Full wwPDB X-ray Structure Validation Report (i)

Dec 7, 2023 – 05:06 am GMT

PDB ID	:	6YRU
Title	:	Crystal structure of FAP in the dark at 100K
Authors	:	Sorigue, D.; Gotthard, G.; Blangy, S.; Nurizzo, D.; Royant, A.; Beisson, F.; Arnoux, P.
Deposited on	:	2020-04-20
Resolution	:	1.78 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

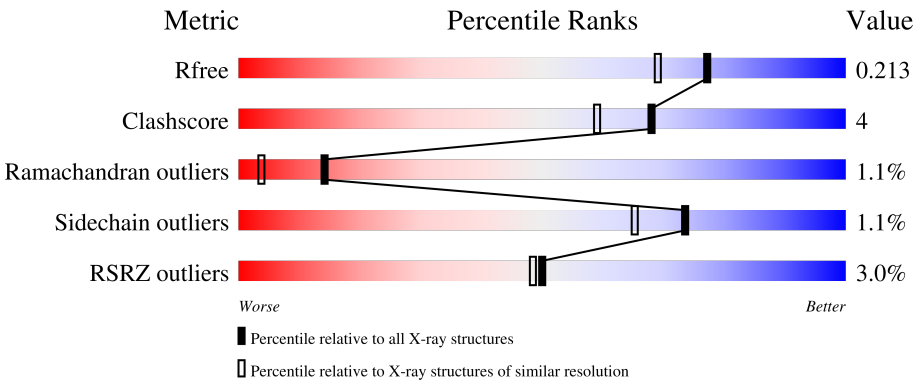
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)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

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.78 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	9185 (1.80-1.76)
Clashscore	141614	10184 (1.80-1.76)
Ramachandran outliers	138981	10051 (1.80-1.76)
Sidechain outliers	138945	10050 (1.80-1.76)
RSRZ outliers	127900	9032 (1.80-1.76)

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	AAA	578	3% 90% 8% ••

2 Entry composition (i)

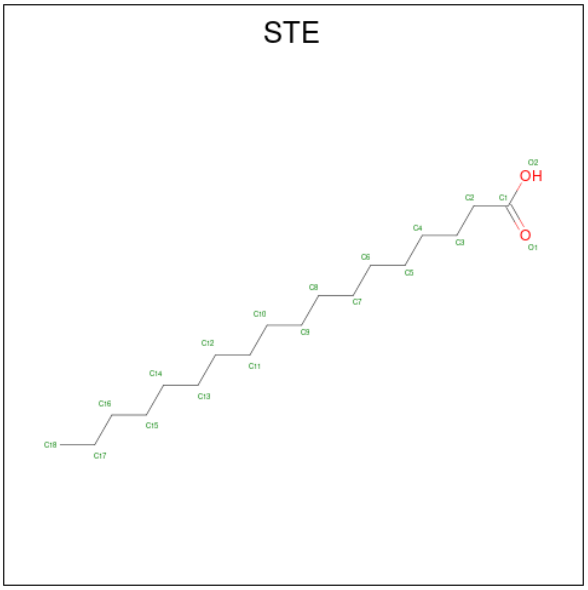
There are 4 unique types of molecules in this entry. The entry contains 4885 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Fatty acid Photodecarboxylase.

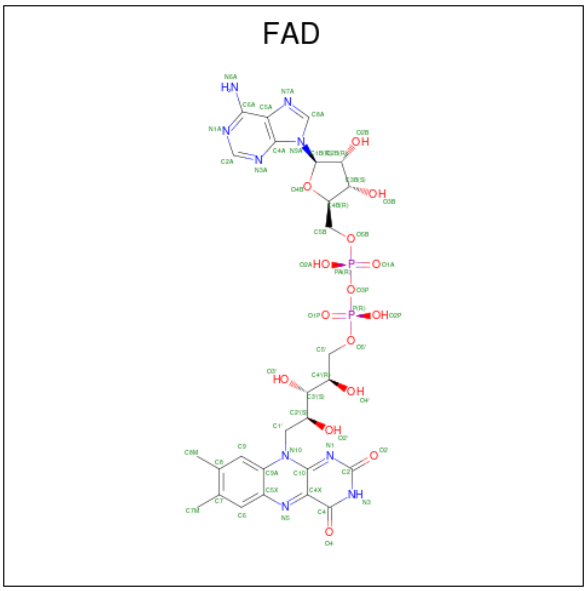
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	AAA	573	4267	2668	764	820	15	0	0	0

- Molecule 2 is STEARIC ACID (three-letter code: STE) (formula: $C_{18}H_{36}O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	AAA	1	20	18	2	0	0
			Total	C	O		
2	AAA	1	20	18	2	0	0

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is water.

