



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 18, 2026 – 03:30 pm BST

PDB ID : 6FAH / pdb_00006fah
Title : Molecular basis of the flavin-based electron-bifurcating caffeyl-CoA reductase reaction
Authors : Demmer, J.K.; Bertsch, J.; Oeppinger, C.; Wohlers, H.; Kayastha, K.; Demmer, U.; Ermler, U.; Mueller, V.
Deposited on : 2017-12-15
Resolution : 3.13 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

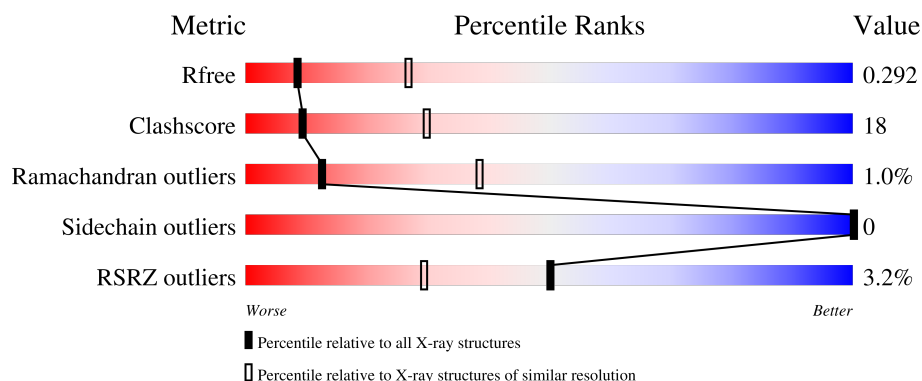
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.13 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}	180053	2351 (3.18-3.10)
Clashscore	190562	2452 (3.18-3.10)
Ramachandran outliers	187476	2324 (3.18-3.10)
Sidechain outliers	187428	2324 (3.18-3.10)
RSRZ outliers	180081	2351 (3.18-3.10)

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 $\leq 5\%$. 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	396	4% 54% 38% 5% ..
1	E	396	7% 58% 34% 6% ..
2	B	262	3% 65% 35%
2	F	262	5% 64% 34% .

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	C	379	 75%24%.
3	D	379	 74%26%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 15985 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 Caffeyl-CoA reductase-Etf complex subunit CarE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	389	Total	C	N	O	S	0	0	0
			2915	1839	479	575	22			
1	E	393	Total	C	N	O	S	0	0	0
			2948	1861	484	581	22			

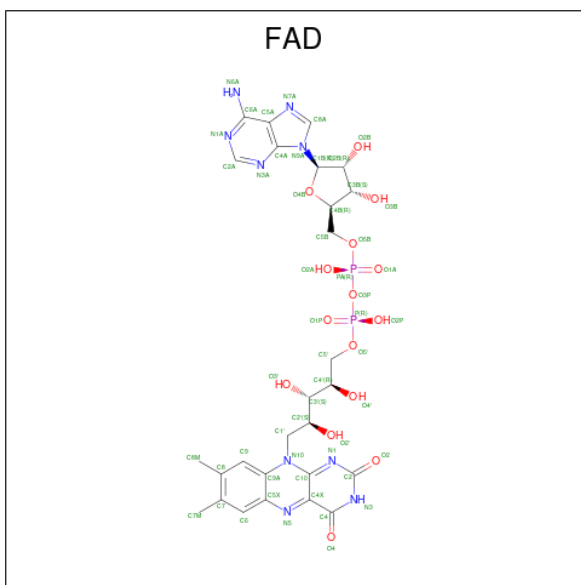
- Molecule 2 is a protein called Caffeyl-CoA reductase-Etf complex subunit CarD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	262	Total	C	N	O	S	0	0	0
			1970	1235	332	387	16			
2	F	262	Total	C	N	O	S	0	0	0
			1970	1235	332	387	16			

- Molecule 3 is a protein called Caffeyl-CoA reductase-Etf complex subunit CarC.

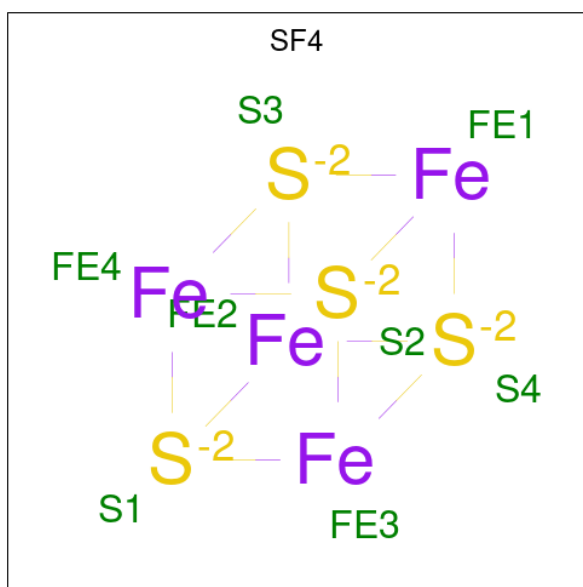
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	379	Total	C	N	O	S	0	1	0
			2903	1841	480	556	26			
3	D	379	Total	C	N	O	S	0	0	0
			2894	1833	480	555	26			

