Full wwPDB X-ray Structure Validation Report (i)

Jan 5, 2024 – 02:07 am GMT

PDB ID	:	5FEX
Title	:	HydE from T. maritima in complex with Se-adenosyl-L-selenocysteine (tfinal of the reaction)
Authors	:	Rohac, R.; Amara, P.; Benjdia, A.; Martin, L.; Ruffie, P.; Favier, A.; Berteau, O.; Mouesca, J.M.; Fontecilla-Camps, J.C.; Nicolet, Y.
Deposited on	:	2015-12-17
Resolution	:	1.32 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 (i) 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 (i)) were used in the production of this report:

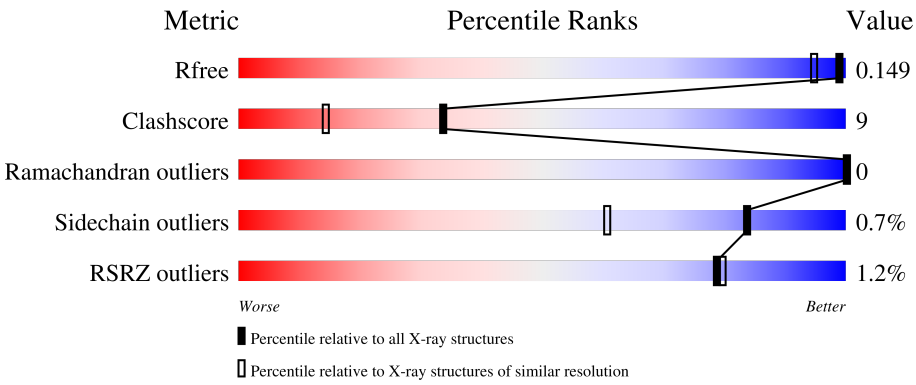
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

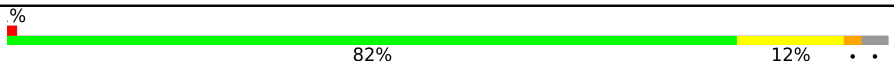
The reported resolution of this entry is 1.32 Å.

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	1611 (1.34-1.30)
Clashscore	141614	1667 (1.34-1.30)
Ramachandran outliers	138981	1615 (1.34-1.30)
Sidechain outliers	138945	1615 (1.34-1.30)
RSRZ outliers	127900	1580 (1.34-1.30)

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 <=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	358	

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
1	OTY	A	114[B]	-	-	X	-

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 3593 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.

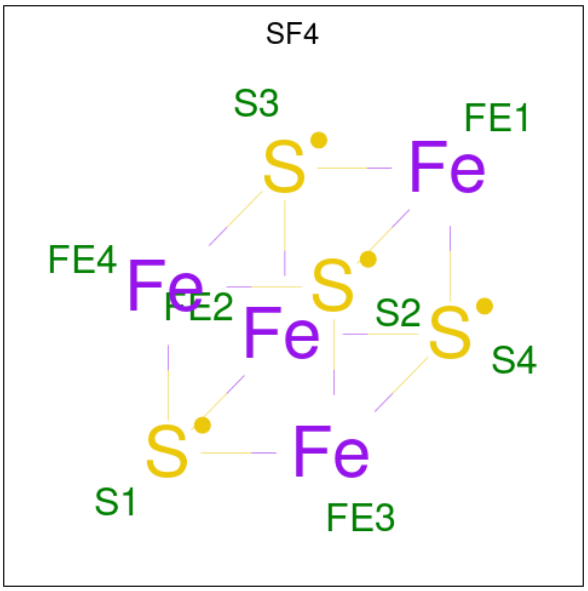
- Molecule 1 is a protein called [FeFe] hydrogenase maturase subunit HydE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	346	2887	1843	496	529	19	0	21	0

There are 14 discrepancies between the modelled and reference sequences:

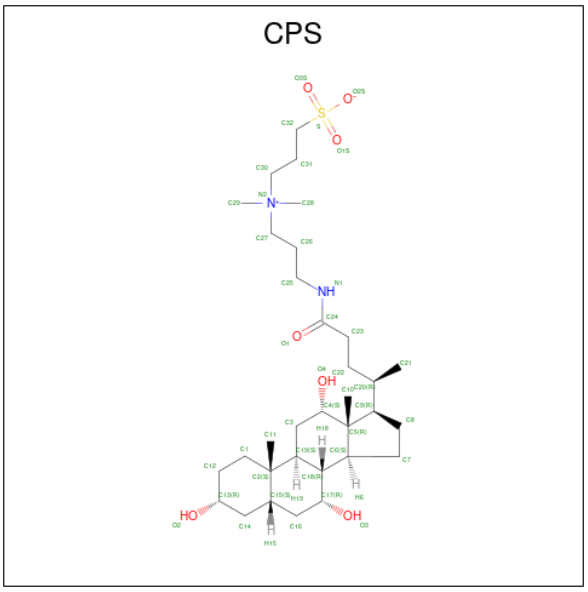
Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	MET	-	initiating methionine	UNP Q9X0Z6
A	-8	TRP	-	expression tag	UNP Q9X0Z6
A	-7	SER	-	expression tag	UNP Q9X0Z6
A	-6	HIS	-	expression tag	UNP Q9X0Z6
A	-5	PRO	-	expression tag	UNP Q9X0Z6
A	-4	GLN	-	expression tag	UNP Q9X0Z6
A	-3	PHE	-	expression tag	UNP Q9X0Z6
A	-2	GLU	-	expression tag	UNP Q9X0Z6
A	-1	LYS	-	expression tag	UNP Q9X0Z6
A	0	ALA	-	expression tag	UNP Q9X0Z6
A	1	SER	-	expression tag	UNP Q9X0Z6
A	311	SER	CYS	engineered mutation	UNP Q9X0Z6
A	319	SER	CYS	engineered mutation	UNP Q9X0Z6
A	322	SER	CYS	engineered mutation	UNP Q9X0Z6

- Molecule 2 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe_4S_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	Fe	S		
2	A	1	8	4	4	0	0

- Molecule 3 is 3-[(3-CHOLAMIDOPROPYL)DIMETHYLAMMONIO]-1-PROPANESULFONATE (three-letter code: CPS) (formula: $C_{32}H_{58}N_2O_7S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
3	A	1	42	32	2	7	1	0	0
			Total	C	N	O	S		
3	A	1	42	32	2	7	1	0	0

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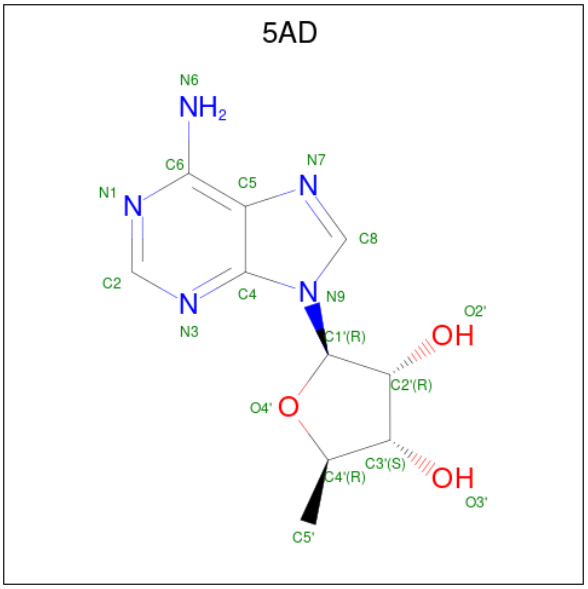
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O 29 24 1 4	0	0
3	A	1	Total C N O 29 24 1 4	0	0
3	A	1	Total C O 50 44 6	0	1

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

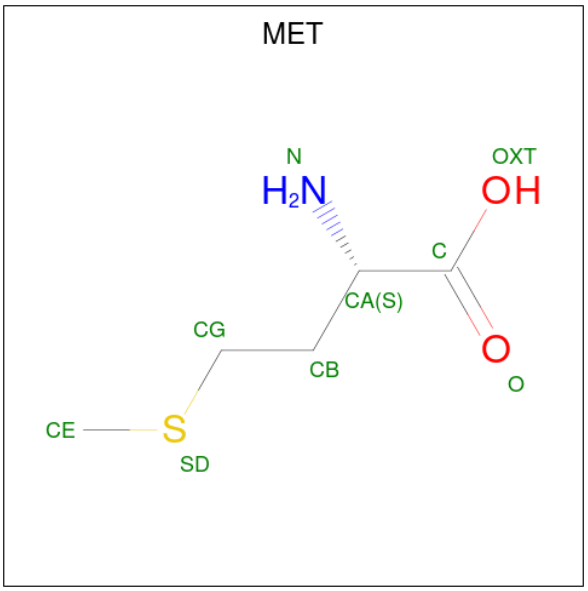
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	3	Total Cl 3 3	0	2

- Molecule 5 is 5'-DEOXYADENOSINE (three-letter code: 5AD) (formula: $C_{10}H_{13}N_5O_3$).