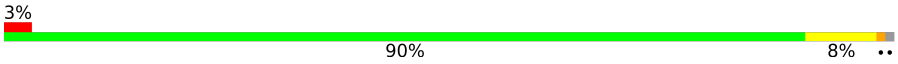
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	AAA	525	525	525	0	0

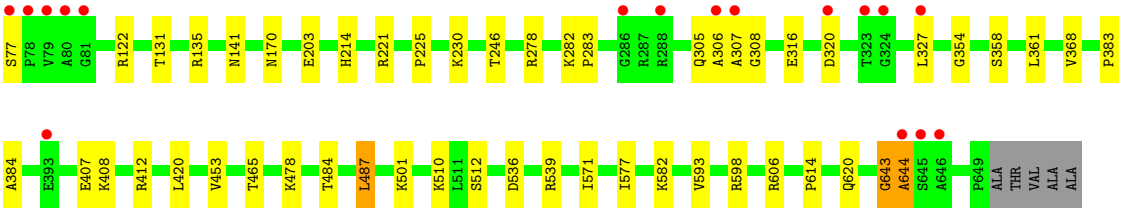
3 Residue-property plots (i)

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: Fatty acid Photodecarboxylase

Chain AAA:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	89.69Å 103.57Å 155.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.50 – 1.78 46.46 – 1.78	Depositor EDS
% Data completeness (in resolution range)	99.2 (46.50-1.78) 99.3 (46.46-1.78)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.50 (at 1.78Å)	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.172 , 0.202 0.181 , 0.213	Depositor DCC
R_{free} test set	3431 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å^2)	29.4	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 42.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	4885	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% 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, STE

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	AAA	0.73	0/4355	0.86	4/5910 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	AAA	221	ARG	NE-CZ-NH1	5.90	123.25	120.30
1	AAA	278	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	AAA	278	ARG	NE-CZ-NH1	5.07	122.83	120.30
1	AAA	606	ARG	NE-CZ-NH2	-5.02	117.79	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	AAA	4267	0	4205	33	0
2	AAA	40	0	70	5	0
3	AAA	53	0	31	3	0
4	AAA	525	0	0	9	4
All	All	4885	0	4306	33	4

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.

All (33) 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:AAA:643:GLY:HA2	1:AAA:644:ALA:CB	2.07	0.85
1:AAA:512:SER:OG	4:AAA:801:HOH:O	1.94	0.85
1:AAA:407:GLU:OE1	4:AAA:802:HOH:O	2.01	0.78
1:AAA:643:GLY:HA2	1:AAA:644:ALA:HB2	1.66	0.78
1:AAA:412:ARG:HD2	4:AAA:1190:HOH:O	1.88	0.72
1:AAA:536:ASP:OD1	1:AAA:539:ARG:NH2	2.26	0.69
1:AAA:484:THR:HG21	2:AAA:701:STE:H81	1.73	0.68
1:AAA:170:ASN:HB2	3:AAA:703:FAD:C5X	2.25	0.66
1:AAA:383:PRO:HB2	1:AAA:571:ILE:HD11	1.81	0.62
1:AAA:643:GLY:HA2	1:AAA:644:ALA:HB3	1.89	0.55
1:AAA:131:THR:OG1	2:AAA:702:STE:H183	2.07	0.54
1:AAA:420:LEU:HD21	2:AAA:702:STE:H62	1.92	0.51
1:AAA:246:THR:HG23	4:AAA:810:HOH:O	2.10	0.51
1:AAA:643:GLY:CA	1:AAA:644:ALA:CB	2.87	0.51
1:AAA:358:SER:HB2	1:AAA:368:VAL:HG21	1.94	0.50
1:AAA:305:GLN:NE2	4:AAA:819:HOH:O	2.44	0.50
1:AAA:320:ASP:OD2	4:AAA:803:HOH:O	2.18	0.50
1:AAA:453:VAL:HG11	2:AAA:701:STE:H121	1.96	0.47
1:AAA:203:GLU:OE1	1:AAA:214:HIS:HA	2.14	0.47
1:AAA:354:GLY:HA2	1:AAA:361:LEU:HD13	1.97	0.47
1:AAA:501:LYS:HG2	1:AAA:510:LYS:HB2	1.97	0.46
1:AAA:170:ASN:HB2	3:AAA:703:FAD:C4X	2.46	0.45
1:AAA:384:ALA:HA	1:AAA:487:LEU:O	2.15	0.45
1:AAA:230:LYS:HG3	4:AAA:1180:HOH:O	2.16	0.44
1:AAA:246:THR:HG22	4:AAA:1205:HOH:O	2.19	0.43
1:AAA:77:SER:OG	4:AAA:804:HOH:O	2.22	0.42
1:AAA:282:LYS:N	1:AAA:283:PRO:CD	2.83	0.42
1:AAA:620:GLN:HB3	3:AAA:703:FAD:C2	2.50	0.42
1:AAA:122:ARG:HG3	2:AAA:702:STE:H31	2.03	0.41
1:AAA:135:ARG:HA	1:AAA:141:ASN:OD1	2.20	0.41
1:AAA:582:LYS:HG3	1:AAA:614:PRO:HA	2.02	0.41
1:AAA:593:VAL:HA	1:AAA:598:ARG:O	2.21	0.41
1:AAA:316:GLU:HA	1:AAA:327:LEU:O	2.20	0.41

