



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 08:41 pm BST

PDB ID : 3ZBN  
Title : Crystal structure of SCP2 thiolase from *Leishmania mexicana*. Complex of the C123A mutant with Coenzyme A.  
Authors : Harijan, R.K.; Kiema, T.-R.; Weiss, M.S.; Michels, P.A.M.; Wierenga, R.K.  
Deposited on : 2012-11-11  
Resolution : 2.45 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

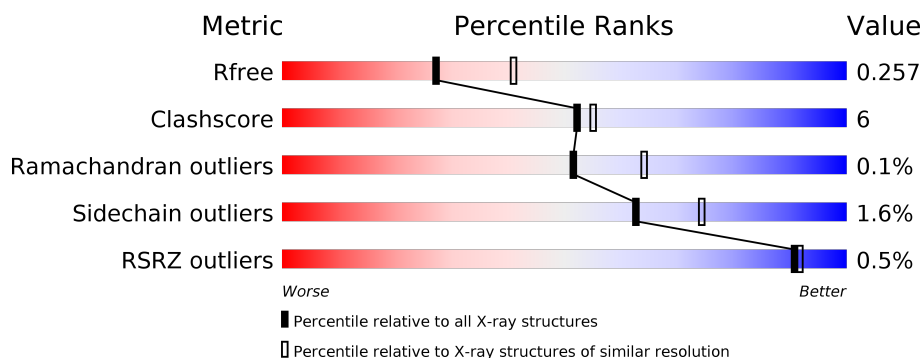
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.45 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	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	457	<div> <div></div> <div>81% 12% 6%</div> </div>
1	B	457	<div> <div></div> <div>83% 10% 6%</div> </div>
1	C	457	<div> <div></div> <div>82% 11% 6%</div> </div>
1	D	457	<div> <div></div> <div>83% 11% 6%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13437 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 3-KETOACYL-COA THIOLASE-LIKE PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	430	Total	C	N	O	S	0	1	0
			3230	2014	566	624	26			
1	B	430	Total	C	N	O	S	0	2	0
			3233	2016	566	625	26			
1	C	430	Total	C	N	O	S	0	1	0
			3227	2012	566	623	26			
1	D	430	Total	C	N	O	S	0	2	0
			3230	2014	566	624	26			

There are 180 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	HIS	-	expression tag	UNP E9AW84
A	-14	HIS	-	expression tag	UNP E9AW84
A	-13	HIS	-	expression tag	UNP E9AW84
A	-12	HIS	-	expression tag	UNP E9AW84
A	-11	HIS	-	expression tag	UNP E9AW84
A	-10	HIS	-	expression tag	UNP E9AW84
A	-9	SER	-	expression tag	UNP E9AW84
A	-8	SER	-	expression tag	UNP E9AW84
A	-7	GLY	-	expression tag	UNP E9AW84
A	-6	LEU	-	expression tag	UNP E9AW84
A	-5	VAL	-	expression tag	UNP E9AW84
A	-4	PRO	-	expression tag	UNP E9AW84
A	-3	ARG	-	expression tag	UNP E9AW84
A	-2	GLY	-	expression tag	UNP E9AW84
A	-1	SER	-	expression tag	UNP E9AW84
A	0	HIS	-	expression tag	UNP E9AW84
A	6	LEU	MET	SEE REMARK 999	UNP E9AW84
A	12	ALA	THR	SEE REMARK 999	UNP E9AW84
A	31	PRO	LYS	SEE REMARK 999	UNP E9AW84
A	34	ILE	VAL	SEE REMARK 999	UNP E9AW84
A	45	LYS	GLN	SEE REMARK 999	UNP E9AW84

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Chain	Residue	Modelled	Actual	Comment	Reference
A	59	GLY	HIS	SEE REMARK 999	UNP E9AW84
A	61	MET	ILE	SEE REMARK 999	UNP E9AW84
A	63	HIS	ASN	SEE REMARK 999	UNP E9AW84
A	64	SER	VAL	SEE REMARK 999	UNP E9AW84
A	65	GLY	SER	SEE REMARK 999	UNP E9AW84
A	67	SER	LYS	SEE REMARK 999	UNP E9AW84
A	69	ARG	LYS	SEE REMARK 999	UNP E9AW84
A	71	GLY	ASP	SEE REMARK 999	UNP E9AW84
A	107	SER	ALA	SEE REMARK 999	UNP E9AW84
A	112	MET	ILE	SEE REMARK 999	UNP E9AW84
A	123	ALA	CYS	engineered mutation	UNP E9AW84
A	134	MET	THR	SEE REMARK 999	UNP E9AW84
A	139	SER	ALA	SEE REMARK 999	UNP E9AW84
A	145	THR	VAL	SEE REMARK 999	UNP E9AW84
A	157	SER	THR	SEE REMARK 999	UNP E9AW84
A	176	GLN	LYS	SEE REMARK 999	UNP E9AW84
A	229	THR	SER	SEE REMARK 999	UNP E9AW84
A	268	VAL	ALA	SEE REMARK 999	UNP E9AW84
A	277	ARG	LYS	SEE REMARK 999	UNP E9AW84
A	327	ILE	VAL	SEE REMARK 999	UNP E9AW84
A	356	GLU	ASP	SEE REMARK 999	UNP E9AW84
A	368	ASP	GLU	SEE REMARK 999	UNP E9AW84
A	398	ILE	VAL	SEE REMARK 999	UNP E9AW84
A	410	GLU	GLY	SEE REMARK 999	UNP E9AW84
B	-15	HIS	-	expression tag	UNP E9AW84
B	-14	HIS	-	expression tag	UNP E9AW84
B	-13	HIS	-	expression tag	UNP E9AW84
B	-12	HIS	-	expression tag	UNP E9AW84
B	-11	HIS	-	expression tag	UNP E9AW84
B	-10	HIS	-	expression tag	UNP E9AW84
B	-9	SER	-	expression tag	UNP E9AW84
B	-8	SER	-	expression tag	UNP E9AW84
B	-7	GLY	-	expression tag	UNP E9AW84
B	-6	LEU	-	expression tag	UNP E9AW84
B	-5	VAL	-	expression tag	UNP E9AW84
B	-4	PRO	-	expression tag	UNP E9AW84
B	-3	ARG	-	expression tag	UNP E9AW84
B	-2	GLY	-	expression tag	UNP E9AW84
B	-1	SER	-	expression tag	UNP E9AW84
B	0	HIS	-	expression tag	UNP E9AW84
B	6	LEU	MET	SEE REMARK 999	UNP E9AW84
B	12	ALA	THR	SEE REMARK 999	UNP E9AW84

