Full wwPDB X-ray Structure Validation Report (i)

Jun 23, 2024 – 04:37 AM EDT

PDB ID	:	6LQ0
Title	:	Crystal Structure of E447A Acyl-CoA Dehydrogenase FadE5 mutant from Mycobacteria smegmatis in complex with C6CoA
Authors	:	Liu, X.; Chen, X.B.
Deposited on	:	2020-01-12
Resolution	:	2.20 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:

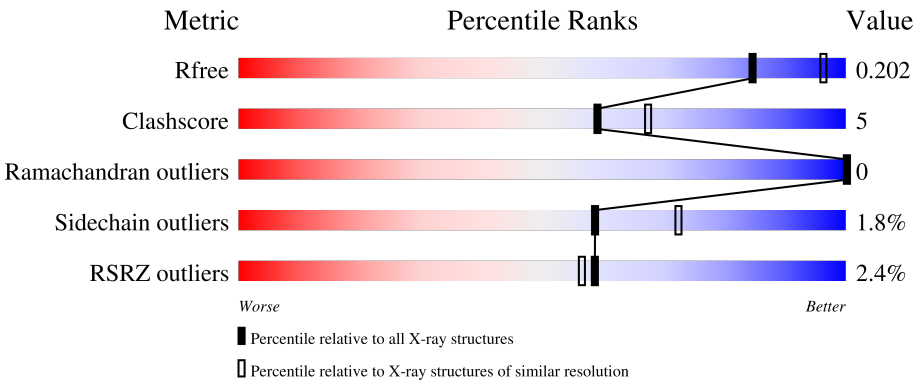
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	611	3% 92% 7% •
1	B	611	2% 92% 7% •

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 19519 atoms, of which 9276 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.

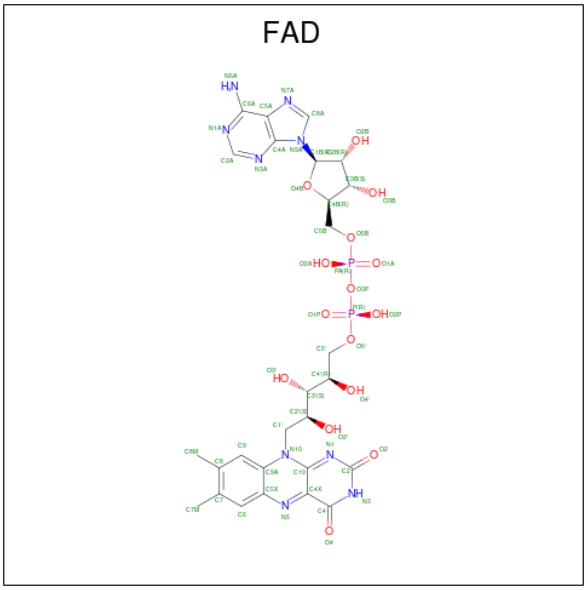
- Molecule 1 is a protein called Acyl-CoA dehydrogenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	607	9227	2948	4581	789	887	22	0	0	0
1	B	607	9217	2946	4574	789	887	21	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	447	ALA	GLU	engineered mutation	UNP A0A0D6G5A8
B	447	ALA	GLU	engineered mutation	UNP A0A0D6G5A8

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O	P		
2	A	1	84	27	31	9	15	2	0	0

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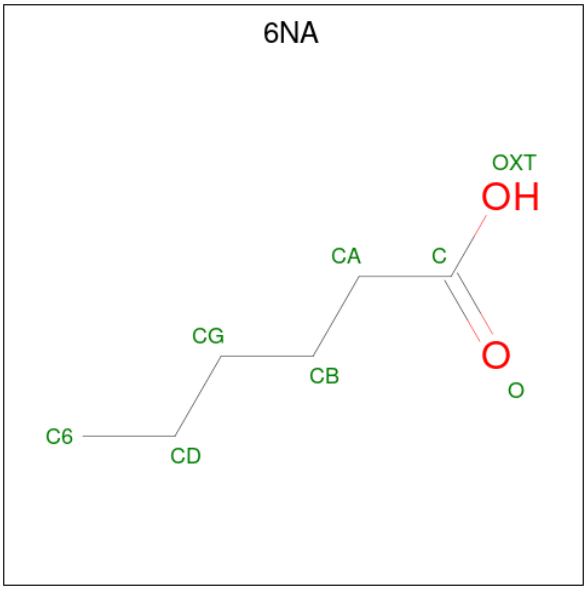
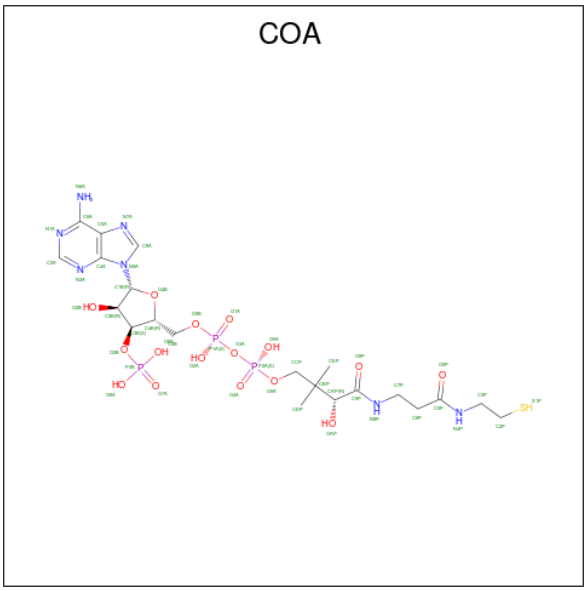
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O	P		
2	B	1	83	27	30	9	15	2	0	0

- Molecule 3 is HEXANOIC ACID (three-letter code: 6NA) (formula: $C_6H_{12}O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
3	A	1	17	6	10	1	0	0
3	B	1	17	6	10	1	0	0

- Molecule 4 is COENZYME A (three-letter code: COA) (formula: $C_{21}H_{36}N_7O_{16}P_3S$).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
			Total	C	H	N	O	P	S		
4	A	1	68	21	20	7	16	3	1	0	0
			Total	C	H	N	O	P	S		
4	B	1	68	21	20	7	16	3	1	0	0

- Molecule 5 is water.

