wwPDB X-ray Structure Validation Summary Report (i)

Jan 5, 2023 – 03:46 pm GMT

PDB ID	:	8APQ
Title	:	CaMct - Mesaconyl-CoA C1:C4 CoA Transferase of Chloroflexus aurantiacus
Authors	:	Pfister, P.; Zarzycki, J.; Erb, T.J.
Deposited on	:	2022-08-10
Resolution	:	2.49 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 (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:

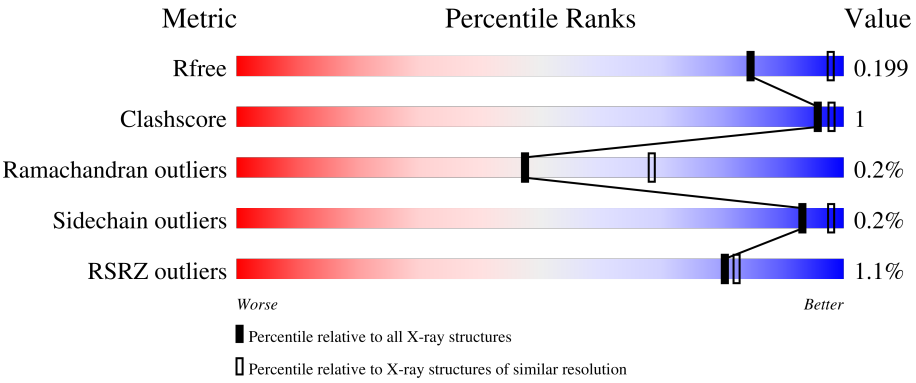
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.31.3
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.3

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 2.49 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	430	91% · 6%
1	B	430	92% · 6%
1	C	430	92% · 6%
1	D	430	92% · 6%
1	E	430	92% · 6%

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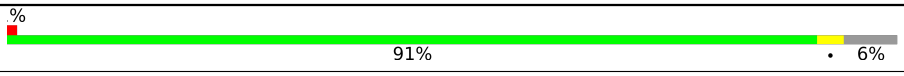
Mol	Chain	Length	Quality of chain
1	F	430	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	MEZ	B	502	-	X	-	X
4	MEZ	E	502	-	-	-	X
4	MEZ	F	502	-	-	-	X

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 20187 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 2-methylfumaryl-CoA isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	406	Total	C	N	O	S	0	0	0
			3134	1989	565	565	15			
1	B	406	Total	C	N	O	S	0	0	0
			3134	1989	565	565	15			
1	C	406	Total	C	N	O	S	0	0	0
			3134	1989	565	565	15			
1	D	406	Total	C	N	O	S	0	0	0
			3133	1989	565	564	15			
1	E	405	Total	C	N	O	S	0	0	0
			3126	1984	564	564	14			
1	F	406	Total	C	N	O	S	0	0	0
			3134	1989	565	565	15			

There are 126 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	initiating methionine	UNP A9WC36
A	-19	GLY	-	expression tag	UNP A9WC36
A	-18	HIS	-	expression tag	UNP A9WC36
A	-17	HIS	-	expression tag	UNP A9WC36
A	-16	HIS	-	expression tag	UNP A9WC36
A	-15	HIS	-	expression tag	UNP A9WC36
A	-14	HIS	-	expression tag	UNP A9WC36
A	-13	HIS	-	expression tag	UNP A9WC36
A	-12	HIS	-	expression tag	UNP A9WC36
A	-11	HIS	-	expression tag	UNP A9WC36
A	-10	HIS	-	expression tag	UNP A9WC36
A	-9	HIS	-	expression tag	UNP A9WC36
A	-8	SER	-	expression tag	UNP A9WC36
A	-7	SER	-	expression tag	UNP A9WC36
A	-6	GLY	-	expression tag	UNP A9WC36
A	-5	HIS	-	expression tag	UNP A9WC36
A	-4	ILE	-	expression tag	UNP A9WC36

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLU	-	expression tag	UNP A9WC36
A	-2	GLY	-	expression tag	UNP A9WC36
A	-1	ARG	-	expression tag	UNP A9WC36
A	0	HIS	-	expression tag	UNP A9WC36
B	-20	MET	-	initiating methionine	UNP A9WC36
B	-19	GLY	-	expression tag	UNP A9WC36
B	-18	HIS	-	expression tag	UNP A9WC36
B	-17	HIS	-	expression tag	UNP A9WC36
B	-16	HIS	-	expression tag	UNP A9WC36
B	-15	HIS	-	expression tag	UNP A9WC36
B	-14	HIS	-	expression tag	UNP A9WC36
B	-13	HIS	-	expression tag	UNP A9WC36
B	-12	HIS	-	expression tag	UNP A9WC36
B	-11	HIS	-	expression tag	UNP A9WC36
B	-10	HIS	-	expression tag	UNP A9WC36
B	-9	HIS	-	expression tag	UNP A9WC36
B	-8	SER	-	expression tag	UNP A9WC36
B	-7	SER	-	expression tag	UNP A9WC36
B	-6	GLY	-	expression tag	UNP A9WC36
B	-5	HIS	-	expression tag	UNP A9WC36
B	-4	ILE	-	expression tag	UNP A9WC36
B	-3	GLU	-	expression tag	UNP A9WC36
B	-2	GLY	-	expression tag	UNP A9WC36
B	-1	ARG	-	expression tag	UNP A9WC36
B	0	HIS	-	expression tag	UNP A9WC36
C	-20	MET	-	initiating methionine	UNP A9WC36
C	-19	GLY	-	expression tag	UNP A9WC36
C	-18	HIS	-	expression tag	UNP A9WC36
C	-17	HIS	-	expression tag	UNP A9WC36
C	-16	HIS	-	expression tag	UNP A9WC36
C	-15	HIS	-	expression tag	UNP A9WC36
C	-14	HIS	-	expression tag	UNP A9WC36
C	-13	HIS	-	expression tag	UNP A9WC36
C	-12	HIS	-	expression tag	UNP A9WC36
C	-11	HIS	-	expression tag	UNP A9WC36
C	-10	HIS	-	expression tag	UNP A9WC36
C	-9	HIS	-	expression tag	UNP A9WC36
C	-8	SER	-	expression tag	UNP A9WC36
C	-7	SER	-	expression tag	UNP A9WC36
C	-6	GLY	-	expression tag	UNP A9WC36
C	-5	HIS	-	expression tag	UNP A9WC36
C	-4	ILE	-	expression tag	UNP A9WC36

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-3	GLU	-	expression tag	UNP A9WC36
C	-2	GLY	-	expression tag	UNP A9WC36
C	-1	ARG	-	expression tag	UNP A9WC36
C	0	HIS	-	expression tag	UNP A9WC36
D	-20	MET	-	initiating methionine	UNP A9WC36
D	-19	GLY	-	expression tag	UNP A9WC36
D	-18	HIS	-	expression tag	UNP A9WC36
D	-17	HIS	-	expression tag	UNP A9WC36
D	-16	HIS	-	expression tag	UNP A9WC36
D	-15	HIS	-	expression tag	UNP A9WC36
D	-14	HIS	-	expression tag	UNP A9WC36
D	-13	HIS	-	expression tag	UNP A9WC36
D	-12	HIS	-	expression tag	UNP A9WC36
D	-11	HIS	-	expression tag	UNP A9WC36
D	-10	HIS	-	expression tag	UNP A9WC36
D	-9	HIS	-	expression tag	UNP A9WC36
D	-8	SER	-	expression tag	UNP A9WC36
D	-7	SER	-	expression tag	UNP A9WC36
D	-6	GLY	-	expression tag	UNP A9WC36
D	-5	HIS	-	expression tag	UNP A9WC36
D	-4	ILE	-	expression tag	UNP A9WC36
D	-3	GLU	-	expression tag	UNP A9WC36
D	-2	GLY	-	expression tag	UNP A9WC36
D	-1	ARG	-	expression tag	UNP A9WC36
D	0	HIS	-	expression tag	UNP A9WC36
E	-20	MET	-	initiating methionine	UNP A9WC36
E	-19	GLY	-	expression tag	UNP A9WC36
E	-18	HIS	-	expression tag	UNP A9WC36
E	-17	HIS	-	expression tag	UNP A9WC36
E	-16	HIS	-	expression tag	UNP A9WC36
E	-15	HIS	-	expression tag	UNP A9WC36
E	-14	HIS	-	expression tag	UNP A9WC36
E	-13	HIS	-	expression tag	UNP A9WC36
E	-12	HIS	-	expression tag	UNP A9WC36
E	-11	HIS	-	expression tag	UNP A9WC36
E	-10	HIS	-	expression tag	UNP A9WC36
E	-9	HIS	-	expression tag	UNP A9WC36
E	-8	SER	-	expression tag	UNP A9WC36
E	-7	SER	-	expression tag	UNP A9WC36
E	-6	GLY	-	expression tag	UNP A9WC36
E	-5	HIS	-	expression tag	UNP A9WC36
E	-4	ILE	-	expression tag	UNP A9WC36

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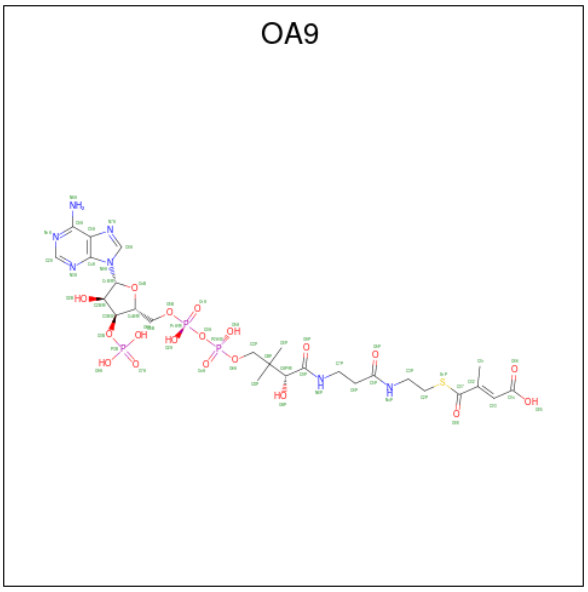
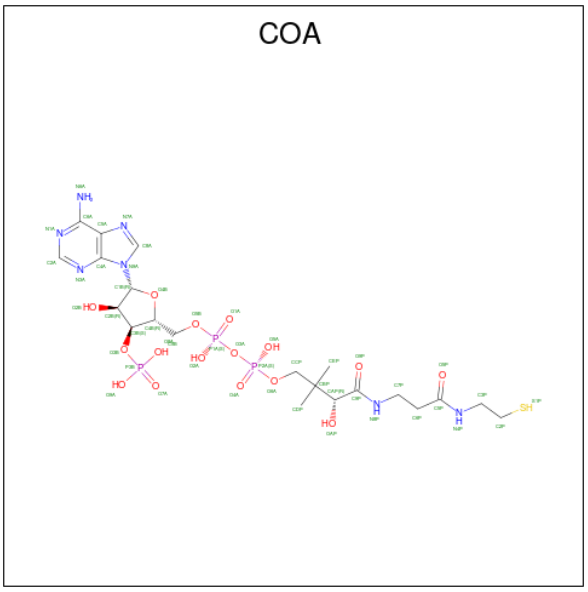
Chain	Residue	Modelled	Actual	Comment	Reference
E	-3	GLU	-	expression tag	UNP A9WC36
E	-2	GLY	-	expression tag	UNP A9WC36
E	-1	ARG	-	expression tag	UNP A9WC36
E	0	HIS	-	expression tag	UNP A9WC36
F	-20	MET	-	initiating methionine	UNP A9WC36
F	-19	GLY	-	expression tag	UNP A9WC36
F	-18	HIS	-	expression tag	UNP A9WC36
F	-17	HIS	-	expression tag	UNP A9WC36
F	-16	HIS	-	expression tag	UNP A9WC36
F	-15	HIS	-	expression tag	UNP A9WC36
F	-14	HIS	-	expression tag	UNP A9WC36
F	-13	HIS	-	expression tag	UNP A9WC36
F	-12	HIS	-	expression tag	UNP A9WC36
F	-11	HIS	-	expression tag	UNP A9WC36
F	-10	HIS	-	expression tag	UNP A9WC36
F	-9	HIS	-	expression tag	UNP A9WC36
F	-8	SER	-	expression tag	UNP A9WC36
F	-7	SER	-	expression tag	UNP A9WC36
F	-6	GLY	-	expression tag	UNP A9WC36
F	-5	HIS	-	expression tag	UNP A9WC36
F	-4	ILE	-	expression tag	UNP A9WC36
F	-3	GLU	-	expression tag	UNP A9WC36
F	-2	GLY	-	expression tag	UNP A9WC36
F	-1	ARG	-	expression tag	UNP A9WC36
F	0	HIS	-	expression tag	UNP A9WC36

