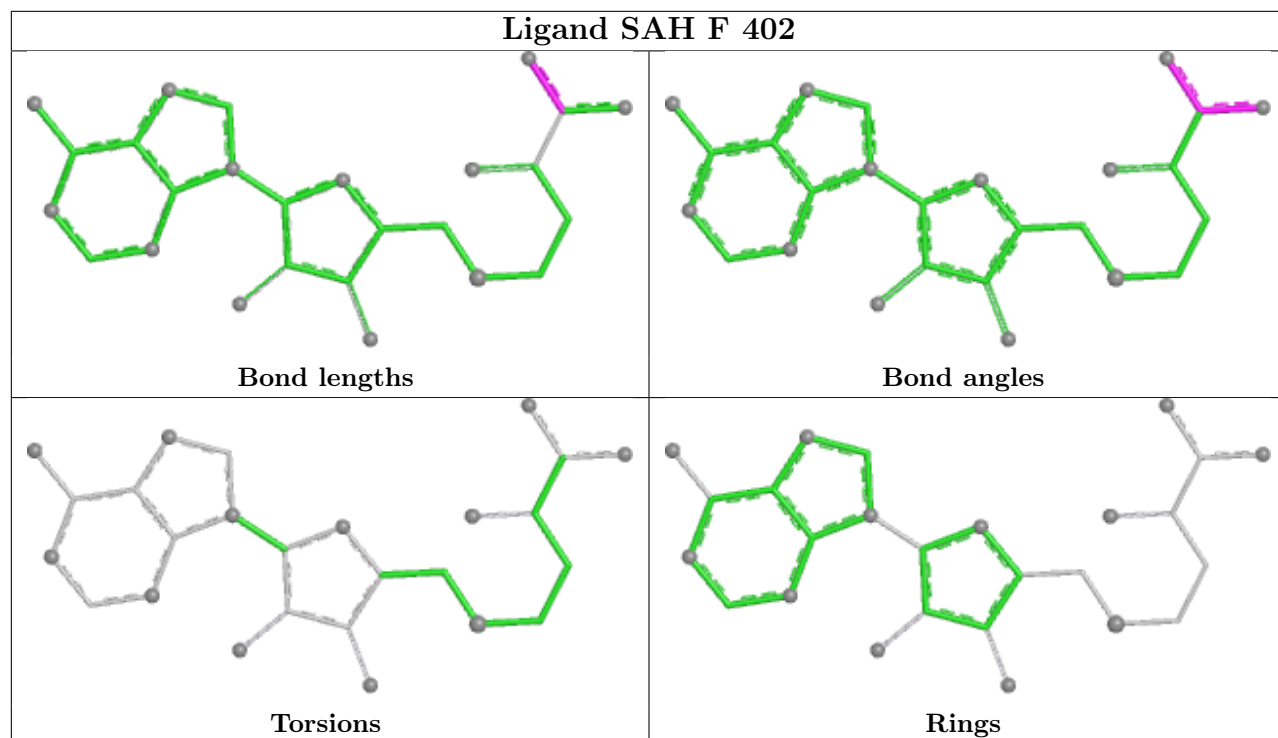
Ligand FPP G 401



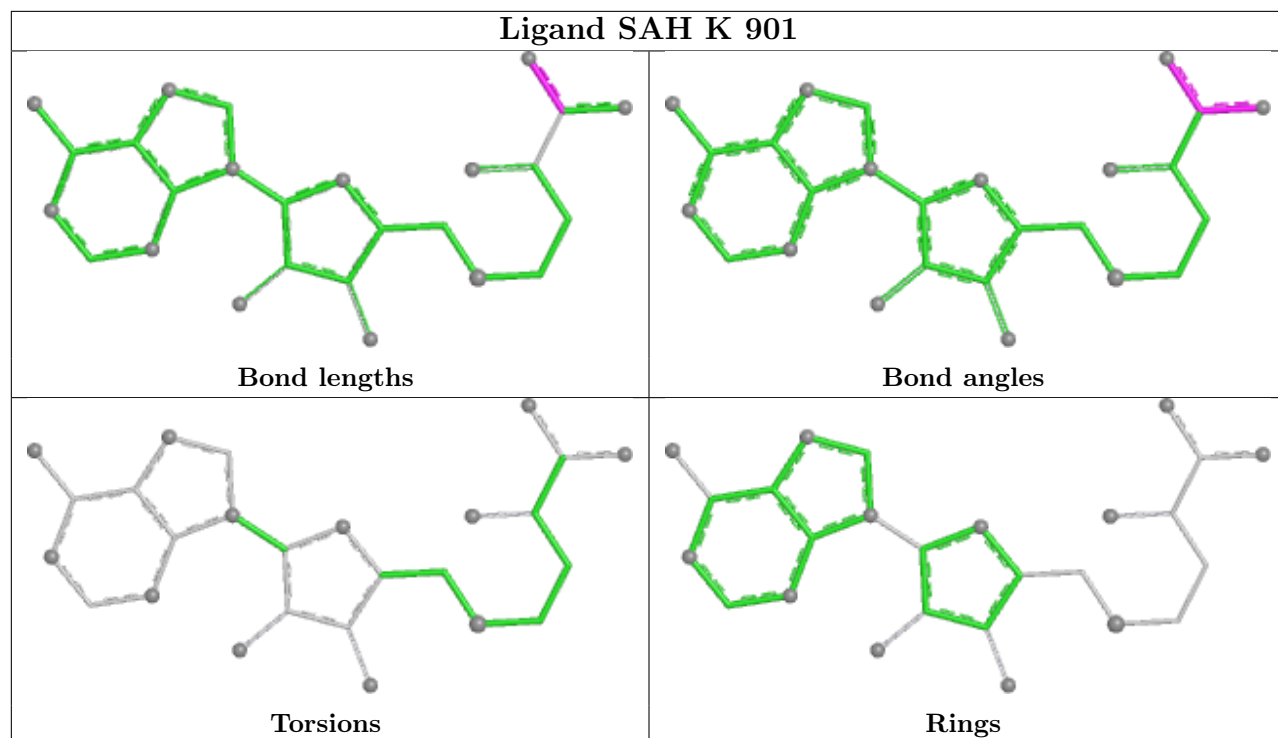
Ligand SAH H 901



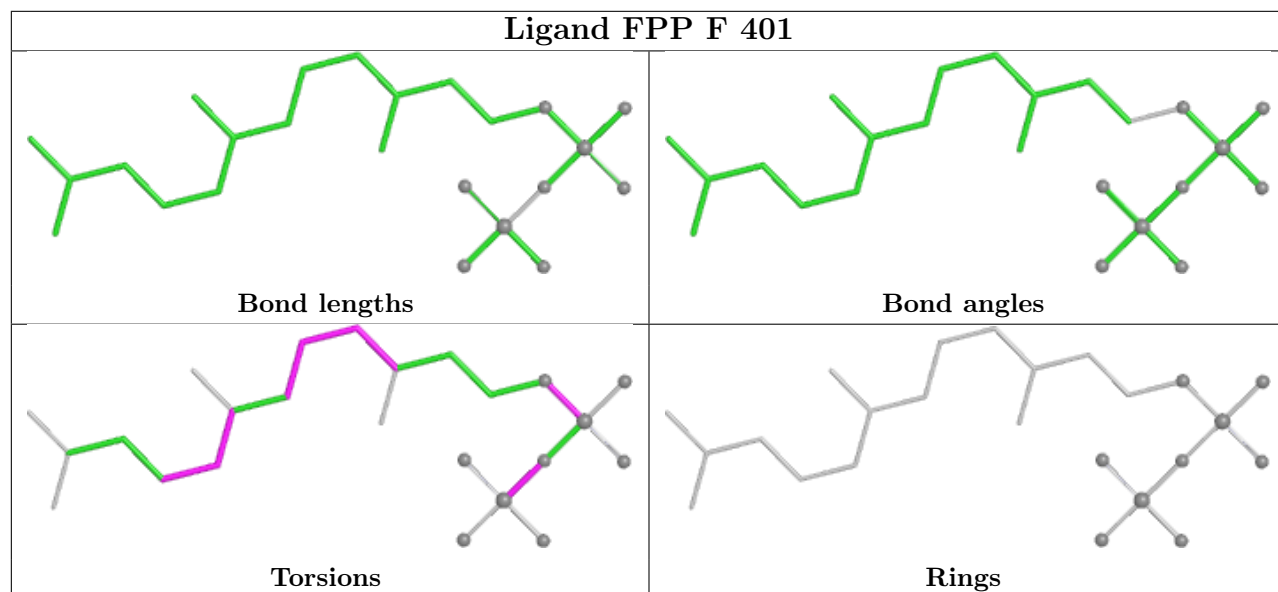
Ligand SAH F 402



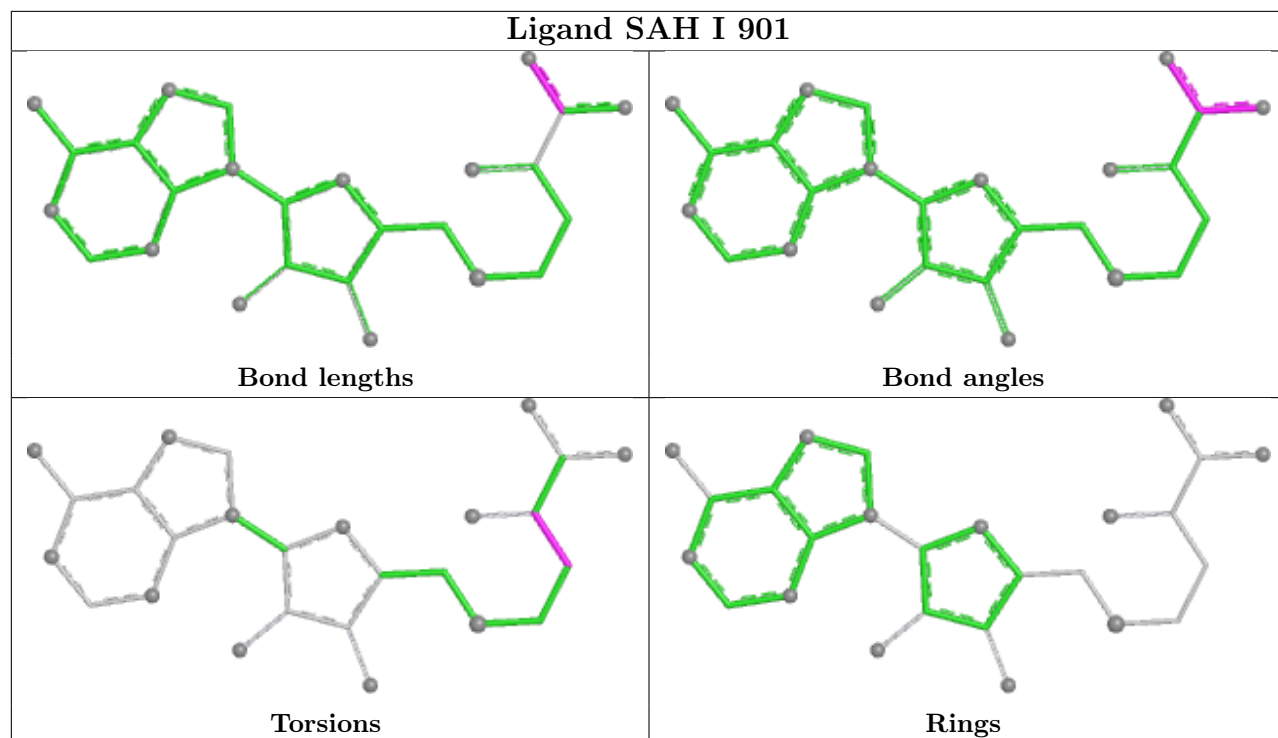
Ligand SAH K 901



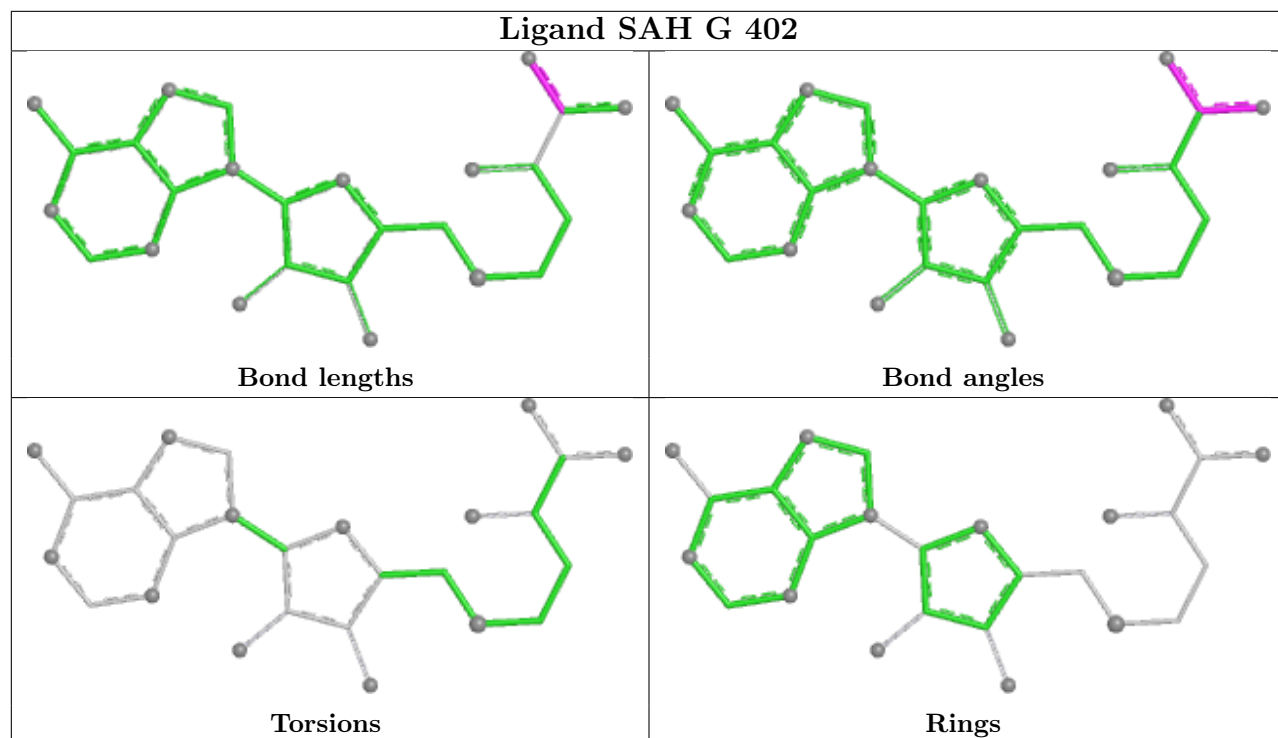
Ligand FPP F 401

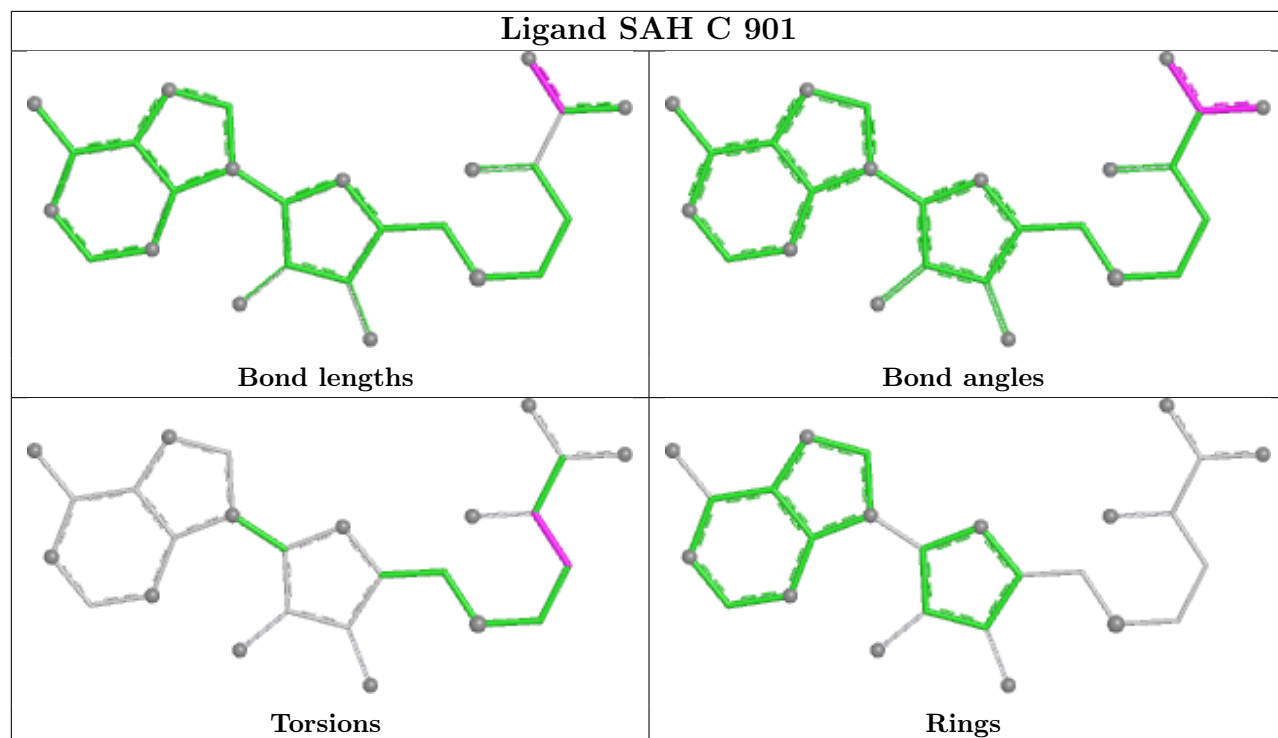


Ligand SAH I 901



Ligand SAH G 402





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	277/303 (91%)	0.11	8 (2%) 53 50	26, 36, 54, 87	0
1	B	276/303 (91%)	0.02	18 (6%) 25 22	23, 29, 59, 100	0
1	C	275/303 (90%)	-0.00	4 (1%) 72 69	23, 32, 56, 92	0
1	D	275/303 (90%)	0.10	9 (3%) 49 46	24, 34, 65, 101	0
1	E	274/303 (90%)	0.13	17 (6%) 26 23	25, 32, 68, 100	0
1	F	278/303 (91%)	-0.09	4 (1%) 73 71	23, 30, 51, 81	0
1	G	274/303 (90%)	0.02	5 (1%) 67 64	23, 35, 54, 90	0