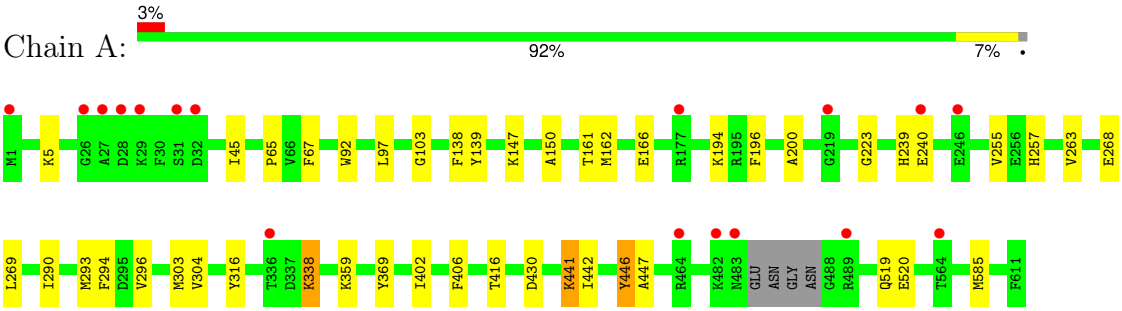
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	363	363	363	0	0
			Total	O		
5	B	375	375	375	0	0

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.

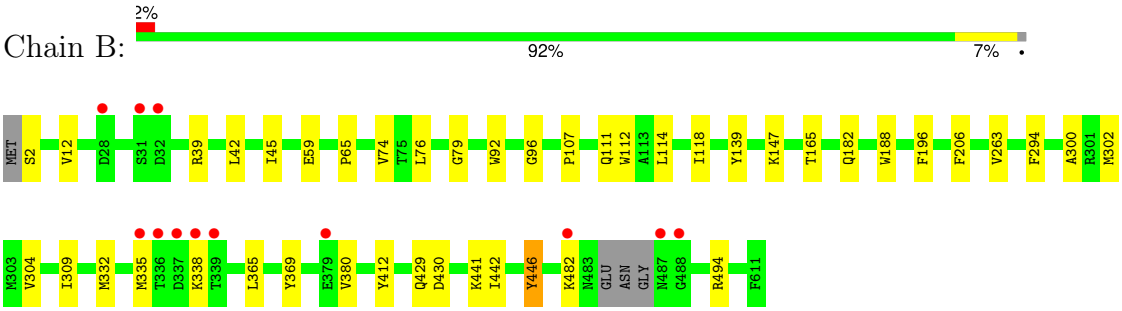
- Molecule 1: Acyl-CoA dehydrogenase

Chain A:



- Molecule 1: Acyl-CoA dehydrogenase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	98.09Å 206.02Å 74.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.63 – 2.20 19.63 – 2.20	Depositor EDS
% Data completeness (in resolution range)	92.5 (19.63-2.20) 92.5 (19.63-2.20)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.54 (at 2.19Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.152 , 0.201 0.154 , 0.202	Depositor DCC
R_{free} test set	3557 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	27.9	Xtriage
Anisotropy	0.347	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.47 , 54.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	19519	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.63% 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: COA, 6NA, FAD

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	0.54	0/4738	0.64	0/6411
1	B	0.58	0/4735	0.68	0/6408
All	All	0.56	0/9473	0.66	0/12819

