wwPDB X-ray Structure Validation Summary Report (i)

Nov 23, 2023 – 02:12 AM JST

PDB ID	:	7Y0B
Title	:	Crystal structure of human short-chain acyl-CoA dehydrogenase
Authors	:	Huang, Y.; Xu, Y.; Li, J.
Deposited on	:	2022-06-04
Resolution	:	2.08 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:

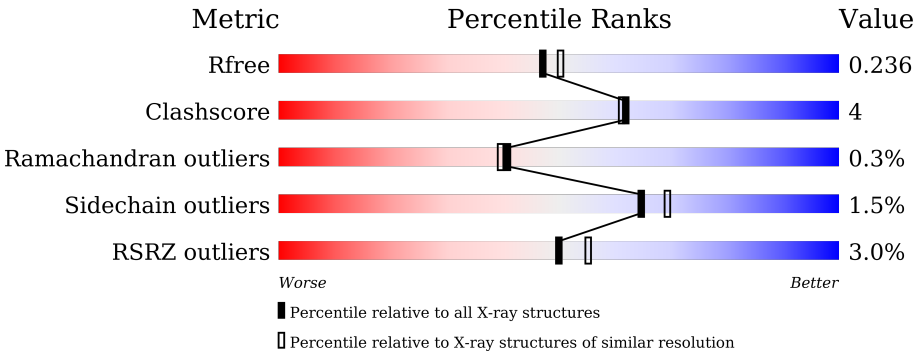
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.08 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	6189 (2.10-2.06)
Clashscore	141614	6738 (2.10-2.06)
Ramachandran outliers	138981	6663 (2.10-2.06)
Sidechain outliers	138945	6664 (2.10-2.06)
RSRZ outliers	127900	6057 (2.10-2.06)

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	398	3% 88% 8%
1	B	398	2% 86% 9%
1	C	398	4% 88% 9%
1	D	398	3% 87% 9%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	7R3	B	502	-	-	-	X

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 12259 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 Short-chain specific acyl-CoA dehydrogenase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	385	Total	C	N	O	S	0	0	0
			2885	1828	495	543	19			
1	B	383	Total	C	N	O	S	0	0	0
			2864	1814	491	539	20			
1	C	387	Total	C	N	O	S	0	2	0
			2903	1842	500	541	20			
1	D	385	Total	C	N	O	S	0	0	0
			2888	1831	494	543	20			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	MET	-	initiating methionine	UNP P16219
A	24	GLY	-	expression tag	UNP P16219
A	413	LEU	-	expression tag	UNP P16219
A	414	GLU	-	expression tag	UNP P16219
A	415	HIS	-	expression tag	UNP P16219
A	416	HIS	-	expression tag	UNP P16219
A	417	HIS	-	expression tag	UNP P16219
A	418	HIS	-	expression tag	UNP P16219
A	419	HIS	-	expression tag	UNP P16219
A	420	HIS	-	expression tag	UNP P16219
B	23	MET	-	initiating methionine	UNP P16219
B	24	GLY	-	expression tag	UNP P16219
B	413	LEU	-	expression tag	UNP P16219
B	414	GLU	-	expression tag	UNP P16219
B	415	HIS	-	expression tag	UNP P16219
B	416	HIS	-	expression tag	UNP P16219
B	417	HIS	-	expression tag	UNP P16219
B	418	HIS	-	expression tag	UNP P16219
B	419	HIS	-	expression tag	UNP P16219
B	420	HIS	-	expression tag	UNP P16219
C	23	MET	-	initiating methionine	UNP P16219

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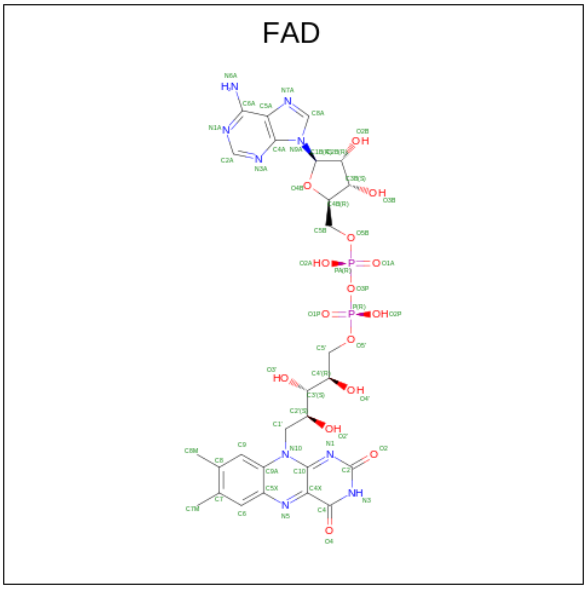
Chain	Residue	Modelled	Actual	Comment	Reference
C	24	GLY	-	expression tag	UNP P16219
C	413	LEU	-	expression tag	UNP P16219
C	414	GLU	-	expression tag	UNP P16219
C	415	HIS	-	expression tag	UNP P16219
C	416	HIS	-	expression tag	UNP P16219
C	417	HIS	-	expression tag	UNP P16219
C	418	HIS	-	expression tag	UNP P16219
C	419	HIS	-	expression tag	UNP P16219
C	420	HIS	-	expression tag	UNP P16219
D	23	MET	-	initiating methionine	UNP P16219
D	24	GLY	-	expression tag	UNP P16219
D	413	LEU	-	expression tag	UNP P16219
D	414	GLU	-	expression tag	UNP P16219
D	415	HIS	-	expression tag	UNP P16219
D	416	HIS	-	expression tag	UNP P16219
D	417	HIS	-	expression tag	UNP P16219
D	418	HIS	-	expression tag	UNP P16219
D	419	HIS	-	expression tag	UNP P16219
D	420	HIS	-	expression tag	UNP P16219

- 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
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

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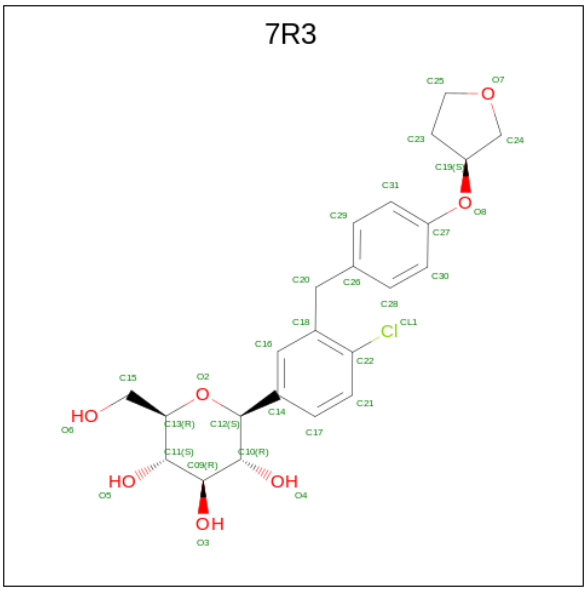
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is (2S,3R,4R,5S,6R)-2-[4-chloranyl-3-[[4-[(3S)-oxolan-3-yl]oxyphenyl]methyl]phenyl]-6-(hydroxymethyl)oxane-3,4,5-triol (three-letter code: 7R3) (formula: $C_{23}H_{27}ClO_7$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	Cl	O	0	0
			31	23	1	7		

- Molecule 4 is water.