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Chain	Residue	Modelled	Actual	Comment	Reference
B	31	PRO	LYS	SEE REMARK 999	UNP E9AW84
B	34	ILE	VAL	SEE REMARK 999	UNP E9AW84
B	45	LYS	GLN	SEE REMARK 999	UNP E9AW84
B	59	GLY	HIS	SEE REMARK 999	UNP E9AW84
B	61	MET	ILE	SEE REMARK 999	UNP E9AW84
B	63	HIS	ASN	SEE REMARK 999	UNP E9AW84
B	64	SER	VAL	SEE REMARK 999	UNP E9AW84
B	65	GLY	SER	SEE REMARK 999	UNP E9AW84
B	67	SER	LYS	SEE REMARK 999	UNP E9AW84
B	69	ARG	LYS	SEE REMARK 999	UNP E9AW84
B	71	GLY	ASP	SEE REMARK 999	UNP E9AW84
B	107	SER	ALA	SEE REMARK 999	UNP E9AW84
B	112	MET	ILE	SEE REMARK 999	UNP E9AW84
B	123	ALA	CYS	engineered mutation	UNP E9AW84
B	134	MET	THR	SEE REMARK 999	UNP E9AW84
B	139	SER	ALA	SEE REMARK 999	UNP E9AW84
B	145	THR	VAL	SEE REMARK 999	UNP E9AW84
B	157	SER	THR	SEE REMARK 999	UNP E9AW84
B	176	GLN	LYS	SEE REMARK 999	UNP E9AW84
B	229	THR	SER	SEE REMARK 999	UNP E9AW84
B	268	VAL	ALA	SEE REMARK 999	UNP E9AW84
B	277	ARG	LYS	SEE REMARK 999	UNP E9AW84
B	327	ILE	VAL	SEE REMARK 999	UNP E9AW84
B	356	GLU	ASP	SEE REMARK 999	UNP E9AW84
B	368	ASP	GLU	SEE REMARK 999	UNP E9AW84
B	398	ILE	VAL	SEE REMARK 999	UNP E9AW84
B	410	GLU	GLY	SEE REMARK 999	UNP E9AW84
C	-15	HIS	-	expression tag	UNP E9AW84
C	-14	HIS	-	expression tag	UNP E9AW84
C	-13	HIS	-	expression tag	UNP E9AW84
C	-12	HIS	-	expression tag	UNP E9AW84
C	-11	HIS	-	expression tag	UNP E9AW84
C	-10	HIS	-	expression tag	UNP E9AW84
C	-9	SER	-	expression tag	UNP E9AW84
C	-8	SER	-	expression tag	UNP E9AW84
C	-7	GLY	-	expression tag	UNP E9AW84
C	-6	LEU	-	expression tag	UNP E9AW84
C	-5	VAL	-	expression tag	UNP E9AW84
C	-4	PRO	-	expression tag	UNP E9AW84
C	-3	ARG	-	expression tag	UNP E9AW84
C	-2	GLY	-	expression tag	UNP E9AW84
C	-1	SER	-	expression tag	UNP E9AW84

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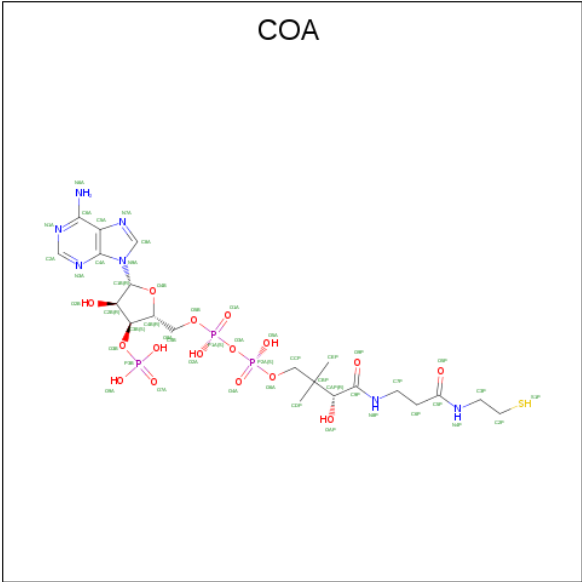
Chain	Residue	Modelled	Actual	Comment	Reference
C	0	HIS	-	expression tag	UNP E9AW84
C	6	LEU	MET	SEE REMARK 999	UNP E9AW84
C	12	ALA	THR	SEE REMARK 999	UNP E9AW84
C	31	PRO	LYS	SEE REMARK 999	UNP E9AW84
C	34	ILE	VAL	SEE REMARK 999	UNP E9AW84
C	45	LYS	GLN	SEE REMARK 999	UNP E9AW84
C	59	GLY	HIS	SEE REMARK 999	UNP E9AW84
C	61	MET	ILE	SEE REMARK 999	UNP E9AW84
C	63	HIS	ASN	SEE REMARK 999	UNP E9AW84
C	64	SER	VAL	SEE REMARK 999	UNP E9AW84
C	65	GLY	SER	SEE REMARK 999	UNP E9AW84
C	67	SER	LYS	SEE REMARK 999	UNP E9AW84
C	69	ARG	LYS	SEE REMARK 999	UNP E9AW84
C	71	GLY	ASP	SEE REMARK 999	UNP E9AW84
C	107	SER	ALA	SEE REMARK 999	UNP E9AW84
C	112	MET	ILE	SEE REMARK 999	UNP E9AW84
C	123	ALA	CYS	engineered mutation	UNP E9AW84
C	134	MET	THR	SEE REMARK 999	UNP E9AW84
C	139	SER	ALA	SEE REMARK 999	UNP E9AW84
C	145	THR	VAL	SEE REMARK 999	UNP E9AW84
C	157	SER	THR	SEE REMARK 999	UNP E9AW84
C	176	GLN	LYS	SEE REMARK 999	UNP E9AW84
C	229	THR	SER	SEE REMARK 999	UNP E9AW84
C	268	VAL	ALA	SEE REMARK 999	UNP E9AW84
C	277	ARG	LYS	SEE REMARK 999	UNP E9AW84
C	327	ILE	VAL	SEE REMARK 999	UNP E9AW84
C	356	GLU	ASP	SEE REMARK 999	UNP E9AW84
C	368	ASP	GLU	SEE REMARK 999	UNP E9AW84
C	398	ILE	VAL	SEE REMARK 999	UNP E9AW84
C	410	GLU	GLY	SEE REMARK 999	UNP E9AW84
D	-15	HIS	-	expression tag	UNP E9AW84
D	-14	HIS	-	expression tag	UNP E9AW84
D	-13	HIS	-	expression tag	UNP E9AW84
D	-12	HIS	-	expression tag	UNP E9AW84
D	-11	HIS	-	expression tag	UNP E9AW84
D	-10	HIS	-	expression tag	UNP E9AW84
D	-9	SER	-	expression tag	UNP E9AW84
D	-8	SER	-	expression tag	UNP E9AW84
D	-7	GLY	-	expression tag	UNP E9AW84
D	-6	LEU	-	expression tag	UNP E9AW84
D	-5	VAL	-	expression tag	UNP E9AW84
D	-4	PRO	-	expression tag	UNP E9AW84

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-3	ARG	-	expression tag	UNP E9AW84
D	-2	GLY	-	expression tag	UNP E9AW84
D	-1	SER	-	expression tag	UNP E9AW84
D	0	HIS	-	expression tag	UNP E9AW84
D	6	LEU	MET	SEE REMARK 999	UNP E9AW84
D	12	ALA	THR	SEE REMARK 999	UNP E9AW84
D	31	PRO	LYS	SEE REMARK 999	UNP E9AW84
D	34	ILE	VAL	SEE REMARK 999	UNP E9AW84
D	45	LYS	GLN	SEE REMARK 999	UNP E9AW84
D	59	GLY	HIS	SEE REMARK 999	UNP E9AW84
D	61	MET	ILE	SEE REMARK 999	UNP E9AW84
D	63	HIS	ASN	SEE REMARK 999	UNP E9AW84
D	64	SER	VAL	SEE REMARK 999	UNP E9AW84
D	65	GLY	SER	SEE REMARK 999	UNP E9AW84
D	67	SER	LYS	SEE REMARK 999	UNP E9AW84
D	69	ARG	LYS	SEE REMARK 999	UNP E9AW84
D	71	GLY	ASP	SEE REMARK 999	UNP E9AW84
D	107	SER	ALA	SEE REMARK 999	UNP E9AW84
D	112	MET	ILE	SEE REMARK 999	UNP E9AW84
D	123	ALA	CYS	engineered mutation	UNP E9AW84
D	134	MET	THR	SEE REMARK 999	UNP E9AW84
D	139	SER	ALA	SEE REMARK 999	UNP E9AW84
D	145	THR	VAL	SEE REMARK 999	UNP E9AW84
D	157	SER	THR	SEE REMARK 999	UNP E9AW84
D	176	GLN	LYS	SEE REMARK 999	UNP E9AW84
D	229	THR	SER	SEE REMARK 999	UNP E9AW84
D	268	VAL	ALA	SEE REMARK 999	UNP E9AW84
D	277	ARG	LYS	SEE REMARK 999	UNP E9AW84
D	327	ILE	VAL	SEE REMARK 999	UNP E9AW84
D	356	GLU	ASP	SEE REMARK 999	UNP E9AW84
D	368	ASP	GLU	SEE REMARK 999	UNP E9AW84
D	398	ILE	VAL	SEE REMARK 999	UNP E9AW84
D	410	GLU	GLY	SEE REMARK 999	UNP E9AW84

- Molecule 2 is COENZYME A (three-letter code: COA) (formula: C<sub>21</sub>H<sub>36</sub>N<sub>7</sub>O<sub>16</sub>P<sub>3</sub>S).