1	H	274/303 (90%)	0.20	17 (6%) 26 23	25, 35, 67, 103	0
1	I	271/303 (89%)	0.39	14 (5%) 33 29	31, 44, 71, 92	0
1	J	276/303 (91%)	0.41	8 (2%) 53 50	30, 43, 74, 104	0
1	K	274/303 (90%)	0.37	19 (6%) 23 20	30, 40, 75, 108	0
1	L	268/303 (88%)	0.33	15 (5%) 30 27	29, 41, 73, 111	0
All	All	3292/3636 (90%)	0.16	138 (4%) 40 37	23, 36, 67, 111	0

The worst 5 of 138 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	37	TYR	5.7
1	L	34	LEU	5.7
1	H	45	PHE	4.9
1	L	262	GLY	4.8
1	H	29	LYS	4.7

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides 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($\text{\AA}^2$)	Q<0.9
6	PE8	E	902	25/25	0.59	0.18	45,50,52,53	0
7	PEG	K	905	7/7	0.61	0.17	52,52,53,53	0
5	PGE	K	903	10/10	0.64	0.18	49,49,51,51	0
8	EDO	E	907	4/4	0.64	0.20	56,57,58,58	0
4	PG4	E	903	13/13	0.69	0.16	47,48,49,50	0
4	PG4	B	902	13/13	0.71	0.18	57,63,64,64	0
7	PEG	E	905	7/7	0.74	0.18	58,61,62,63	0
7	PEG	K	904	7/7	0.75	0.14	50,50,51,52	0
5	PGE	B	903	10/10	0.76	0.16	35,39,42,44	0
2	FPP	F	401	24/24	0.80	0.17	38,49,59,62	8
4	PG4	A	403	13/13	0.81	0.13	44,45,49,50	0
4	PG4	H	902	13/13	0.81	0.15	60,62,63,64	0
2	FPP	G	401	24/24	0.81	0.19	42,53,66,73	8
5	PGE	K	902	10/10	0.84	0.12	45,46,46,47	0
8	EDO	E	908	4/4	0.84	0.19	38,40,40,42	0
5	PGE	D	902	10/10	0.85	0.15	52,53,53,54	0
5	PGE	I	902	10/10	0.85	0.16	57,58,59,59	0
8	EDO	H	903	4/4	0.87	0.18	52,52,52,53	0
4	PG4	F	403	13/13	0.88	0.11	42,43,45,45	0
5	PGE	F	404	10/10	0.88	0.12	40,40,42,42	0
4	PG4	E	904	13/13	0.90	0.12	47,49,54,55	0
2	FPP	A	401	24/24	0.90	0.13	37,45,53,58	8
5	PGE	B	904	10/10	0.91	0.11	52,53,55,55	0
5	PGE	C	902	10/10	0.91	0.12	47,51,52,52	0