All (4) 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)
4:AAA:1099:HOH:O	4:AAA:1262:HOH:O[4_565]	1.97	0.23
4:AAA:1312:HOH:O	4:AAA:1317:HOH:O[7_555]	2.04	0.16
4:AAA:1141:HOH:O	4:AAA:1232:HOH:O[2_665]	2.18	0.02
4:AAA:1274:HOH:O	4:AAA:1274:HOH:O[2_665]	2.19	0.01

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	AAA	571/578 (99%)	547 (96%)	18 (3%)	6 (1%)	14 4

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	AAA	307	ALA
1	AAA	644	ALA
1	AAA	306	ALA
1	AAA	643	GLY
1	AAA	577	ILE
1	AAA	308	GLY

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	AAA	439/441 (100%)	434 (99%)	5 (1%)	73 65

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	225	PRO
1	AAA	408	LYS
1	AAA	465	THR
1	AAA	478	LYS
1	AAA	487	LEU

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

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

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	STE	AAA	701	-	19,19,19	0.76	1 (5%)	19,19,19	1.22	3 (15%)
3	FAD	AAA	703	-	53,58,58	0.78	1 (1%)	68,89,89	0.84	1 (1%)
2	STE	AAA	702	-	19,19,19	0.52	0	19,19,19	1.47	2 (10%)

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	STE	AAA	701	-	-	6/17/17/17	-
3	FAD	AAA	703	-	-	2/30/50/50	0/6/6/6
2	STE	AAA	702	-	-	6/17/17/17	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	AAA	703	FAD	O2'-C2'	2.43	1.48	1.43
2	AAA	701	STE	O2-C1	-2.21	1.23	1.30

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	AAA	701	STE	C11-C10-C9	-2.63	101.09	114.42
2	AAA	702	STE	C3-C2-C1	-2.56	108.01	114.47
2	AAA	701	STE	C6-C5-C4	-2.48	101.83	114.42
2	AAA	702	STE	C5-C4-C3	-2.31	102.68	114.42
2	AAA	701	STE	O2-C1-C2	2.20	121.10	114.03
3	AAA	703	FAD	C5A-C6A-N6A	2.04	123.45	120.35

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	AAA	701	STE	C6-C7-C8-C9
2	AAA	701	STE	C1-C2-C3-C4
2	AAA	702	STE	C6-C7-C8-C9
2	AAA	702	STE	C11-C10-C9-C8
2	AAA	702	STE	C1-C2-C3-C4
2	AAA	701	STE	C12-C13-C14-C15
3	AAA	703	FAD	PA-O3P-P-O5'
2	AAA	701	STE	C11-C10-C9-C8
3	AAA	703	FAD	O4B-C4B-C5B-O5B
2	AAA	701	STE	C14-C15-C16-C17
2	AAA	701	STE	C10-C11-C12-C13
2	AAA	702	STE	O1-C1-C2-C3
2	AAA	702	STE	C12-C13-C14-C15
2	AAA	702	STE	O2-C1-C2-C3

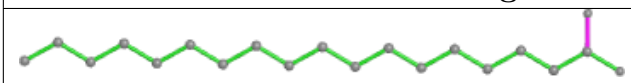
There are no ring outliers.