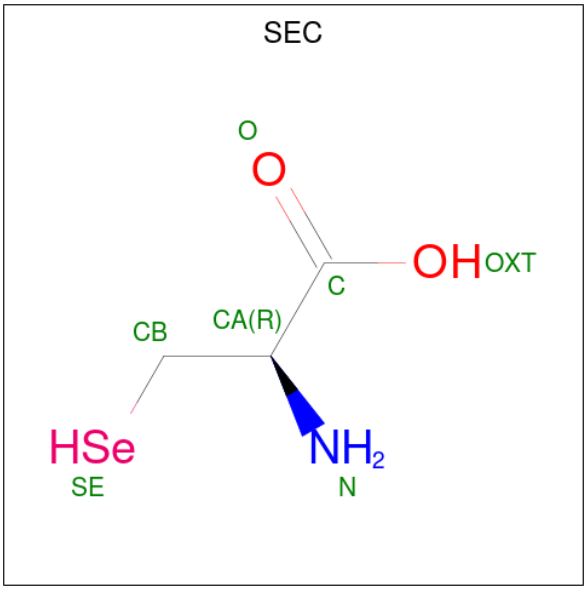
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C N O 18 10 5 3	0	1

- Molecule 6 is METHIONINE (three-letter code: MET) (formula: $C_5H_{11}NO_2S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
6	A	1	9	5	1	2	1	0	1

- Molecule 7 is SELENOCYSTEINE (three-letter code: SEC) (formula: $C_3H_7NO_2Se$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	Se		
7	A	1	7	3	1	2	1	0	1

- Molecule 8 is water.

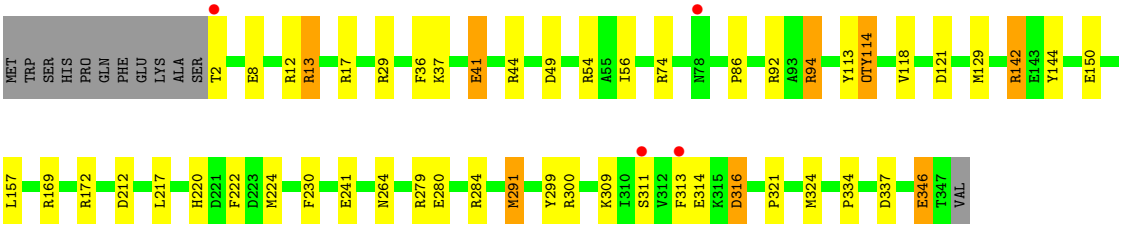
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
8	A	467	469	469	0	7

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: [FeFe] hydrogenase maturase subunit HydE

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	51.12Å 79.46Å 86.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.02 – 1.32 44.02 – 1.32	Depositor EDS
% Data completeness (in resolution range)	99.4 (44.02-1.32) 99.4 (44.02-1.32)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.28 (at 1.32Å)	Xtriage
Refinement program	REFMAC 5.8.0131	Depositor
R, R_{free}	0.113 , 0.147 0.116 , 0.149	Depositor DCC
R_{free} test set	4132 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	12.4	Xtriage
Anisotropy	0.374	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.40 , 57.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.98	EDS
Total number of atoms	3593	wwPDB-VP
Average B, all atoms (Å^2)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.24% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: SF4, SEC, CPS, 5AD, OTY, CL

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	1.22	8/2948 (0.3%)	1.23	25/3982 (0.6%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	41	GLU	CD-OE2	8.46	1.34	1.25
1	A	92	ARG	CZ-NH2	-6.78	1.24	1.33
1	A	346	GLU	CD-OE2	-6.70	1.18	1.25
1	A	280	GLU	CD-OE2	6.60	1.32	1.25
1	A	241	GLU	CD-OE2	-6.11	1.19	1.25
1	A	13	ARG	CZ-NH2	5.90	1.40	1.33
1	A	29	ARG	CZ-NH1	5.82	1.40	1.33
1	A	299	TYR	CE1-CZ	-5.43	1.31	1.38

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	29	ARG	NE-CZ-NH1	11.34	125.97	120.30
1	A	300	ARG	NE-CZ-NH2	-10.76	114.92	120.30
1	A	279	ARG	NE-CZ-NH2	-10.36	115.12	120.30
1	A	316	ASP	CB-CG-OD2	-9.73	109.54	118.30
1	A	13	ARG	NE-CZ-NH1	-9.59	115.51	120.30
1	A	169	ARG	NE-CZ-NH2	-9.51	115.55	120.30
1	A	29	ARG	NE-CZ-NH2	-9.06	115.77	120.30
1	A	172	ARG	NE-CZ-NH2	-8.38	116.11	120.30
1	A	230	PHE	CB-CG-CD2	-7.76	115.37	120.80
1	A	279	ARG	NE-CZ-NH1	7.33	123.96	120.30
1	A	284	ARG	NE-CZ-NH1	6.83	123.72	120.30
1	A	94	ARG	NE-CZ-NH2	6.51	123.56	120.30
1	A	337	ASP	CB-CG-OD1	6.50	124.15	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	74	ARG	NE-CZ-NH1	6.12	123.36	120.30
1	A	54	ARG	NE-CZ-NH2	-6.01	117.29	120.30
1	A	44	ARG	NE-CZ-NH2	-5.79	117.41	120.30
1	A	172	ARG	NE-CZ-NH1	5.63	123.11	120.30
1	A	36	PHE	CB-CG-CD1	5.57	124.70	120.80
1	A	41	GLU	CG-CD-OE2	-5.49	107.32	118.30
1	A	121	ASP	CB-CG-OD2	-5.41	113.43	118.30
1	A	142[A]	ARG	NE-CZ-NH1	-5.34	117.63	120.30
1	A	142[B]	ARG	NE-CZ-NH1	-5.34	117.63	120.30
1	A	49	ASP	CB-CG-OD1	5.32	123.09	118.30
1	A	41	GLU	OE1-CD-OE2	5.08	129.39	123.30
1	A	212	ASP	CB-CG-OD2	-5.07	113.74	118.30

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	2887	0	2891	48	2
2	A	8	0	0	0	0
3	A	192	0	264	14	0
4	A	3	0	0	0	0
5	A	18	0	12	0	0
6	A	9	0	8	0	0
7	A	7	0	3	0	0
8	A	469	0	0	14	2
All	All	3593	0	3178	54	3

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 9.