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



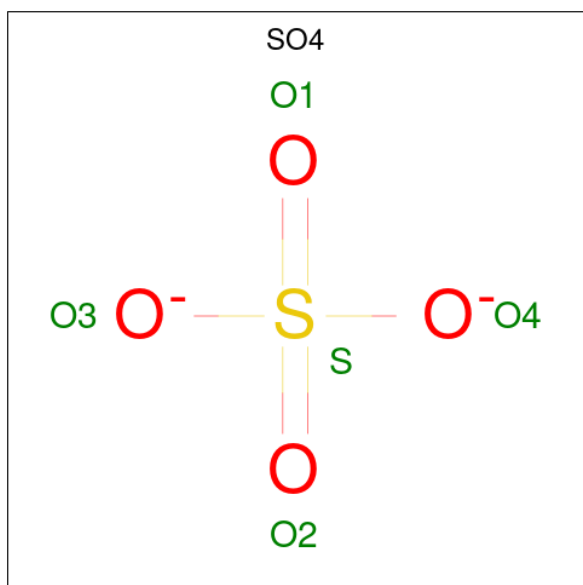
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 53	C 27	N 9	O 15	P 2	0	0
4	B	1	Total 53	C 27	N 9	O 15	P 2	0	0
4	C	1	Total 53	C 27	N 9	O 15	P 2	0	0
4	D	1	Total 53	C 27	N 9	O 15	P 2	0	0
4	E	1	Total 53	C 27	N 9	O 15	P 2	0	0
4	F	1	Total 53	C 27	N 9	O 15	P 2	0	0

- Molecule 5 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	Fe	S	0	0
			8	4	4		
5	A	1	Total	Fe	S	0	0
			8	4	4		
5	E	1	Total	Fe	S	0	0
			8	4	4		
5	E	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).

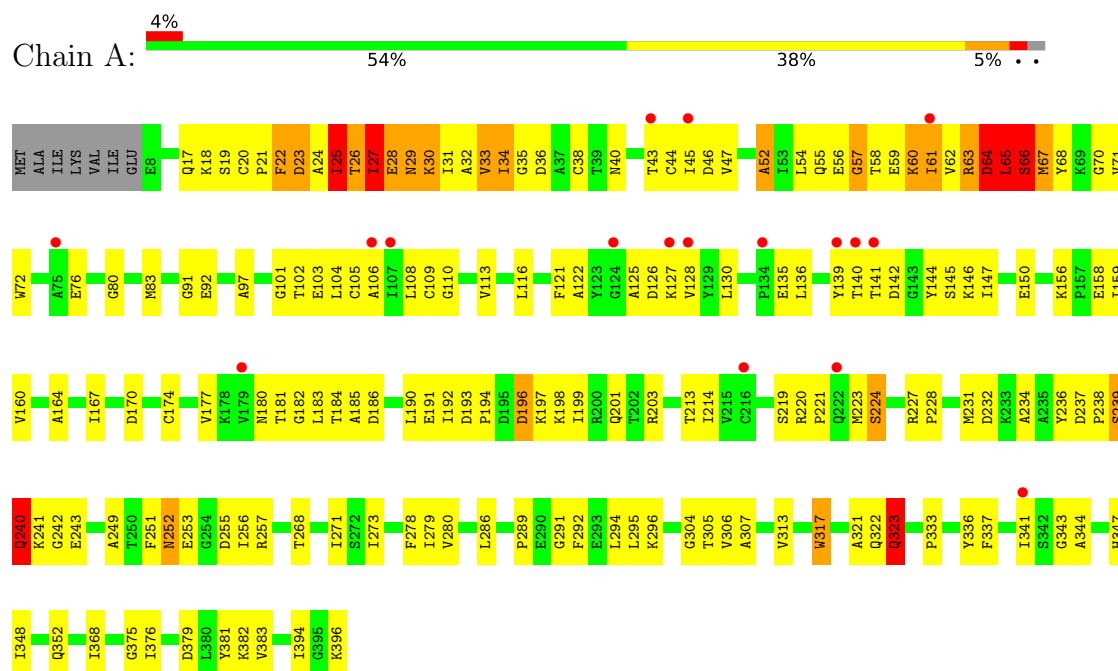


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	C	1	Total	O	S	0	0
			5	4	1		
6	C	1	Total	O	S	0	0
			5	4	1		
6	C	1	Total	O	S	0	0
			5	4	1		
6	C	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		

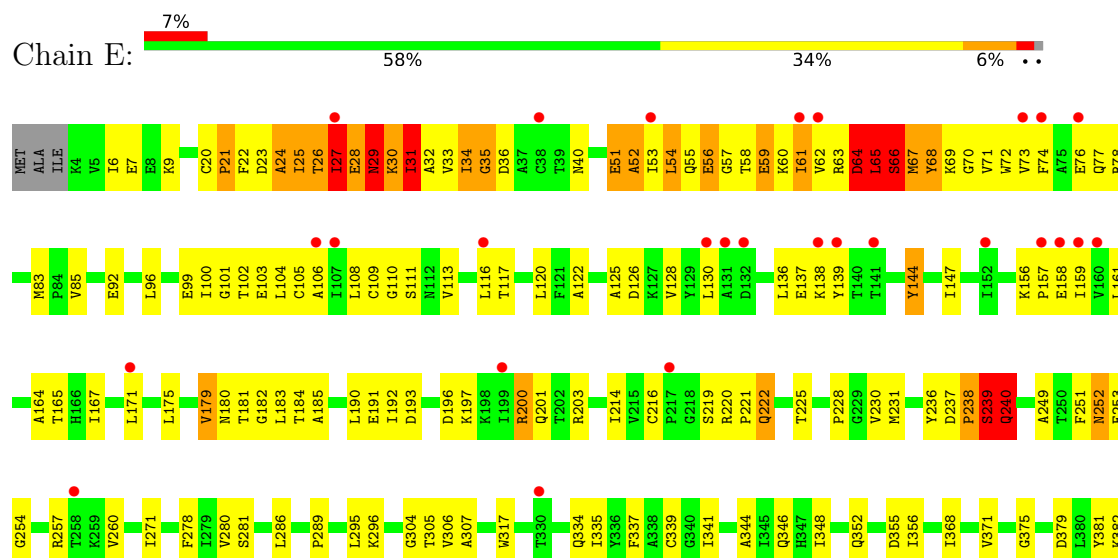
3 Residue-property plots

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: Caffeyl-CoA reductase-Etf complex subunit CarE