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4646	4581	4579	41	0
1	B	4643	4574	4569	50	0
2	A	53	31	31	8	0
2	B	53	30	31	14	0
3	A	7	10	11	1	0
3	B	7	10	11	0	0
4	A	48	20	31	2	0
4	B	48	20	30	3	0
5	A	363	0	0	5	0
5	B	375	0	0	4	0
All	All	10243	9276	9293	89	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (89) 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)
4:B:703:COA:O4B	4:B:703:COA:C1B	1.63	1.19
1:A:442:ILE:CD1	2:A:701:FAD:HM83	1.79	1.11
1:A:442:ILE:HD11	2:A:701:FAD:HM83	1.13	1.09
1:B:442:ILE:HD11	2:B:701:FAD:HM83	1.52	0.92
1:A:442:ILE:HD11	2:A:701:FAD:C8M	2.04	0.78
1:B:442:ILE:CD1	2:B:701:FAD:HM83	2.15	0.77
1:A:304:VAL:HG22	1:A:446:TYR:HE1	1.53	0.73
1:B:442:ILE:CD1	2:B:701:FAD:C7M	2.68	0.71
1:B:2:SER:N	5:B:801:HOH:O	2.24	0.70
1:B:494:ARG:NH2	5:B:802:HOH:O	2.25	0.70
1:A:442:ILE:HG13	1:A:446:TYR:CE2	2.26	0.70
1:B:442:ILE:CD1	2:B:701:FAD:HM71	2.23	0.68
1:B:304:VAL:HG22	1:B:446:TYR:HE1	1.57	0.68
1:A:442:ILE:HG13	1:A:446:TYR:CD2	2.28	0.67
1:B:442:ILE:HD11	2:B:701:FAD:C8M	2.26	0.65
1:A:416:THR:HG23	1:B:441:LYS:HE2	1.77	0.65
1:B:45:ILE:HD12	1:B:92:TRP:CE3	2.34	0.62
1:A:519:GLN:OE1	5:A:801:HOH:O	2.16	0.60
1:A:442:ILE:HD13	2:A:701:FAD:HM83	1.83	0.58
1:B:442:ILE:HD12	2:B:701:FAD:HM71	1.86	0.58
1:A:139:TYR:CZ	1:A:147:LYS:HD3	2.40	0.57
1:B:107:PRO:HB3	1:B:380:VAL:HG21	1.88	0.56
1:B:304:VAL:HG22	1:B:446:TYR:CE1	2.41	0.55
1:B:65:PRO:HG3	1:B:76:LEU:HD23	1.88	0.54
1:A:223:GLY:CA	1:B:338:LYS:HD2	2.38	0.54
2:A:701:FAD:H9	2:A:701:FAD:H2'	1.90	0.54
1:B:338:LYS:O	1:B:338:LYS:HG3	2.08	0.54
1:B:442:ILE:CD1	1:B:446:TYR:CD2	2.92	0.53
1:A:196:PHE:O	2:A:701:FAD:C4X	2.57	0.53
1:B:442:ILE:HD11	2:B:701:FAD:C7M	2.37	0.53
1:A:223:GLY:HA2	1:B:338:LYS:HD2	1.92	0.52
1:B:442:ILE:O	1:B:442:ILE:HG12	2.09	0.52
2:B:701:FAD:C2	4:B:703:COA:S1P	2.97	0.52
1:B:196:PHE:O	2:B:701:FAD:C4X	2.58	0.52
1:A:442:ILE:CG1	1:A:446:TYR:CE2	2.94	0.51
1:A:223:GLY:HA3	1:B:338:LYS:CD	2.41	0.50
1:A:268:GLU:HG2	5:A:826:HOH:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:412:TYR:CE1	1:B:441:LYS:HD2	2.47	0.50
1:A:166:GLU:HG2	1:A:194:LYS:HD3	1.94	0.50
1:A:257:HIS:HB3	1:B:429:GLN:HE21	1.77	0.49
1:A:239:HIS:HD2	5:A:1096:HOH:O	1.95	0.49
1:B:309:ILE:HG12	1:B:365:LEU:HB3	1.94	0.49
1:B:112:TRP:CE2	1:B:302:MET:HG2	2.47	0.49
2:A:701:FAD:H2'	2:A:701:FAD:C9	2.43	0.48
1:A:442:ILE:CG1	1:A:446:TYR:CD2	2.96	0.48
1:A:161:THR:HG21	1:A:200:ALA:HA	1.95	0.47
1:B:65:PRO:HD3	1:B:263:VAL:HA	1.96	0.47
1:B:196:PHE:O	2:B:701:FAD:N5	2.47	0.47
1:A:304:VAL:CG2	1:A:446:TYR:HE1	2.22	0.47
1:B:304:VAL:CG2	1:B:446:TYR:HE1	2.27	0.47
1:B:139:TYR:CZ	1:B:147:LYS:HD3	2.49	0.47
1:A:138:PHE:CD2	1:A:150:ALA:HB2	2.50	0.47
1:A:223:GLY:HA3	1:B:338:LYS:HG2	1.97	0.47
1:B:45:ILE:HD11	1:B:118:ILE:HD11	1.97	0.46
1:B:65:PRO:HB3	1:B:74:VAL:HG13	1.95	0.46
1:B:332:MET:HA	1:B:335:MET:HG3	1.97	0.46
1:A:304:VAL:HG22	1:A:446:TYR:CE1	2.42	0.46
1:A:441:LYS:CE	5:A:875:HOH:O	2.63	0.45
1:B:59:GLU:OE1	1:B:79:GLY:HA3	2.16	0.45
1:B:442:ILE:HD13	1:B:446:TYR:CE2	2.51	0.45
1:B:442:ILE:HD13	1:B:446:TYR:CD2	2.51	0.45
1:A:200:ALA:HB1	1:A:269:LEU:HD13	2.00	0.44
1:A:303:MET:CE	3:A:702:6NA:H6C3	2.46	0.44
1:A:67:PHE:HB2	1:A:255:VAL:HG23	1.98	0.44
1:A:441:LYS:HE3	5:A:875:HOH:O	2.17	0.44
1:A:402:ILE:O	1:A:406:PHE:HB3	2.18	0.43
1:A:45:ILE:HD12	1:A:92:TRP:CE3	2.54	0.43
1:B:300:ALA:O	1:B:304:VAL:HG23	2.18	0.43
1:B:442:ILE:CD1	2:B:701:FAD:HM73	2.46	0.43
2:A:701:FAD:C2	4:A:703:COA:S1P	3.07	0.43
1:A:447:ALA:HA	4:A:703:COA:H32	2.01	0.43
1:B:42:LEU:HD23	1:B:114:LEU:HD23	2.01	0.43
1:A:338:LYS:CG	5:B:806:HOH:O	2.67	0.43
1:A:65:PRO:HD3	1:A:263:VAL:HA	2.01	0.42
1:A:97:LEU:O	1:A:103:GLY:HA3	2.20	0.42
1:B:65:PRO:HG3	1:B:76:LEU:CD2	2.49	0.42
1:B:12:VAL:HG21	1:B:39:ARG:HH11	1.85	0.42
1:B:442:ILE:HD11	2:B:701:FAD:C8	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:GLY:HA3	1:B:338:LYS:CG	2.51	0.41
1:A:316:TYR:CZ	1:A:359:LYS:HB2	2.55	0.41
1:A:162:MET:HE3	1:A:296:VAL:HG12	2.01	0.41
1:A:223:GLY:CA	1:B:338:LYS:CD	2.99	0.41
1:B:96:GLY:HA3	1:B:111:GLN:OE1	2.20	0.41
4:B:703:COA:O4B	4:B:703:COA:N9A	2.40	0.41
1:A:290:ILE:HD12	1:A:290:ILE:HA	1.98	0.40
1:B:165:THR:OG1	2:B:701:FAD:H1'1	2.21	0.40
1:B:441:LYS:HE3	5:B:883:HOH:O	2.21	0.40
1:B:442:ILE:HD11	2:B:701:FAD:C7	2.51	0.40
1:B:182:GLN:HB2	1:B:188:TRP:CZ3	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	603/611 (99%)	592 (98%)	11 (2%)	0	100	100
1	B	603/611 (99%)	593 (98%)	10 (2%)	0	100	100
All	All	1206/1222 (99%)	1185 (98%)	21 (2%)	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	474/478 (99%)	463 (98%)	11 (2%)	50	63
1	B	473/478 (99%)	467 (99%)	6 (1%)	69	81
All	All	947/956 (99%)	930 (98%)	17 (2%)	59	72