- Molecule 2 is Mesaconyl Coenzme A (three-letter code: OA9) (formula: $C_{26}H_{40}N_7O_{19}P_3S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			56	26	7	19	3	1		

- Molecule 3 is COENZYME A (three-letter code: COA) (formula: $C_{21}H_{36}N_7O_{16}P_3S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		
3	C	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

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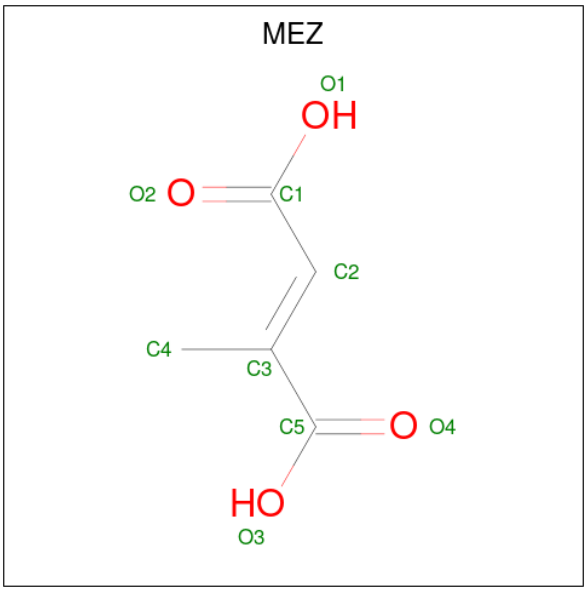
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	1	Total C N O P S 48 21 7 16 3 1	0	0
3	E	1	Total C N O P S 48 21 7 16 3 1	0	0
3	F	1	Total C N O P S 48 21 7 16 3 1	0	0

- Molecule 4 is (2E)-2-METHYLBUT-2-ENEDIOIC ACID (three-letter code: MEZ) (formula: $C_5H_6O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C O 8 5 3	0	0
4	C	1	Total C O 8 5 3	0	0
4	D	1	Total C O 9 5 4	0	0
4	E	1	Total C O 8 5 3	0	0
4	F	1	Total C O 8 5 3	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	173	Total O 173 173	0	0

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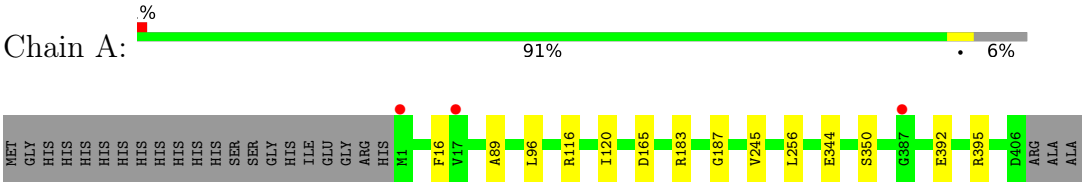
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	151	Total O 151 151	0	0
5	C	172	Total O 172 172	0	0
5	D	177	Total O 177 177	0	0
5	E	185	Total O 185 185	0	0
5	F	197	Total O 197 197	0	0

3 Residue-property plots (i)

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

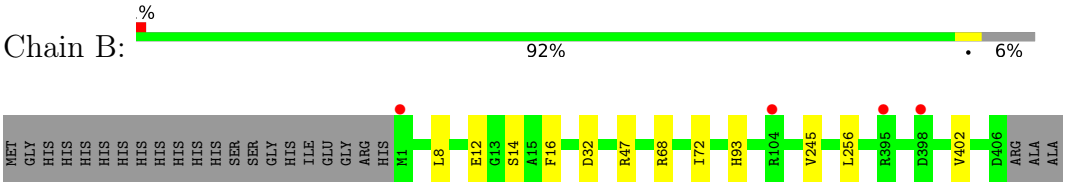
- Molecule 1: 2-methylfumaryl-CoA isomerase

Chain A:



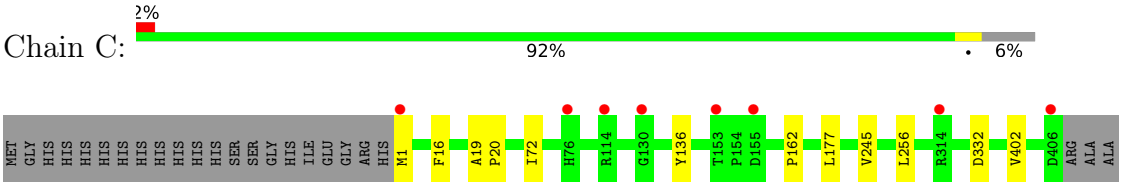
- Molecule 1: 2-methylfumaryl-CoA isomerase

Chain B:



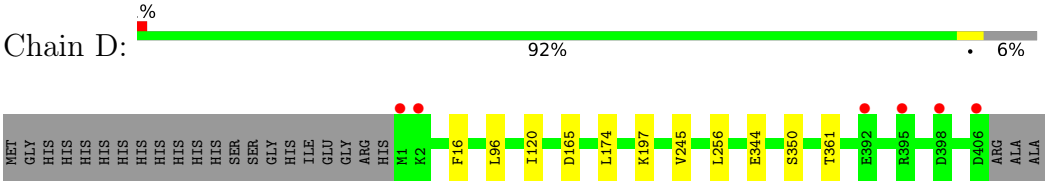
- Molecule 1: 2-methylfumaryl-CoA isomerase

Chain C:



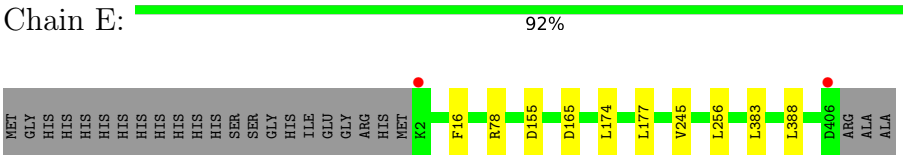
- Molecule 1: 2-methylfumaryl-CoA isomerase

Chain D:



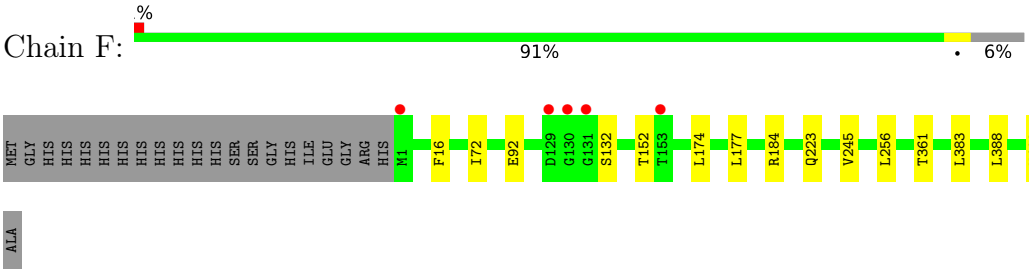
- Molecule 1: 2-methylfumaryl-CoA isomerase

Chain E:



- Molecule 1: 2-methylfumaryl-CoA isomerase

Chain F:


%
91%
6%
MET GLY HIS HIS HIS HIS HIS HIS HIS HIS HIS HIS SER SER GLY HIS ILE GLU GLY ARG HIS M1 F16 I72 E92 D129 G130 G131 S132 T152 T153 L174 L177 R184 Q223 V245 L256 T361 L383 L388 V402 D406 ARG ALA
ALA

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	193.84Å 193.84Å 251.97Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.69 – 2.49 29.69 – 2.49	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.69-2.49) 99.8 (29.69-2.49)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.67 (at 2.48Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.185 , 0.200 0.185 , 0.199	Depositor DCC
R_{free} test set	1998 reflections (1.05%)	wwPDB-VP
Wilson B-factor (Å^2)	47.1	Xtriage
Anisotropy	0.276	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 33.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.010 for -h,-k,l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	20187	wwPDB-VP
Average B, all atoms (Å^2)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.04% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: COA, MEZ, OA9

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	0.24	0/3205	0.51	0/4350
1	B	0.25	0/3205	0.50	0/4350
1	C	0.24	0/3205	0.51	0/4350
1	D	0.24	0/3204	0.51	0/4348
1	E	0.24	0/3197	0.52	1/4340 (0.0%)
1	F	0.25	0/3205	0.50	0/4350
All	All	0.24	0/19221	0.51	1/26088 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	165	ASP	CB-CG-OD2	5.20	122.98	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3134	0	3141	6	0
1	B	3134	0	3141	6	0
1	C	3134	0	3141	8	0
1	D	3133	0	3141	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	3126	0	3129	4	1
1	F	3134	0	3141	10	0
2	A	56	0	0	0	0
3	B	48	0	32	0	0
3	C	48	0	32	0	0
3	D	48	0	31	1	0
3	E	48	0	32	0	0
3	F	48	0	32	1	0
4	B	8	0	4	1	0
4	C	8	0	4	2	0
4	D	9	0	4	1	0
4	E	8	0	4	1	0
4	F	8	0	4	1	0
5	A	173	0	0	0	0
5	B	151	0	0	0	0
5	C	172	0	0	3	0
5	D	177	0	0	1	1
5	E	185	0	0	0	0
5	F	197	0	0	3	0
All	All	20187	0	19013	41	1

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 1.