All (54) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:PRO:HB3	8:A:864[B]:HOH:O	1.21	1.26
1:A:113:TYR:CE2	1:A:114[B]:OTY:CE1	2.19	1.25
1:A:17:ARG:NH1	8:A:501:HOH:O	1.69	1.22
1:A:113:TYR:HE2	1:A:114[B]:OTY:CE1	1.53	1.17
1:A:113:TYR:CE2	1:A:114[B]:OTY:H9	1.76	1.17
1:A:113:TYR:HE2	1:A:114[B]:OTY:H9	1.05	1.06
1:A:56[B]:ILE:HD11	1:A:291[B]:MET:HE2	1.49	0.94
1:A:324[B]:MET:CE	3:A:406[B]:CPS:C11	2.48	0.91
1:A:324[B]:MET:HE2	3:A:406[B]:CPS:H11B	1.53	0.90
1:A:321:PRO:HB3	3:A:406[A]:CPS:H11A	1.53	0.89
1:A:113:TYR:CD2	1:A:114[B]:OTY:CE1	2.59	0.85
1:A:314:GLU:O	8:A:502:HOH:O	1.98	0.82
1:A:324[B]:MET:HE2	3:A:406[B]:CPS:C11	2.09	0.81
1:A:113:TYR:CD2	1:A:114[B]:OTY:CD1	2.67	0.78
1:A:157:LEU:HD23	1:A:224[A]:MET:HE2	1.68	0.74
1:A:324[B]:MET:HE3	3:A:406[B]:CPS:C11	2.16	0.73
1:A:309:LYS:CD	8:A:651:HOH:O	2.38	0.72
1:A:217:LEU:HD11	8:A:511:HOH:O	1.92	0.69
1:A:2:THR:HG22	8:A:924:HOH:O	1.94	0.67
1:A:8:GLU:CG	1:A:12:ARG:CZ	2.74	0.66
1:A:334:PRO:HA	8:A:622[B]:HOH:O	1.97	0.65
1:A:324[B]:MET:CE	3:A:406[B]:CPS:H11B	2.17	0.65
1:A:334:PRO:CB	8:A:622[B]:HOH:O	2.47	0.63
1:A:157:LEU:CD2	1:A:224[A]:MET:HE2	2.28	0.62
1:A:114[B]:OTY:CE1	1:A:118:VAL:HG11	2.29	0.62
1:A:334:PRO:CA	8:A:622[B]:HOH:O	2.48	0.61
1:A:13:ARG:NH1	1:A:346:GLU:OE1	2.33	0.60
1:A:157:LEU:HD23	1:A:224[A]:MET:CE	2.31	0.60
1:A:157:LEU:CD2	1:A:224[A]:MET:CE	2.81	0.58
1:A:314:GLU:OE1	8:A:503:HOH:O	2.17	0.58
1:A:8:GLU:CG	1:A:12:ARG:NH2	2.67	0.57
1:A:324[B]:MET:CE	3:A:406[B]:CPS:H11	2.35	0.56
1:A:324[B]:MET:HE3	3:A:406[B]:CPS:H11	1.88	0.55
1:A:8:GLU:O	1:A:12:ARG:HG3	2.07	0.55
1:A:56[B]:ILE:CD1	1:A:291[B]:MET:HE2	2.29	0.54
1:A:94:ARG:HG3	1:A:129[B]:MET:SD	2.47	0.54
1:A:113:TYR:HD2	1:A:114[B]:OTY:CD1	2.18	0.54
1:A:86:PRO:HD3	1:A:114[B]:OTY:OH	2.09	0.52
1:A:316:ASP:HB3	8:A:834:HOH:O	2.09	0.51
3:A:406[A]:CPS:H4	3:A:406[A]:CPS:H21A	1.92	0.50
1:A:114[B]:OTY:CD1	1:A:118:VAL:HG11	2.42	0.49
1:A:150:GLU:CD	8:A:505:HOH:O	2.50	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:TYR:CE2	1:A:114[B]:OTY:CD1	2.89	0.48
1:A:220:HIS:HB2	1:A:222:PHE:CE1	2.50	0.46
3:A:406[B]:CPS:H21A	3:A:406[B]:CPS:H4	1.98	0.46
1:A:142[A]:ARG:NH1	8:A:506:HOH:O	2.33	0.46
3:A:403:CPS:H4	3:A:403:CPS:H21A	1.99	0.44
1:A:13:ARG:HH11	1:A:13:ARG:HD2	1.66	0.43
1:A:224[A]:MET:HA	1:A:264:ASN:O	2.17	0.43
3:A:405:CPS:H21A	3:A:405:CPS:H4	2.01	0.42
1:A:37[B]:LYS:CE	8:A:644:HOH:O	2.67	0.41
3:A:404:CPS:H21A	3:A:404:CPS:H4	2.01	0.41
3:A:402:CPS:H21B	3:A:402:CPS:C24	2.50	0.41
1:A:113:TYR:HE2	1:A:114[B]:OTY:CZ	2.21	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:313:PHE:O	8:A:768:HOH:O[3_755]	1.99	0.21
8:A:769:HOH:O	8:A:877:HOH:O[2_655]	2.05	0.15
1:A:41:GLU:OE2	1:A:144:TYR:OH[2_654]	2.10	0.10

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	363/358 (101%)	351 (97%)	12 (3%)	0	100 100

There are no Ramachandran outliers to report.

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	309/312 (99%)	306 (99%)	3 (1%)	76 47

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	291[A]	MET
1	A	291[B]	MET
1	A	311	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\#\|Z\| > 2$
1	OTY	A	114[A]	1	12,13,14	1.55	2 (16%)	14,17,19	1.57	4 (28%)
1	OTY	A	114[B]	1	11,12,14	1.52	1 (9%)	12,15,19	3.02	6 (50%)

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
1	OTY	A	114[A]	1	-	0/5/6/8	0/1/1/1
1	OTY	A	114[B]	1	-	4/5/6/8	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	114[B]	OTY	CB-CG	3.71	1.60	1.51
1	A	114[A]	OTY	CD1-CG	-2.63	1.36	1.40
1	A	114[A]	OTY	CB-CG	2.54	1.54	1.51

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	114[B]	OTY	CG-CB-CA	-6.32	101.30	114.10
1	A	114[B]	OTY	CD2-CE2-CZ	4.42	124.72	119.88
1	A	114[B]	OTY	CE1-CD1-CG	4.21	126.82	121.03
1	A	114[B]	OTY	CD2-CG-CD1	-3.66	112.42	118.17
1	A	114[B]	OTY	CB-CG-CD1	3.42	127.70	120.91
1	A	114[A]	OTY	OD1-CD1-CG	-3.22	110.63	118.89
1	A	114[A]	OTY	CB-CA-C	-2.59	106.61	111.47
1	A	114[A]	OTY	OD1-CD1-CE1	2.52	126.22	119.46
1	A	114[B]	OTY	CE2-CZ-CE1	-2.25	115.97	119.77
1	A	114[A]	OTY	CE1-CD1-CG	2.08	123.15	120.70

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	114[B]	OTY	C-CA-CB-CG
1	A	114[B]	OTY	CA-CB-CG-CD1
1	A	114[B]	OTY	CA-CB-CG-CD2
1	A	114[B]	OTY	N-CA-CB-CG

There are no ring outliers.

1 monomer is involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	114[B]	OTY	12	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 13 ligands modelled in this entry, 3 are monoatomic - leaving 10 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	$\#\|Z\| > 2$
7	SEC	A	412[B]	5	3,6,6	1.41	0	3,7,7	0.77	0
3	CPS	A	406[B]	-	28,28,45	1.08	3 (10%)	46,46,70	1.40	6 (13%)
3	CPS	A	402	-	45,45,45	1.57	6 (13%)	69,70,70	1.31	9 (13%)
3	CPS	A	404	-	32,32,45	1.18	2 (6%)	51,51,70	1.24	5 (9%)
2	SF4	A	401	1,6	0,12,12	-	-	-		
6	MET	A	411[B]	2	7,8,8	0.73	0	7,9,9	0.59	0
3	CPS	A	403	-	45,45,45	1.26	3 (6%)	69,70,70	1.64	13 (18%)
5	5AD	A	410[B]	7	17,20,20	1.31	2 (11%)	15,30,30	1.56	3 (20%)
3	CPS	A	406[A]	-	28,28,45	1.41	4 (14%)	46,46,70	2.04	15 (32%)
3	CPS	A	405	-	32,32,45	1.19	5 (15%)	51,51,70	1.06	3 (5%)

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
7	SEC	A	412[B]	5	-	0/4/6/6	-
3	CPS	A	406[B]	-	-	0/4/69/90	0/4/4/4
3	CPS	A	402	-	-	0/25/90/90	0/4/4/4
3	CPS	A	404	-	-	0/9/74/90	0/4/4/4
2	SF4	A	401	1,6	-	-	0/6/5/5
6	MET	A	411[B]	2	-	1/8/8/8	-
3	CPS	A	403	-	-	0/25/90/90	0/4/4/4
5	5AD	A	410[B]	7	-	0/0/20/20	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	CPS	A	406[A]	-	-	0/4/69/90	0/4/4/4
3	CPS	A	405	-	-	0/9/74/90	0/4/4/4

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	CPS	C32-S	-6.99	1.67	1.77
3	A	403	CPS	C32-S	-3.90	1.72	1.77
3	A	406[A]	CPS	C5-C9	-3.80	1.49	1.55
5	A	410[B]	5AD	C3'-C4'	3.63	1.57	1.52
3	A	404	CPS	O3-C17	3.29	1.50	1.43
3	A	405	CPS	C3-C4	3.08	1.58	1.53
3	A	406[A]	CPS	C14-C15	2.91	1.58	1.53
3	A	406[B]	CPS	O3-C17	2.84	1.49	1.43
3	A	404	CPS	O4-C4	2.78	1.48	1.43
3	A	402	CPS	O3-C17	2.75	1.49	1.43
3	A	406[A]	CPS	C18-C17	-2.59	1.49	1.53
3	A	405	CPS	O4-C4	2.51	1.47	1.43
3	A	406[A]	CPS	C11-C2	2.31	1.58	1.54
3	A	402	CPS	C1-C2	2.30	1.58	1.54
3	A	403	CPS	O3S-S	-2.29	1.38	1.45
3	A	405	CPS	C2-C19	2.28	1.60	1.56
3	A	405	CPS	C5-C6	-2.28	1.51	1.55
3	A	402	CPS	C2-C19	2.27	1.60	1.56
3	A	402	CPS	O1-C24	2.25	1.27	1.23
3	A	402	CPS	C10-C5	-2.17	1.50	1.54
3	A	406[B]	CPS	C18-C17	-2.14	1.49	1.53
3	A	406[B]	CPS	C3-C19	2.11	1.57	1.53
5	A	410[B]	5AD	C4-N3	-2.06	1.32	1.35
3	A	405	CPS	O1-C24	-2.06	1.17	1.24
3	A	403	CPS	C18-C19	2.04	1.57	1.53