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

- Molecule 3 is water.

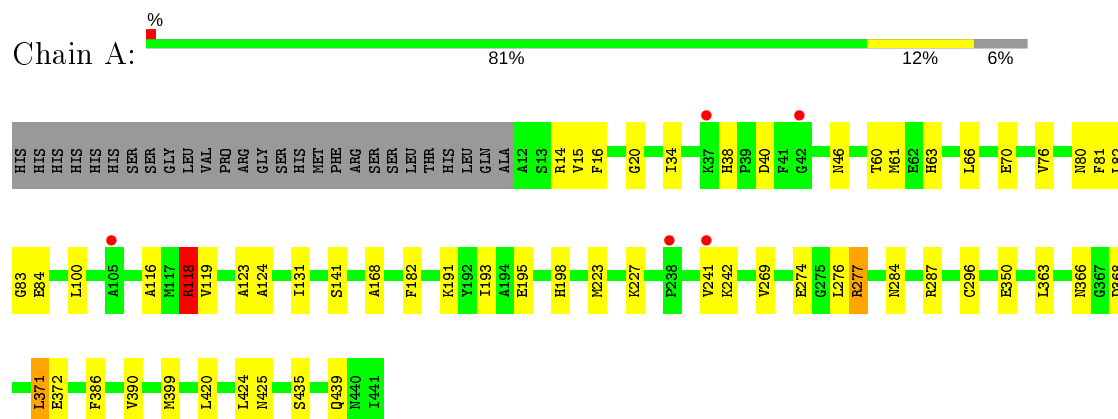
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	76	Total	O	0	0
			76	76		
3	B	118	Total	O	0	0
			118	118		
3	C	85	Total	O	0	0
			85	85		
3	D	94	Total	O	0	0
			94	94		



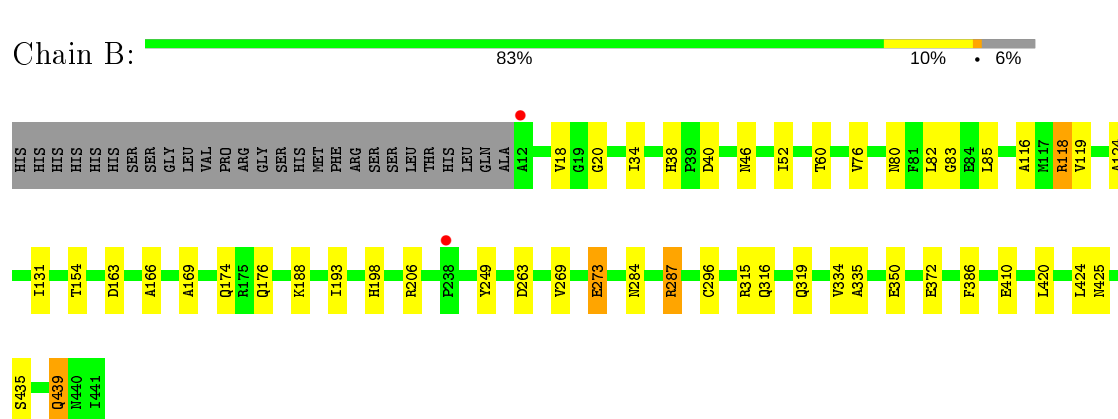
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

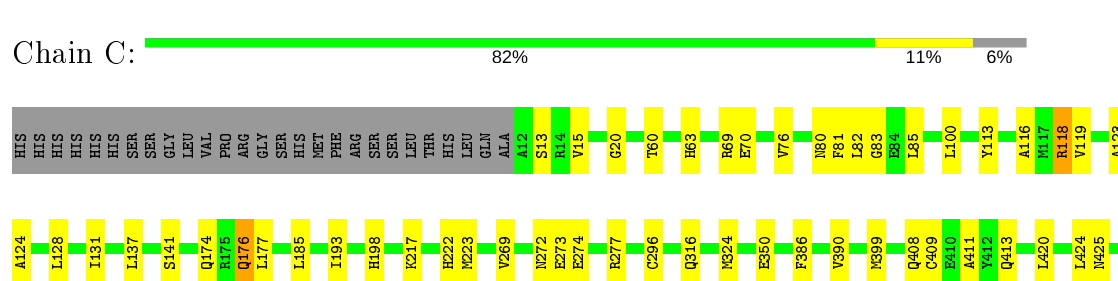
#### • Molecule 1: 3-KETOACYL-COA THIOLASE-LIKE PROTEIN



#### • Molecule 1: 3-KETOACYL-COA THIOLASE-LIKE PROTEIN

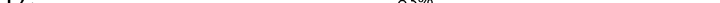


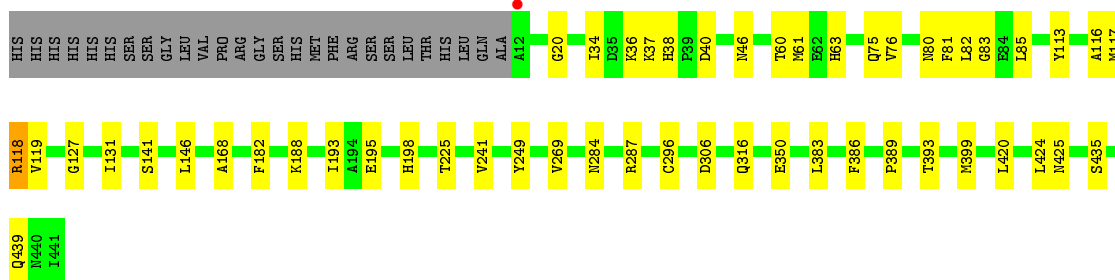
#### • Molecule 1: 3-KETOACYL-COA THIOLASE-LIKE PROTEIN





- Molecule 1: 3-KETOACYL-COA THIOLASE-LIKE PROTEIN

Chain D:  83% 11% 6%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.26Å 89.83Å 126.24Å 90.00° 105.80° 90.00°	Depositor
Resolution (Å)	121.47 – 2.45 60.74 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.7 (121.47-2.45) 99.8 (60.74-2.45)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.60 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, $R_{free}$	0.209 , 0.254 0.212 , 0.257	Depositor DCC
$R_{free}$ test set	3195 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	16.5	Xtriage
Anisotropy	0.382	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 31.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.009 for h,-k,-h-l	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	13437	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.01% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

### 5.1 Standard geometry

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

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	0.65	0/3286	0.79	4/4429 (0.1%)
1	B	0.70	0/3292	0.79	4/4437 (0.1%)
1	C	0.67	0/3283	0.76	1/4426 (0.0%)
1	D	0.65	0/3289	0.77	1/4433 (0.0%)
All	All	0.67	0/13150	0.78	10/17725 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	223	MET	CG-SD-CE	6.76	111.01	100.20
1	A	14	ARG	N-CA-C	6.67	129.00	111.00
1	B	118	ARG	NE-CZ-NH1	6.60	123.60	120.30
1	A	277	ARG	NE-CZ-NH1	6.16	123.38	120.30
1	B	206	ARG	NE-CZ-NH1	5.68	123.14	120.30
1	B	287	ARG	CG-CD-NE	5.53	123.41	111.80
1	A	118	ARG	NE-CZ-NH2	-5.50	117.55	120.30
1	D	75	GLN	CB-CA-C	-5.43	99.53	110.40
1	B	118	ARG	NE-CZ-NH2	-5.38	117.61	120.30
1	C	118	ARG	NE-CZ-NH2	-5.21	117.70	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	389	PRO	Peptide

## 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	3230	0	3208	37	1
1	B	3233	0	3213	32	0
1	C	3227	0	3207	39	0
1	D	3230	0	3212	36	0
2	A	48	0	32	4	0
2	B	48	0	32	3	0
2	D	48	0	32	4	0
3	A	76	0	0	2	0
3	B	118	0	0	5	1
3	C	85	0	0	4	0
3	D	94	0	0	3	0
All	All	13437	0	12936	144	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 6.