3	SAH	J	901	26/26	0.92	0.10	40,45,46,46	0
8	EDO	E	906	4/4	0.93	0.09	46,46,47,47	0
3	SAH	I	901	26/26	0.93	0.09	40,43,45,46	0
3	SAH	C	901	26/26	0.94	0.08	27,31,33,33	0
3	SAH	K	901	26/26	0.95	0.07	32,33,34,34	0
3	SAH	L	901	26/26	0.95	0.07	34,36,39,39	0
3	SAH	D	901	26/26	0.95	0.08	30,35,36,36	0
3	SAH	H	901	26/26	0.95	0.07	29,30,31,31	0

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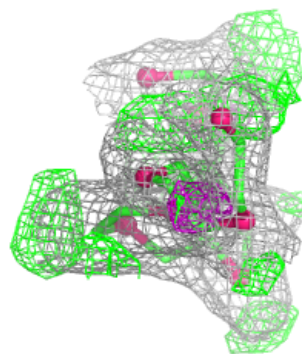
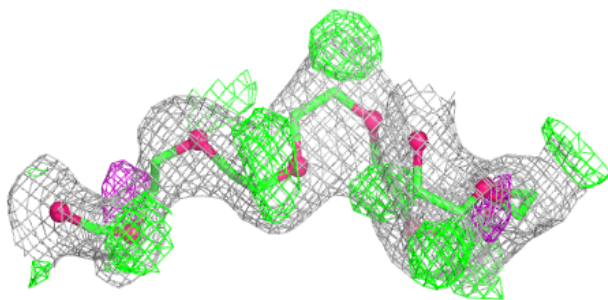
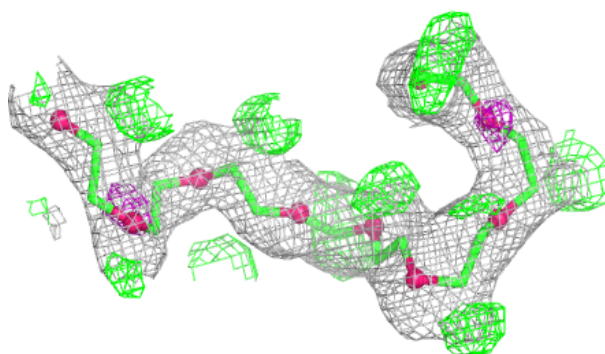
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SAH	F	402	26/26	0.96	0.07	23,24,25,25	0
3	SAH	A	402	26/26	0.96	0.06	27,33,34,35	0
3	SAH	B	901	26/26	0.96	0.06	25,26,26,26	0
3	SAH	E	901	26/26	0.96	0.06	27,29,30,30	0
3	SAH	G	402	26/26	0.97	0.06	30,33,33,34	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.