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	403	CPS	O2S-S-C32	4.38	112.85	105.77
3	A	406[A]	CPS	C7-C6-C5	-4.22	99.41	103.55
3	A	406[A]	CPS	C3-C19-C2	-4.11	109.49	113.73
3	A	406[A]	CPS	C9-C5-C6	4.10	104.22	100.09
3	A	406[A]	CPS	C12-C1-C2	3.95	119.56	112.78
3	A	403	CPS	C25-N1-C24	-3.88	115.63	122.84
3	A	406[A]	CPS	C5-C9-C20	-3.84	114.90	119.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	403	CPS	O1-C24-C23	-3.80	115.07	122.02
3	A	406[B]	CPS	C14-C15-C2	3.77	116.66	112.66
3	A	406[A]	CPS	C22-C20-C9	-3.70	102.52	112.42
3	A	403	CPS	C14-C15-C2	3.67	116.56	112.66
3	A	402	CPS	C14-C13-C12	-3.52	106.35	110.55
3	A	403	CPS	O3S-S-C32	3.50	111.13	106.92
3	A	403	CPS	C23-C22-C20	-3.48	108.16	114.52
5	A	410[B]	5AD	C5'-C4'-C3'	3.22	119.08	115.70
3	A	403	CPS	C9-C5-C4	3.13	120.52	117.67
3	A	406[B]	CPS	C1-C2-C19	3.07	116.18	111.35
3	A	406[A]	CPS	C8-C9-C5	-3.05	100.56	103.55
3	A	402	CPS	C16-C15-C14	2.85	114.48	111.19
3	A	402	CPS	C10-C5-C6	2.85	115.67	111.21
3	A	406[A]	CPS	C19-C18-C17	2.84	115.27	111.88
5	A	410[B]	5AD	C1'-N9-C4	-2.83	121.68	126.64
3	A	406[B]	CPS	C10-C5-C9	-2.81	106.81	111.21
3	A	403	CPS	C29-N2-C27	2.79	116.55	109.46
3	A	406[A]	CPS	C22-C20-C21	-2.72	102.98	110.59
3	A	406[A]	CPS	C11-C2-C19	2.72	114.93	111.18
3	A	405	CPS	C3-C19-C2	-2.69	110.96	113.73
3	A	406[A]	CPS	C19-C2-C15	-2.67	104.82	108.58
3	A	404	CPS	C9-C5-C4	-2.64	115.25	117.67
3	A	406[B]	CPS	C12-C1-C2	2.61	117.26	112.78
3	A	402	CPS	O2-C13-C14	-2.58	104.71	109.85
3	A	404	CPS	C8-C9-C5	-2.58	101.02	103.55
3	A	406[B]	CPS	C9-C5-C4	2.55	120.00	117.67
3	A	402	CPS	C23-C22-C20	2.54	119.16	114.52
3	A	402	CPS	C29-N2-C27	-2.53	103.02	109.46
3	A	402	CPS	C29-N2-C30	2.50	115.82	109.46
3	A	405	CPS	C1-C2-C15	2.47	111.42	107.77
3	A	404	CPS	C16-C17-C18	2.38	114.03	111.48
5	A	410[B]	5AD	O3'-C3'-C4'	-2.35	104.70	110.47
3	A	406[B]	CPS	C11-C2-C1	-2.30	104.55	108.26
3	A	403	CPS	C10-C5-C9	-2.26	107.67	111.21
3	A	404	CPS	O2-C13-C14	2.26	114.35	109.85
3	A	403	CPS	C21-C20-C22	2.22	113.84	110.36
3	A	406[A]	CPS	C21-C20-C9	2.21	118.34	112.42
3	A	403	CPS	O2S-S-O1S	-2.18	105.94	111.27
3	A	402	CPS	C3-C4-C5	-2.16	109.03	111.24
3	A	402	CPS	C16-C17-C18	2.13	113.75	111.48
3	A	406[A]	CPS	C16-C17-C18	-2.12	109.21	111.48
3	A	404	CPS	C10-C5-C6	2.04	114.41	111.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	403	CPS	C19-C18-C6	2.04	112.51	109.71
3	A	406[A]	CPS	C15-C16-C17	2.04	116.71	114.46
3	A	403	CPS	O1-C24-N1	2.03	126.85	123.01
3	A	405	CPS	C23-C22-C20	2.03	118.23	114.52
3	A	406[A]	CPS	C19-C18-C6	2.03	112.49	109.71

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	411[B]	MET	C-CA-CB-CG

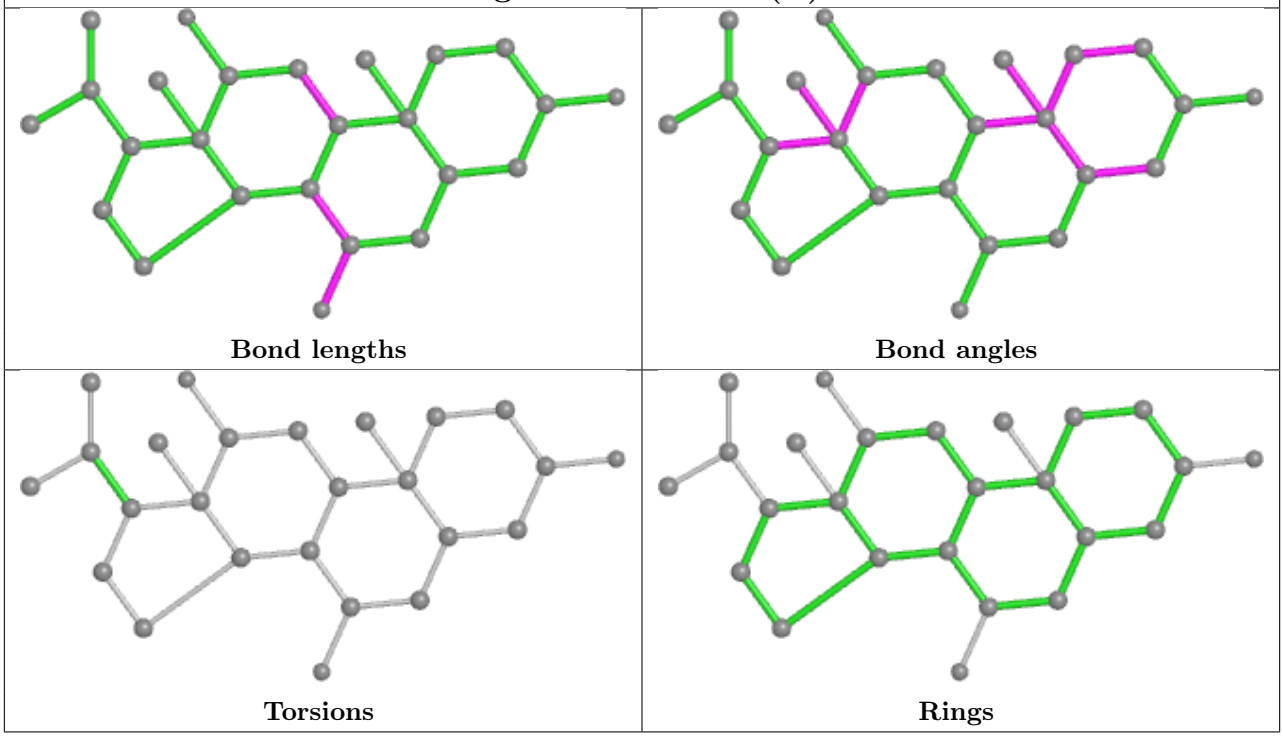
There are no ring outliers.

6 monomers are involved in 14 short contacts:

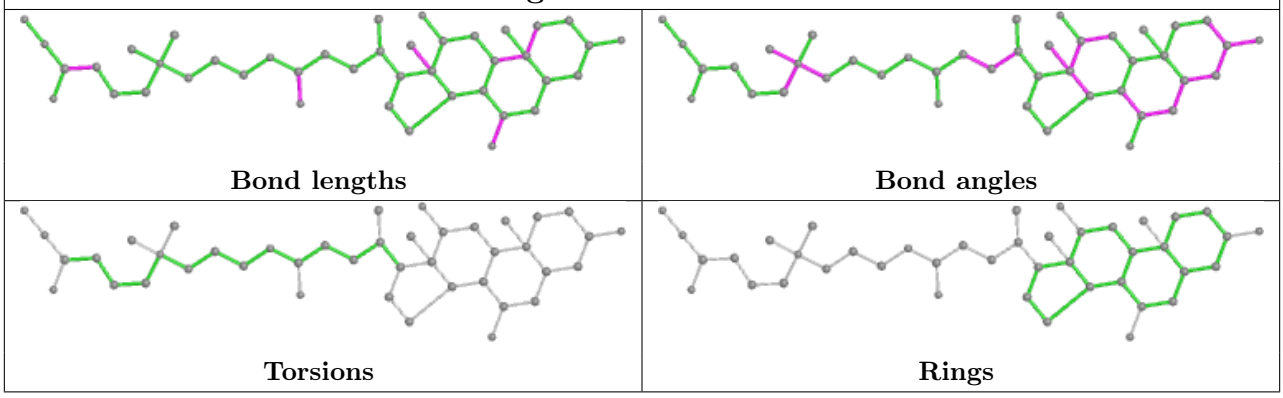
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	406[B]	CPS	8	0
3	A	402	CPS	1	0
3	A	404	CPS	1	0
3	A	403	CPS	1	0
3	A	406[A]	CPS	2	0
3	A	405	CPS	1	0

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 then 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.