All (144) 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:61:MET:HE2	1:A:70:GLU:CG	1.87	1.03
1:A:61:MET:HE2	1:A:70:GLU:HG3	1.37	1.02
2:A:1442:COA:O9P	2:A:1442:COA:H131	1.72	0.87
1:B:18:VAL:O	1:B:287:ARG:NH1	2.07	0.86
1:A:274:GLU:OE2	1:A:277:ARG:NH2	2.09	0.86
1:A:123:ALA:HA	1:A:390:VAL:HG13	1.61	0.81
1:D:80:ASN:HD21	1:D:83:GLY:H	1.25	0.79
1:C:80:ASN:HD21	1:C:83:GLY:H	1.30	0.79
1:A:61:MET:CE	1:A:70:GLU:HG3	2.13	0.79
2:B:1442:COA:S1P	3:B:2100:HOH:O	2.48	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:ASN:HD21	1:A:83:GLY:H	1.38	0.71
1:B:34:ILE:H	1:B:46:ASN:HD21	1.39	0.70
1:B:80:ASN:HD21	1:B:83:GLY:H	1.39	0.69
2:A:1442:COA:CDP	2:A:1442:COA:O9P	2.43	0.67
1:C:119:VAL:HG12	1:D:119:VAL:HG12	1.75	0.67
1:A:193:ILE:HG23	1:A:198:HIS:HB3	1.78	0.65
1:C:193:ILE:HG23	1:C:198:HIS:HB3	1.78	0.64
1:C:176:GLN:NE2	1:C:177:LEU:HD22	2.12	0.64
1:A:420:LEU:HD21	1:A:439:GLN:HE21	1.64	0.63
1:D:193:ILE:HG23	1:D:198:HIS:HB3	1.81	0.63
1:D:61:MET:CE	1:D:146:LEU:HD13	2.30	0.62
2:D:1442:COA:O9P	2:D:1442:COA:CDP	2.47	0.62
1:A:424:LEU:HD13	1:A:435:SER:HB2	1.82	0.61
1:B:273:GLU:OE1	3:B:2001:HOH:O	2.16	0.61
1:D:80:ASN:HD21	1:D:83:GLY:N	1.97	0.61
1:D:195:GLU:HG3	3:D:2044:HOH:O	2.00	0.60
1:D:60:THR:HG22	1:D:269:VAL:HG23	1.84	0.59
1:D:80:ASN:ND2	1:D:83:GLY:H	2.00	0.58
1:B:176:GLN:HG3	3:B:2046:HOH:O	2.03	0.58
1:B:193:ILE:HG23	1:B:198:HIS:HB3	1.87	0.57
1:A:70:GLU:HG2	1:A:100:LEU:HD21	1.85	0.57
1:D:118:ARG:HG3	1:D:119:VAL:N	2.17	0.56
1:C:424:LEU:HD13	1:C:435:SER:HB2	1.87	0.56
1:B:316:GLN:HE22	1:C:69:ARG:HH11	1.52	0.56
1:B:131:ILE:CD1	1:B:296:CYS:HB2	2.36	0.56
1:D:34:ILE:H	1:D:46:ASN:HD21	1.54	0.56
2:A:1442:COA:H52A	2:A:1442:COA:N3A	2.22	0.54
2:D:1442:COA:H133	2:D:1442:COA:O9P	2.06	0.54
1:C:123:ALA:HA	1:C:390:VAL:CG1	2.38	0.54
1:A:241:VAL:HG12	1:A:242:LYS:N	2.23	0.54
1:C:123:ALA:HA	1:C:390:VAL:HG13	1.87	0.54
1:B:315:ARG:HE	1:B:319:GLN:NE2	2.06	0.54
1:A:61:MET:HE2	1:A:70:GLU:CB	2.39	0.53
1:A:61:MET:HE3	1:A:66:LEU:CB	2.38	0.53
1:C:222:HIS:CD2	1:C:223:MET:CE	2.91	0.53
1:C:420:LEU:HD12	3:C:2073:HOH:O	2.07	0.53
1:A:70:GLU:CG	1:A:100:LEU:HD21	2.37	0.53
1:D:63:HIS:O	1:D:287:ARG:NH2	2.42	0.52
1:A:372:GLU:HA	1:A:372:GLU:OE1	2.09	0.52
1:D:306:ASP:HB3	3:D:2046:HOH:O	2.10	0.52
1:B:420:LEU:HD23	1:B:439:GLN:HG2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:128:LEU:HD22	1:D:117:MET:HB3	1.92	0.51
2:B:1442:COA:O9P	2:B:1442:COA:H141	2.09	0.51
1:C:131:ILE:CD1	1:C:296:CYS:HB2	2.41	0.51
1:B:410:GLU:HB2	3:B:2071:HOH:O	2.10	0.51
1:B:316:GLN:NE2	1:C:69:ARG:HH11	2.09	0.51
1:A:118:ARG:HG3	1:A:119:VAL:N	2.23	0.51
2:A:1442:COA:S1P	3:A:2026:HOH:O	2.59	0.51
1:A:63:HIS:O	1:A:287:ARG:NH2	2.43	0.51
1:B:82:LEU:HD22	1:B:85:LEU:HD12	1.93	0.51
1:B:372:GLU:HG2	3:B:2109:HOH:O	2.11	0.50
1:D:76:VAL:O	1:D:116:ALA:HA	2.12	0.50
1:A:61:MET:HE3	1:A:66:LEU:HB3	1.94	0.50
1:B:198:HIS:HE1	1:B:350:GLU:OE1	1.94	0.50
1:C:324:MET:HG3	3:C:2071:HOH:O	2.12	0.49
1:C:76:VAL:O	1:C:116:ALA:HA	2.13	0.49
1:B:60:THR:HG22	1:B:269:VAL:HG23	1.94	0.49
1:C:222:HIS:CD2	1:C:223:MET:HE2	2.48	0.49
1:C:316:GLN:HG2	1:D:113:TYR:HB3	1.95	0.49
1:A:61:MET:CE	1:A:66:LEU:HB3	2.43	0.48
1:A:34:ILE:H	1:A:46:ASN:HD21	1.61	0.48
1:A:119:VAL:HG12	1:B:119:VAL:HG12	1.96	0.48
1:C:222:HIS:CD2	1:C:223:MET:HE3	2.48	0.48
1:C:60:THR:HG22	1:C:269:VAL:HG23	1.95	0.48
1:C:272:ASN:HB2	3:C:2001:HOH:O	2.14	0.48
1:D:36:LYS:HG3	1:D:37:LYS:HG3	1.95	0.47
1:B:76:VAL:O	1:B:116:ALA:HA	2.14	0.47
1:C:80:ASN:HD21	1:C:83:GLY:N	2.07	0.47
1:C:81:PHE:CD2	1:C:82:LEU:HG	2.50	0.47
1:B:20:GLY:HA2	1:B:60:THR:HG23	1.97	0.47
1:A:76:VAL:O	1:A:116:ALA:HA	2.15	0.47
1:C:274:GLU:OE1	1:C:277:ARG:NH1	2.45	0.47
1:B:163:ASP:O	1:B:166:ALA:HB3	2.15	0.47
1:D:241:VAL:HG21	2:D:1442:COA:C8A	2.45	0.47
1:D:82:LEU:HD22	1:D:85:LEU:HD12	1.97	0.47
1:C:408:GLN:NE2	3:C:2081:HOH:O	2.47	0.46
1:B:80:ASN:HD21	1:B:83:GLY:N	2.10	0.46
1:D:61:MET:HE3	1:D:146:LEU:HD13	1.98	0.46
1:B:124:ALA:HA	1:B:425:ASN:ND2	2.31	0.45
1:D:424:LEU:HD13	1:D:435:SER:HB2	1.98	0.45
2:D:1442:COA:S1P	3:D:2020:HOH:O	2.54	0.45
1:A:16:PHE:CD2	1:A:276:LEU:HD11	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:HIS:CD2	1:A:40:ASP:H	2.35	0.44
1:B:154:THR:HG23	1:B:263:ASP:OD1	2.17	0.44
1:A:168:ALA:HB3	1:A:182:PHE:CE2	2.52	0.44
1:A:80:ASN:HD21	1:A:83:GLY:N	2.10	0.44
1:A:198:HIS:HE1	1:A:350:GLU:OE1	2.00	0.44
1:C:124:ALA:HA	1:C:425:ASN:ND2	2.32	0.44
1:A:124:ALA:HA	1:A:425:ASN:ND2	2.32	0.44
1:B:38:HIS:HD2	1:B:40:ASP:H	1.66	0.44
1:C:20:GLY:HA3	1:C:399:MET:SD	2.57	0.44
2:B:1442:COA:O9P	2:B:1442:COA:CEP	2.65	0.43
1:D:383:LEU:HD11	1:D:393:THR:HG21	1.99	0.43
1:D:131:ILE:CD1	1:D:296:CYS:HB2	2.48	0.43
1:D:198:HIS:HE1	1:D:350:GLU:OE1	2.01	0.43
1:A:191:LYS:NZ	1:A:195:GLU:OE2	2.40	0.43
1:C:85:LEU:HD23	1:D:85:LEU:HD21	2.00	0.43
1:B:38:HIS:CD2	1:B:40:ASP:H	2.36	0.43
1:C:174:GLN:NE2	1:C:185:LEU:HD21	2.33	0.43
1:C:176:GLN:HE22	1:C:177:LEU:HD22	1.79	0.43
1:A:363:LEU:HD22	1:A:368:ASP:HB2	2.00	0.43
1:C:85:LEU:HD21	1:D:85:LEU:HD23	1.99	0.43
1:C:113:TYR:CG	1:D:316:GLN:HG3	2.54	0.43
1:A:131:ILE:CD1	1:A:296:CYS:HB2	2.49	0.43
1:D:188:LYS:HA	1:D:249:TYR:CD1	2.54	0.43
1:D:420:LEU:HD21	1:D:439:GLN:HE21	1.84	0.43
1:A:371:LEU:HD23	3:A:2072:HOH:O	2.19	0.43
1:A:61:MET:HE3	1:A:66:LEU:HB2	2.00	0.43
1:C:198:HIS:HE1	1:C:350:GLU:OE1	2.01	0.43
1:B:80:ASN:ND2	1:B:83:GLY:H	2.12	0.42
1:A:81:PHE:CD2	1:A:82:LEU:HG	2.54	0.42
1:B:169:ALA:HB1	1:B:174:GLN:OE1	2.20	0.42
1:A:60:THR:HG22	1:A:269:VAL:HG23	2.01	0.42
1:B:420:LEU:CD2	1:B:439:GLN:HG2	2.49	0.42
1:C:63:HIS:O	1:C:63:HIS:CD2	2.72	0.42
1:D:131:ILE:HD13	1:D:131:ILE:HG21	1.81	0.42
1:D:168:ALA:HB3	1:D:182:PHE:CE2	2.55	0.42
1:D:20:GLY:HA3	1:D:399:MET:SD	2.60	0.41
1:C:131:ILE:HD11	1:C:434:VAL:HG13	2.02	0.41
1:D:63:HIS:O	1:D:63:HIS:CD2	2.73	0.41
1:B:424:LEU:CD1	1:B:435:SER:HB2	2.50	0.41
1:B:316:GLN:NE2	1:C:69:ARG:NH1	2.67	0.41
1:C:217:LYS:O	1:C:411:ALA:HB3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:38:HIS:CD2	1:D:40:ASP:H	2.38	0.41
1:C:70:GLU:HG2	1:C:100:LEU:HD21	2.01	0.41
1:C:15:VAL:HG13	1:C:137:LEU:HD13	2.01	0.41
1:B:334:VAL:HG22	1:B:335:ALA:N	2.36	0.41
1:D:81:PHE:CD2	1:D:82:LEU:HG	2.56	0.41
1:B:188:LYS:HA	1:B:249:TYR:CD1	2.56	0.41
1:C:409:CYS:HB2	1:C:413:GLN:HA	2.02	0.41
1:D:383:LEU:HD11	1:D:393:THR:CG2	2.51	0.41
1:D:127:GLY:HA3	1:D:425:ASN:HD22	1.86	0.40
1:A:20:GLY:HA3	1:A:399:MET:SD	2.62	0.40
1:A:84[A]:GLU:OE2	1:A:118:ARG:NH2	2.46	0.40