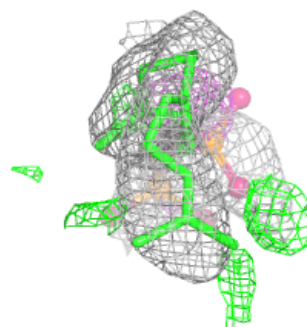
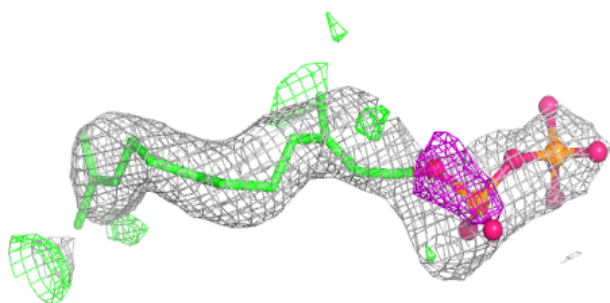
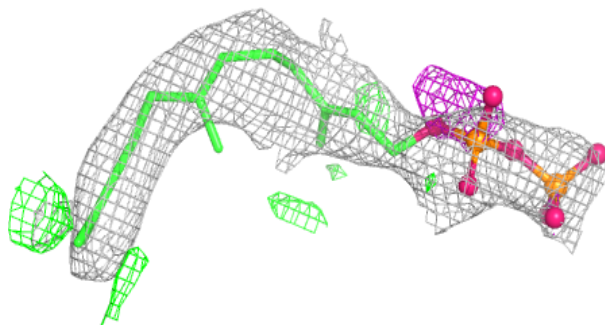
Electron density around PE8 E 902:

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

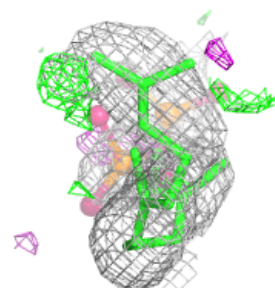
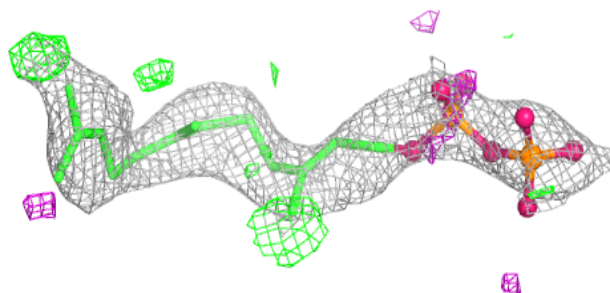
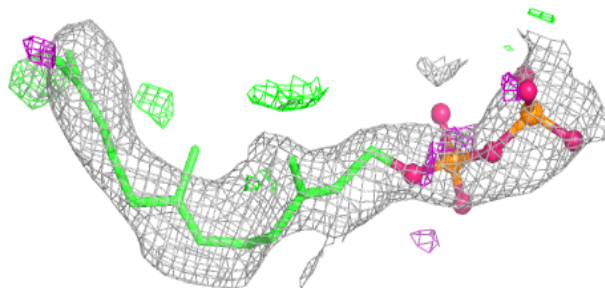


Electron density around FPP F 401:

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

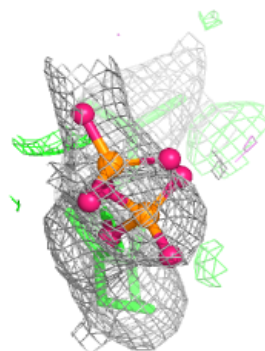
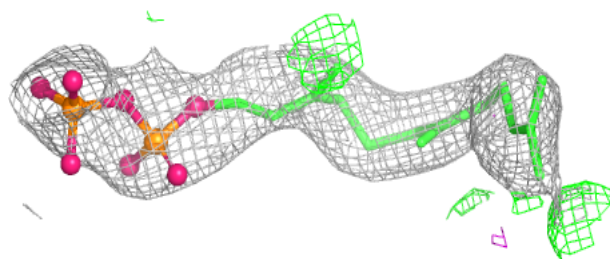
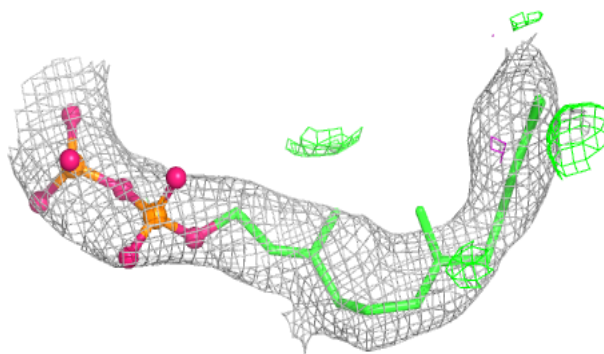
**Electron density around FPP G 401:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

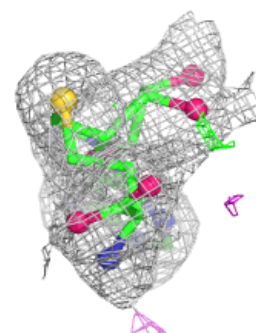
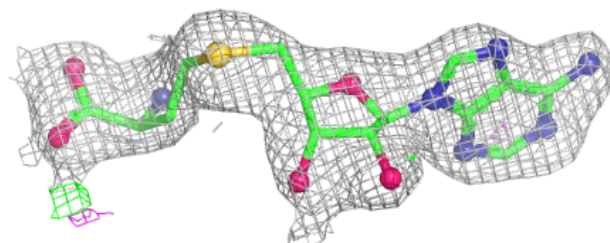
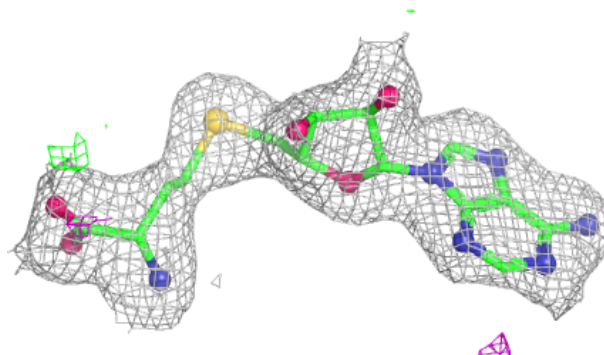


Electron density around FPP A 401:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