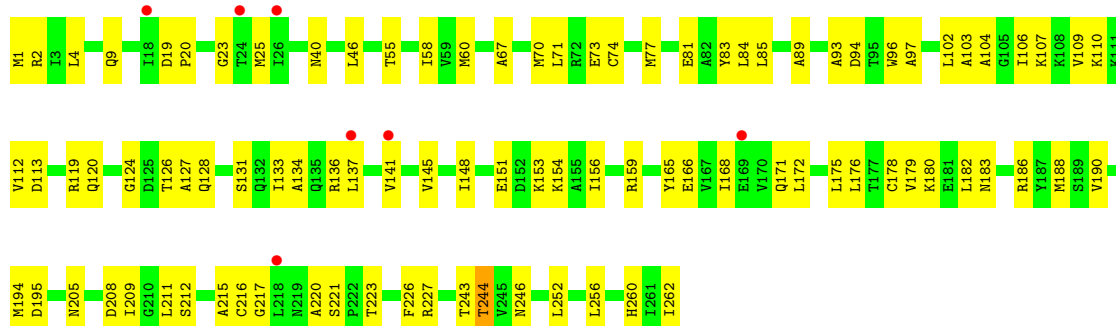


- Molecule 1: Caffeyl-CoA reductase-Etf complex subunit CarE

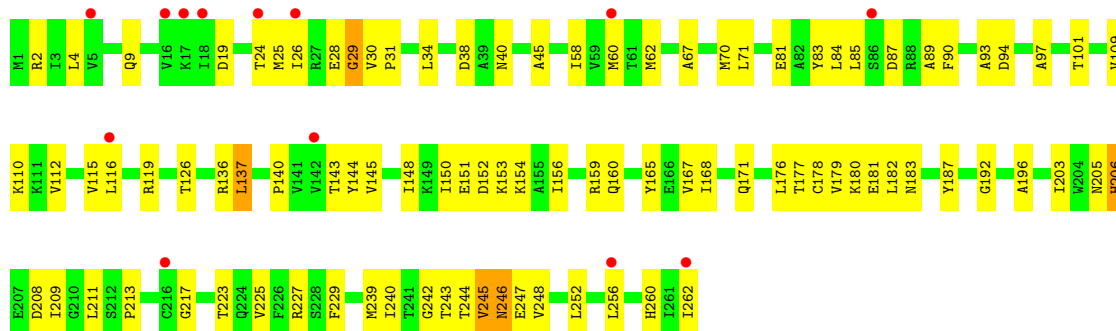




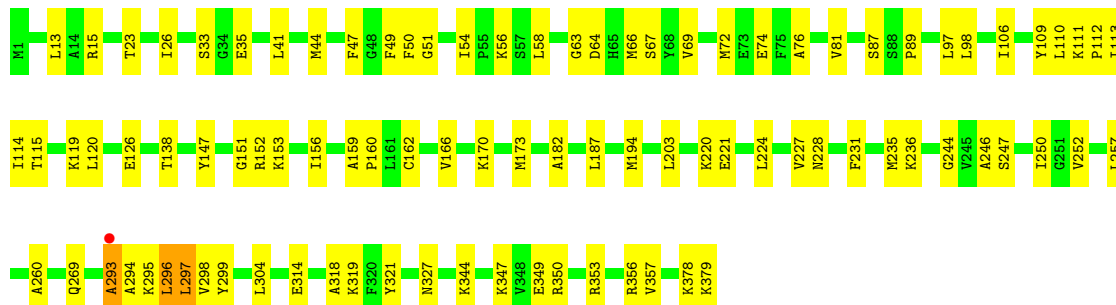
• Molecule 2: Caffeyl-CoA reductase-Etf complex subunit CarD



• Molecule 2: Caffeyl-CoA reductase-Etf complex subunit CarD

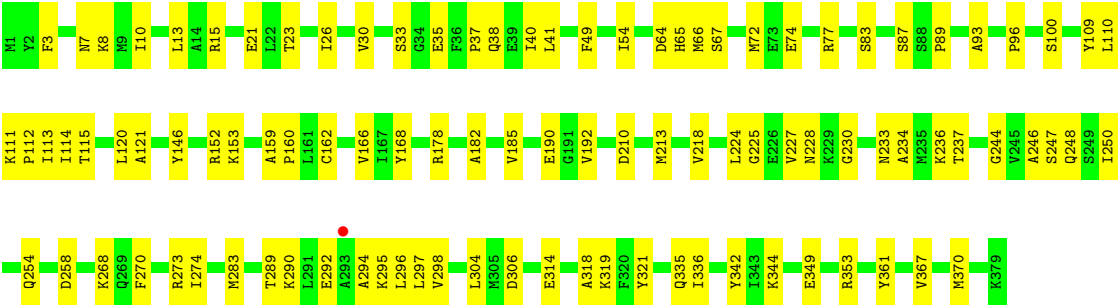


• Molecule 3: Caffeyl-CoA reductase-Etf complex subunit CarC



• Molecule 3: Caffeyl-CoA reductase-Etf complex subunit CarC





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	213.74Å 144.26Å 102.84Å 90.00° 98.93° 90.00°	Depositor
Resolution (Å)	49.52 – 3.13 49.52 – 3.13	Depositor EDS
% Data completeness (in resolution range)	75.2 (49.52-3.13) 75.2 (49.52-3.13)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 3.12Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.234 , 0.286 0.240 , 0.292	Depositor DCC
R_{free} test set	2057 reflections (3.79%)	wwPDB-VP
Wilson B-factor (Å ²)	85.3	Xtriage
Anisotropy	0.131	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 115.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	15985	wwPDB-VP
Average B, all atoms (Å ²)	112.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.35% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: SO4, SF4, 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	0.82	13/2958 (0.4%)	1.64	86/4004 (2.1%)
1	E	0.73	8/2991 (0.3%)	1.58	91/4048 (2.2%)
2	B	0.40	0/1999	0.72	4/2710 (0.1%)
2	F	0.59	4/1999 (0.2%)	1.24	13/2710 (0.5%)
3	C	0.58	0/2948	0.68	5/3947 (0.1%)
3	D	0.56	0/2935	0.70	1/3929 (0.0%)
All	All	0.64	25/15830 (0.2%)	1.19	200/21348 (0.9%)