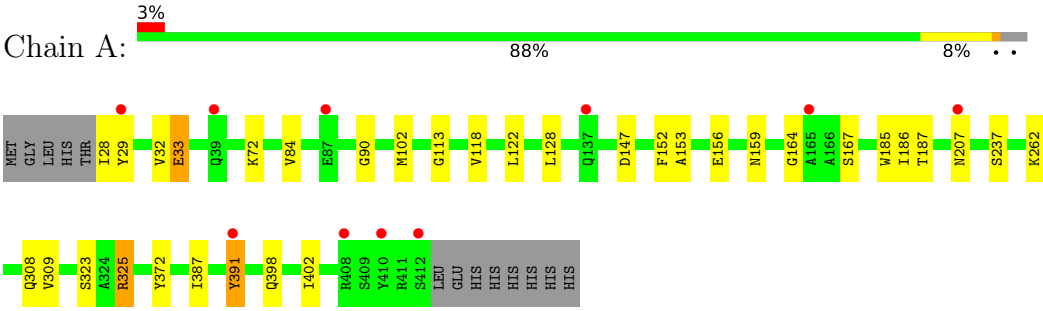
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	138	Total	O	0	0
			138	138		
4	B	75	Total	O	0	0
			75	75		
4	C	132	Total	O	0	0
			132	132		
4	D	131	Total	O	0	0
			131	131		

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.

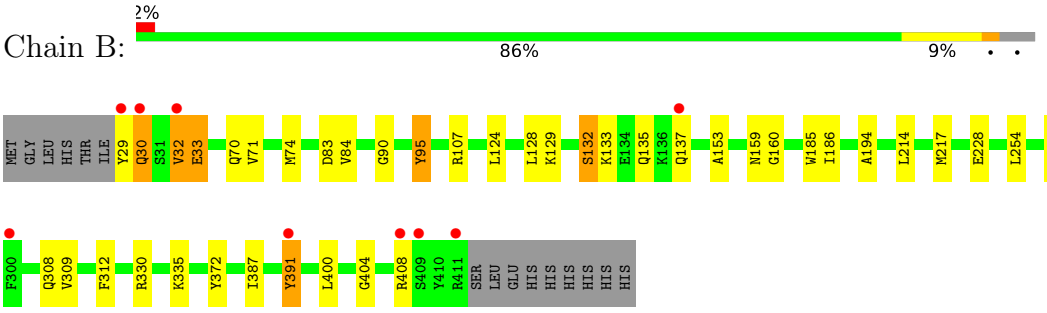
• Molecule 1: Short-chain specific acyl-CoA dehydrogenase, mitochondrial

Chain A:



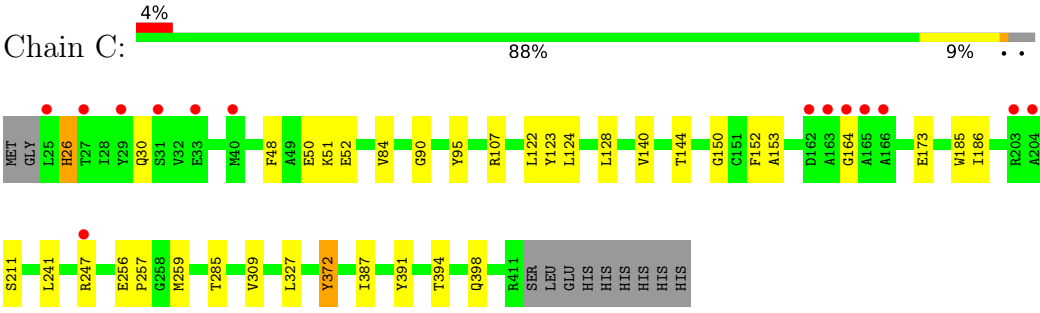
• Molecule 1: Short-chain specific acyl-CoA dehydrogenase, mitochondrial

Chain B:



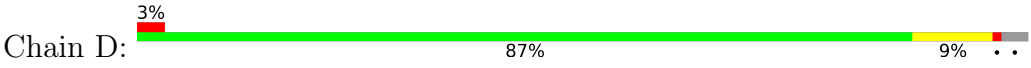
• Molecule 1: Short-chain specific acyl-CoA dehydrogenase, mitochondrial

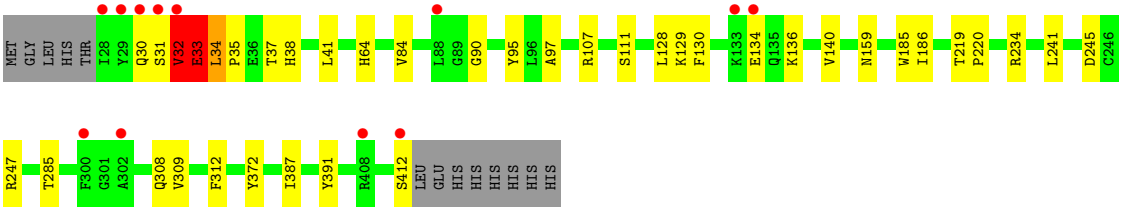
Chain C:



• Molecule 1: Short-chain specific acyl-CoA dehydrogenase, mitochondrial

Chain D:




MET GLY LEU HIS THR I28 Y29 Q30 S31 V32 E33 L34 P35 E36 T37 H38 L41 H64 V84 L88 G89 G90 Y95 L96 A97 R107 S111 L128 K129 F130 K133 E134 Q135 K136 V140 N159 W185 I186 T219 P220 R234 L241 D245 C246
R247 T285 F300 G301 A302 Q308 V309 F312 Y372 I387 Y391 R408 S412 LEU GLU HIS HIS HIS HIS HIS HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	80.99Å 142.08Å 154.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.17 – 2.08 39.17 – 2.08	Depositor EDS
% Data completeness (in resolution range)	96.8 (39.17-2.08) 96.8 (39.17-2.08)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.11 (at 2.08Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.197 , 0.237 0.196 , 0.236	Depositor DCC
R_{free} test set	2004 reflections (1.92%)	wwPDB-VP
Wilson B-factor (Å^2)	17.9	Xtriage
Anisotropy	0.826	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 48.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	12259	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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: FAD, 7R3