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:A:227:LYS:NZ	3:B:2114:HOH:O[1_655]	2.19	0.01

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	429/457 (94%)	416 (97%)	12 (3%)	1 (0%)	47	57
1	B	430/457 (94%)	418 (97%)	12 (3%)	0	100	100
1	C	429/457 (94%)	417 (97%)	11 (3%)	1 (0%)	47	57
1	D	430/457 (94%)	418 (97%)	12 (3%)	0	100	100
All	All	1718/1828 (94%)	1669 (97%)	47 (3%)	2 (0%)	51	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	13	SER
1	A	15	VAL

### 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	339/362 (94%)	333 (98%)	6 (2%)	59	71
1	B	340/362 (94%)	334 (98%)	6 (2%)	59	71
1	C	339/362 (94%)	334 (98%)	5 (2%)	65	76
1	D	340/362 (94%)	335 (98%)	5 (2%)	65	76
All	All	1358/1448 (94%)	1336 (98%)	22 (2%)	62	74

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	118	ARG
1	A	141	SER
1	A	284	ASN
1	A	366	ASN
1	A	371	LEU
1	A	386	PHE
1	B	52	ILE
1	B	118	ARG
1	B	273	GLU
1	B	284	ASN
1	B	386	PHE
1	B	439	GLN
1	C	118	ARG
1	C	141	SER
1	C	176	GLN
1	C	273	GLU
1	C	386	PHE
1	D	118	ARG
1	D	141	SER
1	D	225	THR

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Mol	Chain	Res	Type
1	D	284	ASN
1	D	386	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	HIS
1	A	46	ASN
1	A	58	GLN
1	A	75	GLN
1	A	80	ASN
1	A	172	GLN
1	A	198	HIS
1	A	359	HIS
1	A	425	ASN
1	B	38	HIS
1	B	46	ASN
1	B	63	HIS
1	B	80	ASN
1	B	198	HIS
1	B	316	GLN
1	B	319	GLN
1	B	359	HIS
1	B	425	ASN
1	C	38	HIS
1	C	63	HIS
1	C	80	ASN
1	C	172	GLN
1	C	176	GLN
1	C	198	HIS
1	C	408	GLN
1	C	425	ASN
1	D	38	HIS
1	D	46	ASN
1	D	63	HIS
1	D	80	ASN
1	D	172	GLN
1	D	198	HIS
1	D	425	ASN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