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	240	GLU
1	A	293	MET
1	A	294	PHE
1	A	338	LYS
1	A	369	TYR
1	A	430	ASP
1	A	441	LYS
1	A	446	TYR
1	A	520	GLU
1	A	585	MET
1	B	206	PHE
1	B	294	PHE
1	B	369	TYR
1	B	430	ASP
1	B	446	TYR
1	B	482	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	455	GLN
1	B	429	GLN
1	B	519	GLN

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)

6 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	6NA	A	702	4	5,6,7	0.87	0	4,5,7	0.85	0
2	FAD	A	701	-	54,58,58	0.62	0	71,89,89	0.58	1 (1%)
4	COA	B	703	3	43,50,50	3.78	12 (27%)	56,75,75	2.01	13 (23%)
2	FAD	B	701	-	54,58,58	0.87	1 (1%)	71,89,89	0.61	1 (1%)
4	COA	A	703	3	43,50,50	3.69	14 (32%)	56,75,75	1.75	9 (16%)
3	6NA	B	702	4	5,6,7	0.94	0	4,5,7	0.71	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	6NA	A	702	4	-	1/4/4/5	-
2	FAD	A	701	-	-	2/30/50/50	0/6/6/6
4	COA	B	703	3	-	8/44/64/64	0/3/3/3
2	FAD	B	701	-	-	2/30/50/50	0/6/6/6
4	COA	A	703	3	-	9/44/64/64	0/3/3/3
3	6NA	B	702	4	-	0/4/4/5	-

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	703	COA	O4B-C1B	17.20	1.63	1.40
4	A	703	COA	O4B-C1B	16.84	1.63	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	703	COA	P2A-O3A	8.03	1.68	1.59
4	A	703	COA	P2A-O3A	7.14	1.67	1.59
4	B	703	COA	P1A-O3A	6.92	1.67	1.59
4	A	703	COA	O4B-C4B	-6.62	1.30	1.45
4	B	703	COA	O4B-C4B	-6.46	1.30	1.45
4	A	703	COA	P1A-O3A	6.43	1.66	1.59
4	B	703	COA	C9P-N8P	5.90	1.47	1.33
4	A	703	COA	C5P-N4P	5.81	1.47	1.33
4	A	703	COA	C9P-N8P	5.44	1.46	1.33
4	B	703	COA	C5P-N4P	5.25	1.45	1.33
2	B	701	FAD	P-O3P	4.84	1.64	1.59
4	A	703	COA	P3B-O3B	4.63	1.67	1.59
4	B	703	COA	P3B-O3B	4.57	1.67	1.59
4	B	703	COA	O2B-C2B	4.31	1.53	1.43
4	A	703	COA	O2B-C2B	4.12	1.53	1.43
4	A	703	COA	C6A-N6A	3.16	1.45	1.34
4	A	703	COA	C2A-N3A	3.05	1.36	1.32
4	B	703	COA	C6A-N6A	3.04	1.45	1.34
4	B	703	COA	C2A-N3A	3.03	1.36	1.32
4	B	703	COA	P2A-O6A	2.29	1.68	1.59
4	A	703	COA	O5P-C5P	-2.20	1.18	1.23
4	B	703	COA	O3B-C3B	-2.17	1.36	1.44
4	A	703	COA	P1A-O5B	2.16	1.67	1.59
4	A	703	COA	P2A-O6A	2.03	1.67	1.59
4	A	703	COA	O3B-C3B	-2.01	1.37	1.44

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	703	COA	N3A-C2A-N1A	-6.83	119.39	128.67
4	B	703	COA	N3A-C2A-N1A	-6.39	120.00	128.67
4	B	703	COA	C5A-C6A-N6A	5.54	128.76	120.31
4	A	703	COA	C5A-C6A-N6A	5.18	128.21	120.31
4	B	703	COA	O4B-C1B-N9A	-4.83	102.34	108.75
4	B	703	COA	C3B-C2B-C1B	4.12	108.96	99.89
4	A	703	COA	C3B-C2B-C1B	3.85	108.37	99.89
4	B	703	COA	N6A-C6A-N1A	-3.59	110.67	118.33
4	B	703	COA	C4B-O4B-C1B	-3.53	106.69	109.92
4	A	703	COA	O4B-C1B-N9A	-3.45	104.17	108.75
4	A	703	COA	N6A-C6A-N1A	-3.24	111.42	118.33
4	B	703	COA	CDP-CBP-CCP	3.20	113.50	108.22
4	B	703	COA	C2P-C3P-N4P	-2.99	105.52	112.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	703	COA	C3P-N4P-C5P	-2.94	117.36	122.82
4	B	703	COA	O5P-C5P-N4P	-2.82	117.49	123.03
4	A	703	COA	C6P-C7P-N8P	-2.81	106.02	112.00
4	A	703	COA	CDP-CBP-CAP	2.70	113.37	108.77
4	B	703	COA	C6P-C7P-N8P	-2.69	106.28	112.00
4	A	703	COA	P3B-O3B-C3B	-2.58	116.53	123.43
2	A	701	FAD	C5A-C6A-N6A	2.45	124.04	120.31
4	A	703	COA	C4B-O4B-C1B	-2.38	107.75	109.92
2	B	701	FAD	C5A-C6A-N6A	2.24	123.72	120.31
4	B	703	COA	O5P-C5P-C6P	2.22	126.04	122.02
4	B	703	COA	CEP-CBP-CCP	2.18	111.83	108.22