The worst 5 of 41 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:C:332:ASP:OD2	5:C:601:HOH:O	2.13	0.66
1:F:72:ILE:HD13	1:F:402:VAL:HG12	1.77	0.66
1:D:245:VAL:HG21	1:D:256:LEU:HD13	1.81	0.63
1:F:245:VAL:HG21	1:F:256:LEU:HD13	1.85	0.59
1:E:245:VAL:HG21	1:E:256:LEU:HD13	1.86	0.58

All (1) 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:E:78:ARG:NH2	5:D:644:HOH:O[2_564]	2.17	0.03

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	404/430 (94%)	394 (98%)	9 (2%)	1 (0%)	47 68
1	B	404/430 (94%)	394 (98%)	9 (2%)	1 (0%)	47 68
1	C	404/430 (94%)	395 (98%)	8 (2%)	1 (0%)	47 68
1	D	404/430 (94%)	393 (97%)	10 (2%)	1 (0%)	47 68
1	E	403/430 (94%)	391 (97%)	11 (3%)	1 (0%)	47 68
1	F	404/430 (94%)	392 (97%)	11 (3%)	1 (0%)	47 68
All	All	2423/2580 (94%)	2359 (97%)	58 (2%)	6 (0%)	47 68

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	16	PHE
1	C	16	PHE
1	D	16	PHE
1	E	16	PHE
1	F	16	PHE

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	323/342 (94%)	322 (100%)	1 (0%)	92 97
1	B	323/342 (94%)	323 (100%)	0	100 100
1	C	323/342 (94%)	323 (100%)	0	100 100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	D	323/342 (94%)	322 (100%)	1 (0%)	92 97
1	E	322/342 (94%)	321 (100%)	1 (0%)	92 97
1	F	323/342 (94%)	323 (100%)	0	100 100
All	All	1937/2052 (94%)	1934 (100%)	3 (0%)	93 98

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

Mol	Chain	Res	Type
1	A	165	ASP
1	D	165	ASP
1	E	155	ASP

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)

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

11 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	$\#\|Z\|>2$
3	COA	F	501	-	41,50,50	1.33	4 (9%)	52,75,75	1.88	11 (21%)
3	COA	D	501	1	41,50,50	1.40	5 (12%)	52,75,75	1.60	8 (15%)
3	COA	B	501	-	41,50,50	1.28	3 (7%)	52,75,75	1.59	8 (15%)
4	MEZ	E	502	1	7,7,8	2.06	2 (28%)	7,8,10	4.13	2 (28%)
2	OA9	A	500	-	50,58,58	1.90	13 (26%)	62,86,86	2.19	13 (20%)
4	MEZ	D	502	-	8,8,8	1.45	1 (12%)	10,10,10	1.37	1 (10%)
3	COA	E	501	-	41,50,50	1.30	4 (9%)	52,75,75	1.55	8 (15%)
4	MEZ	C	502	1	7,7,8	1.96	2 (28%)	7,8,10	2.14	2 (28%)
4	MEZ	F	502	1	7,7,8	1.39	1 (14%)	7,8,10	2.68	2 (28%)
4	MEZ	B	502	1	7,7,8	2.80	4 (57%)	7,8,10	6.52	3 (42%)
3	COA	C	501	-	41,50,50	1.25	4 (9%)	52,75,75	1.74	11 (21%)

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
3	COA	F	501	-	-	6/44/64/64	0/3/3/3
3	COA	D	501	1	-	6/44/64/64	0/3/3/3
3	COA	B	501	-	-	8/44/64/64	0/3/3/3
4	MEZ	E	502	1	-	4/6/6/8	-
2	OA9	A	500	-	-	23/52/75/75	0/3/3/3
4	MEZ	D	502	-	-	2/8/8/8	-
3	COA	E	501	-	-	10/44/64/64	0/3/3/3
4	MEZ	C	502	1	-	1/7/7/8	-
4	MEZ	F	502	1	-	1/7/7/8	-
4	MEZ	B	502	1	-	2/6/6/8	-
3	COA	C	501	-	-	11/44/64/64	0/3/3/3

The worst 5 of 43 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	OA9	C07-S1P	5.20	1.91	1.77
4	B	502	MEZ	C5-C3	4.86	1.52	1.46
2	A	500	OA9	P2A-O6A	4.53	1.77	1.59
4	E	502	MEZ	C5-C3	4.12	1.51	1.46
4	B	502	MEZ	O2-C1	4.11	1.33	1.23

The worst 5 of 69 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	502	MEZ	O4-C5-C3	-16.02	108.87	125.16
2	A	500	OA9	C2P-S1P-C07	10.47	112.95	99.80
4	E	502	MEZ	O4-C5-C3	-9.70	115.29	125.16
4	F	502	MEZ	O2-C1-C2	-6.55	115.69	123.66
3	F	501	COA	P2A-O3A-P1A	-6.13	111.78	132.83

There are no chirality outliers.

5 of 74 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	500	OA9	O08-C07-S1P-C2P
2	A	500	OA9	S1P-C2P-C3P-N4P
2	A	500	OA9	C3P-C2P-S1P-C07
2	A	500	OA9	N8P-C9P-CAP-OAP
2	A	500	OA9	C9P-CAP-CBP-CEP

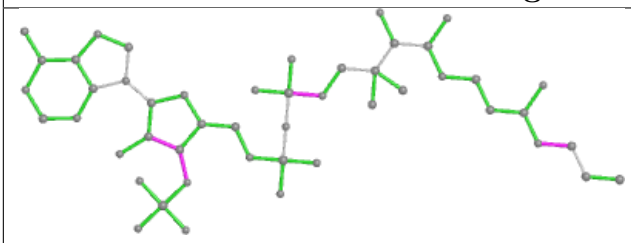
There are no ring outliers.

7 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	501	COA	1	0
3	D	501	COA	1	0
4	E	502	MEZ	1	0
4	D	502	MEZ	1	0
4	C	502	MEZ	2	0
4	F	502	MEZ	1	0
4	B	502	MEZ	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 COA F 501



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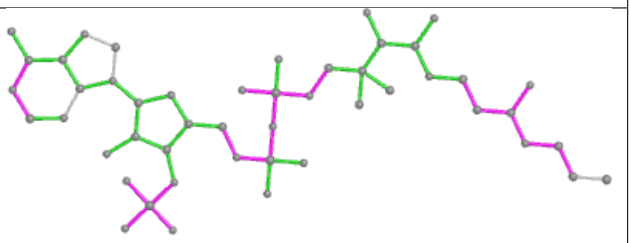

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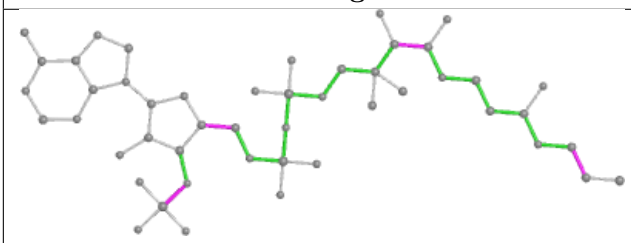

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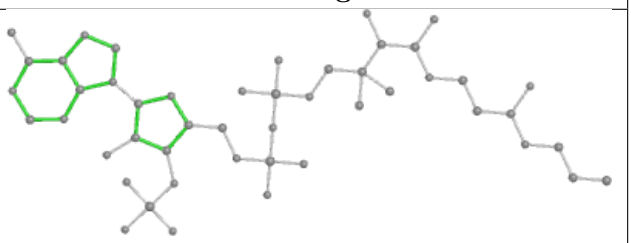
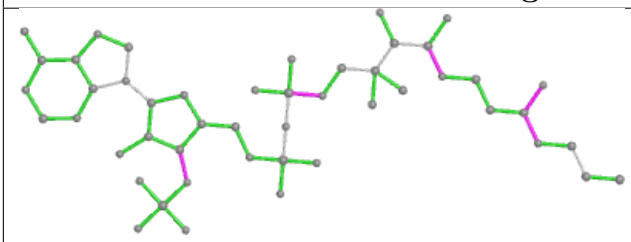


Rings

Ligand COA D 501



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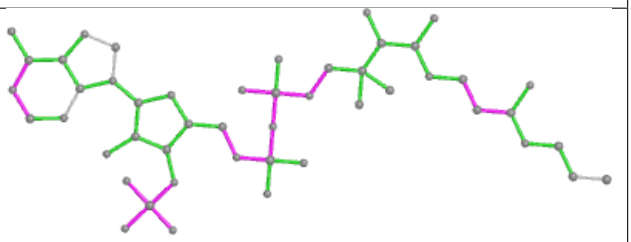

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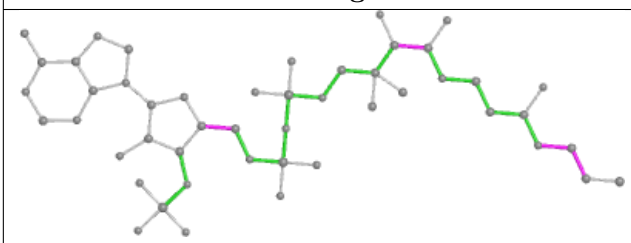

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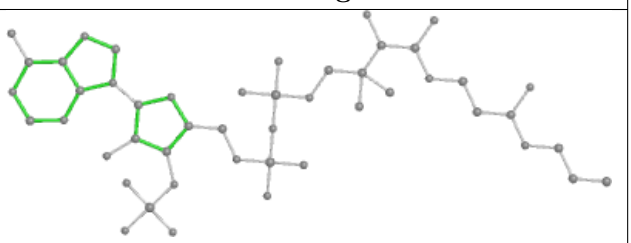

Rings

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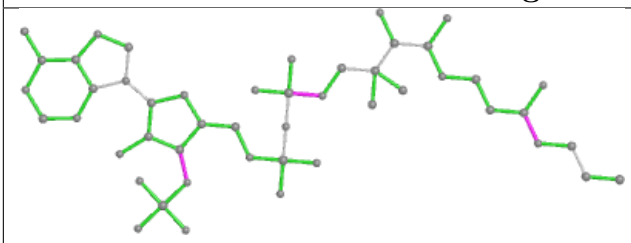

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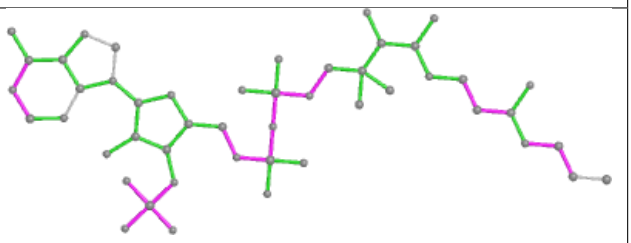