3 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
2	COA	A	1442	-	41,50,50	1.01	3 (7%)	52,75,75	1.34	6 (11%)
2	COA	D	1442	-	41,50,50	1.16	4 (9%)	52,75,75	1.74	12 (23%)
2	COA	B	1442	-	41,50,50	1.18	6 (14%)	52,75,75	2.05	13 (25%)

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
2	COA	A	1442	-	-	23/44/64/64	0/3/3/3
2	COA	D	1442	-	-	20/44/64/64	0/3/3/3
2	COA	B	1442	-	-	9/44/64/64	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1442	COA	O4B-C1B	2.94	1.45	1.41
2	B	1442	COA	C5A-C4A	2.77	1.48	1.40
2	A	1442	COA	C5A-C4A	2.63	1.47	1.40
2	B	1442	COA	C3P-N4P	2.56	1.52	1.46
2	B	1442	COA	C5P-N4P	2.53	1.39	1.33
2	A	1442	COA	O4B-C1B	2.53	1.44	1.41
2	D	1442	COA	C5P-N4P	2.53	1.39	1.33
2	D	1442	COA	C5A-C4A	2.48	1.47	1.40
2	B	1442	COA	C2P-S1P	2.23	1.88	1.80
2	D	1442	COA	C2A-N3A	2.15	1.35	1.32
2	A	1442	COA	P3B-O3B	2.15	1.63	1.59
2	B	1442	COA	C4A-N3A	-2.04	1.32	1.35
2	B	1442	COA	C2A-N1A	2.01	1.37	1.33

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1442	COA	O5P-C5P-C6P	-6.03	110.99	122.02
2	B	1442	COA	C6P-C5P-N4P	5.05	124.92	116.42
2	A	1442	COA	N3A-C2A-N1A	-4.92	121.00	128.68
2	D	1442	COA	C2P-C3P-N4P	4.89	123.47	112.31
2	B	1442	COA	N3A-C2A-N1A	-4.52	121.61	128.68
2	D	1442	COA	O5P-C5P-C6P	-4.47	113.85	122.02
2	D	1442	COA	N3A-C2A-N1A	-4.32	121.93	128.68
2	B	1442	COA	O6A-CCP-CBP	4.24	117.37	110.55
2	D	1442	COA	C6P-C5P-N4P	3.76	122.76	116.42
2	D	1442	COA	C6P-C7P-N8P	3.41	118.79	111.90
2	B	1442	COA	CAP-C9P-N8P	3.33	123.21	116.58
2	B	1442	COA	C2A-N1A-C6A	3.27	124.35	118.75
2	B	1442	COA	C3B-C2B-C1B	3.11	106.78	99.89
2	A	1442	COA	P2A-O3A-P1A	-2.99	122.56	132.83
2	B	1442	COA	C6P-C7P-N8P	2.84	117.64	111.90
2	B	1442	COA	C2P-C3P-N4P	2.79	118.67	112.31
2	A	1442	COA	C3B-C2B-C1B	2.73	105.94	99.89
2	D	1442	COA	O4B-C4B-C5B	2.71	118.29	109.37
2	D	1442	COA	O4B-C4B-C3B	-2.71	99.07	104.87
2	A	1442	COA	C2A-N1A-C6A	2.68	123.33	118.75
2	B	1442	COA	P2A-O3A-P1A	-2.55	124.08	132.83
2	B	1442	COA	N6A-C6A-N1A	2.53	123.83	118.57
2	B	1442	COA	C2B-C3B-C4B	-2.46	98.86	103.22
2	D	1442	COA	CDP-CBP-CAP	2.44	113.05	108.82
2	B	1442	COA	C4A-C5A-N7A	2.35	111.85	109.40
2	A	1442	COA	O5P-C5P-C6P	-2.33	117.75	122.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1442	COA	P2A-O3A-P1A	-2.30	124.93	132.83
2	D	1442	COA	C2A-N1A-C6A	2.24	122.58	118.75
2	A	1442	COA	C6P-C5P-N4P	2.22	120.16	116.42
2	D	1442	COA	O9A-P3B-O8A	2.12	115.74	107.64
2	D	1442	COA	C3P-N4P-C5P	2.04	126.63	122.84

There are no chirality outliers.

All (52) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1442	COA	C3B-O3B-P3B-O7A
2	A	1442	COA	C5B-O5B-P1A-O1A
2	A	1442	COA	OAP-CAP-CBP-CCP
2	A	1442	COA	C9P-CAP-CBP-CCP
2	A	1442	COA	OAP-CAP-CBP-CDP
2	A	1442	COA	C9P-CAP-CBP-CDP
2	A	1442	COA	O9P-C9P-CAP-CBP
2	A	1442	COA	N8P-C9P-CAP-CBP
2	A	1442	COA	N8P-C9P-CAP-OAP
2	D	1442	COA	O4B-C4B-C5B-O5B
2	D	1442	COA	C5B-O5B-P1A-O1A
2	D	1442	COA	C5B-O5B-P1A-O2A
2	D	1442	COA	CCP-O6A-P2A-O3A
2	D	1442	COA	OAP-CAP-CBP-CCP
2	D	1442	COA	C9P-CAP-CBP-CCP
2	D	1442	COA	OAP-CAP-CBP-CDP
2	D	1442	COA	C9P-CAP-CBP-CDP
2	D	1442	COA	OAP-CAP-CBP-CEP
2	D	1442	COA	C9P-CAP-CBP-CEP
2	D	1442	COA	O9P-C9P-CAP-OAP
2	D	1442	COA	N8P-C9P-CAP-OAP
2	B	1442	COA	CCP-O6A-P2A-O3A
2	B	1442	COA	CCP-O6A-P2A-O4A
2	B	1442	COA	CCP-O6A-P2A-O5A
2	B	1442	COA	CDP-CBP-CCP-O6A
2	B	1442	COA	CEP-CBP-CCP-O6A
2	B	1442	COA	CAP-CBP-CCP-O6A
2	B	1442	COA	S1P-C2P-C3P-N4P
2	D	1442	COA	C3B-C4B-C5B-O5B
2	A	1442	COA	C2B-C3B-O3B-P3B
2	B	1442	COA	C2B-C3B-O3B-P3B
2	A	1442	COA	C4B-C3B-O3B-P3B