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	FAD	C2'-C1'-N10-C10
2	B	701	FAD	C2'-C1'-N10-C10
4	B	703	COA	CDP-CBP-CCP-O6A
4	B	703	COA	CEP-CBP-CCP-O6A
4	B	703	COA	CAP-CBP-CCP-O6A
4	A	703	COA	O9P-C9P-CAP-OAP
4	B	703	COA	O9P-C9P-CAP-OAP
4	A	703	COA	N8P-C9P-CAP-OAP
2	A	701	FAD	C1'-C2'-C3'-O3'
2	B	701	FAD	C1'-C2'-C3'-O3'
4	A	703	COA	P2A-O3A-P1A-O1A
4	B	703	COA	P2A-O3A-P1A-O1A
4	A	703	COA	CEP-CBP-CCP-O6A
3	A	702	6NA	C6-CD-CG-CB
4	A	703	COA	S1P-C2P-C3P-N4P
4	B	703	COA	S1P-C2P-C3P-N4P
4	A	703	COA	C5B-O5B-P1A-O1A
4	A	703	COA	C2P-C3P-N4P-C5P
4	B	703	COA	N8P-C9P-CAP-OAP
4	A	703	COA	CAP-CBP-CCP-O6A
4	A	703	COA	P2A-O3A-P1A-O2A
4	B	703	COA	P2A-O3A-P1A-O2A

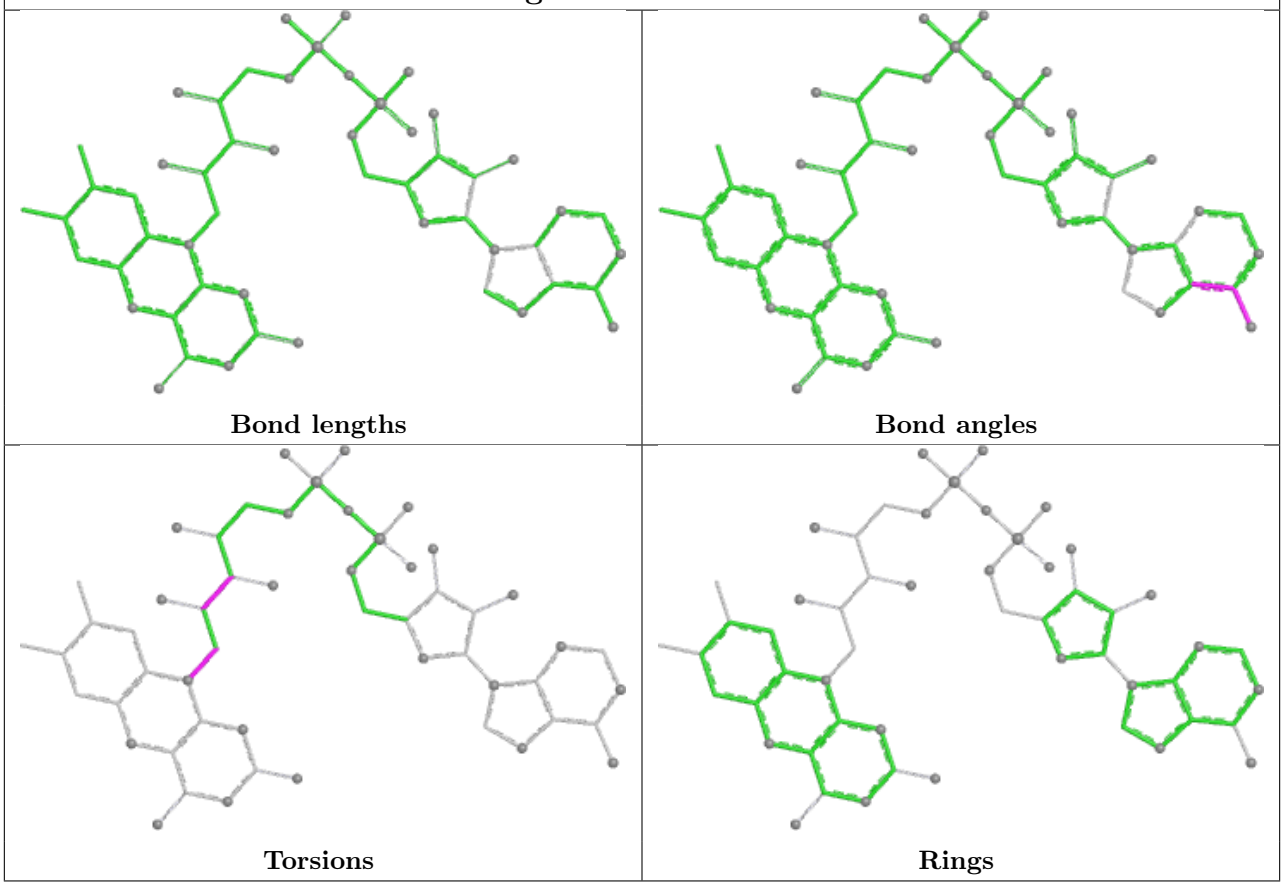
There are no ring outliers.

5 monomers are involved in 26 short contacts:

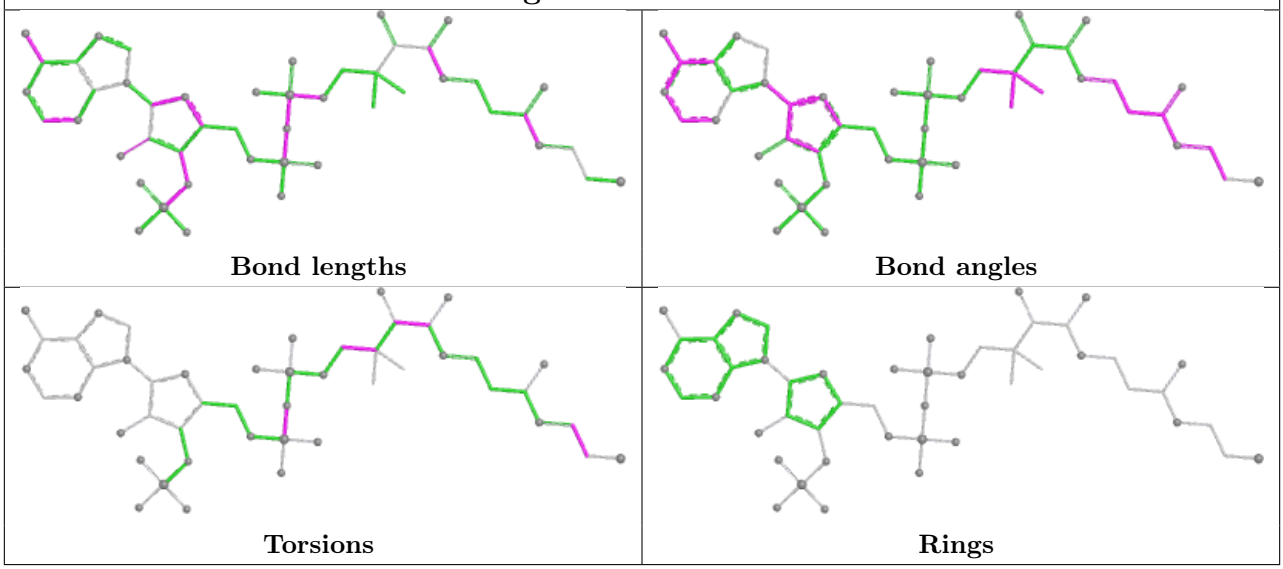
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	702	6NA	1	0
2	A	701	FAD	8	0
4	B	703	COA	3	0
2	B	701	FAD	14	0
4	A	703	COA	2	0

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 FAD A 701



Ligand COA B 703



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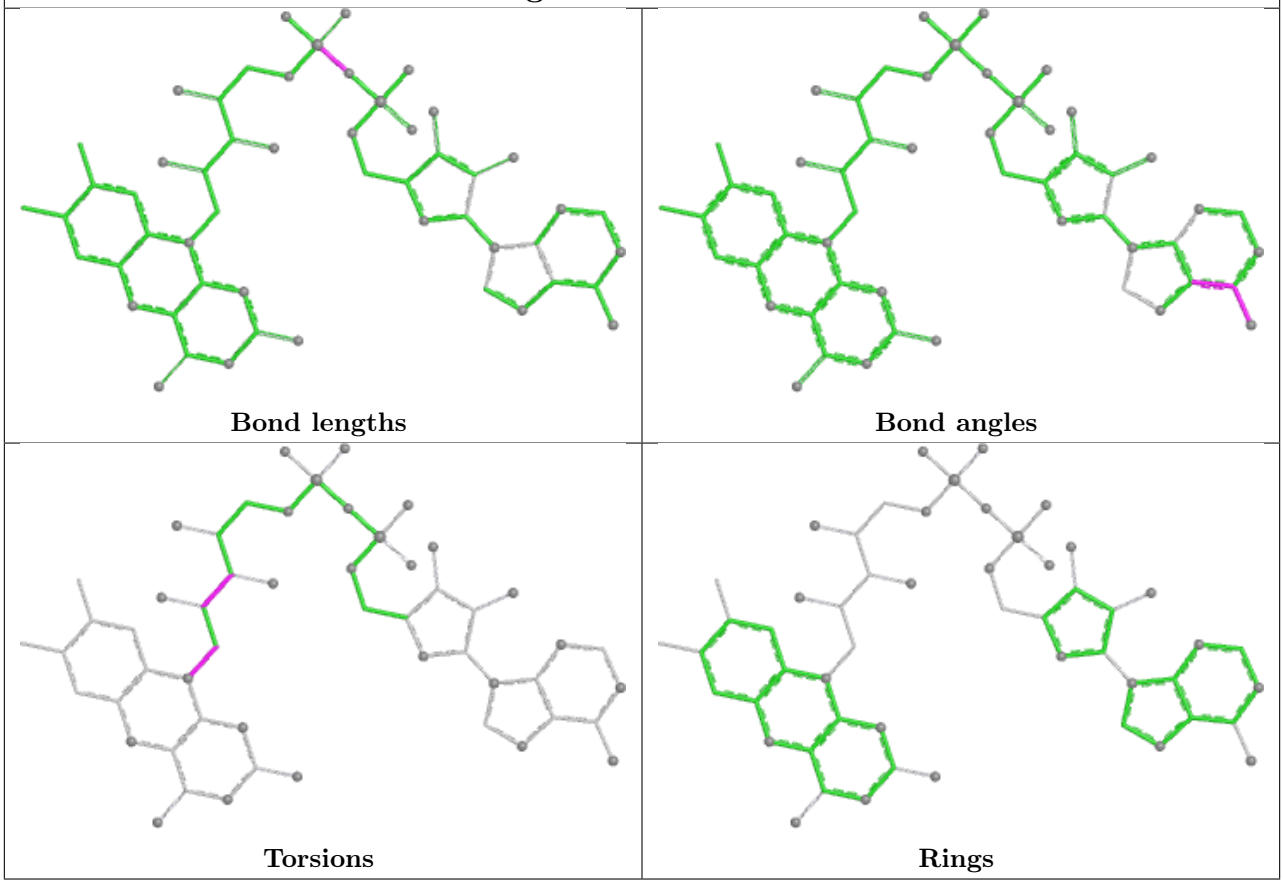
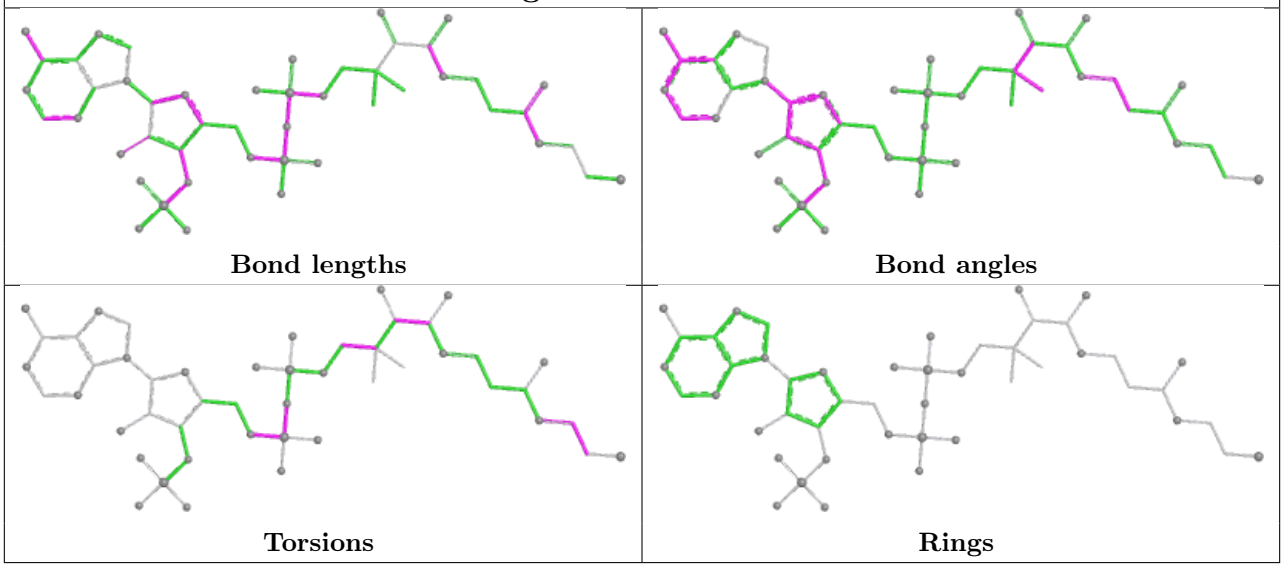