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	A	0	5
1	E	0	6
All	All	0	11

The worst 5 of 25 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	246	ASN	CG-ND2	11.68	1.57	1.33
1	A	323	GLN	CD-OE1	-11.08	1.02	1.23
1	A	323	GLN	CD-NE2	8.14	1.50	1.33
1	E	35	GLY	CA-C	-7.10	1.41	1.51
1	A	240	GLN	CG-CD	6.97	1.69	1.52

The worst 5 of 200 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	246	ASN	OD1-CG-ND2	-30.02	92.58	122.60
2	F	246	ASN	CB-CG-OD1	29.59	179.98	120.80
1	A	240	GLN	OE1-CD-NE2	-21.01	101.59	122.60
2	F	246	ASN	CB-CG-ND2	-19.31	87.43	116.40
1	A	323	GLN	OE1-CD-NE2	-18.90	103.70	122.60

There are no chirality outliers.

5 of 11 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	240	GLN	Sidechain,Mainchain
1	A	26	THR	Mainchain
1	A	323	GLN	Sidechain
1	A	52	ALA	Mainchain
1	E	29	ASN	Sidechain

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	2915	0	2934	153	0
1	E	2948	0	2973	151	0
2	B	1970	0	2005	69	0
2	F	1970	0	2005	86	0
3	C	2903	0	2952	68	0
3	D	2894	0	2943	62	1
4	A	53	0	31	3	0
4	B	53	0	31	5	0
4	C	53	0	31	5	0
4	D	53	0	31	4	0
4	E	53	0	31	3	0
4	F	53	0	31	8	0
5	A	16	0	0	1	0
5	E	16	0	0	0	0
6	C	25	0	0	0	0
6	D	10	0	0	0	0
All	All	15985	0	15998	574	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 18.

The worst 5 of 574 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:E:65:LEU:HD21	1:E:68:TYR:OH	1.42	1.19
1:A:201:GLN:HE22	1:A:224:SER:HB3	1.05	1.16
1:A:26:THR:HG23	1:A:31:ILE:HA	1.31	1.12
1:E:77:GLN:NE2	1:E:109:CYS:O	1.81	1.10
1:E:99:GLU:HB3	1:E:192:ILE:HD11	1.31	1.07

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)
3:D:321:TYR:OH	3:D:321:TYR:OH[2_555]	2.00	0.20

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	387/396 (98%)	332 (86%)	45 (12%)	10 (3%)	4	18
1	E	391/396 (99%)	333 (85%)	50 (13%)	8 (2%)	6	23
2	B	260/262 (99%)	247 (95%)	13 (5%)	0	100	100
2	F	260/262 (99%)	245 (94%)	14 (5%)	1 (0%)	30	59
3	C	378/379 (100%)	359 (95%)	19 (5%)	0	100	100
3	D	377/379 (100%)	358 (95%)	18 (5%)	1 (0%)	36	64
All	All	2053/2074 (99%)	1874 (91%)	159 (8%)	20 (1%)	12	38

5 of 20 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	25	ILE
1	A	27	ILE
1	A	52	ALA
1	A	64	ASP
1	A	65	LEU

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	315/321 (98%)	315 (100%)	0	100	100
1	E	319/321 (99%)	319 (100%)	0	100	100
2	B	216/216 (100%)	216 (100%)	0	100	100
2	F	216/216 (100%)	216 (100%)	0	100	100
3	C	307/306 (100%)	307 (100%)	0	100	100
3	D	306/306 (100%)	306 (100%)	0	100	100
All	All	1679/1686 (100%)	1679 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 15 such sidechains are listed below:

Mol	Chain	Res	Type
3	D	368	GLN
2	F	206	HIS
1	E	180	ASN
2	F	246	ASN
1	E	240	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

17 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	$\# Z > 2$
6	SO4	C	406	-	4,4,4	0.25	0	6,6,6	0.09	0
5	SF4	E	402	1	0,12,12	-	-	-		
6	SO4	C	403	-	4,4,4	0.31	0	6,6,6	0.18	0
6	SO4	D	402	-	4,4,4	0.21	0	6,6,6	0.21	0
4	FAD	B	301	-	56,58,58	0.57	1 (1%)	81,89,89	0.43	1 (1%)
4	FAD	F	301	-	56,58,58	0.54	1 (1%)	81,89,89	0.46	1 (1%)
5	SF4	A	403	1	0,12,12	-	-	-		
4	FAD	A	401	-	56,58,58	0.56	1 (1%)	81,89,89	0.55	2 (2%)
5	SF4	E	403	1	0,12,12	-	-	-		
4	FAD	D	401	-	56,58,58	0.74	2 (3%)	81,89,89	0.58	1 (1%)
6	SO4	D	403	-	4,4,4	0.19	0	6,6,6	0.11	0
4	FAD	C	401	-	56,58,58	0.84	2 (3%)	81,89,89	0.50	1 (1%)
4	FAD	E	401	-	56,58,58	0.52	1 (1%)	81,89,89	0.45	0
5	SF4	A	402	1	0,12,12	-	-	-		
6	SO4	C	404	-	4,4,4	0.20	0	6,6,6	0.27	0
6	SO4	C	405	-	4,4,4	0.23	0	6,6,6	0.08	0
6	SO4	C	402	-	4,4,4	0.20	0	6,6,6	0.09	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
5	SF4	E	402	1	-	-	0/6/5/5
4	FAD	B	301	-	-	8/34/50/50	0/6/6/6
4	FAD	F	301	-	-	9/34/50/50	0/6/6/6
5	SF4	A	403	1	-	-	0/6/5/5
4	FAD	A	401	-	-	9/34/50/50	0/6/6/6
5	SF4	E	403	1	-	-	0/6/5/5
4	FAD	D	401	-	-	10/34/50/50	0/6/6/6
4	FAD	C	401	-	-	8/34/50/50	0/6/6/6
4	FAD	E	401	-	-	7/34/50/50	0/6/6/6
5	SF4	A	402	1	-	-	0/6/5/5