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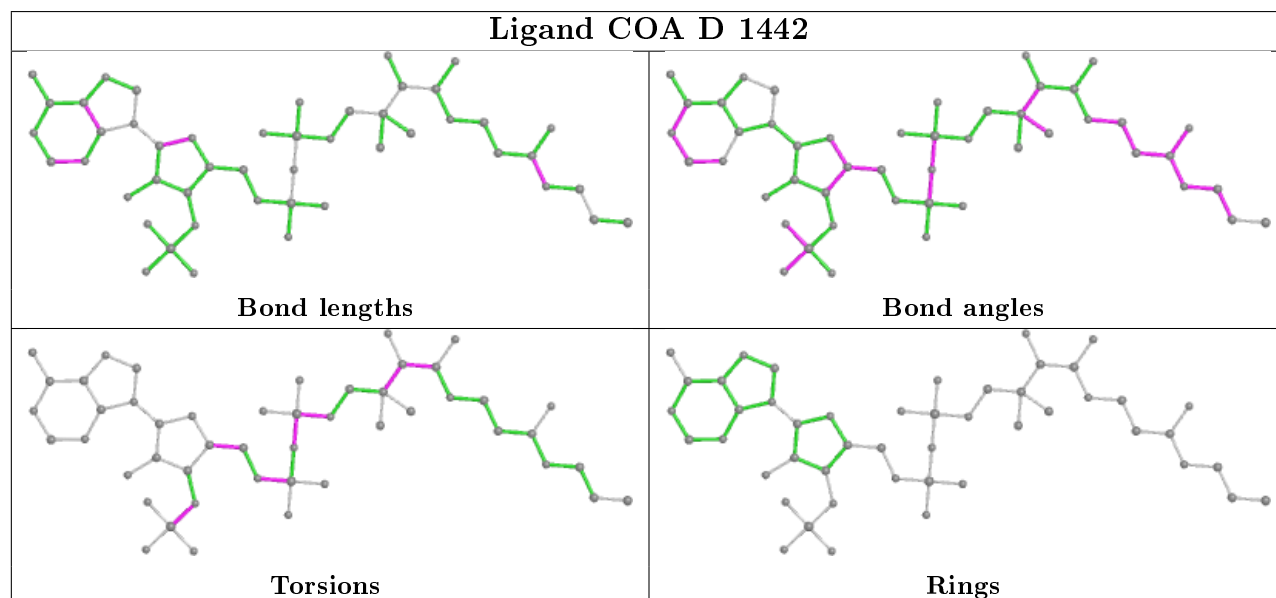
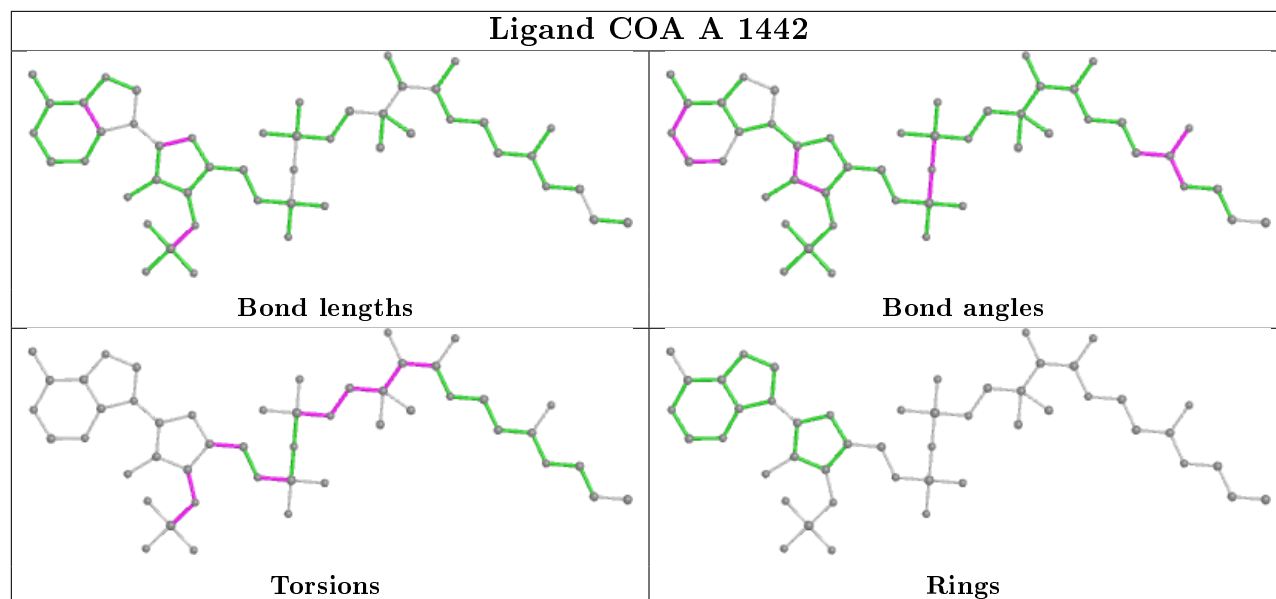
Mol	Chain	Res	Type	Atoms
2	B	1442	COA	C4B-C3B-O3B-P3B
2	A	1442	COA	O9P-C9P-CAP-OAP
2	A	1442	COA	OAP-CAP-CBP-CEP
2	D	1442	COA	O9P-C9P-CAP-CBP
2	D	1442	COA	N8P-C9P-CAP-CBP
2	A	1442	COA	C3B-O3B-P3B-O8A
2	A	1442	COA	C5B-O5B-P1A-O3A
2	D	1442	COA	P1A-O3A-P2A-O5A
2	A	1442	COA	C5B-O5B-P1A-O2A
2	D	1442	COA	CCP-O6A-P2A-O4A
2	D	1442	COA	CCP-O6A-P2A-O5A
2	A	1442	COA	CAP-CBP-CCP-O6A
2	A	1442	COA	CEP-CBP-CCP-O6A
2	A	1442	COA	CDP-CBP-CCP-O6A
2	A	1442	COA	C9P-CAP-CBP-CEP
2	D	1442	COA	C3B-O3B-P3B-O9A
2	D	1442	COA	C5B-O5B-P1A-O3A
2	A	1442	COA	O4B-C4B-C5B-O5B
2	A	1442	COA	CBP-CCP-O6A-P2A
2	A	1442	COA	CCP-O6A-P2A-O4A

There are no ring outliers.

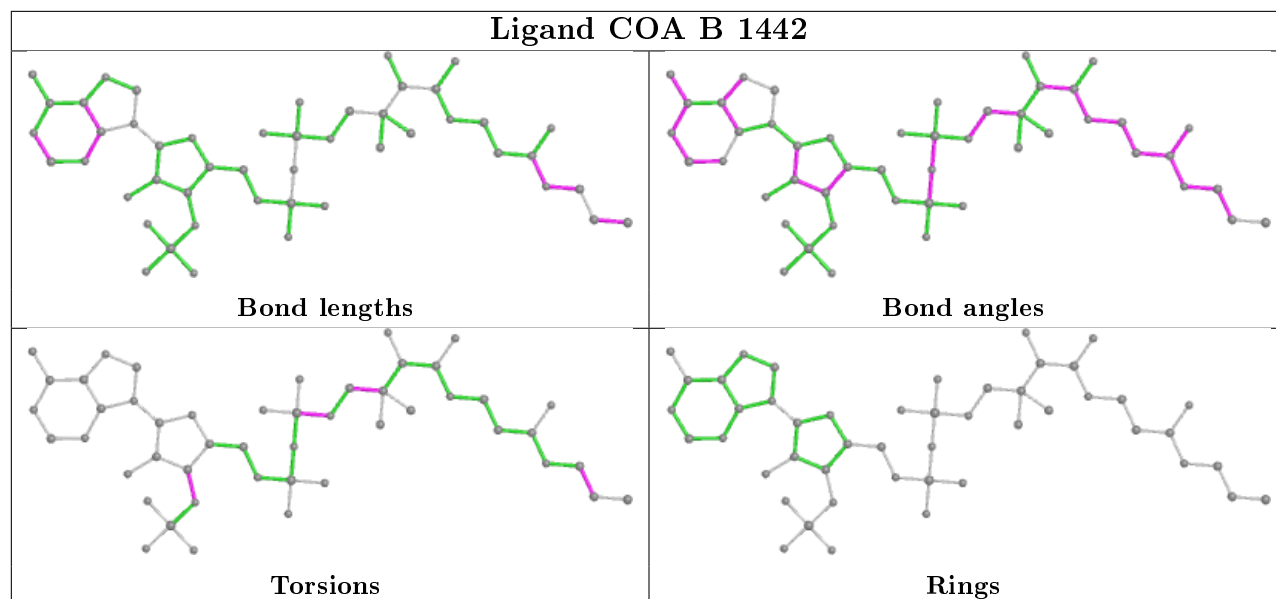
3 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1442	COA	4	0
2	D	1442	COA	4	0
2	B	1442	COA	3	0