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	0.52	3/2936 (0.1%)	0.58	2/3970 (0.1%)
1	B	0.48	0/2915	0.61	3/3942 (0.1%)
1	C	0.58	2/2968 (0.1%)	0.60	1/4014 (0.0%)
1	D	0.65	1/2939 (0.0%)	0.67	5/3973 (0.1%)
All	All	0.56	6/11758 (0.1%)	0.62	11/15899 (0.1%)

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	156	GLU	CD-OE2	-7.19	1.17	1.25
1	D	111	SER	CB-OG	-6.77	1.33	1.42
1	A	164	GLY	C-O	-6.25	1.13	1.23
1	C	164	GLY	C-O	-5.88	1.14	1.23
1	A	167	SER	CB-OG	-5.19	1.35	1.42

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	32	VAL	N-CA-C	-8.10	89.13	111.00
1	D	32	VAL	CA-CB-CG1	6.56	120.73	110.90
1	C	207	ASN	N-CA-CB	-6.54	98.82	110.60
1	D	32	VAL	CB-CA-C	6.47	123.69	111.40
1	D	34	LEU	N-CA-C	6.42	128.32	111.00

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	2885	0	2908	26	0
1	B	2864	0	2870	33	0
1	C	2903	0	2926	26	0
1	D	2888	0	2912	33	0
2	A	53	0	31	5	0
2	B	53	0	31	7	0
2	C	53	0	29	4	0
2	D	53	0	31	5	0
3	B	31	0	0	1	0
4	A	138	0	0	2	0
4	B	75	0	0	0	0
4	C	132	0	0	1	0
4	D	131	0	0	0	0
All	All	12259	0	11738	104	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 4.

The worst 5 of 104 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:B:330:ARG:HD2	1:D:31:SER:HB2	1.51	0.92
1:A:387:ILE:HD11	2:A:601:FAD:HM83	1.61	0.83
1:B:160:GLY:HA3	2:B:501:FAD:H5'2	1.66	0.78
1:A:187:THR:HG22	2:A:601:FAD:O4	1.84	0.78
1:B:330:ARG:HD2	1:D:31:SER:CB	2.13	0.77

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	383/398 (96%)	365 (95%)	17 (4%)	1 (0%)	41	39
1	B	381/398 (96%)	367 (96%)	13 (3%)	1 (0%)	41	39
1	C	387/398 (97%)	374 (97%)	12 (3%)	1 (0%)	41	39
1	D	383/398 (96%)	366 (96%)	16 (4%)	1 (0%)	41	39
All	All	1534/1592 (96%)	1472 (96%)	58 (4%)	4 (0%)	41	39

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	32	VAL
1	A	33	GLU
1	B	33	GLU
1	C	26	HIS

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	293/309 (95%)	289 (99%)	4 (1%)	67	72
1	B	288/309 (93%)	282 (98%)	6 (2%)	53	57
1	C	294/309 (95%)	292 (99%)	2 (1%)	84	87
1	D	293/309 (95%)	287 (98%)	6 (2%)	55	59
All	All	1168/1236 (94%)	1150 (98%)	18 (2%)	65	69

5 of 18 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	95	TYR
1	D	412	SER
1	D	391	TYR

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Mol	Chain	Res	Type
1	B	159	ASN
1	D	33	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	43	GLN
1	B	137	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)

5 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	FAD	C	501	-	53,58,58	2.21	20 (37%)	68,89,89	1.74	15 (22%)
3	7R3	B	502	-	34,34,34	5.56	26 (76%)	45,48,48	2.54	14 (31%)
2	FAD	A	601	-	53,58,58	0.54	0	68,89,89	0.61	1 (1%)
2	FAD	B	501	-	53,58,58	0.62	0	68,89,89	1.11	7 (10%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	FAD	D	601	-	53,58,58	2.16	21 (39%)	68,89,89	1.50	9 (13%)

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	FAD	C	501	-	-	4/30/50/50	0/6/6/6
3	7R3	B	502	-	-	2/14/41/41	0/4/4/4
2	FAD	A	601	-	-	3/30/50/50	0/6/6/6
2	FAD	B	501	-	-	9/30/50/50	0/6/6/6
2	FAD	D	601	-	-	4/30/50/50	0/6/6/6

The worst 5 of 67 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	502	7R3	C16-C14	10.66	1.55	1.39
3	B	502	7R3	C16-C18	9.73	1.56	1.39
3	B	502	7R3	C31-C29	9.00	1.55	1.38
3	B	502	7R3	C22-C18	8.98	1.53	1.39
3	B	502	7R3	C30-C28	8.68	1.54	1.38

The worst 5 of 46 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	502	7R3	O2-C12-C14	8.33	122.46	108.00
3	B	502	7R3	C17-C14-C12	-4.99	111.36	120.64
3	B	502	7R3	C13-O2-C12	-4.78	103.45	112.14
2	C	501	FAD	N3A-C2A-N1A	-4.41	121.78	128.68
2	C	501	FAD	C4-C4X-N5	4.37	124.45	118.23

There are no chirality outliers.

5 of 22 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	FAD	C5B-O5B-PA-O1A
2	A	601	FAD	C5B-O5B-PA-O2A
2	B	501	FAD	C1'-C2'-C3'-C4'
2	B	501	FAD	C5'-O5'-P-O3P

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Mol	Chain	Res	Type	Atoms
3	B	502	7R3	O2-C13-C15-O6

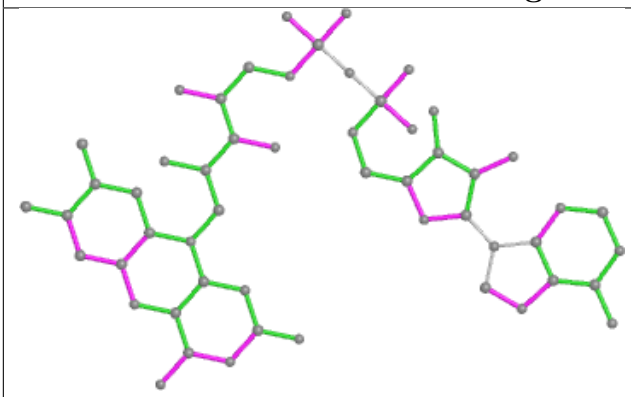
There are no ring outliers.