3 monomers are involved in 8 short contacts:

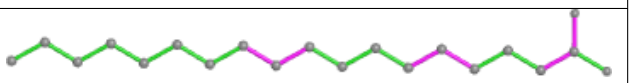
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	AAA	701	STE	2	0
3	AAA	703	FAD	3	0
2	AAA	702	STE	3	0

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 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.

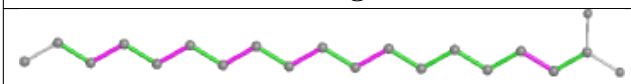
Ligand STE AAA 701



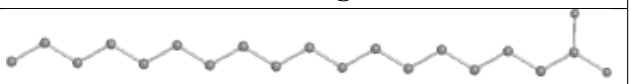
Bond lengths



Bond angles

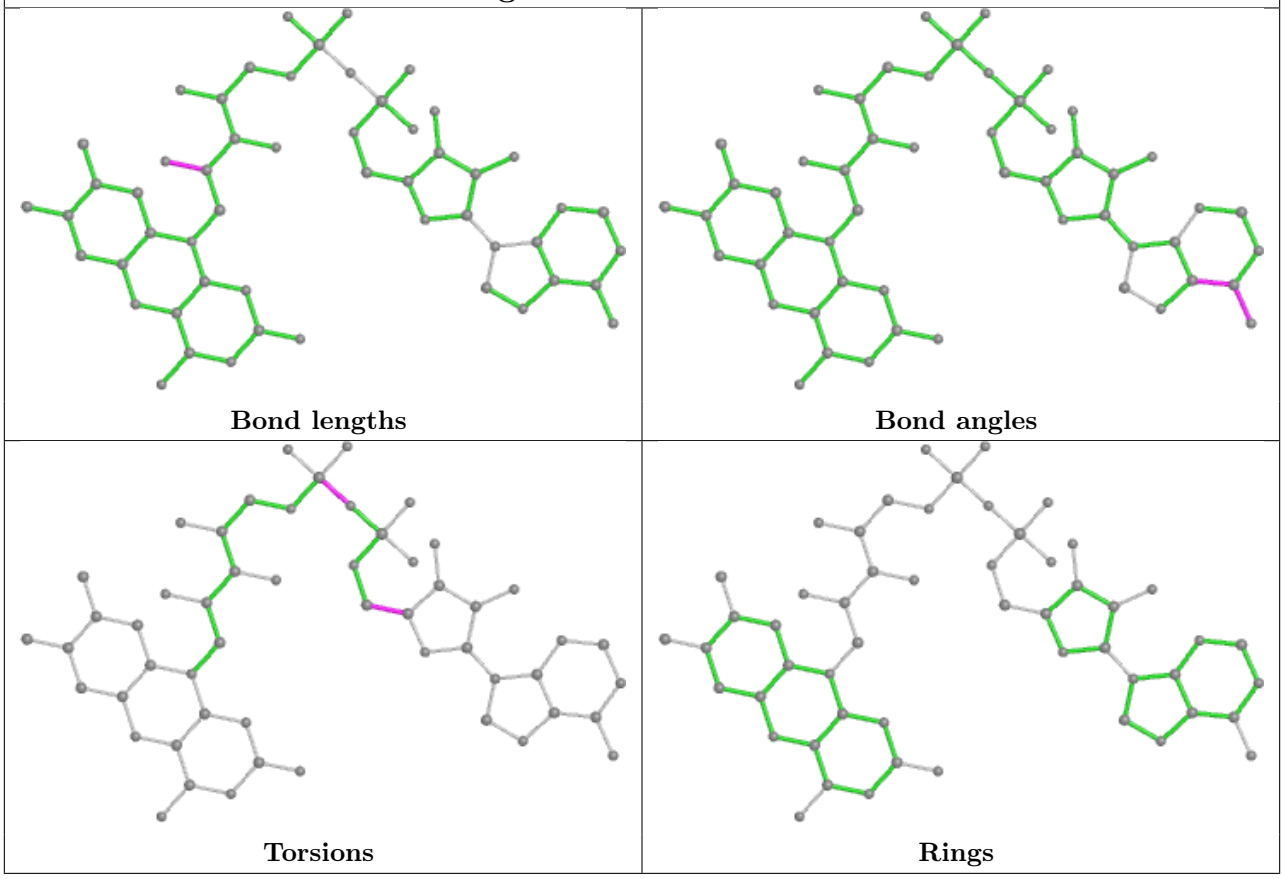


Torsions

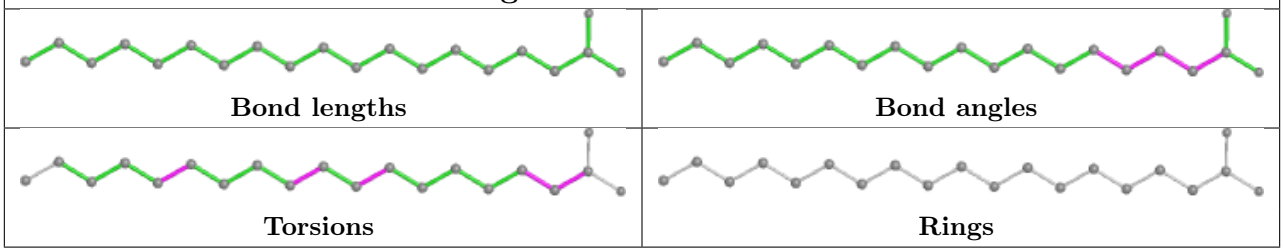


Rings

Ligand FAD AAA 703



Ligand STE AAA 702



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	AAA	573/578 (99%)	0.10	17 (2%) 50 48	21, 29, 51, 94	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	307	ALA	9.1