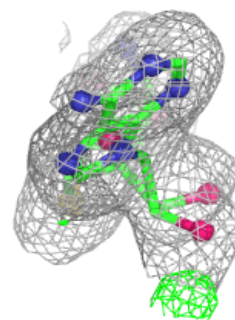
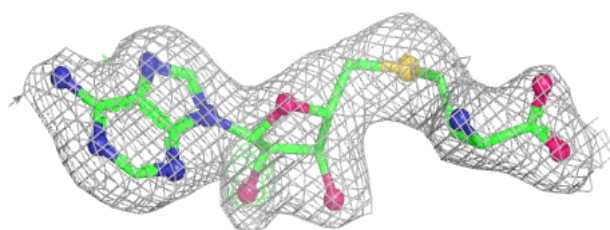
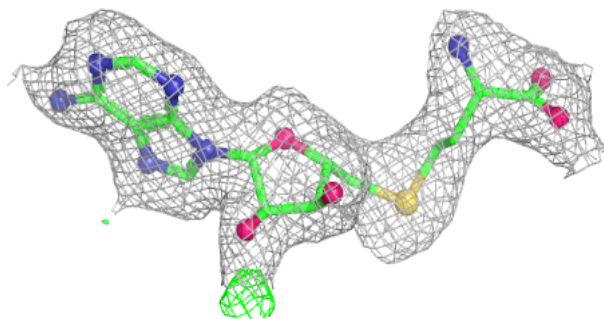
**Electron density around SAH J 901:**

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

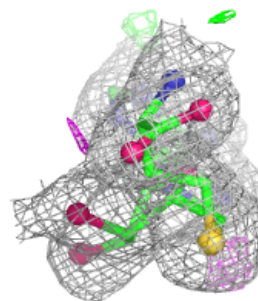
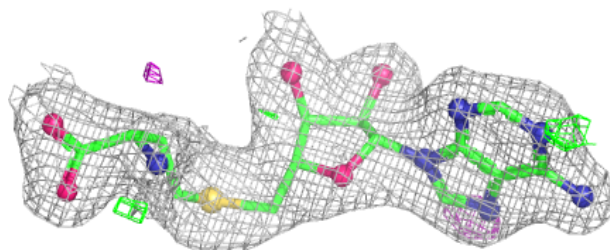
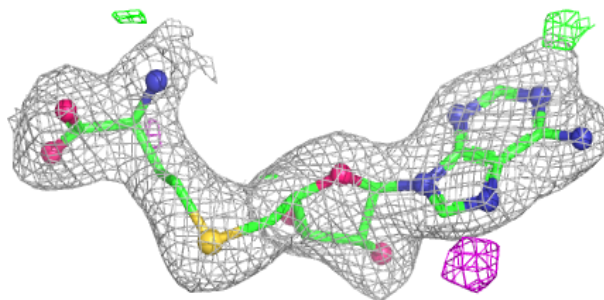


Electron density around SAH I 901:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

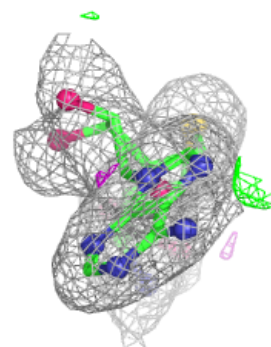
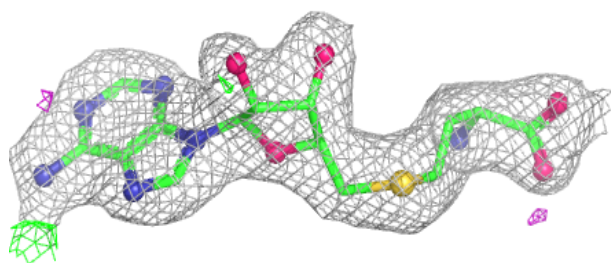
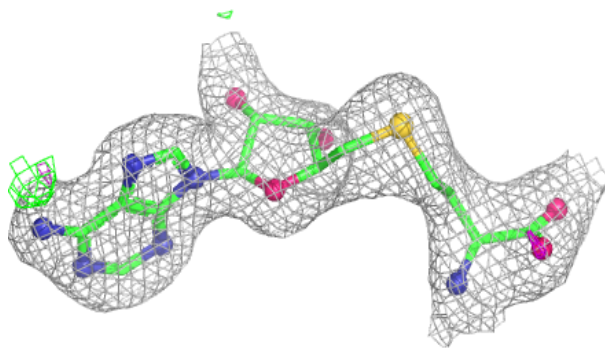
**Electron density around SAH C 901:**

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

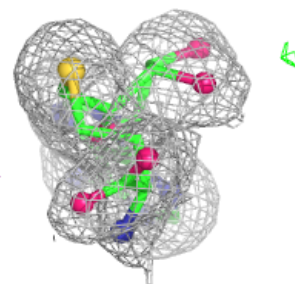
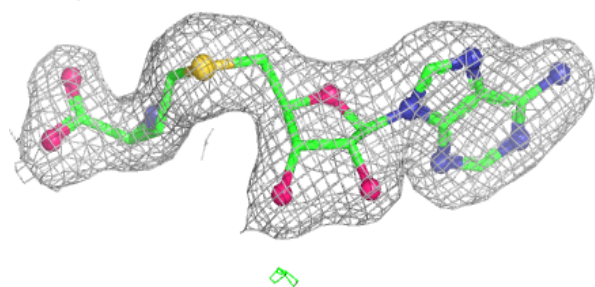
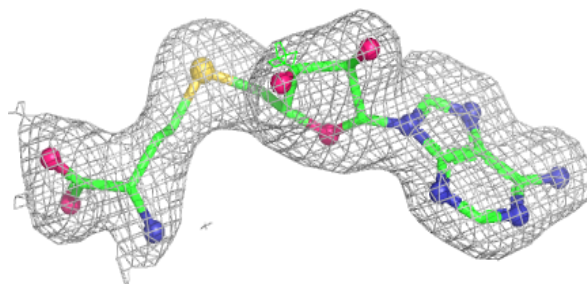


Electron density around SAH K 901:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