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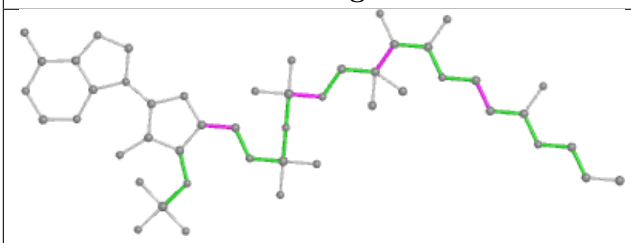

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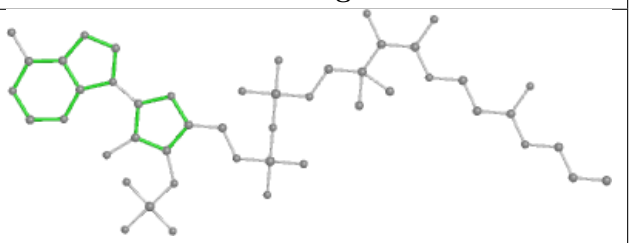
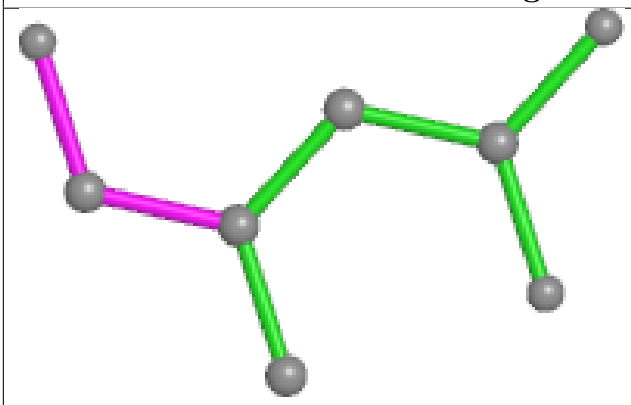


Rings

Ligand MEZ E 502



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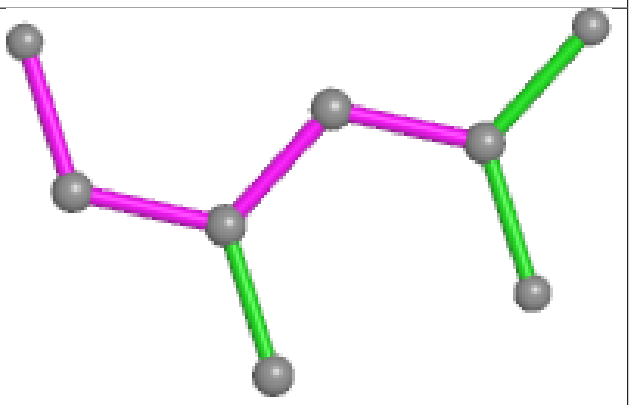

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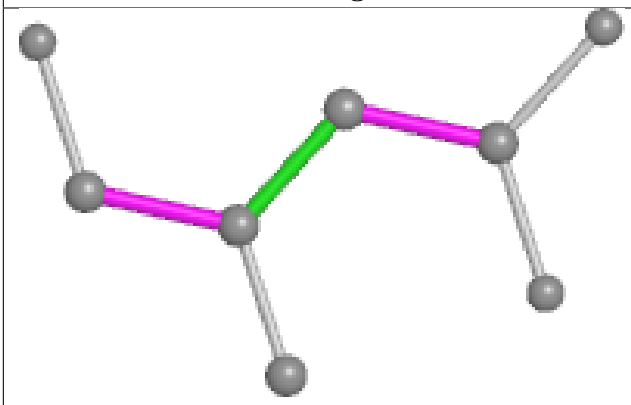

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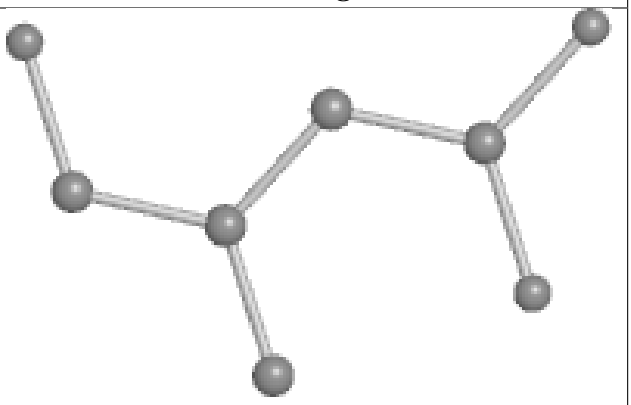
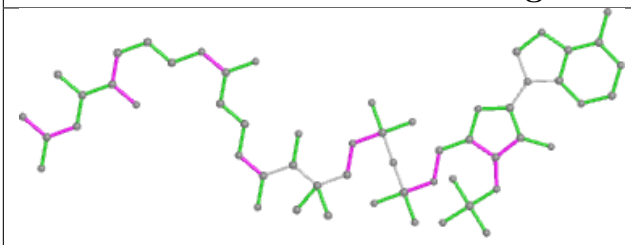


Rings

Ligand OA9 A 500



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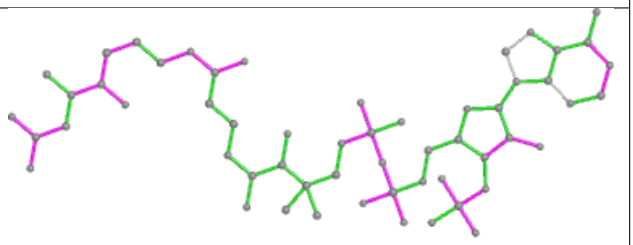

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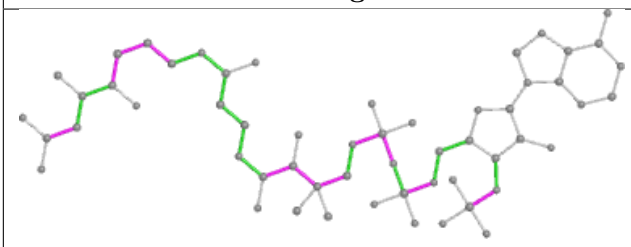

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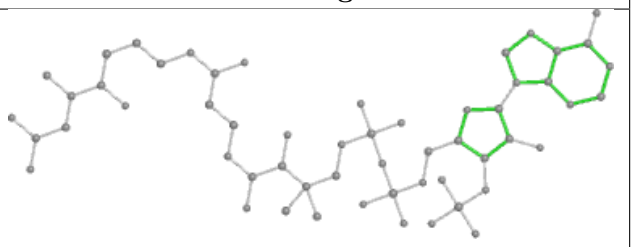

Rings

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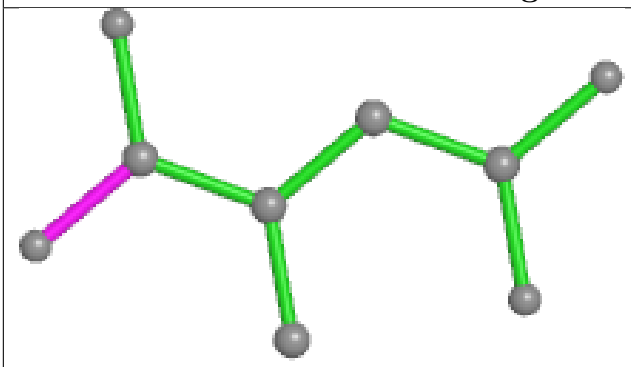

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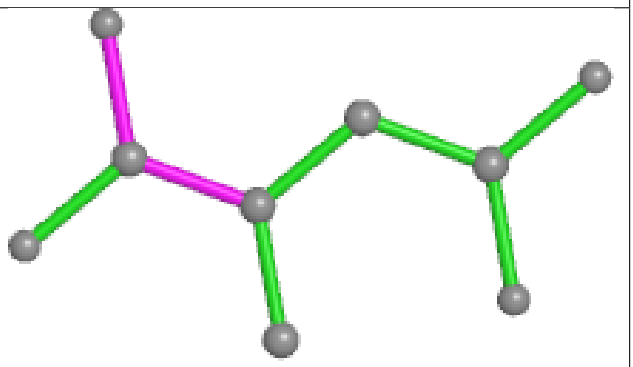

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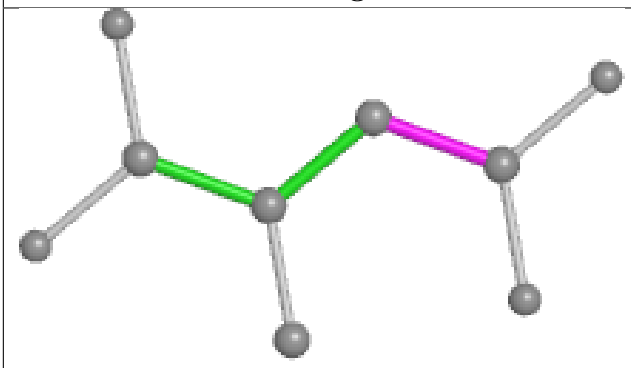

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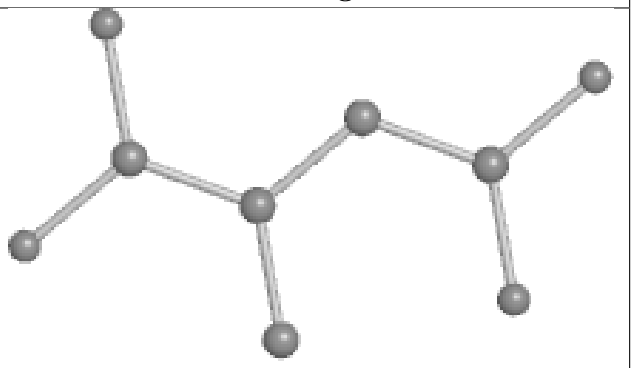
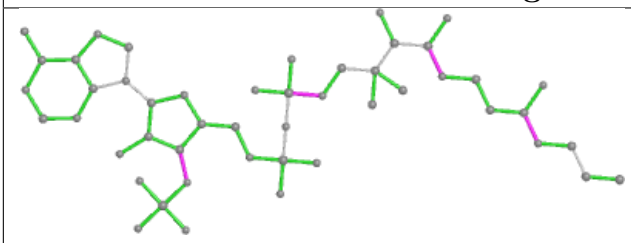


Rings

Ligand COA E 501



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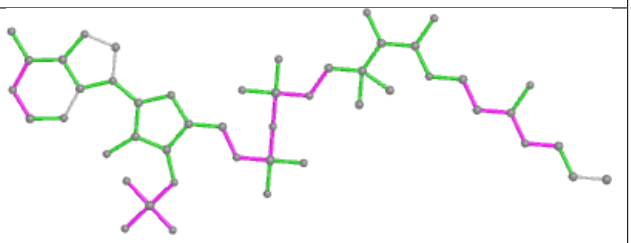