1	AAA	306	ALA	5.7
1	AAA	80	ALA	4.6
1	AAA	288	ARG	4.3
1	AAA	323	THR	3.5
1	AAA	79	VAL	3.4
1	AAA	81	GLY	3.0
1	AAA	644	ALA	2.9
1	AAA	646	ALA	2.9
1	AAA	286	GLY	2.8
1	AAA	78	PRO	2.7
1	AAA	393	GLU	2.6
1	AAA	77	SER	2.6
1	AAA	327	LEU	2.3
1	AAA	645	SER	2.2
1	AAA	320	ASP	2.1
1	AAA	324	GLY	2.1

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)

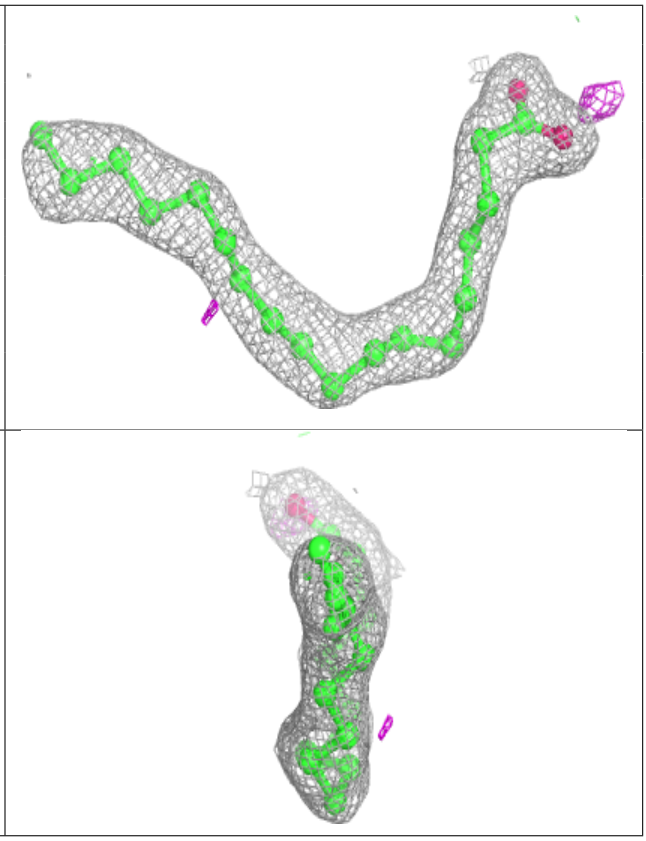
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	STE	AAA	702	20/20	0.90	0.13	37,45,52,52	0
2	STE	AAA	701	20/20	0.93	0.13	24,30,35,38	0
3	FAD	AAA	703	53/53	0.97	0.11	20,23,29,31	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.