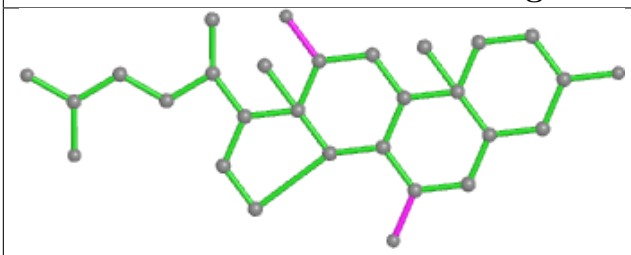
Ligand CPS A 406 (B)



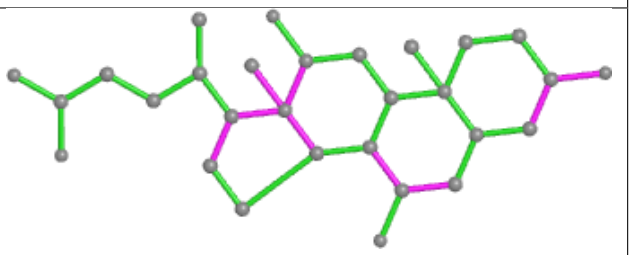
Ligand CPS A 402



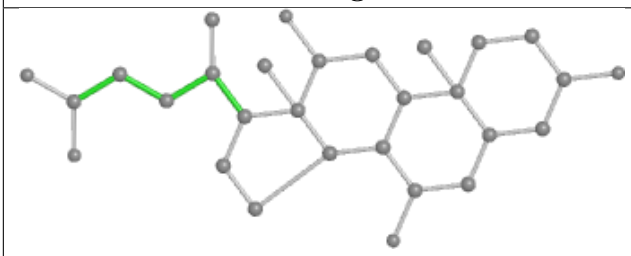
Ligand CPS A 404



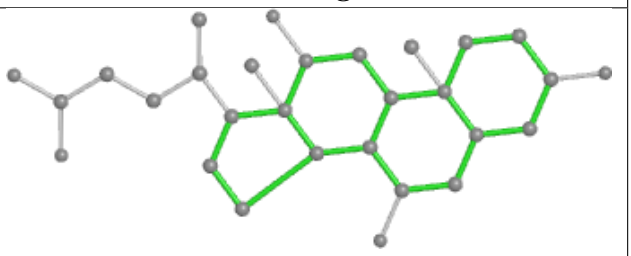
Bond lengths



Bond angles

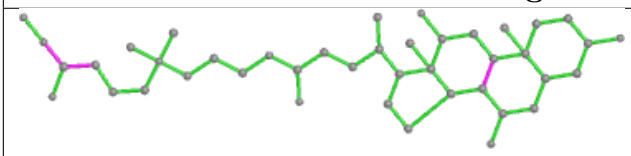


Torsions



Rings

Ligand CPS A 403



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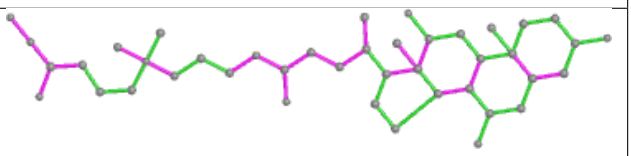
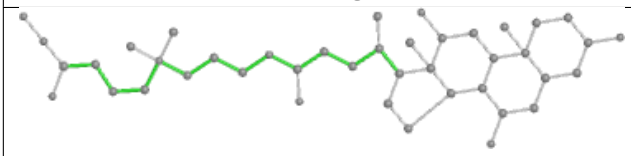
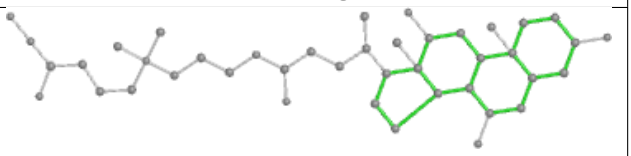


Bond angles

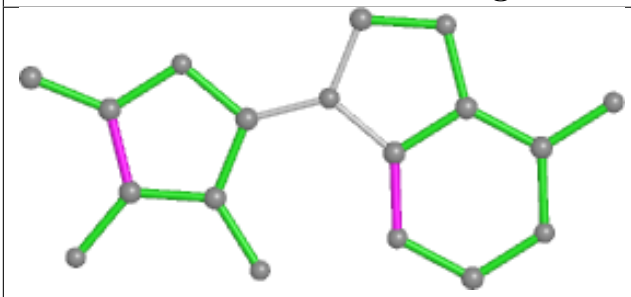


Torsions

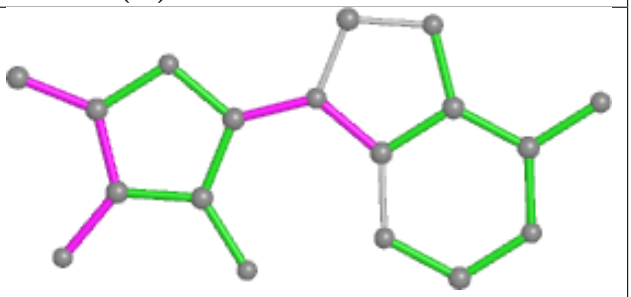


Rings

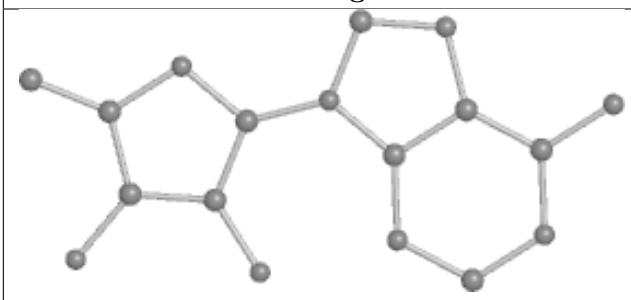
Ligand 5AD A 410 (B)