5 monomers are involved in 22 short contacts:

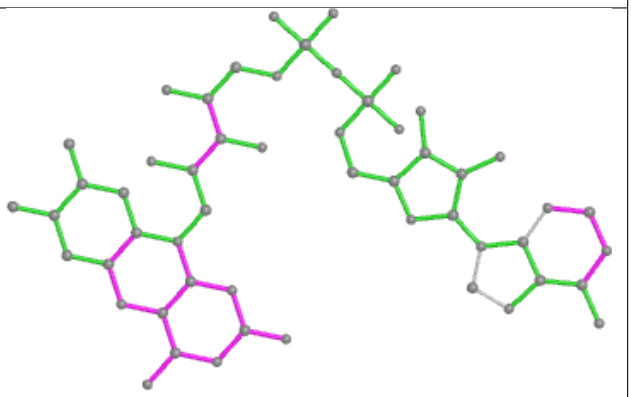
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	501	FAD	4	0
3	B	502	7R3	1	0
2	A	601	FAD	5	0
2	B	501	FAD	7	0
2	D	601	FAD	5	0

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.

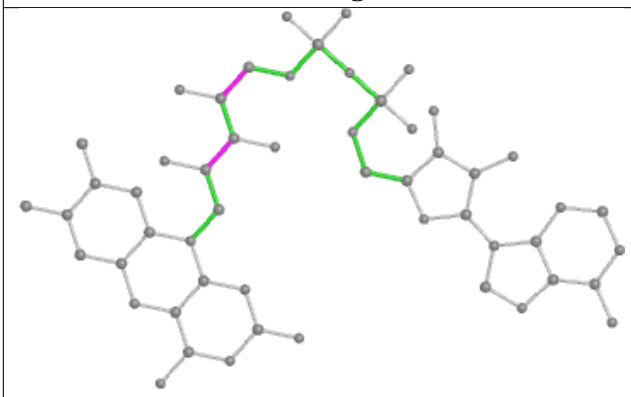
Ligand FAD C 501



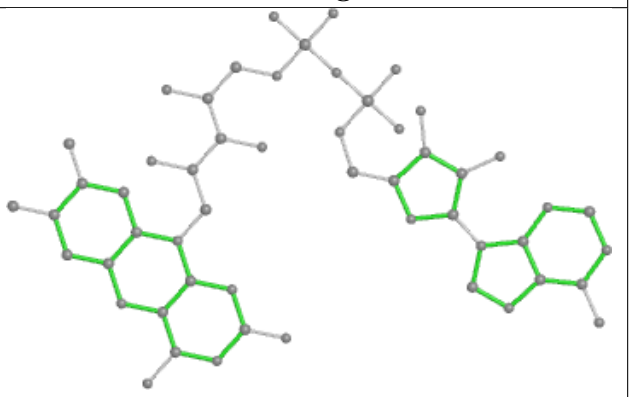
Bond lengths



Bond angles

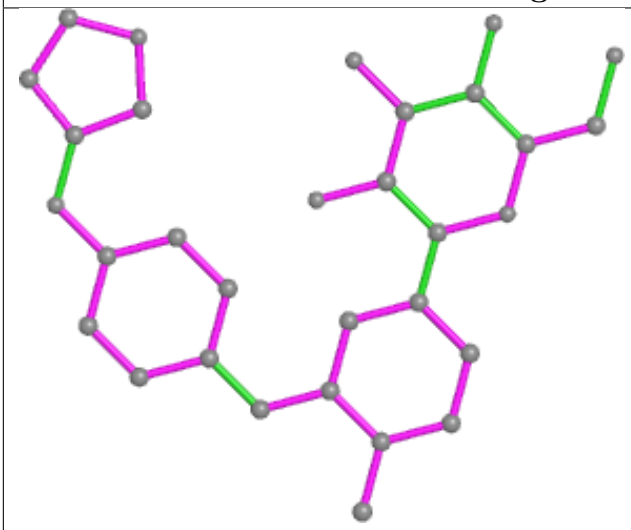


Torsions

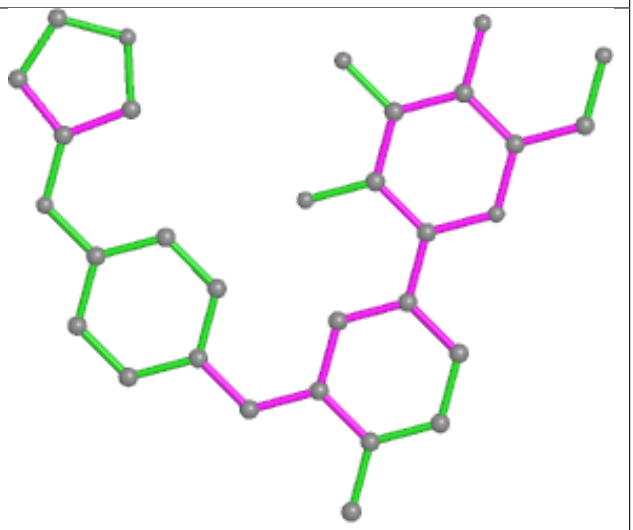


Rings

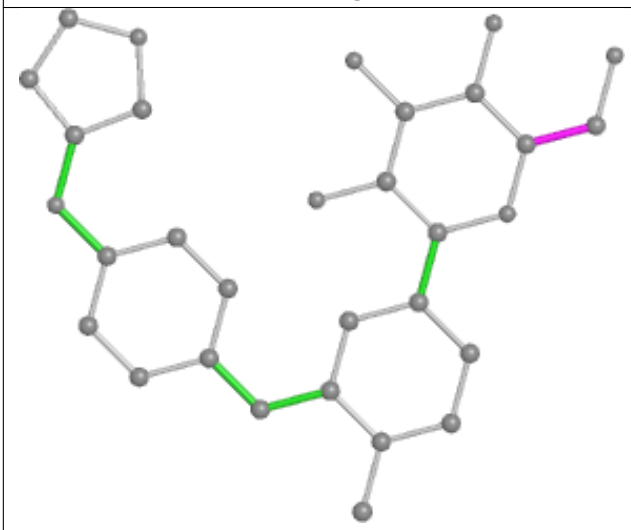
Ligand 7R3 B 502



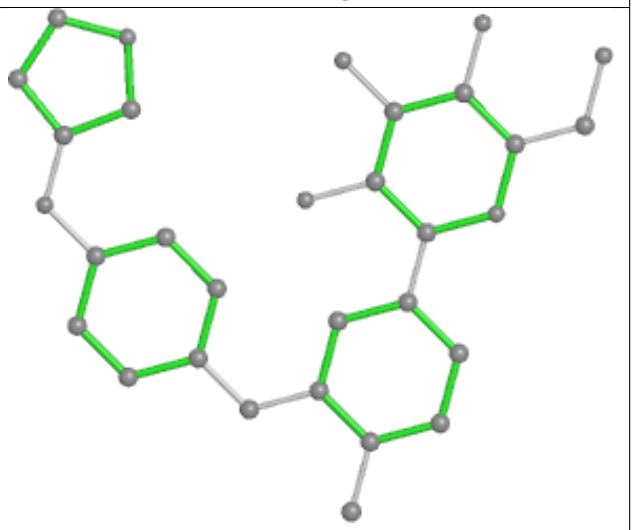