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	401	FAD	P-O2P	-3.32	1.39	1.55
4	C	401	FAD	PA-O5B	-3.10	1.46	1.59
4	D	401	FAD	PA-O5B	-2.92	1.47	1.59
4	D	401	FAD	P-O2P	-2.89	1.41	1.55
4	B	301	FAD	P-O2P	-2.59	1.43	1.55

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	401	FAD	P-O3P-PA	-3.84	119.63	132.83
4	C	401	FAD	P-O3P-PA	-2.57	123.99	132.83
4	B	301	FAD	P-O3P-PA	-2.43	124.48	132.83
4	F	301	FAD	P-O3P-PA	-2.39	124.62	132.83
4	A	401	FAD	P-O3P-PA	-2.12	125.55	132.83

There are no chirality outliers.

5 of 51 torsion outliers are listed below:

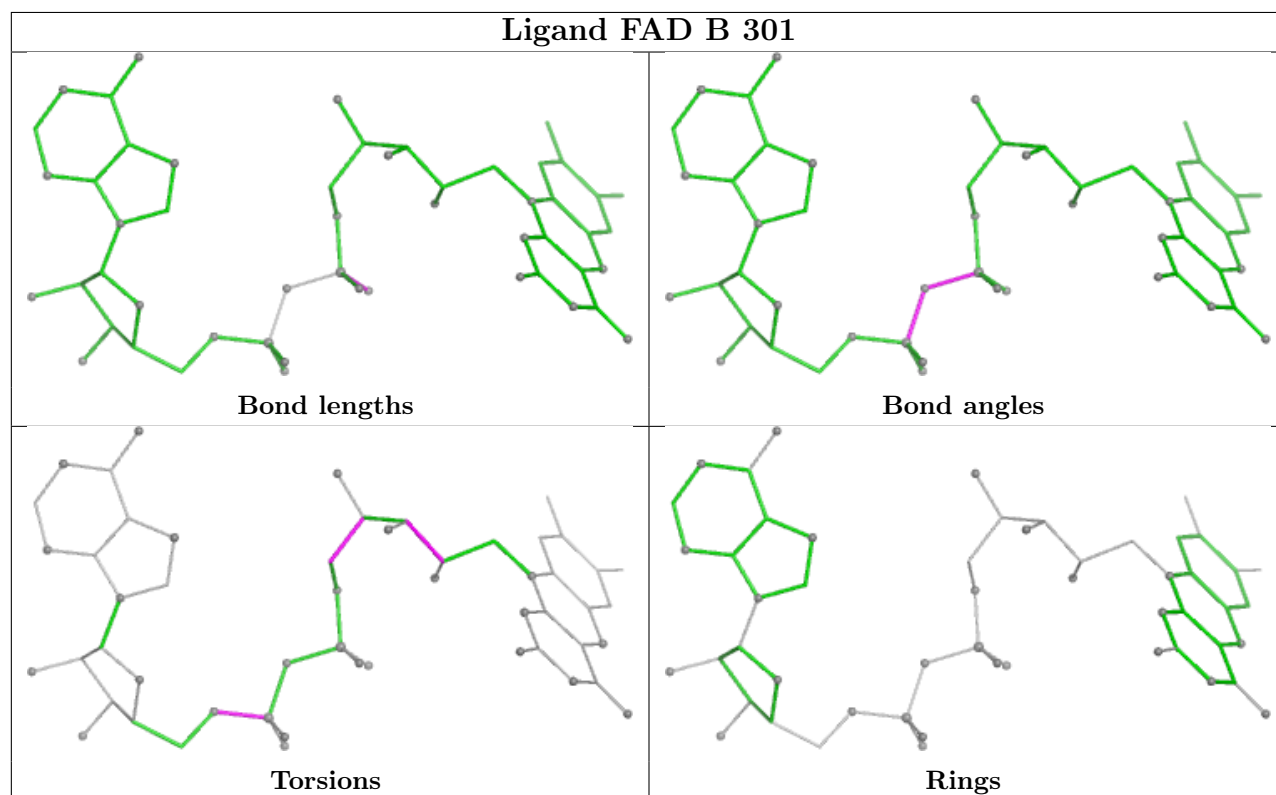
Mol	Chain	Res	Type	Atoms
4	A	401	FAD	C5B-O5B-PA-O1A
4	A	401	FAD	C5B-O5B-PA-O2A
4	B	301	FAD	C5B-O5B-PA-O1A
4	B	301	FAD	C5B-O5B-PA-O3P
4	B	301	FAD	C1'-C2'-C3'-O3'

There are no ring outliers.