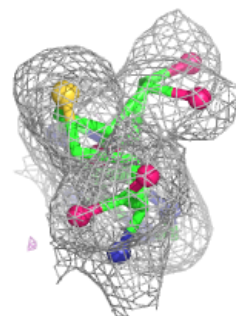
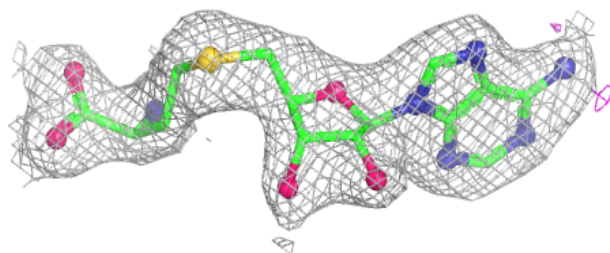
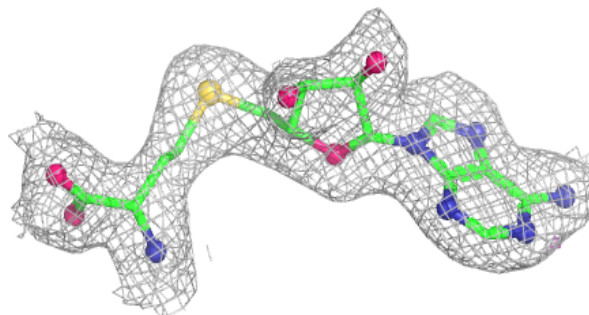
**Electron density around SAH L 901:**

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and green (positive)

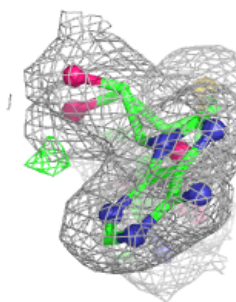
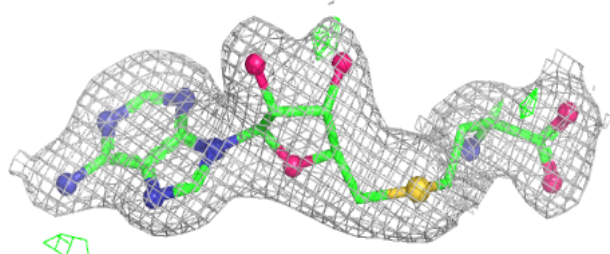
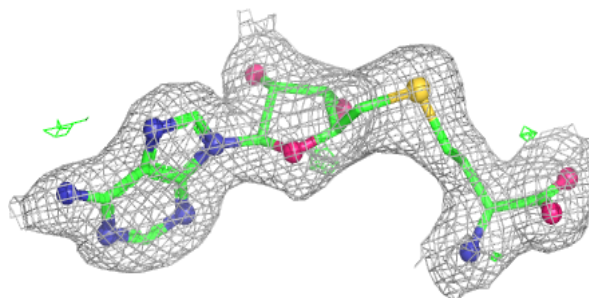


Electron density around SAH D 901:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

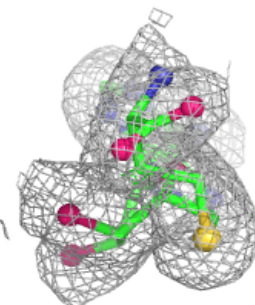
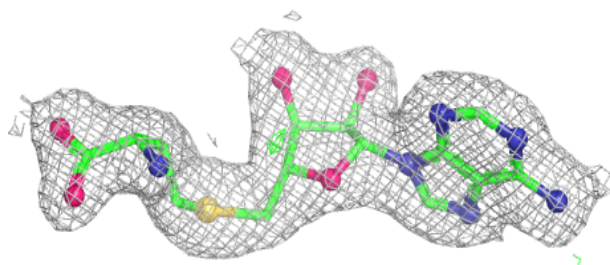
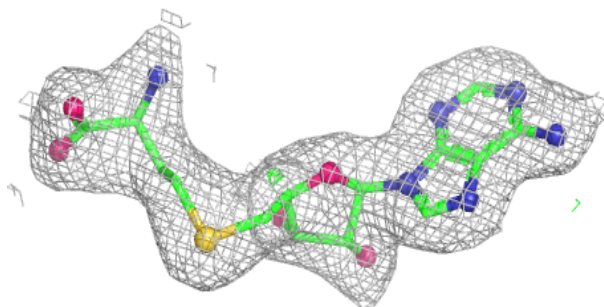
**Electron density around SAH H 901:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

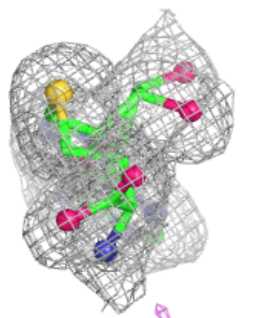
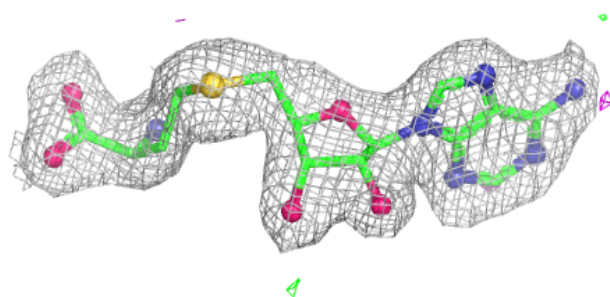
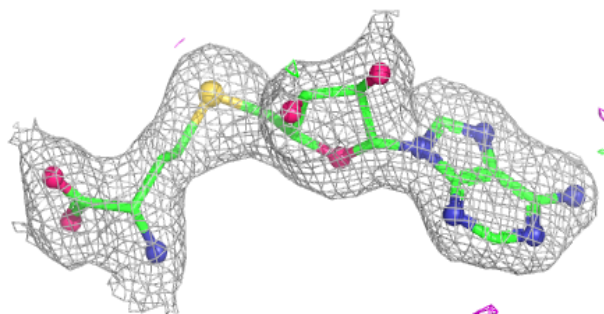


Electron density around SAH F 402:

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

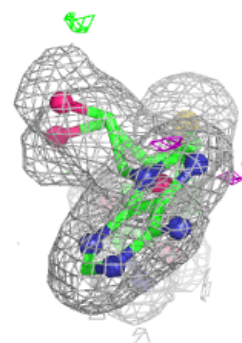
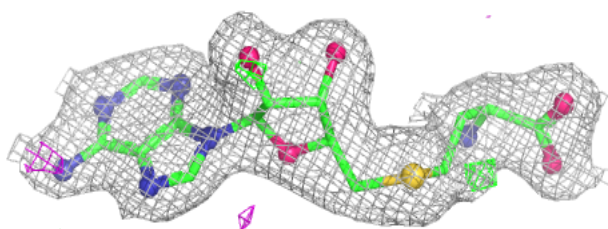
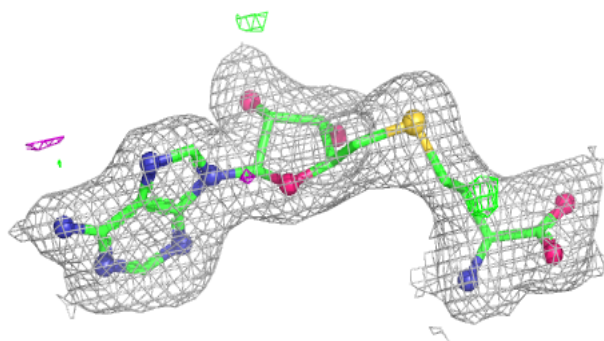
**Electron density around SAH A 402:**

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

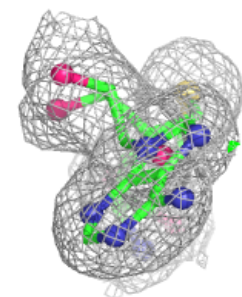
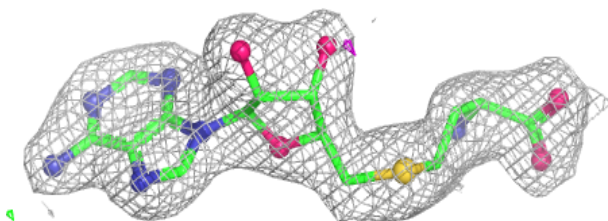
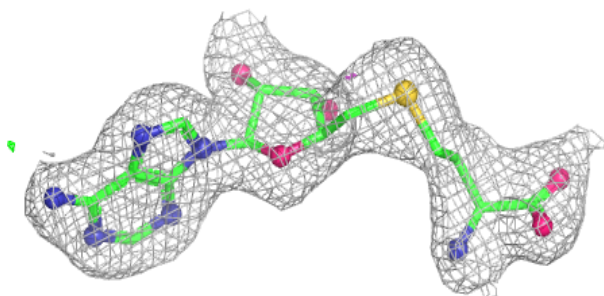


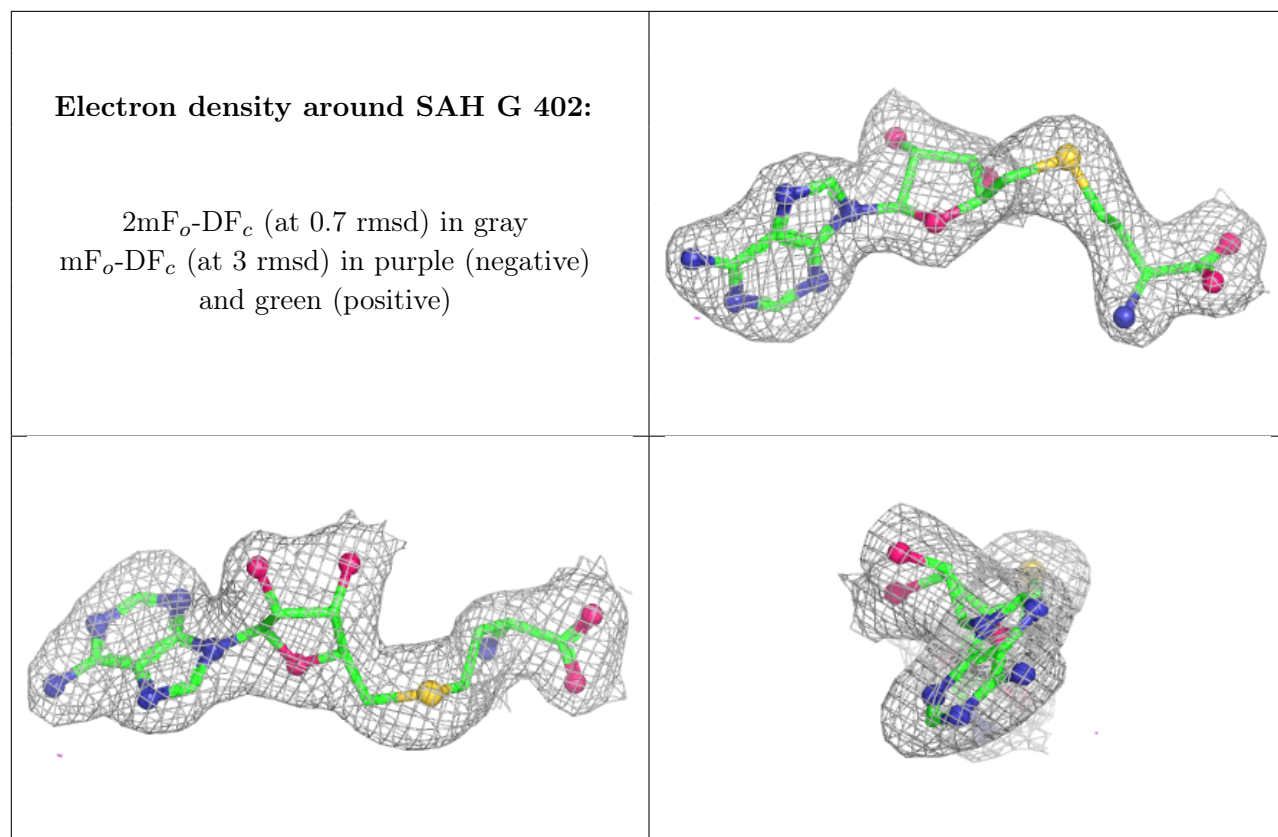
Electron density around SAH B 901:

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

**Electron density around SAH E 901:**

$2mF_o-DF_c$ (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.