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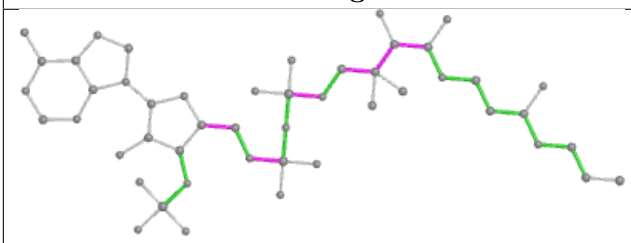

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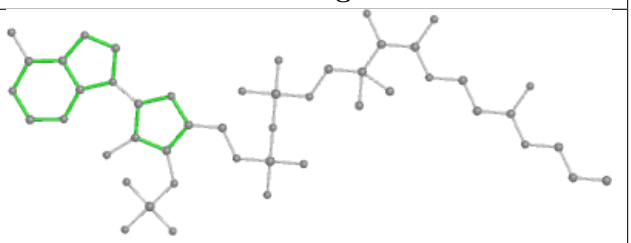

Rings

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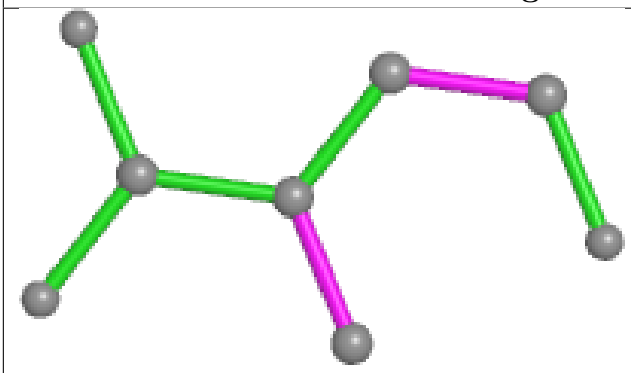

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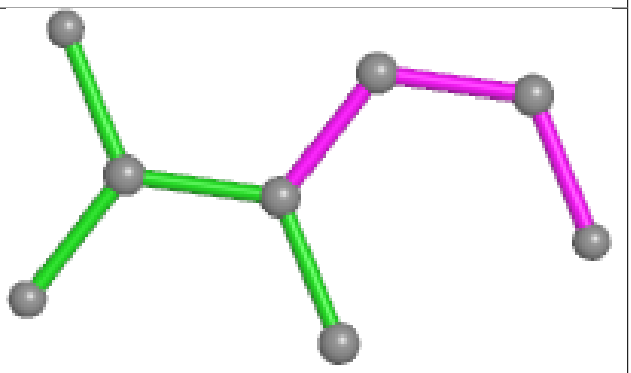
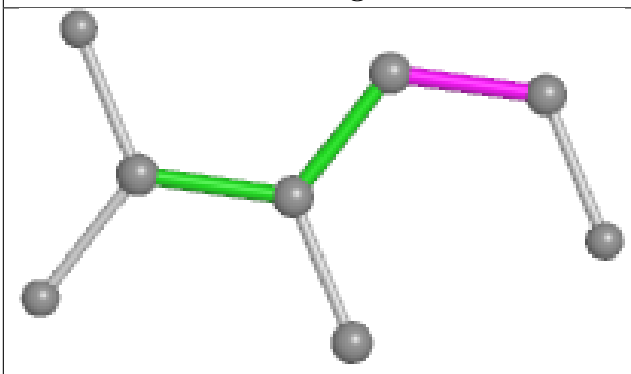


Bond angles



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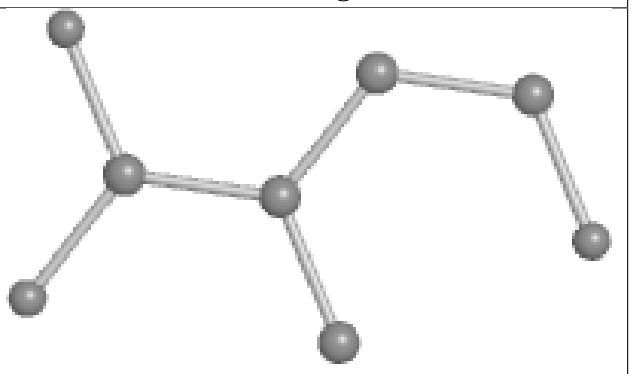
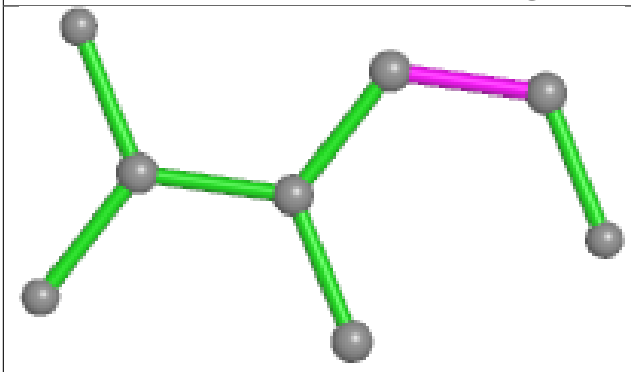
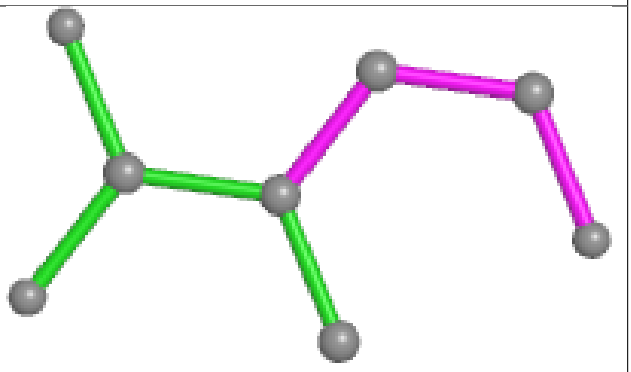