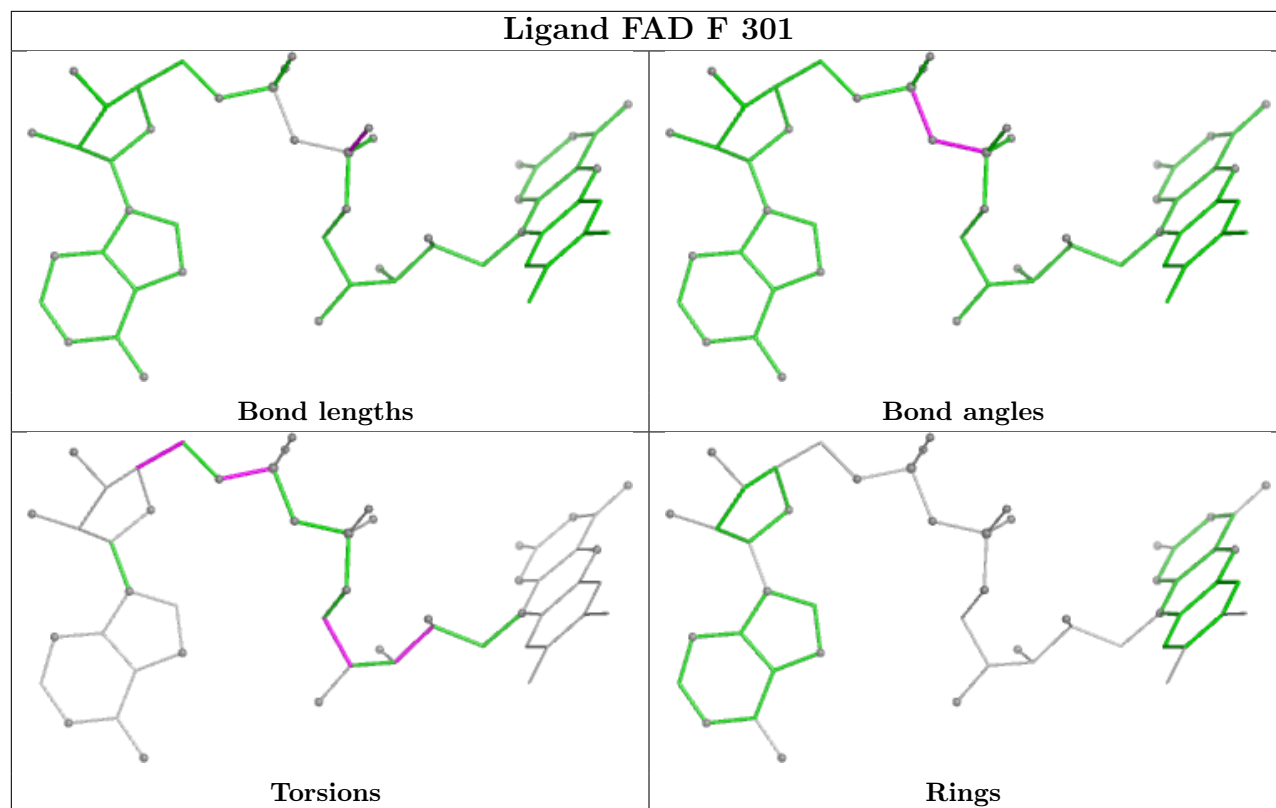
7 monomers are involved in 29 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	301	FAD	5	0
4	F	301	FAD	8	0
5	A	403	SF4	1	0
4	A	401	FAD	3	0
4	D	401	FAD	4	0
4	C	401	FAD	5	0
4	E	401	FAD	3	0

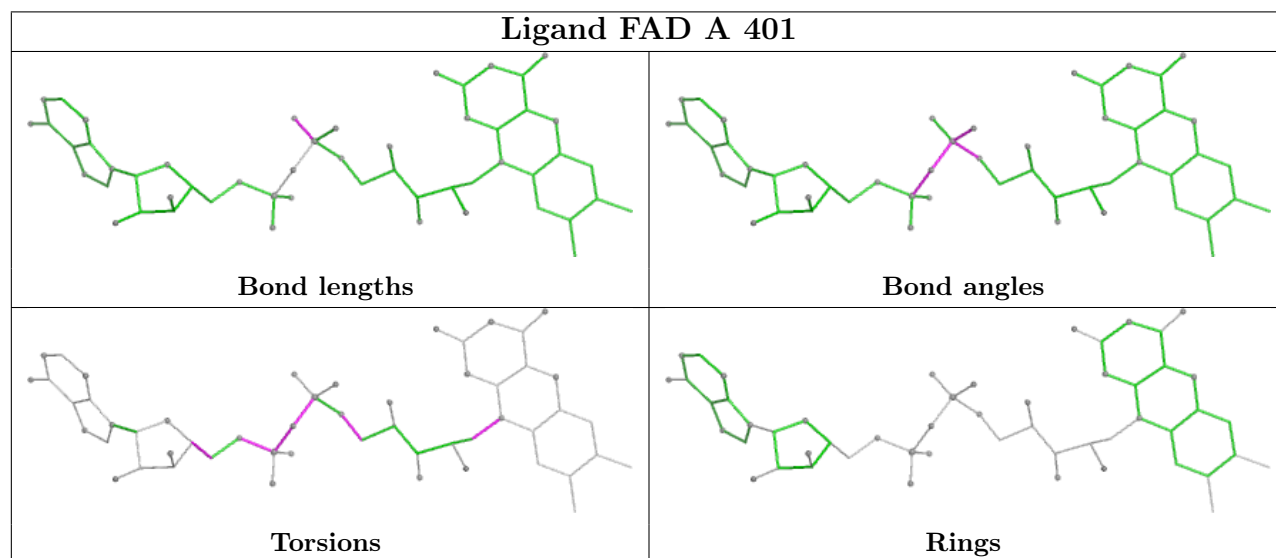
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