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







## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	430/457 (94%)	0.03	5 (1%) 79 77	7, 19, 43, 73	0
1	B	430/457 (94%)	-0.23	2 (0%) 91 92	5, 12, 29, 55	0
1	C	430/457 (94%)	-0.08	0 100 100	5, 17, 44, 73	0
1	D	430/457 (94%)	-0.11	1 (0%) 95 95	5, 16, 37, 66	0
All	All	1720/1828 (94%)	-0.10	8 (0%) 91 92	5, 16, 40, 73	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	12	ALA	4.2
1	D	12	ALA	3.2
1	A	42	GLY	3.2
1	A	37	LYS	2.6
1	A	105	ALA	2.3
1	A	238	PRO	2.1
1	A	241	VAL	2.0
1	B	238	PRO	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

### 6.3 Carbohydrates [i](#)

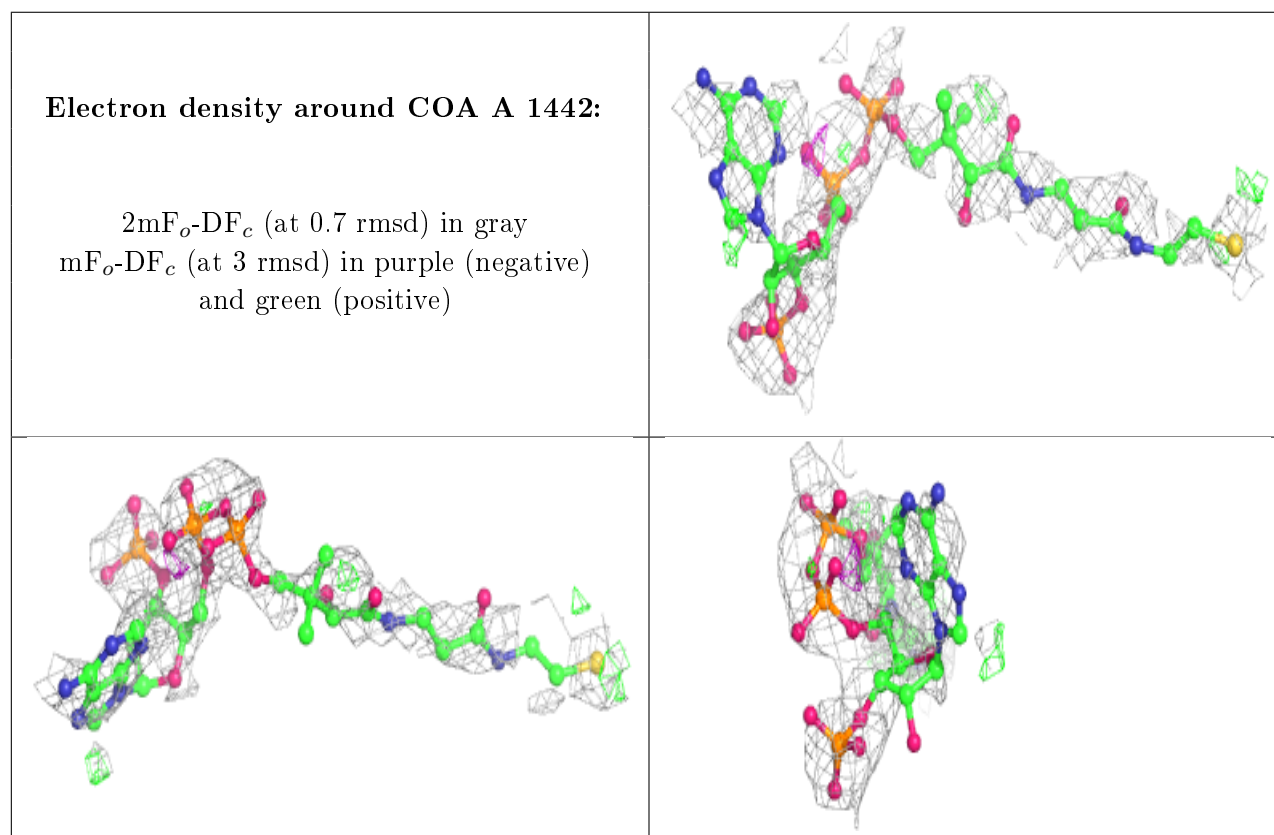
There are no carbohydrates in this entry.

## 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, 95<sup>th</sup> 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.

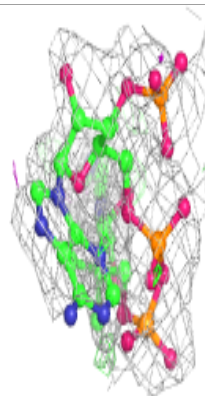
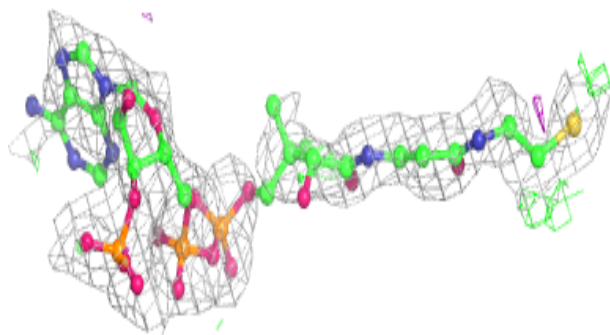
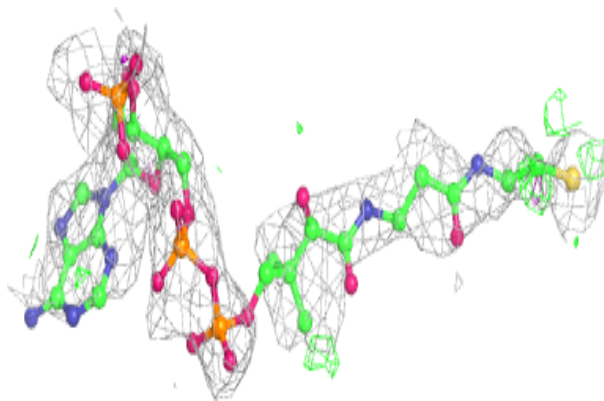
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	COA	A	1442	48/48	0.74	0.36	63,79,85,86	0
2	COA	D	1442	48/48	0.79	0.23	34,59,74,81	0
2	COA	B	1442	48/48	0.94	0.15	11,15,20,24	0

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.

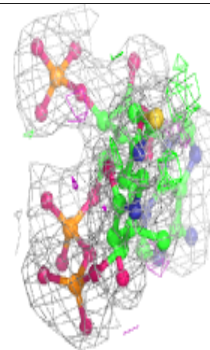
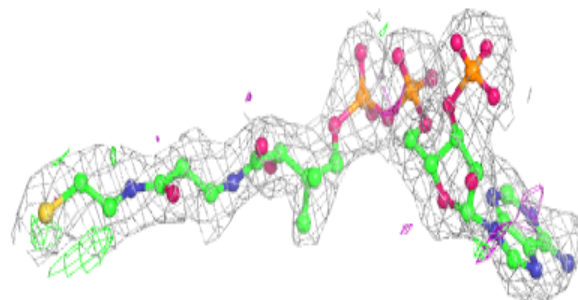
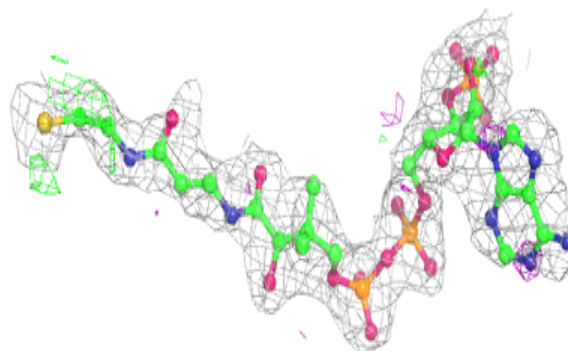


**Electron density around COA D 1442:**

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

**Electron density around COA B 1442:**

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



## 6.5 Other polymers

There are no such residues in this entry.