Bond lengths



Bond angles

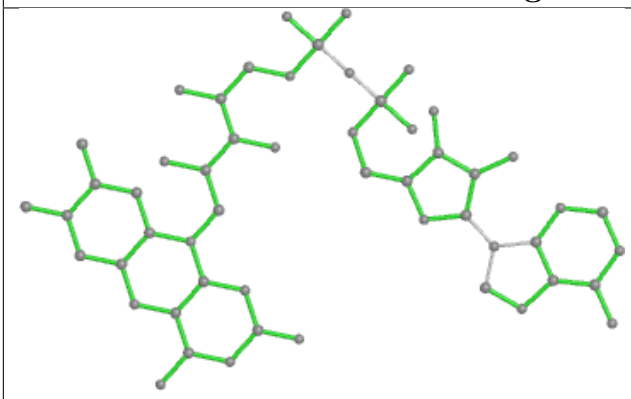


Torsions

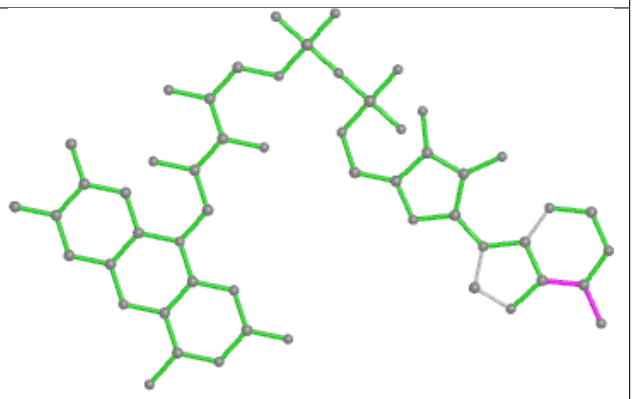


Rings

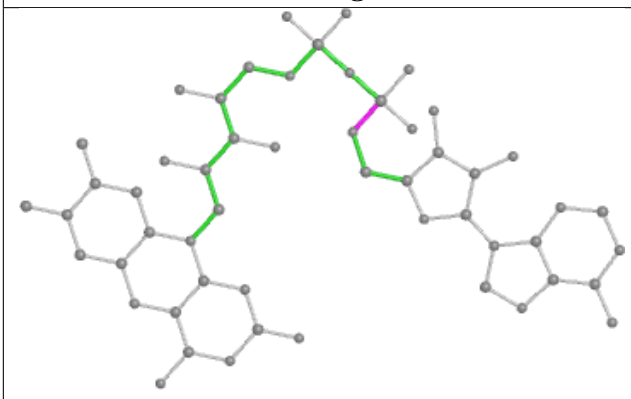
Ligand FAD A 601



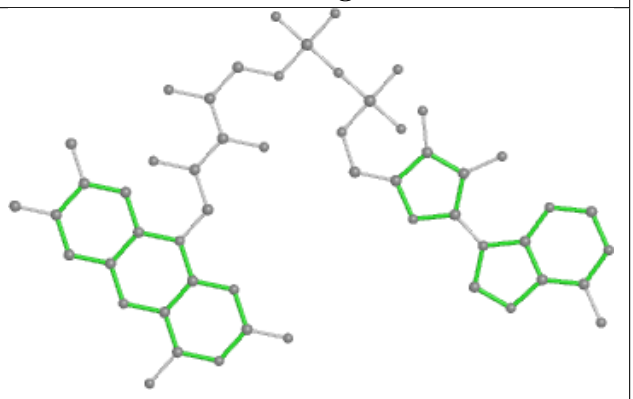
Bond lengths



Bond angles

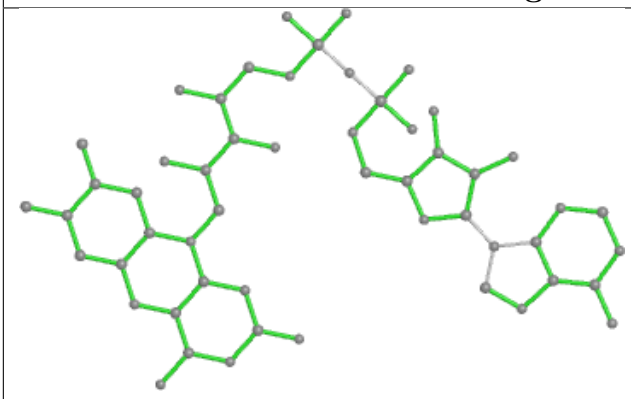


Torsions

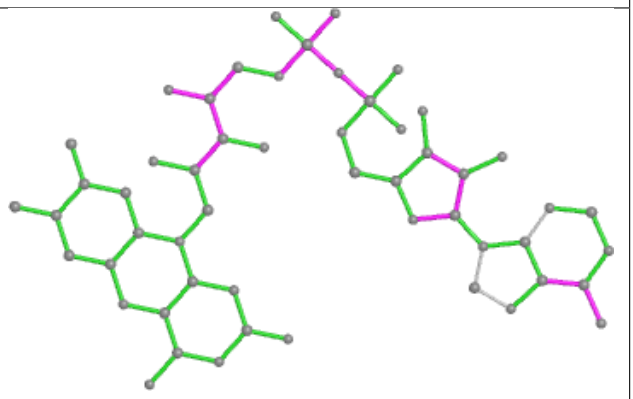


Rings

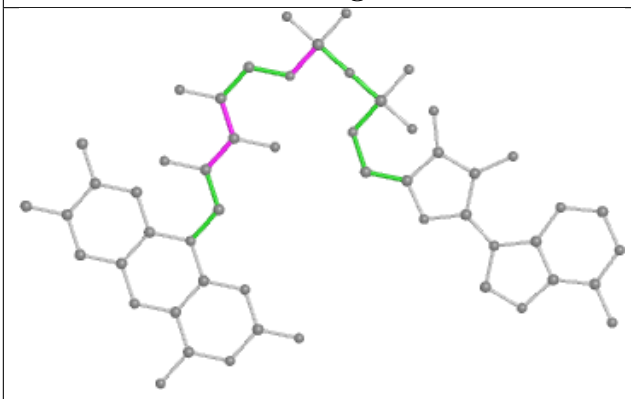
Ligand FAD B 501



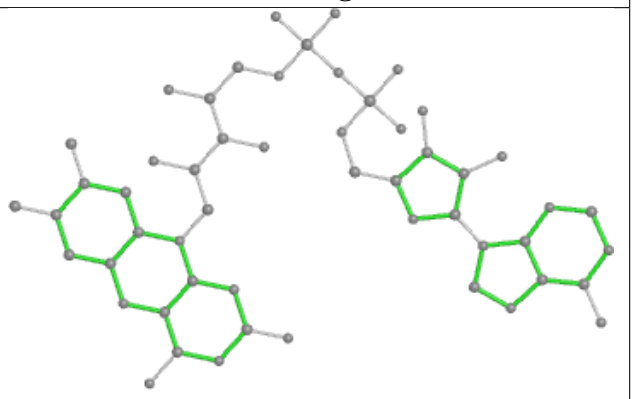
Bond lengths



Bond angles

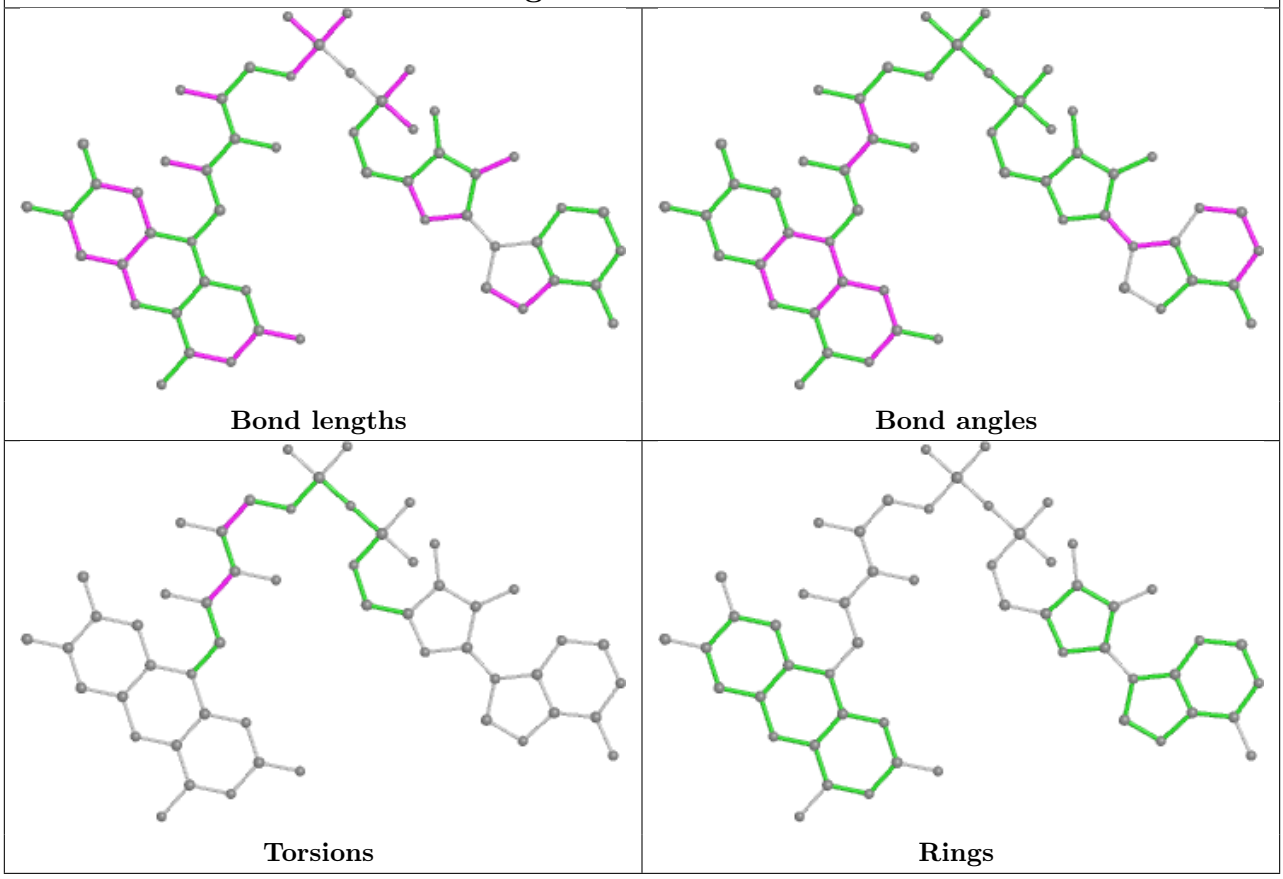


Torsions



Rings

Ligand FAD D 601



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	385/398 (96%)	0.09	10 (2%) 56 60	12, 18, 33, 44	0
1	B	383/398 (96%)	0.06	9 (2%) 60 64	12, 18, 30, 49	0
1	C	387/398 (97%)	0.08	15 (3%) 39 44	12, 20, 36, 46	0
1	D	385/398 (96%)	0.13	12 (3%) 49 54	12, 20, 34, 53	0
All	All	1540/1592 (96%)	0.09	46 (2%) 50 55	12, 19, 33, 53	0

The worst 5 of 46 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	412	SER	4.1
1	D	32	VAL	4.1
1	C	204	ALA	4.1
1	C	165	ALA	4.0
1	C	163	ALA	3.7

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

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

6.3 Carbohydrates

There are no monosaccharides in this entry.