Rings

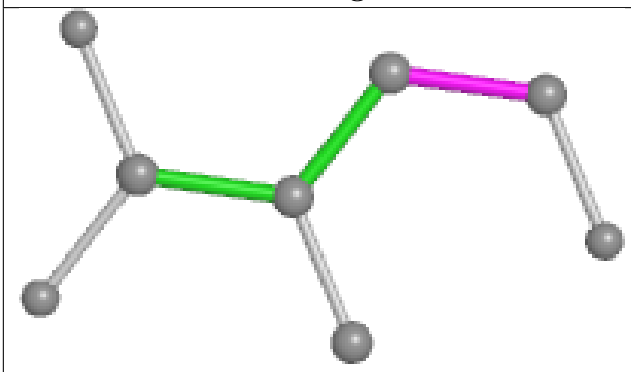
Ligand MEZ F 502



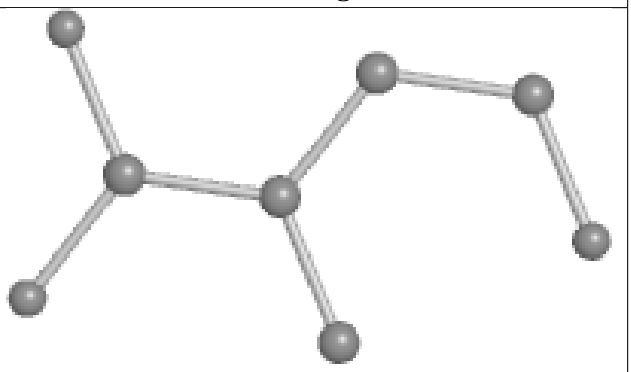
Bond lengths



Bond angles

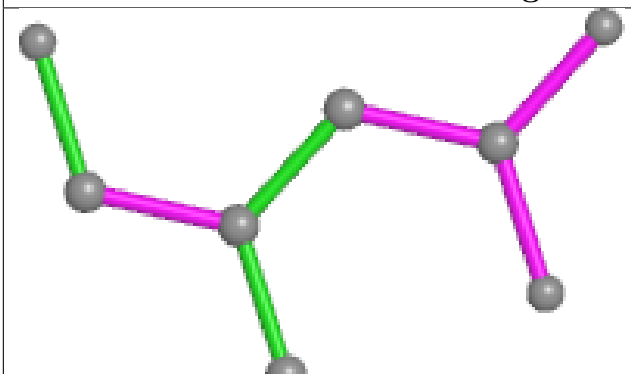


Torsions

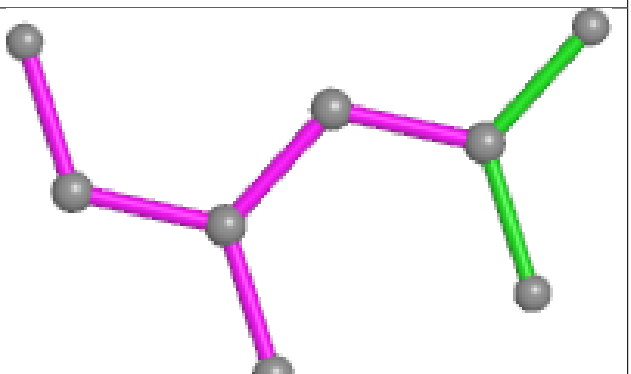


Rings

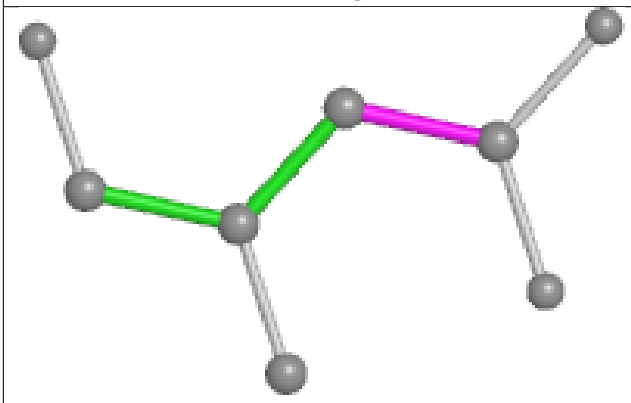
Ligand MEZ B 502



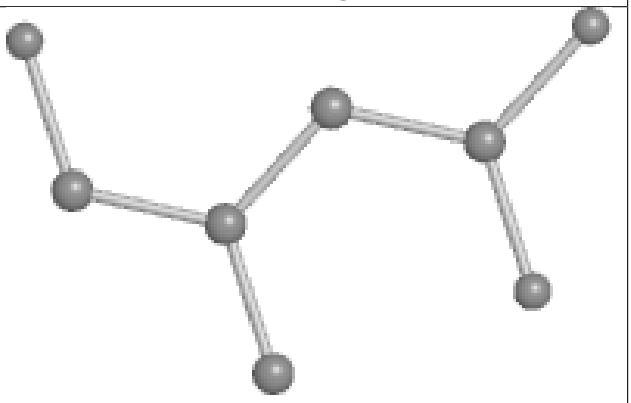
Bond lengths



Bond angles

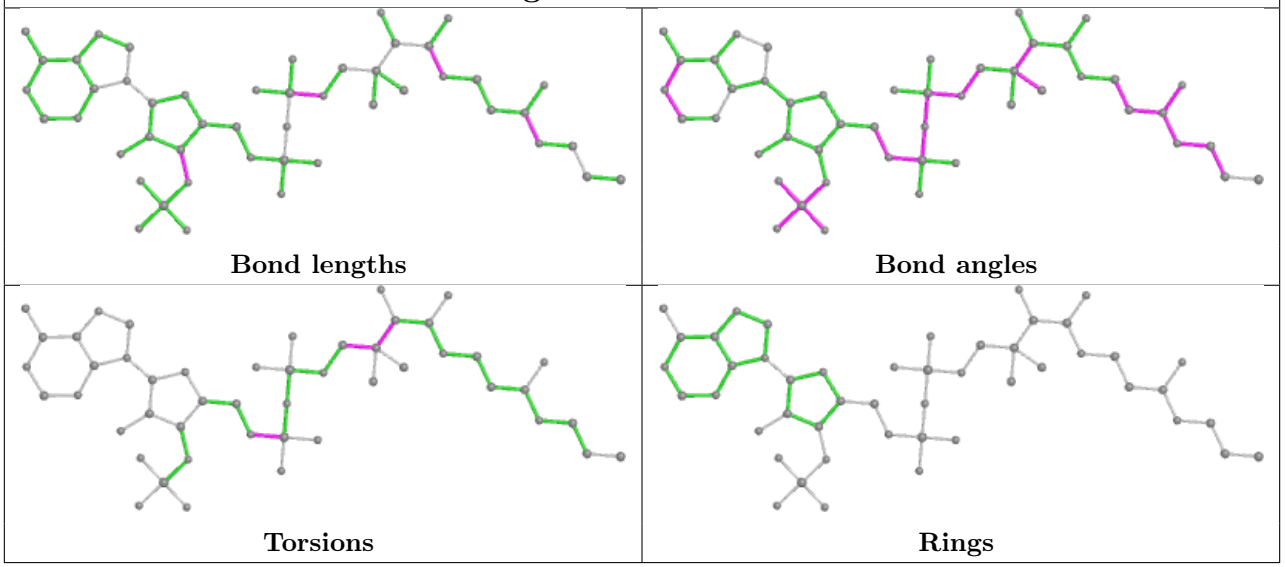


Torsions



Rings

Ligand COA C 501



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

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, 95^{th} 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	406/430 (94%)	-0.40	3 (0%)	87	89	38, 49, 69, 102	0
1	B	406/430 (94%)	-0.36	4 (0%)	82	84	38, 50, 82, 106	0
1	C	406/430 (94%)	-0.30	8 (1%)	65	68	38, 47, 70, 105	0
1	D	406/430 (94%)	-0.42	6 (1%)	73	75	39, 46, 65, 94	0
1	E	405/430 (94%)	-0.47	2 (0%)	91	91	38, 45, 61, 90	0
1	F	406/430 (94%)	-0.38	5 (1%)	79	80	38, 46, 67, 114	0
All	All	2435/2580 (94%)	-0.39	28 (1%)	80	82	38, 47, 70, 114	0

The worst 5 of 28 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	130	GLY	4.3
1	F	131	GLY	3.4
1	A	1	MET	3.4
1	C	1	MET	3.4
1	F	1	MET	3.2

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 monosaccharides in this entry.

6.4 Ligands (i)

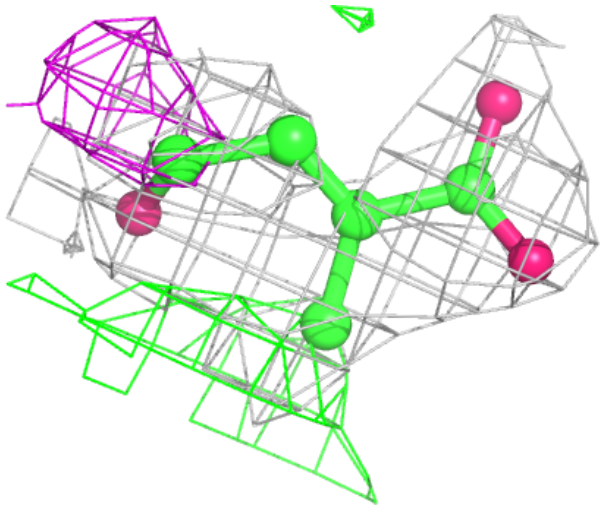
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

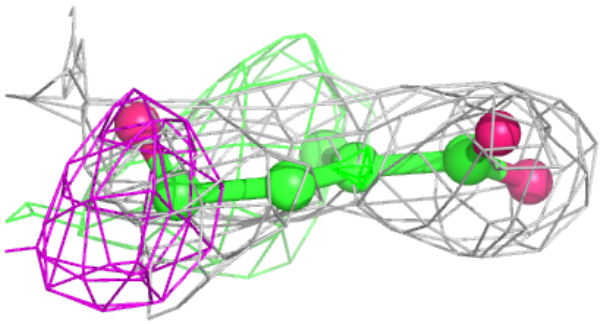
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	MEZ	F	502	8/9	0.67	0.54	49,56,60,62	8
4	MEZ	B	502	8/9	0.69	0.51	46,57,63,66	8
3	COA	F	501	48/48	0.71	0.29	69,82,105,107	48
4	MEZ	E	502	8/9	0.72	0.47	38,53,56,57	8
3	COA	B	501	48/48	0.78	0.29	79,93,114,118	48
4	MEZ	C	502	8/9	0.79	0.26	67,74,76,78	0
4	MEZ	D	502	9/9	0.83	0.40	52,54,59,60	9
3	COA	C	501	48/48	0.85	0.20	65,76,100,102	48
3	COA	E	501	48/48	0.87	0.19	53,62,84,86	48
2	OA9	A	500	56/56	0.93	0.15	44,51,70,73	0
3	COA	D	501	48/48	0.94	0.13	38,47,64,68	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 MEZ F 502:

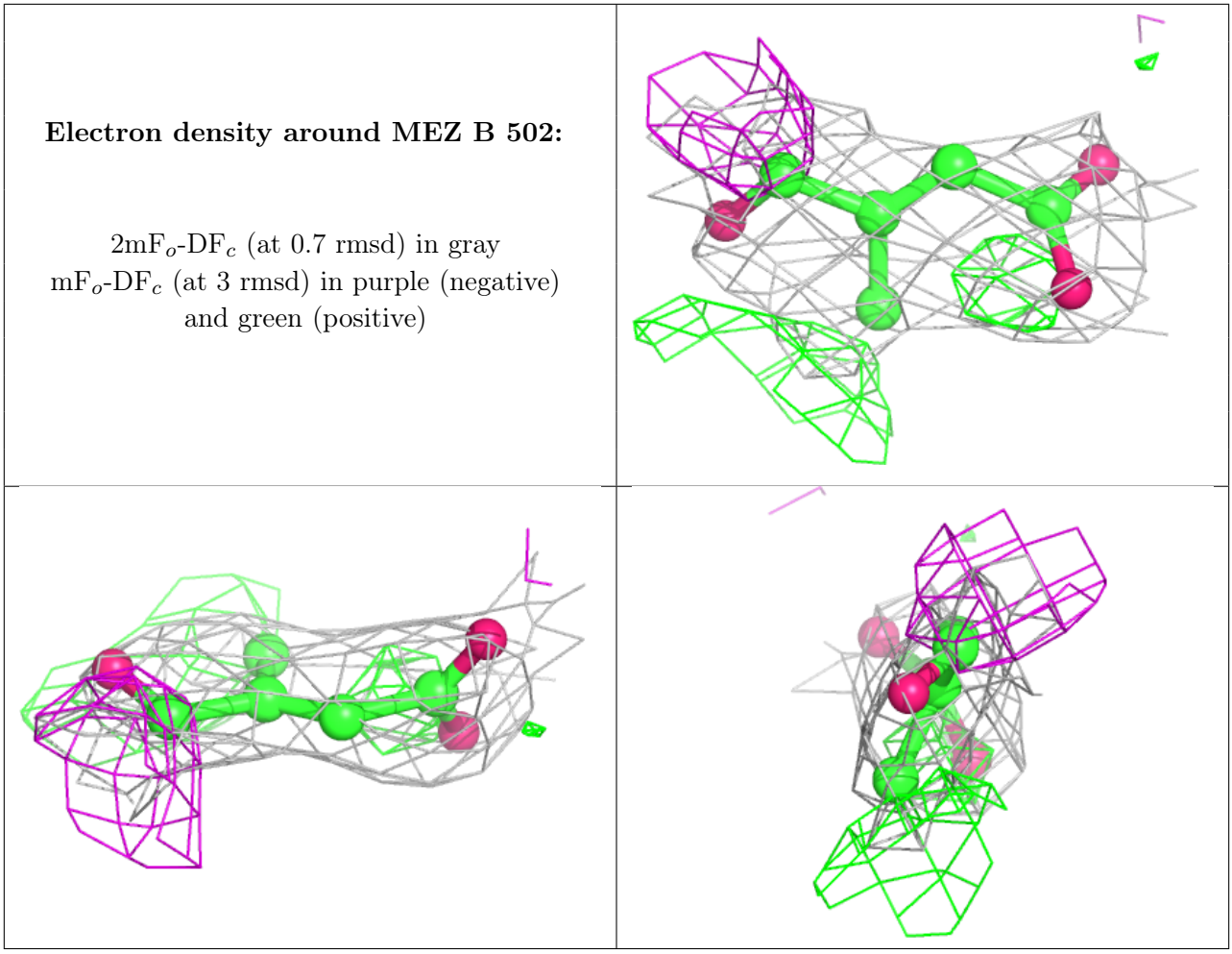
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)





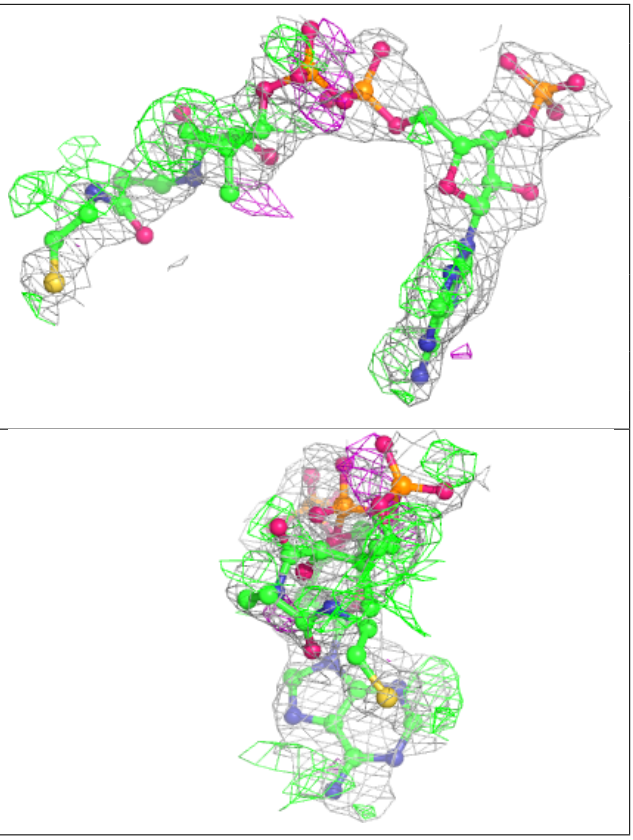
Electron density around MEZ B 502:

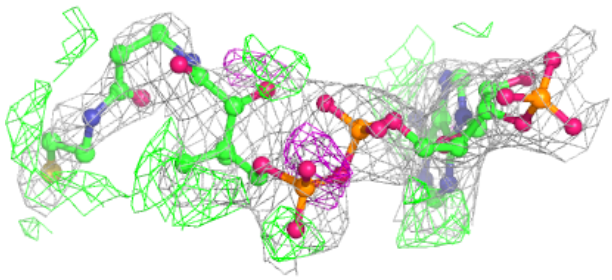
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



Electron density around COA F 501:

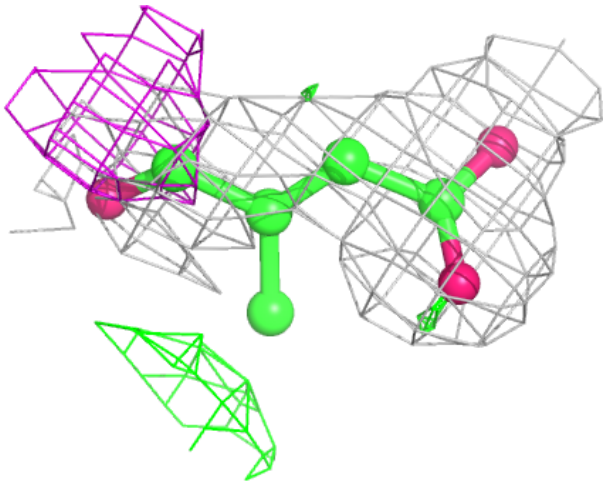
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

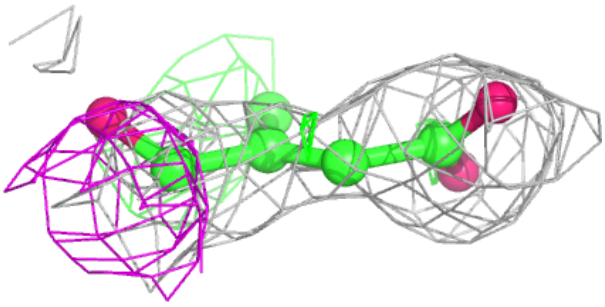


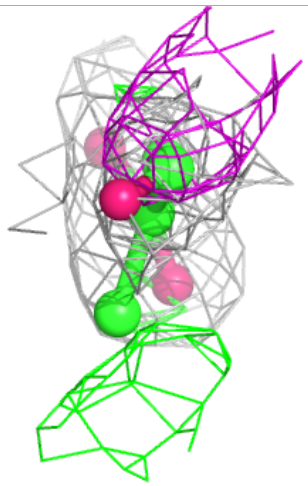


Electron density around MEZ E 502:

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

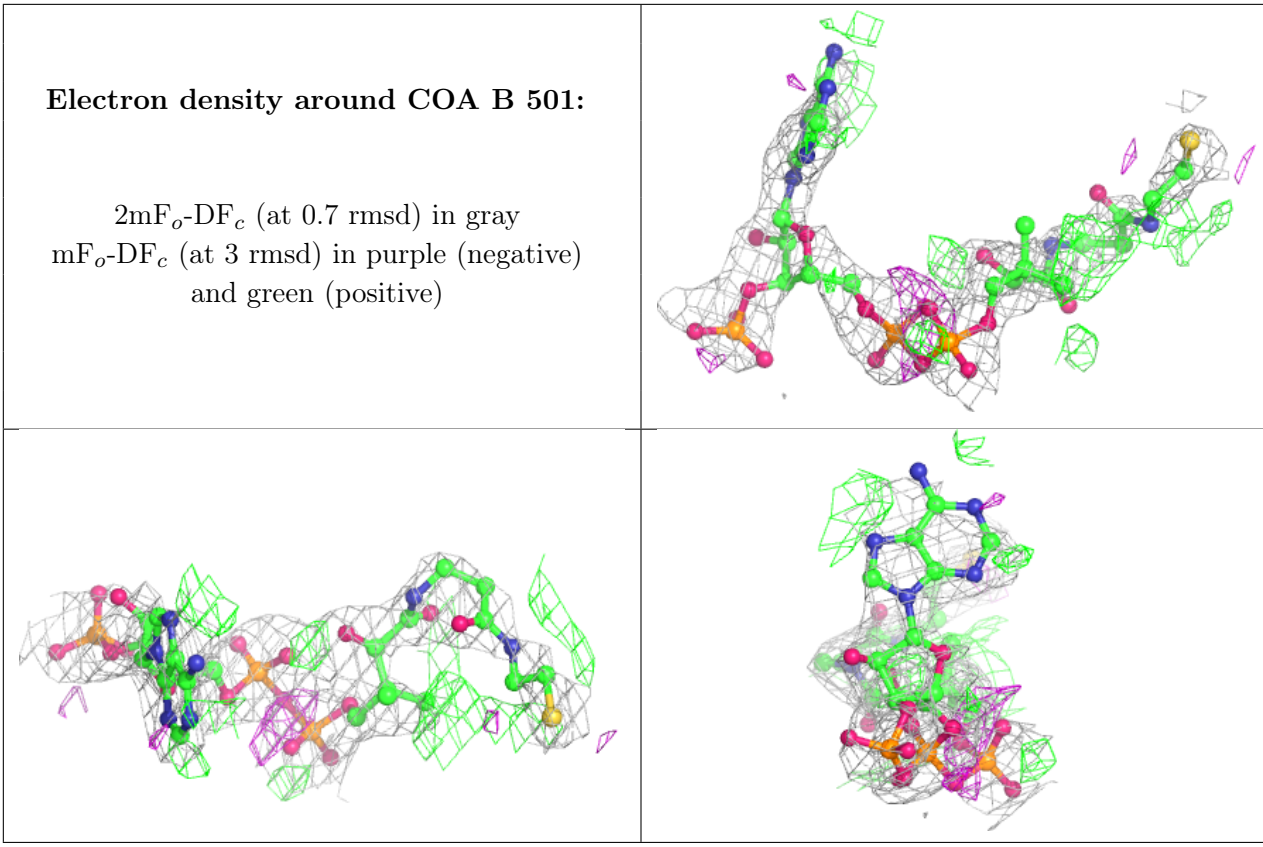






Electron density around COA B 501:

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



Electron density around MEZ C 502:

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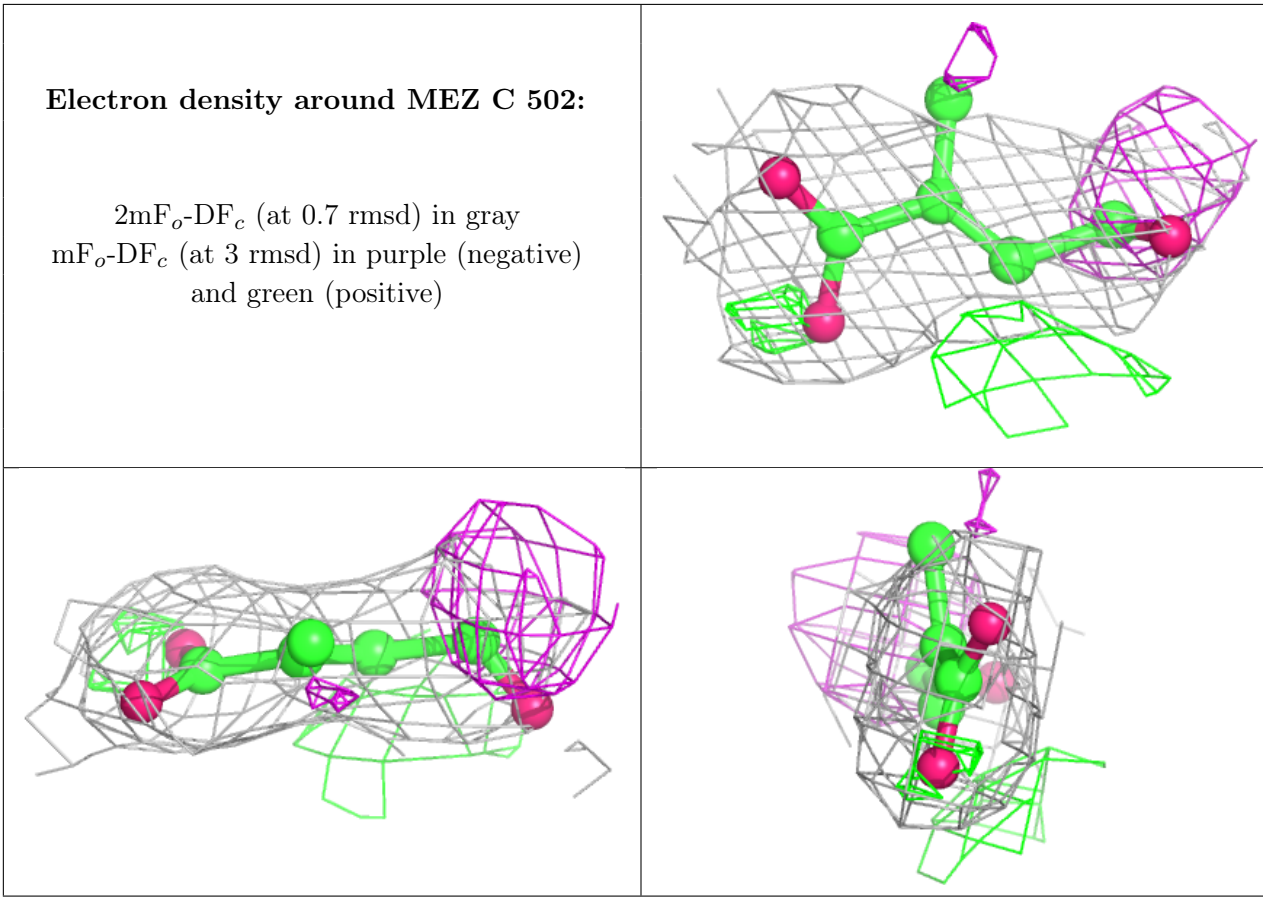
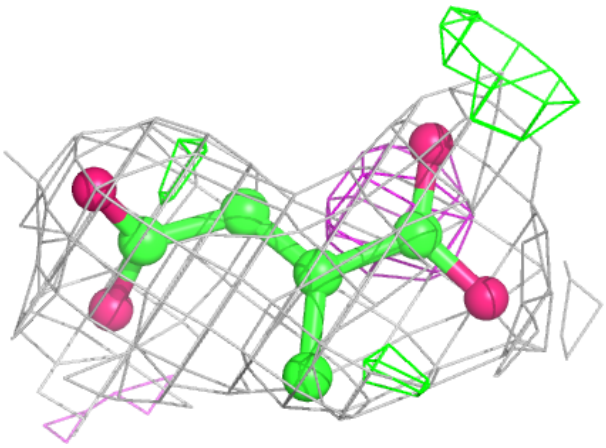
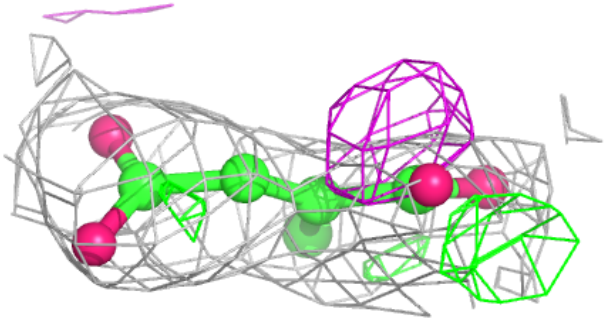
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