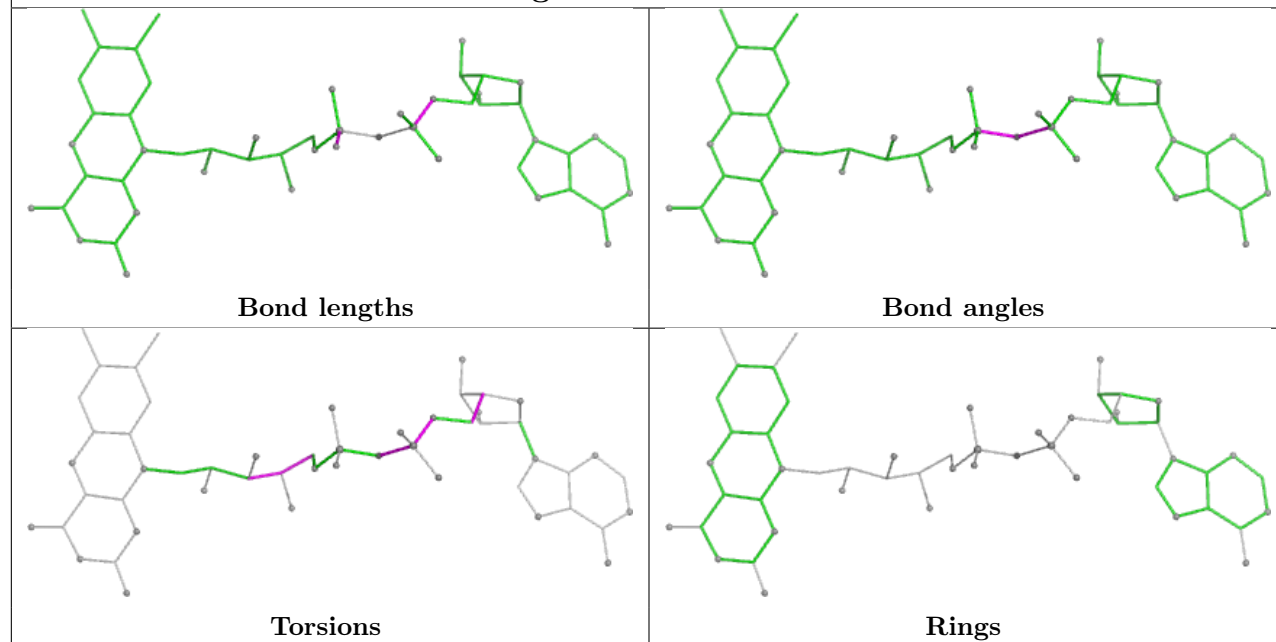
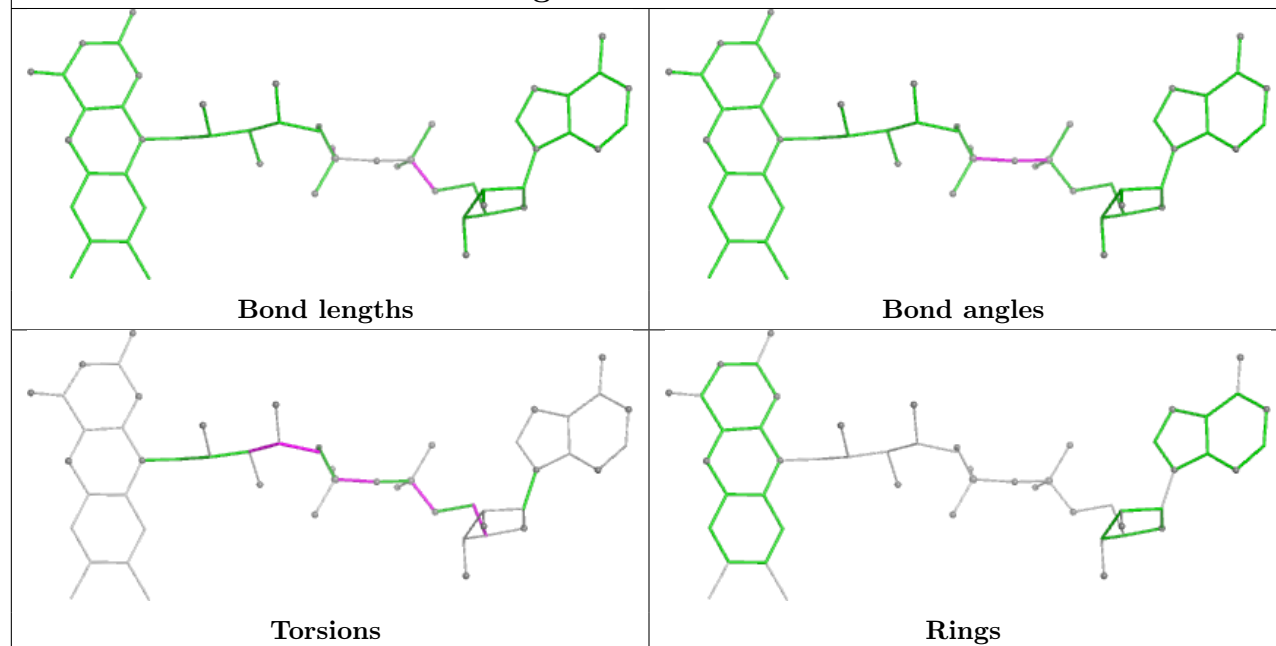