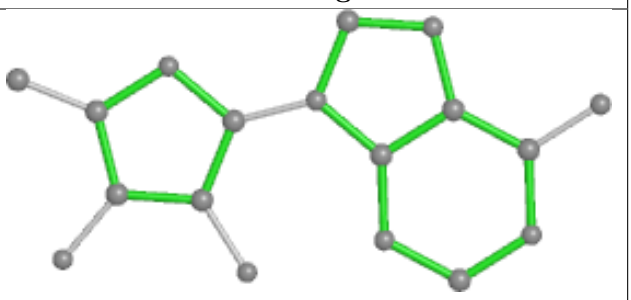
Bond lengths



Bond angles

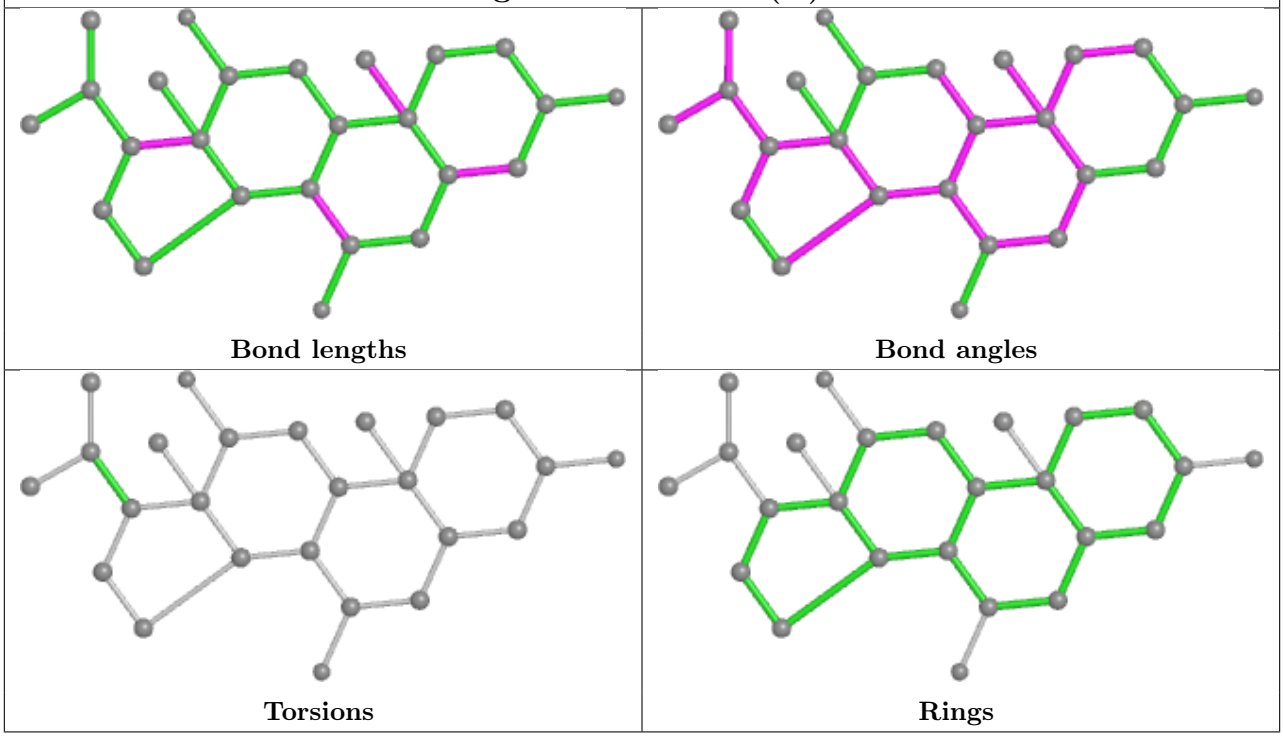


Torsions

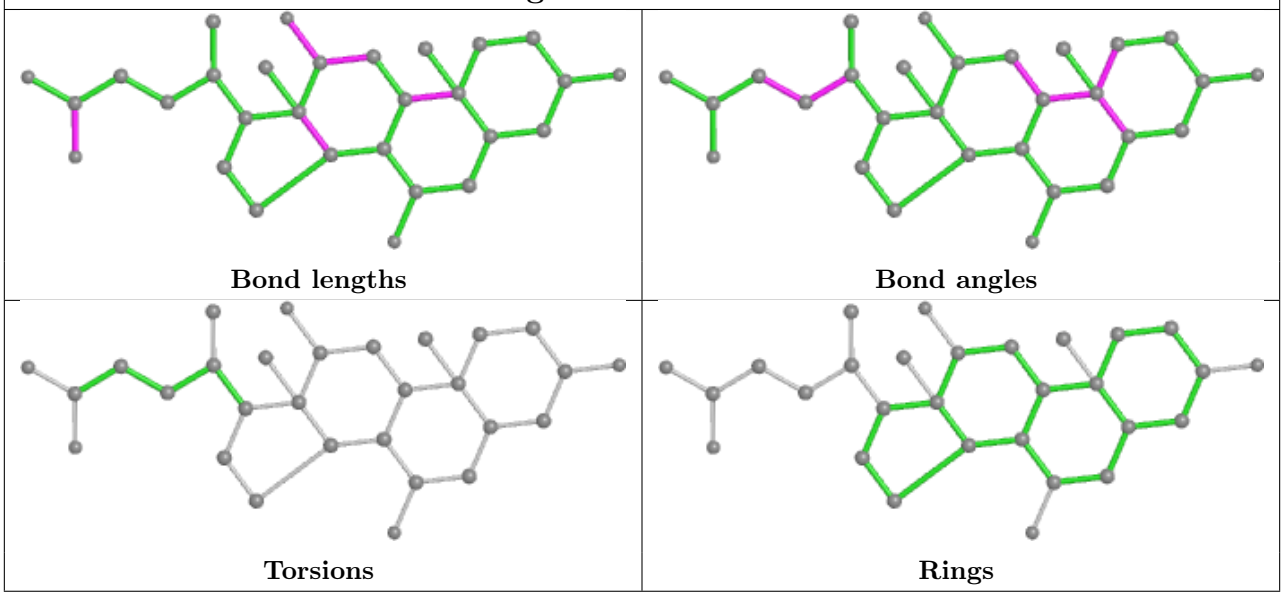


Rings

Ligand CPS A 406 (A)



Ligand CPS A 405



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	345/358 (96%)	-0.41	4 (1%) 79 80	9, 14, 30, 45	7 (2%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	311	SER	3.0
1	A	78[A]	ASN	2.6
1	A	2	THR	2.6
1	A	313	PHE	2.4

6.2 Non-standard residues in protein, DNA, RNA chains (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
1	OTY	A	114[A]	13/14	0.98	0.07	11,12,15,17	1
1	OTY	A	114[B]	12/14	0.98	0.07	7,13,14,14	12

6.3 Carbohydrates (i)

There are no monosaccharides 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,

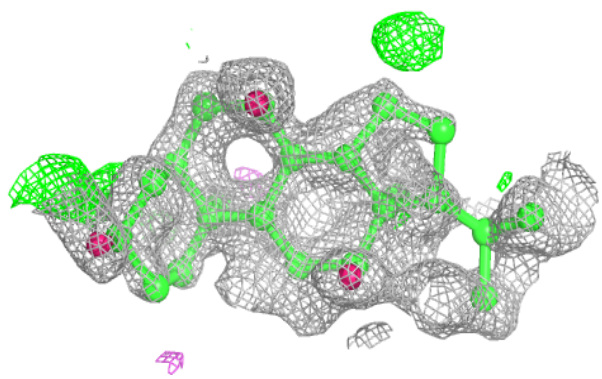
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

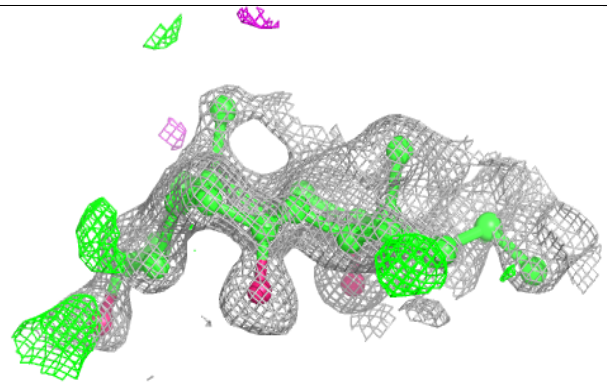
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	CPS	A	406[A]	25/42	0.89	0.13	10,11,13,15	25
3	CPS	A	406[B]	25/42	0.89	0.13	18,21,29,32	25
3	CPS	A	405	29/42	0.90	0.09	15,18,25,32	0
3	CPS	A	403	42/42	0.94	0.08	12,16,29,36	5
3	CPS	A	404	29/42	0.96	0.06	13,15,43,52	0
3	CPS	A	402	42/42	0.96	0.07	10,13,24,28	6
5	5AD	A	410[B]	18/18	0.96	0.07	14,15,18,18	0
6	MET	A	411[B]	9/9	0.98	0.07	14,15,20,21	0
4	CL	A	408[B]	1/1	0.99	0.05	21,21,21,21	1
7	SEC	A	412[B]	7/7	0.99	0.06	13,15,16,18	7
4	CL	A	407	1/1	1.00	0.04	16,16,16,16	0
2	SF4	A	401	8/8	1.00	0.04	12,13,13,14	0
4	CL	A	409[A]	1/1	1.00	0.03	34,34,34,34	1

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 CPS A 406 (A):