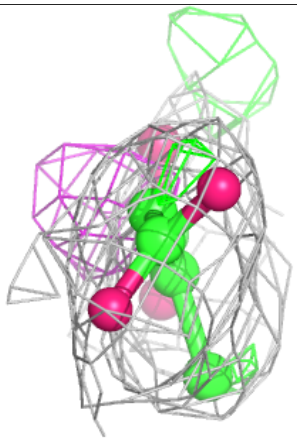


Electron density around MEZ D 502:

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

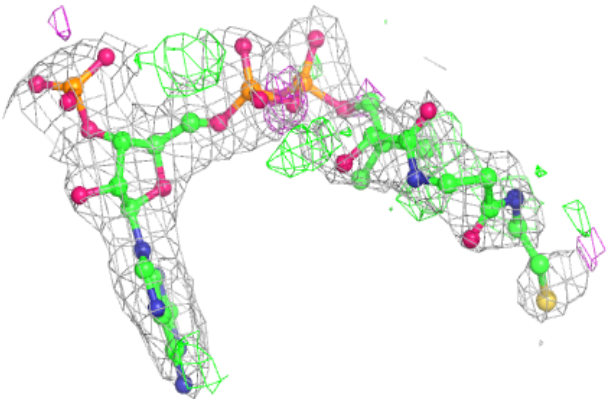


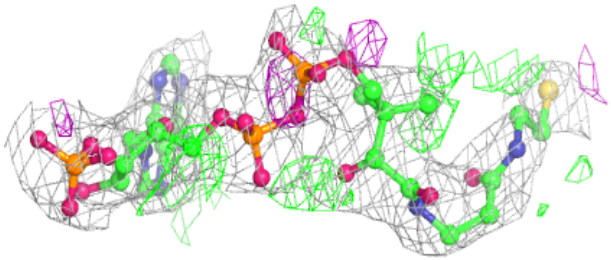


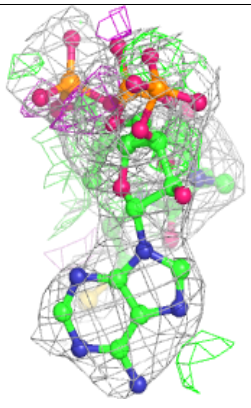


Electron density around COA C 501:

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







Electron density around COA E 501:

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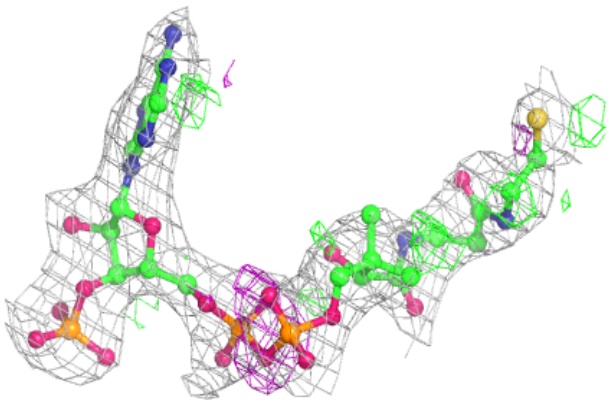
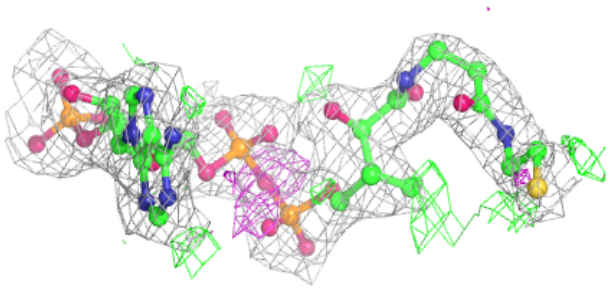
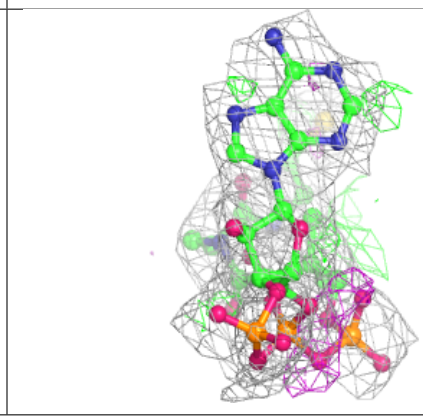
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)







Electron density around OA9 A 500:

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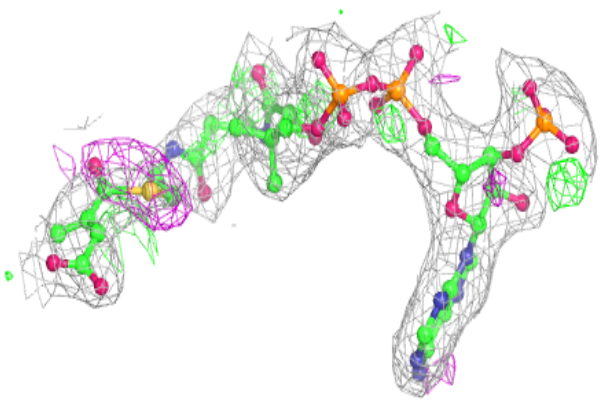
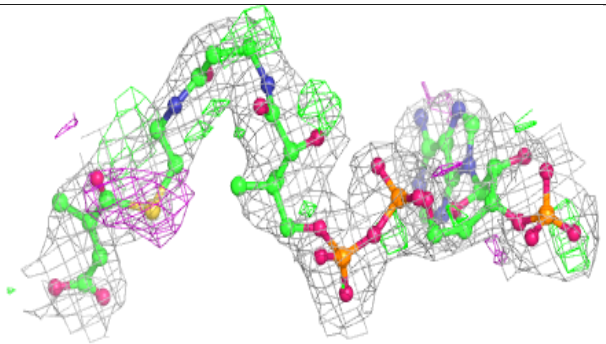
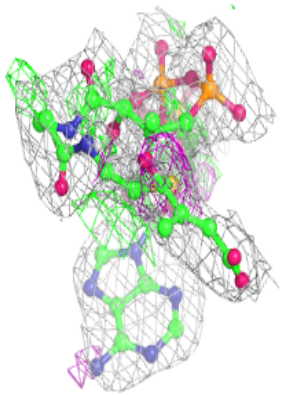
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

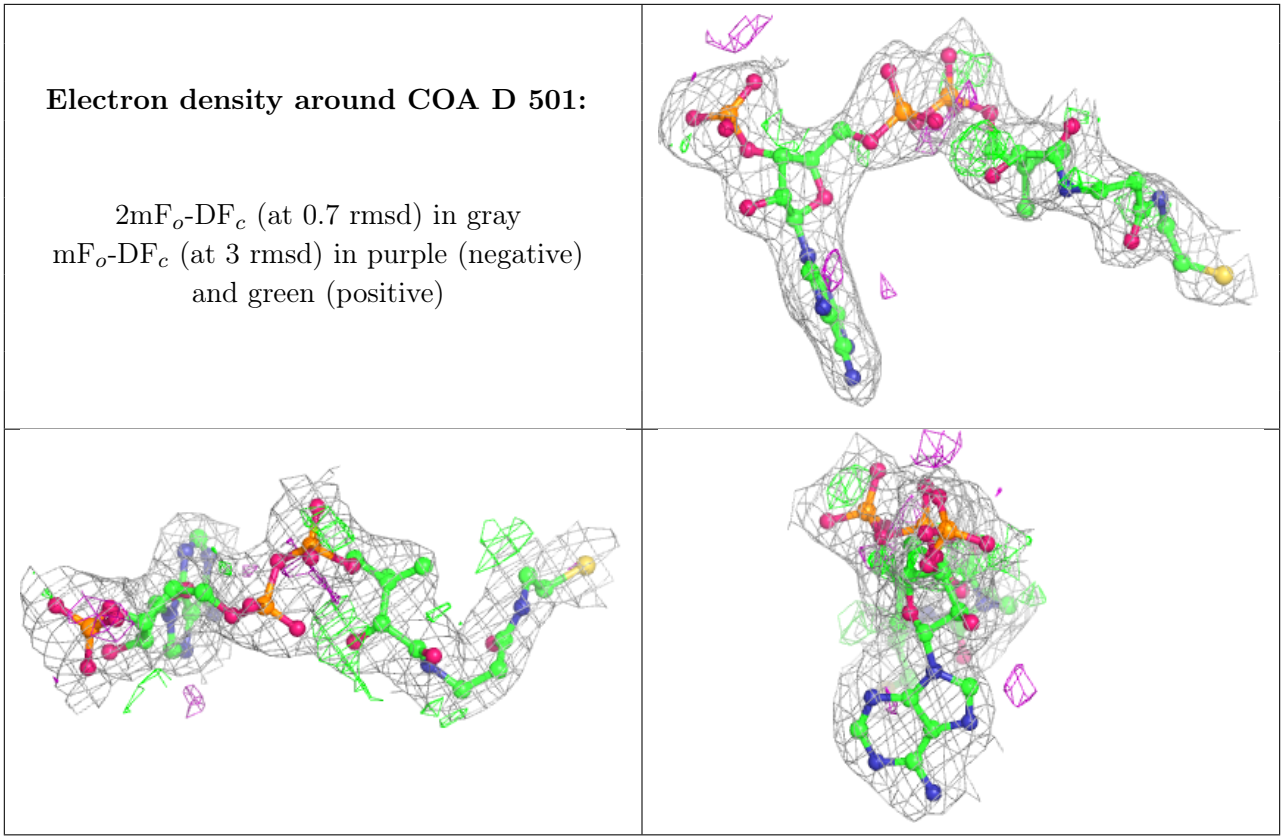






Electron density around COA D 501:

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



6.5 Other polymers (i)

There are no such residues in this entry.