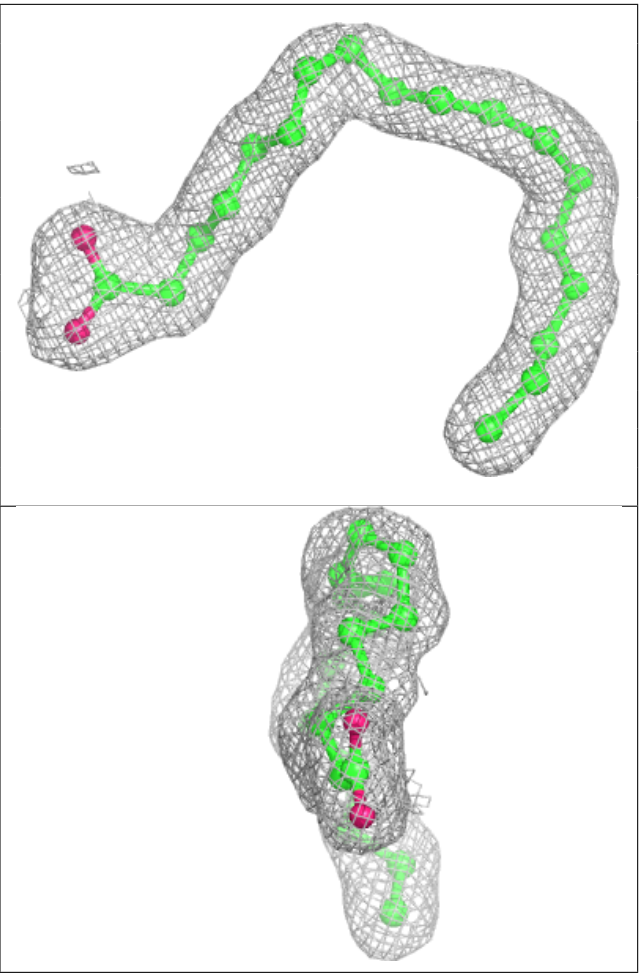
Electron density around STE AAA 702:

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



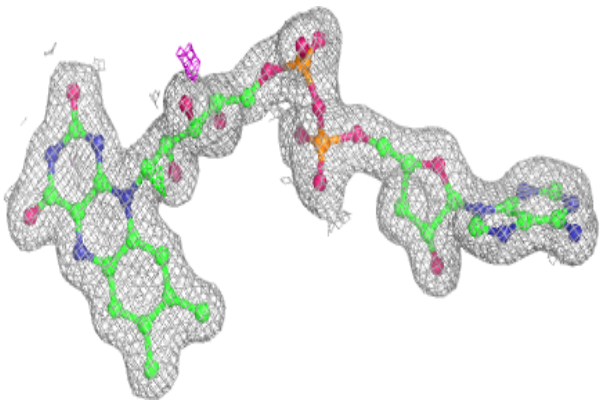
Electron density around STE AAA 701:

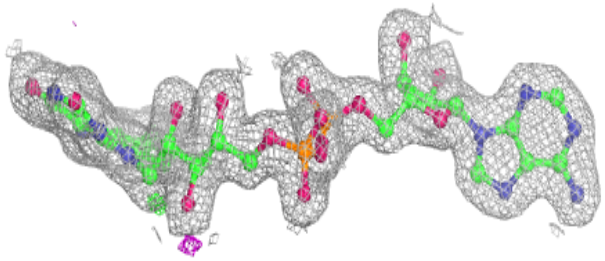
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

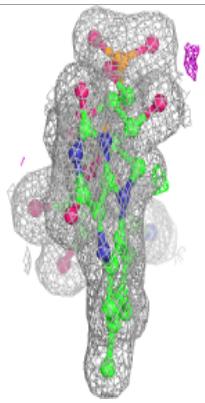


Electron density around FAD AAA 703:

$2mF_o$-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.