6.4 Ligands

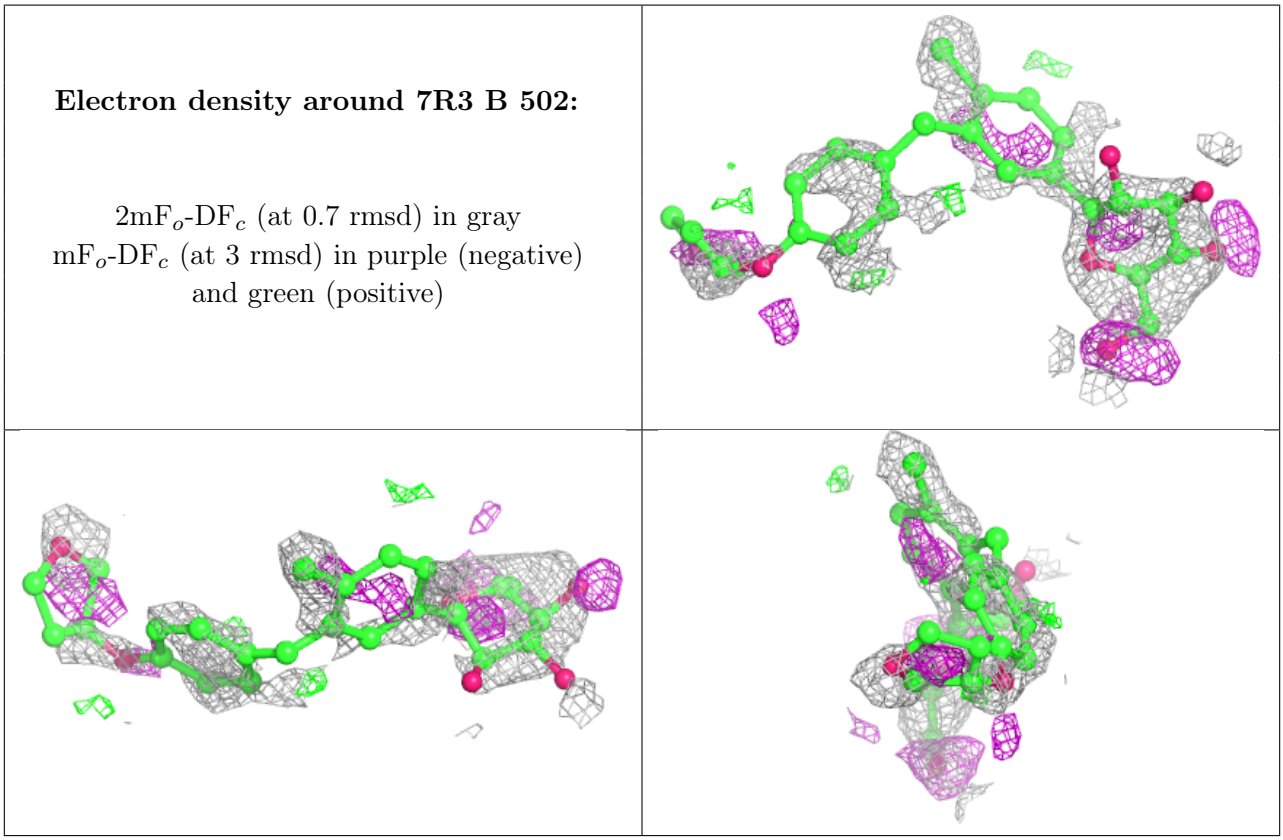
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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
3	7R3	B	502	31/31	0.66	0.59	26,46,55,65	0
2	FAD	B	501	53/53	0.93	0.16	9,17,26,36	0
2	FAD	C	501	53/53	0.94	0.12	13,17,21,24	0
2	FAD	A	601	53/53	0.95	0.12	11,15,19,21	0
2	FAD	D	601	53/53	0.96	0.11	9,16,21,23	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.

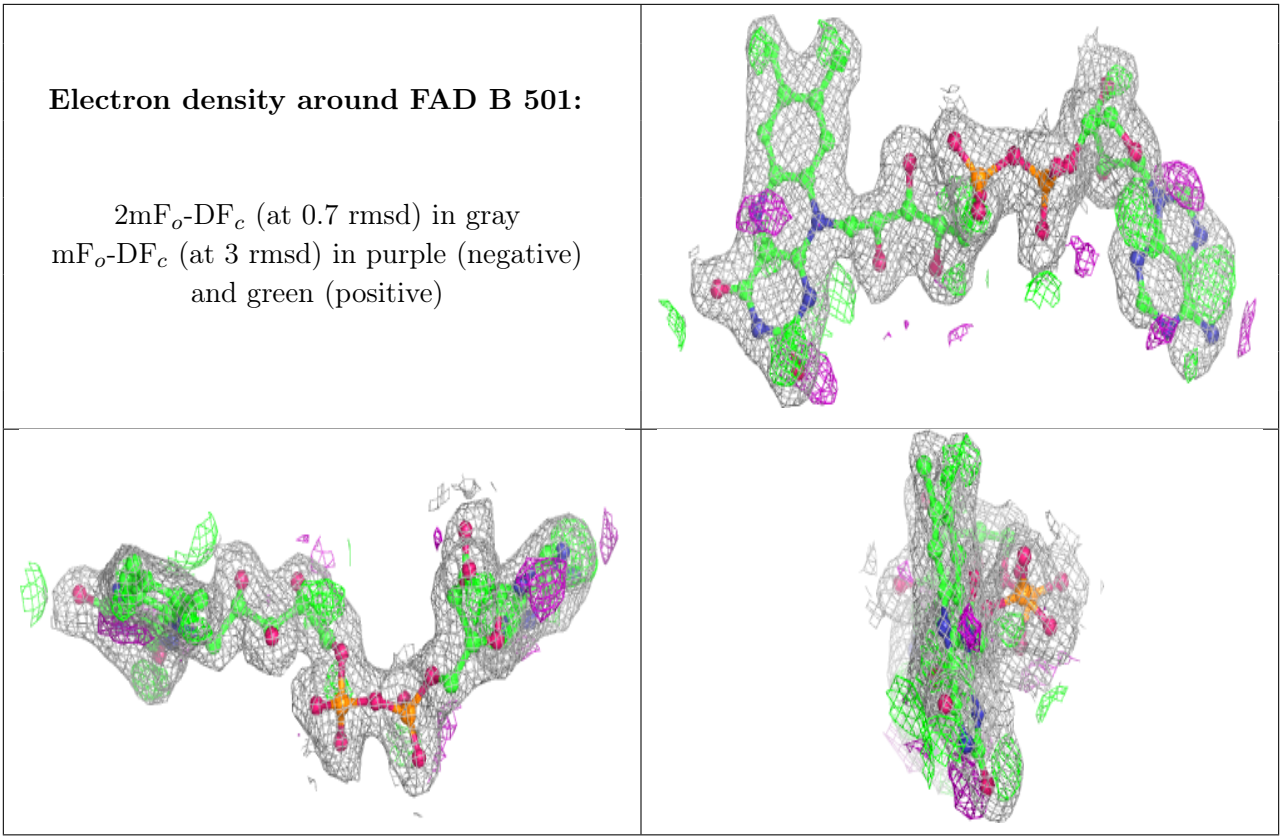
Electron density around 7R3 B 502:

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



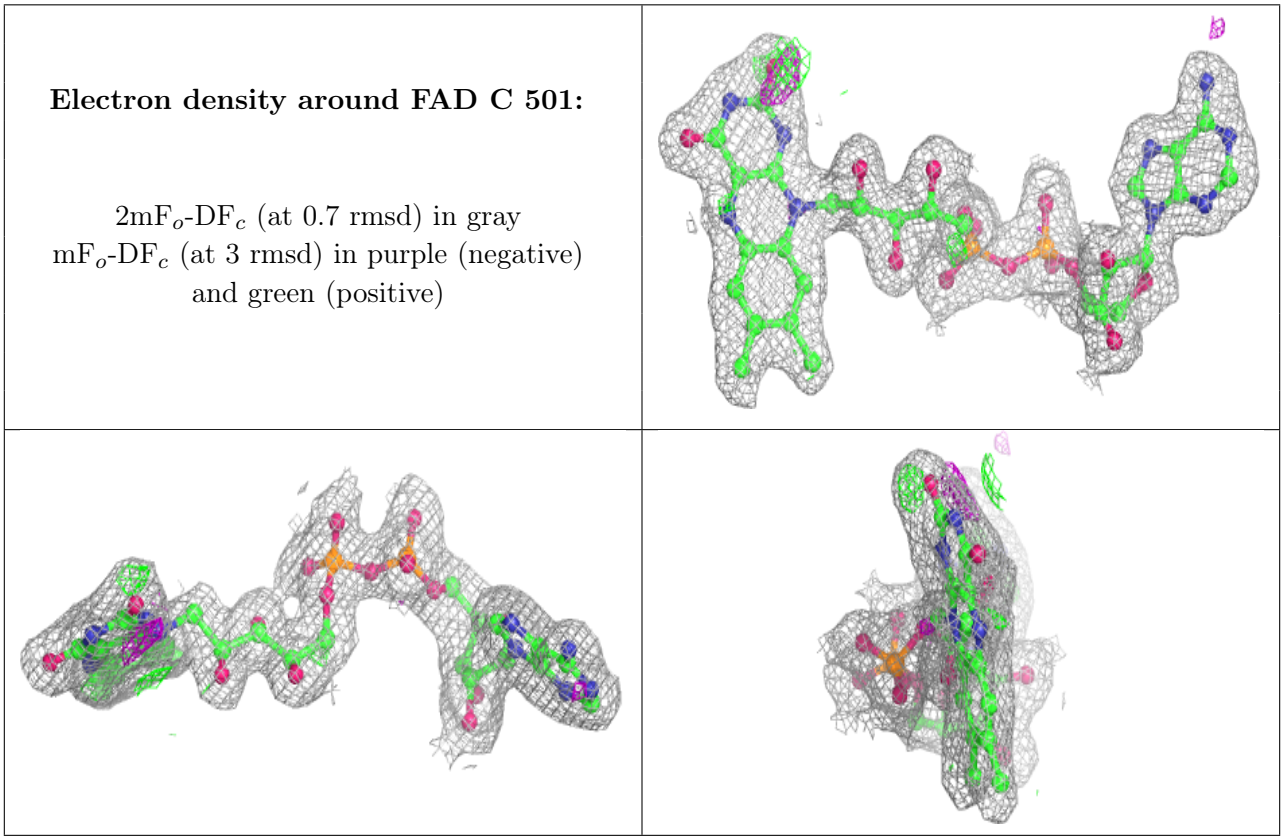
Electron density around FAD B 501:

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



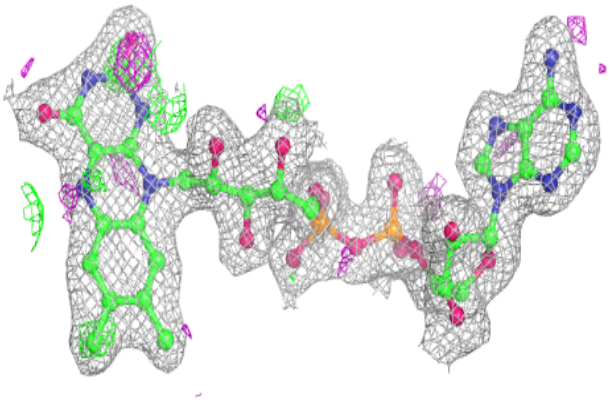
Electron density around FAD C 501:

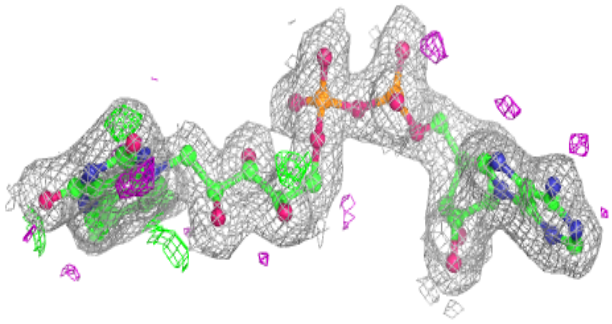
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

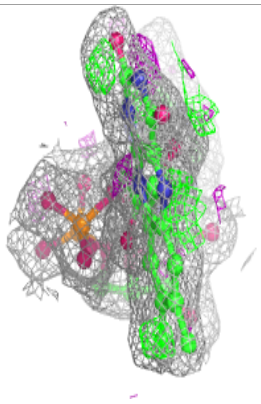


Electron density around FAD A 601:

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

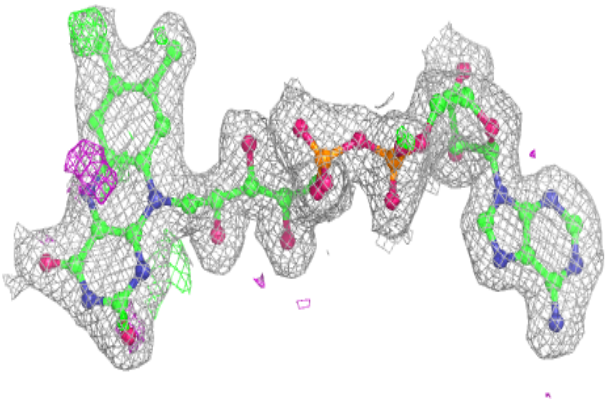


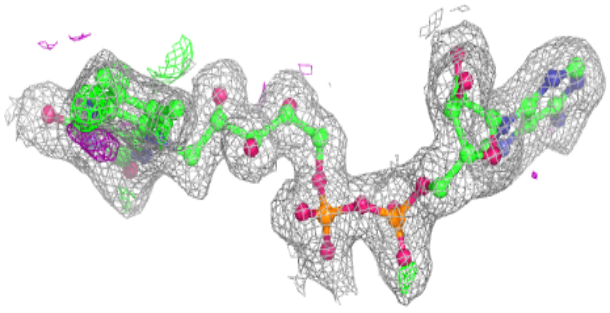


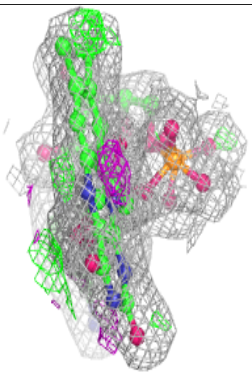


Electron density around FAD D 601:

$2mF_o$-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.