Ligand COA A 703



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	607/611 (99%)	-0.30	17 (2%)	53	51	20, 30, 47, 69	0
1	B	607/611 (99%)	-0.36	12 (1%)	65	63	20, 28, 44, 70	0
All	All	1214/1222 (99%)	-0.33	29 (2%)	59	56	20, 29, 46, 70	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	7.1
1	B	488	GLY	4.8
1	B	336	THR	4.7
1	A	27	ALA	4.1
1	B	32	ASP	3.8
1	A	32	ASP	3.8
1	B	28	ASP	3.5
1	A	28	ASP	3.4
1	A	489	ARG	3.3
1	A	31	SER	3.2
1	B	338	LYS	3.0
1	B	339	THR	2.9
1	B	31	SER	2.9
1	A	564	THR	2.7
1	A	26	GLY	2.7
1	B	335	MET	2.6
1	A	29	LYS	2.5
1	B	337	ASP	2.4
1	A	246	GLU	2.4
1	A	483	ASN	2.4
1	A	336	THR	2.3
1	A	464	ARG	2.2
1	A	177	ARG	2.2
1	A	240	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	487	ASN	2.1
1	A	219	GLY	2.1
1	A	482	LYS	2.1
1	B	379	GLU	2.0
1	B	482	LYS	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 monosaccharides in this entry.

6.4 Ligands (i)

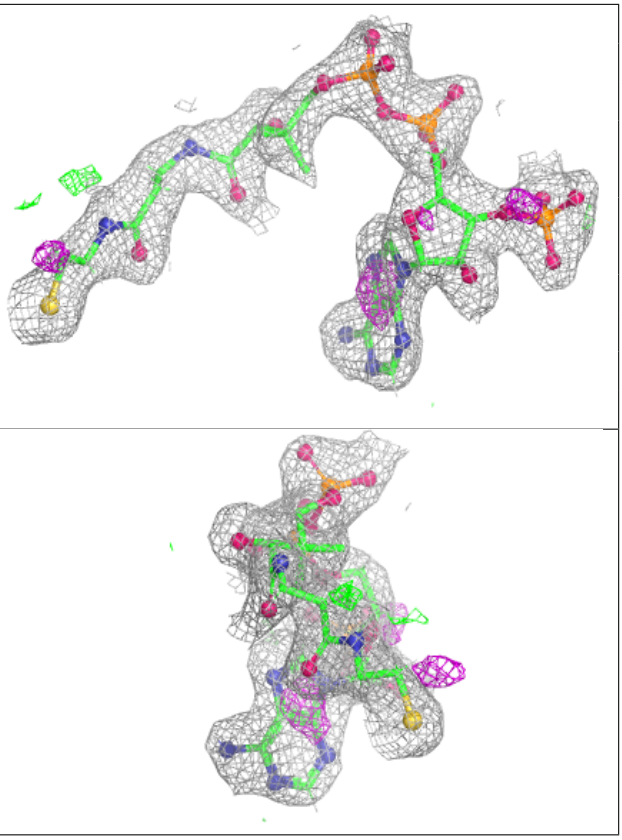
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

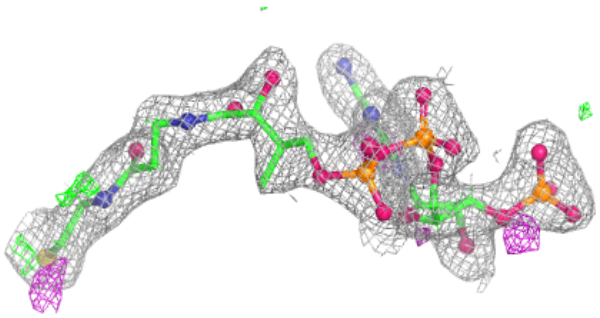
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	COA	B	703	48/48	0.87	0.18	34,45,57,63	0
3	6NA	B	702	7/8	0.89	0.17	39,49,55,55	0
4	COA	A	703	48/48	0.92	0.17	35,46,56,61	0
3	6NA	A	702	7/8	0.93	0.15	37,44,54,54	0
2	FAD	B	701	53/53	0.94	0.11	21,30,39,47	0
2	FAD	A	701	53/53	0.95	0.11	23,33,46,54	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 B 703:

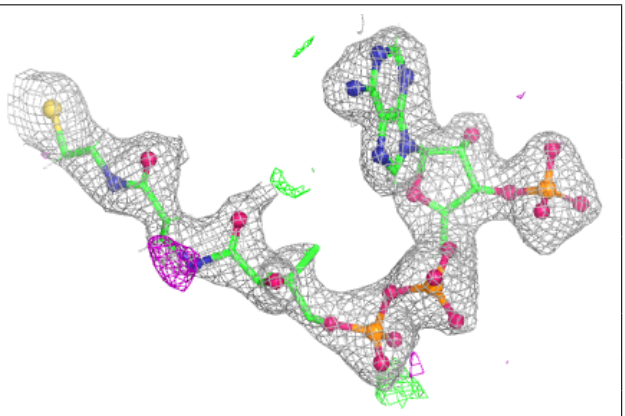
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

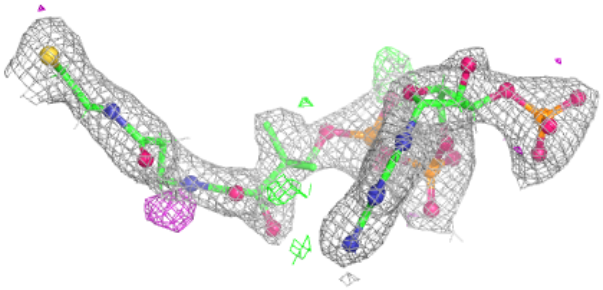


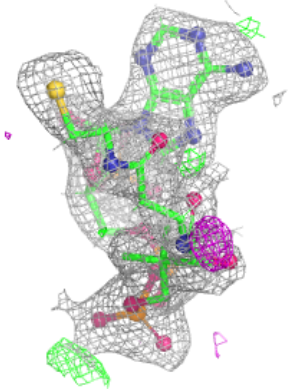


Electron density around COA A 703:

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

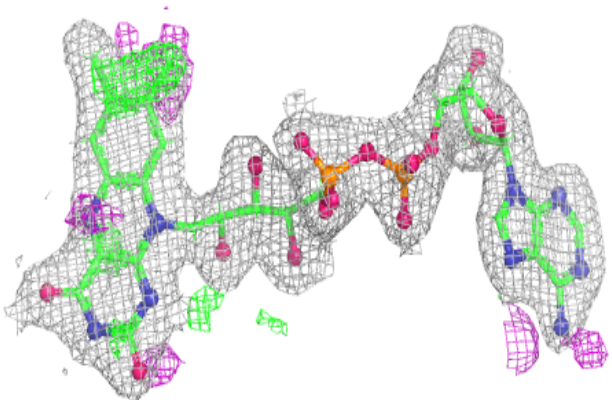


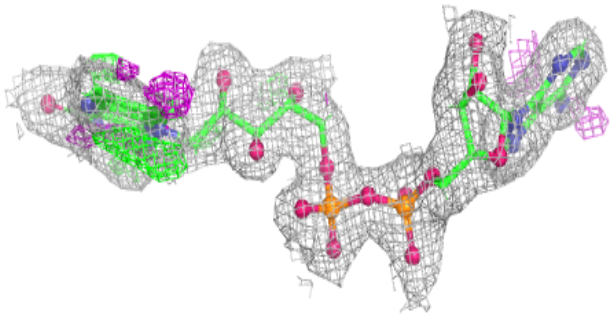


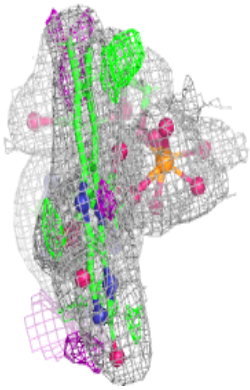


Electron density around FAD B 701:

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

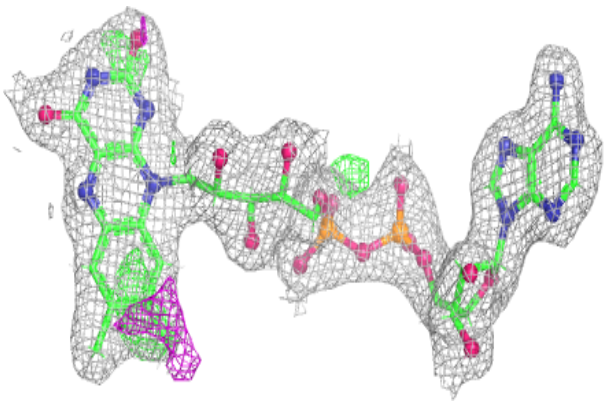


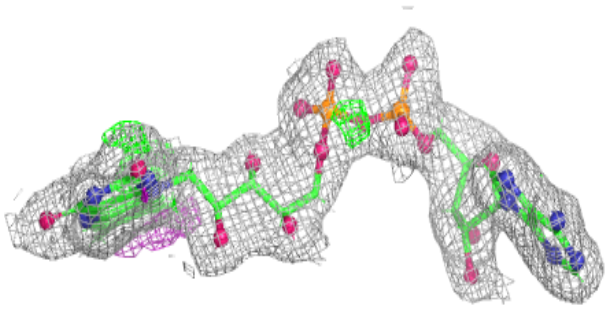


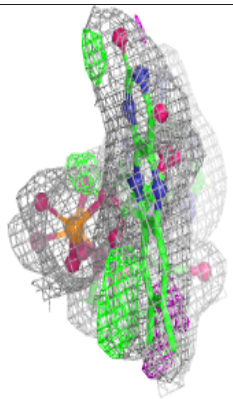


Electron density around FAD A 701:

$2mF_o$-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.