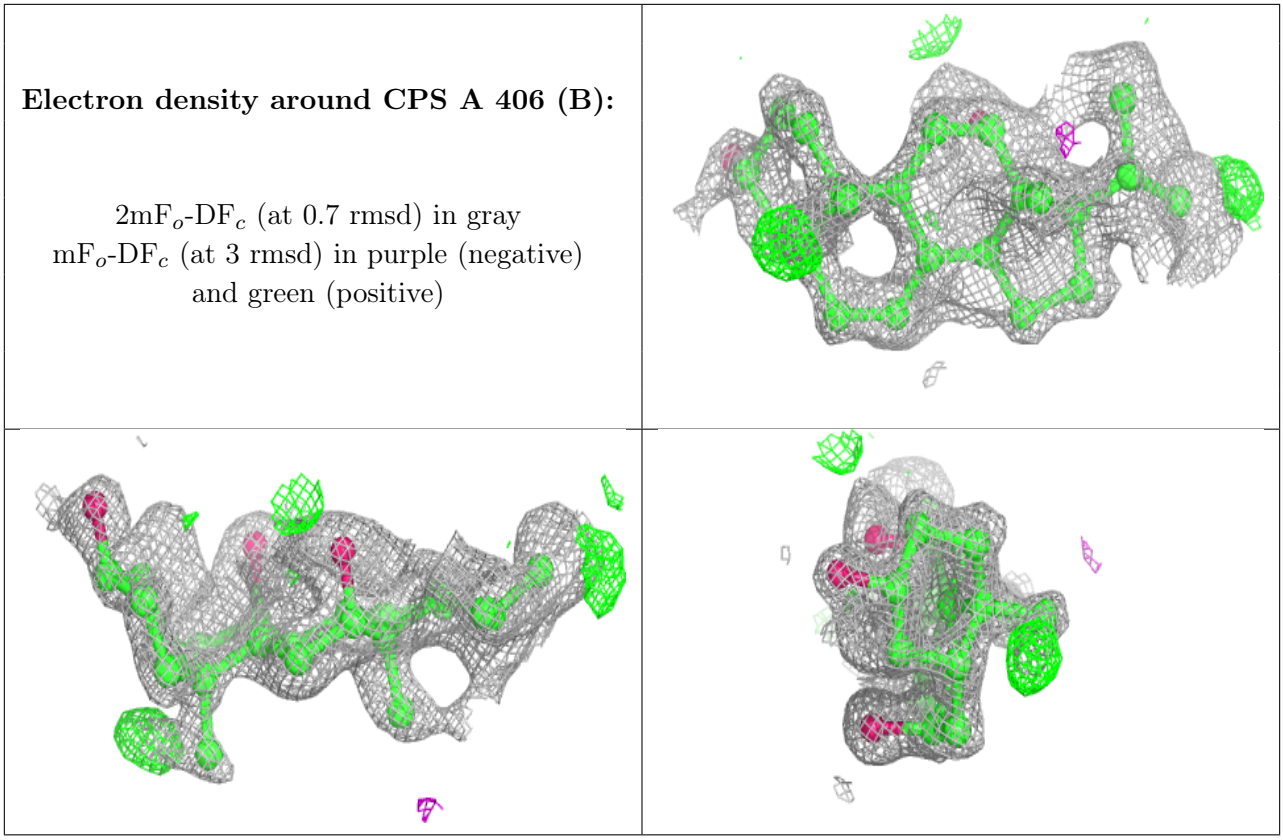
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)





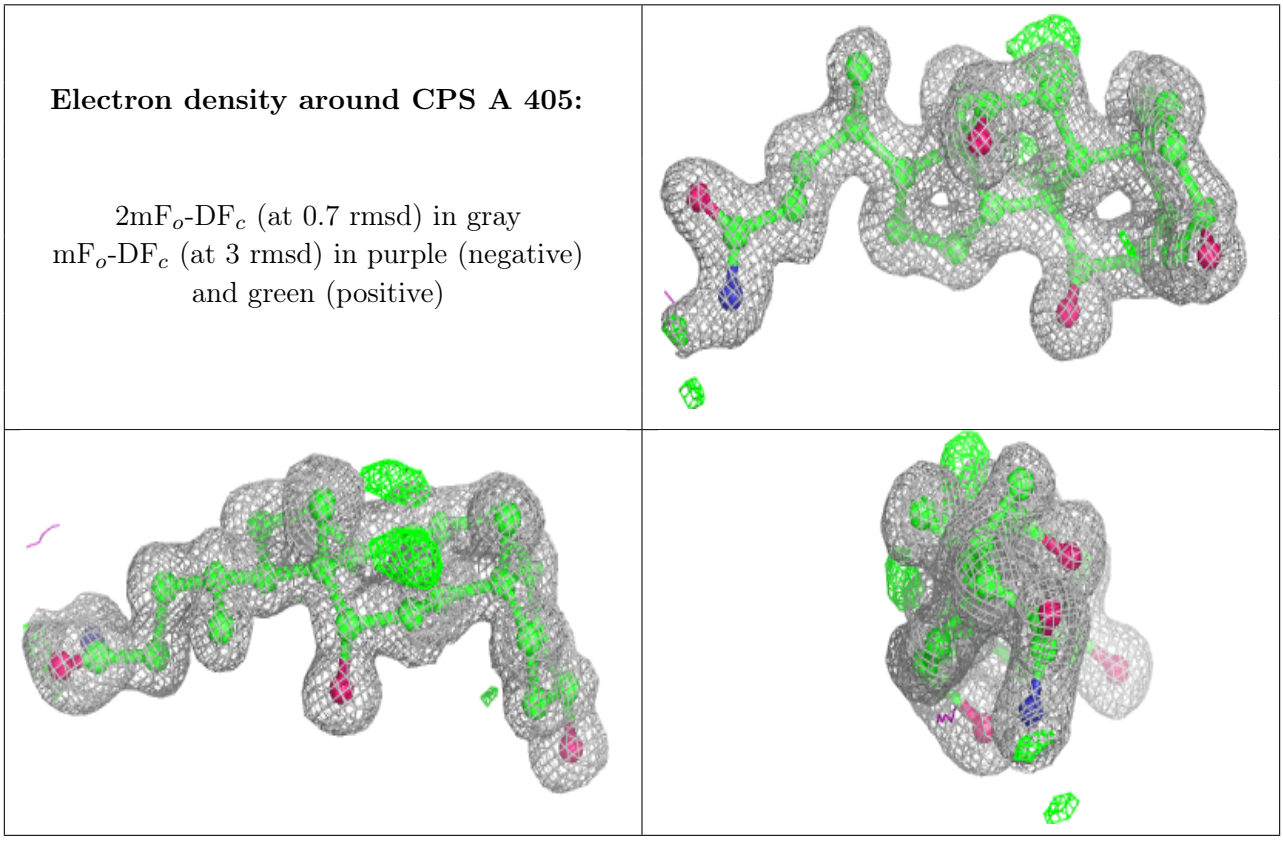
Electron density around CPS A 406 (B):

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



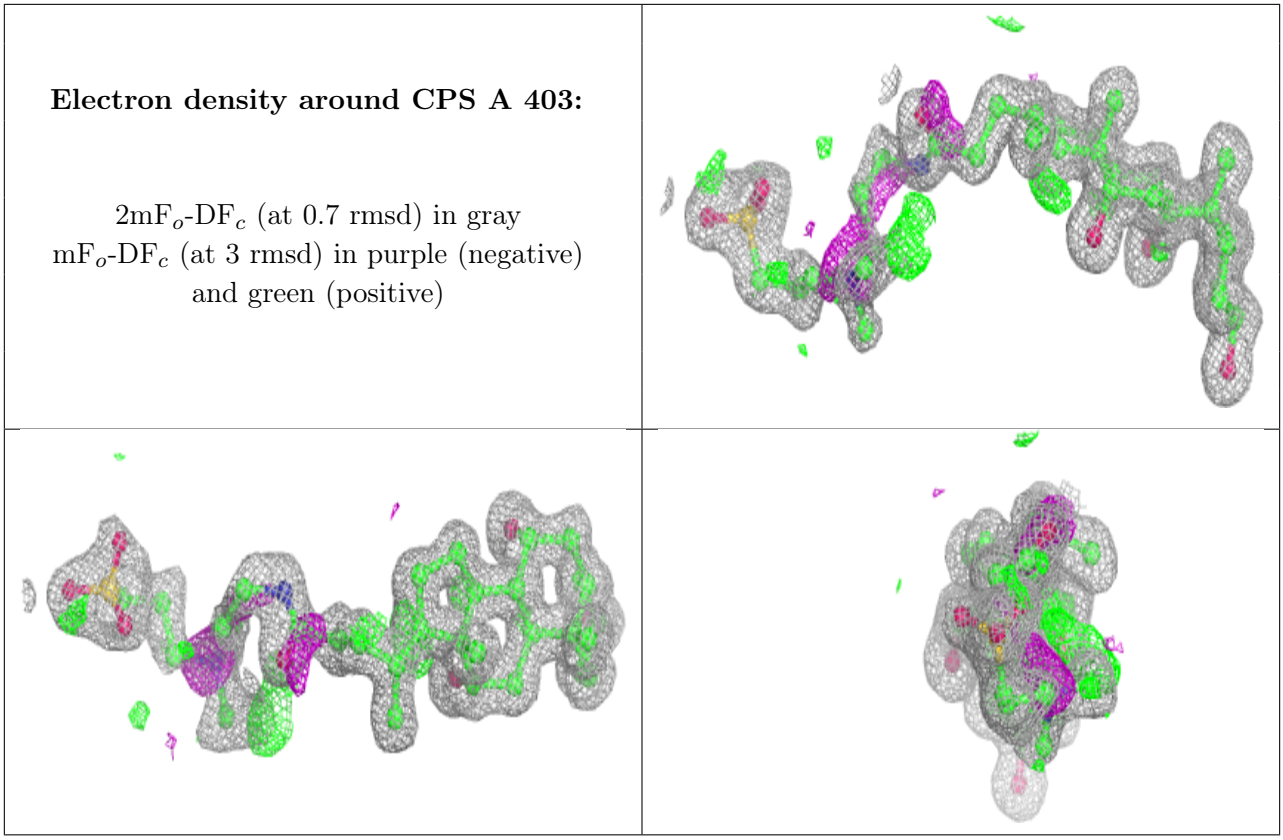
Electron density around CPS A 405:

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



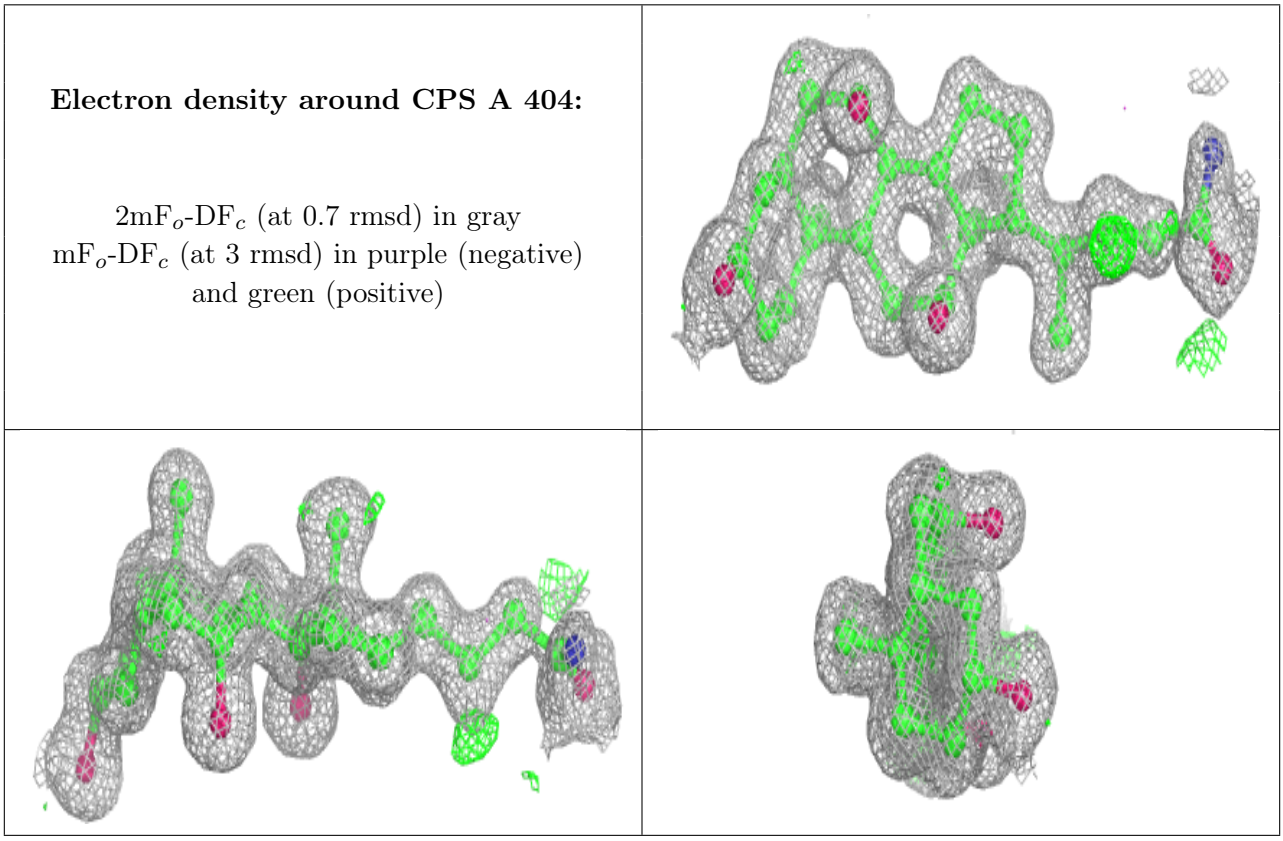
Electron density around CPS A 403:

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



Electron density around CPS A 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 CPS A 402:

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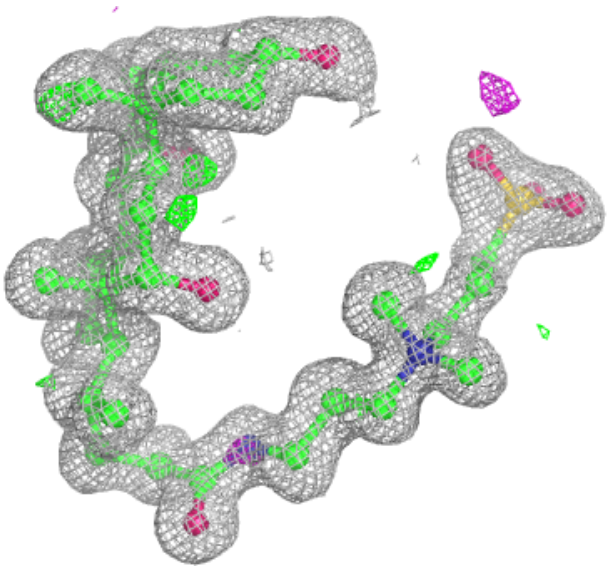
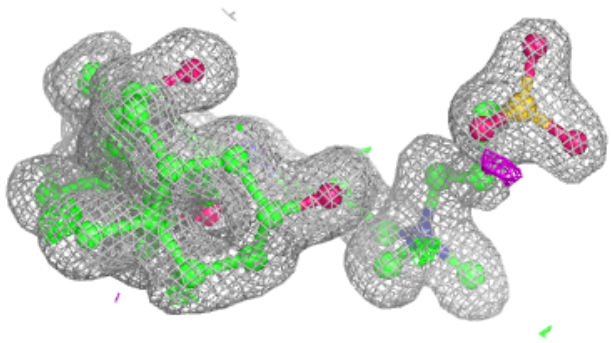
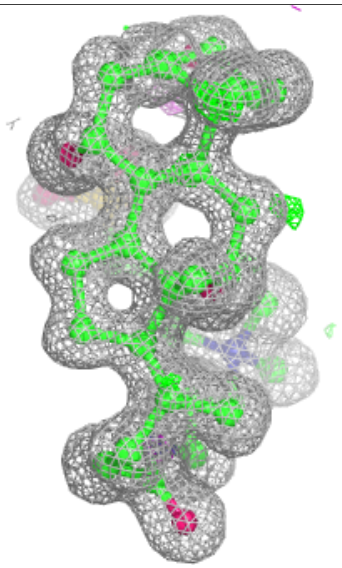
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

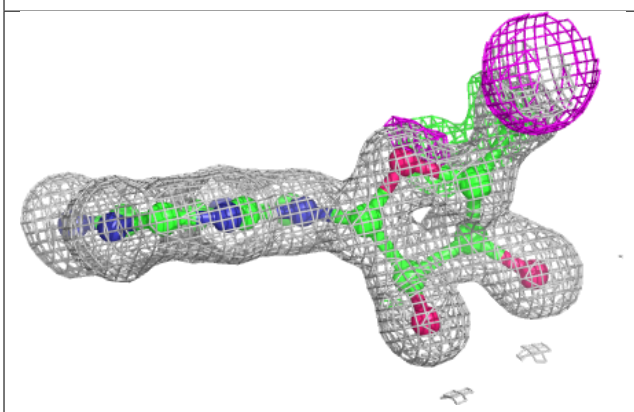


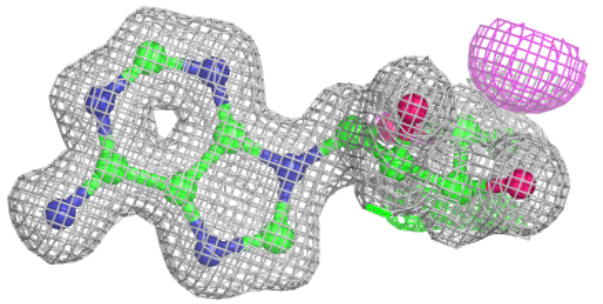


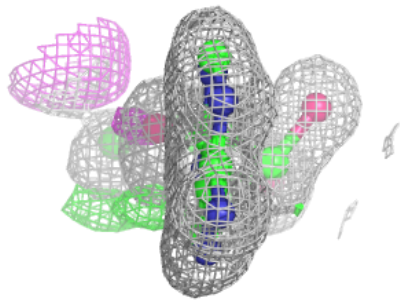


Electron density around 5AD A 410 (B):

$2mF_o$-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.