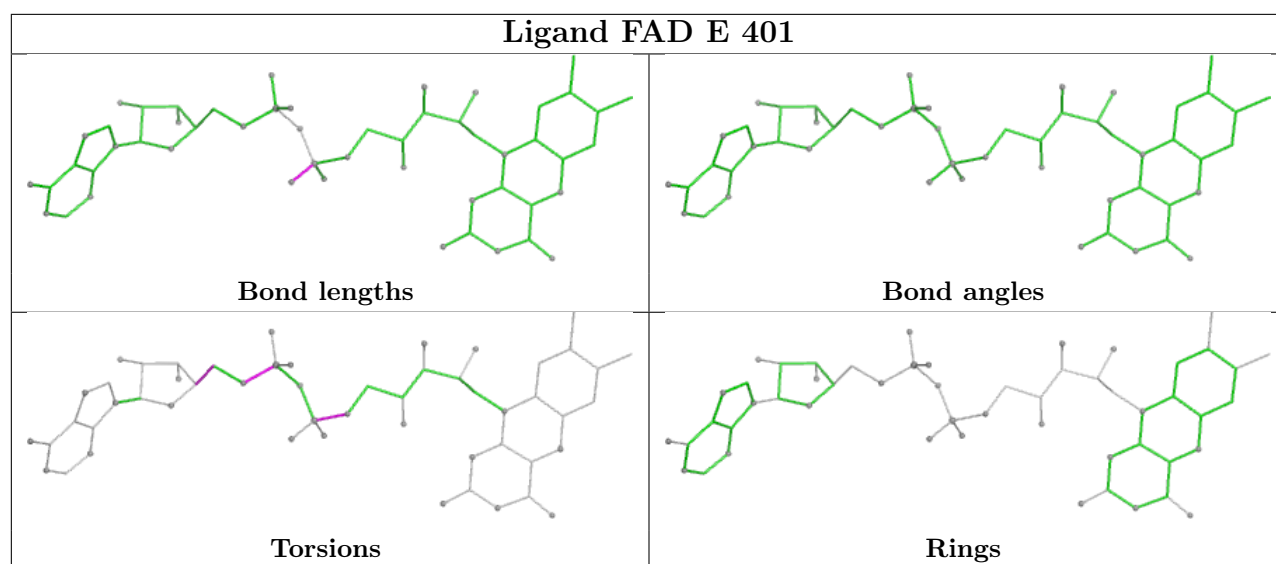
Ligand FAD F 301



Ligand FAD A 401



Ligand FAD D 401**Ligand FAD C 401**



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	389/396 (98%)	0.62	17 (4%)	39 20	33, 133, 234, 270	0
1	E	393/396 (99%)	0.79	27 (6%)	23 11	52, 156, 253, 314	0
2	B	262/262 (100%)	0.44	7 (2%)	56 34	64, 118, 177, 230	0
2	F	262/262 (100%)	0.65	13 (4%)	34 17	83, 134, 213, 260	0
3	C	379/379 (100%)	-0.10	1 (0%)	90 80	26, 62, 103, 128	1 (0%)
3	D	379/379 (100%)	-0.14	1 (0%)	90 80	22, 67, 104, 148	0
All	All	2064/2074 (99%)	0.36	66 (3%)	50 29	22, 101, 218, 314	1 (0%)

The worst 5 of 66 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	293	ALA	6.2
1	E	106	ALA	5.5
3	C	293	ALA	5.2
1	E	130	LEU	5.1
1	E	107	ILE	4.4

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 oligosaccharides in this entry.

6.4 Ligands ⓘ

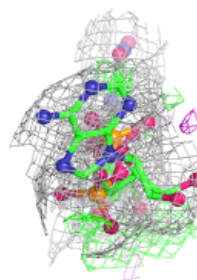
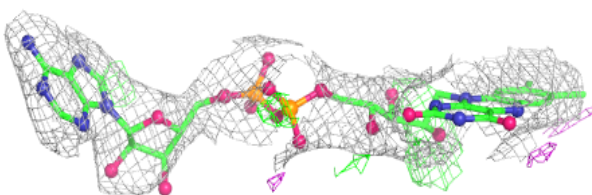
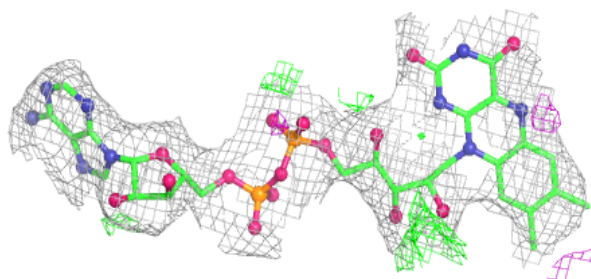
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	SF4	E	402	8/8	0.73	0.10	347,351,352,352	0
6	SO4	D	403	5/5	0.73	0.32	304,305,306,311	0
6	SO4	C	402	5/5	0.78	0.34	134,140,146,147	0
5	SF4	A	402	8/8	0.81	0.08	257,267,273,273	0
6	SO4	C	404	5/5	0.82	0.21	100,117,124,128	0
6	SO4	C	405	5/5	0.86	0.24	139,143,145,146	0
4	FAD	E	401	53/53	0.86	0.11	41,79,98,111	0
6	SO4	D	402	5/5	0.87	0.27	168,170,172,173	0
6	SO4	C	406	5/5	0.89	0.22	129,131,136,138	0
4	FAD	A	401	53/53	0.91	0.11	32,68,85,104	0
5	SF4	E	403	8/8	0.91	0.07	238,241,245,249	0
5	SF4	A	403	8/8	0.92	0.06	237,240,244,247	0
4	FAD	F	301	53/53	0.92	0.10	67,109,135,144	0
4	FAD	B	301	53/53	0.92	0.09	44,95,122,134	0
6	SO4	C	403	5/5	0.93	0.14	90,95,105,108	0
4	FAD	D	401	53/53	0.94	0.10	10,41,67,79	0
4	FAD	C	401	53/53	0.95	0.09	13,39,58,61	0

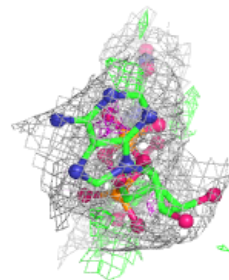
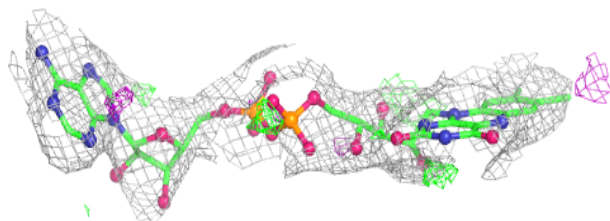
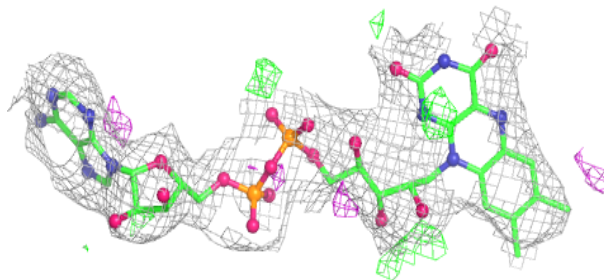
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around FAD E 401:

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

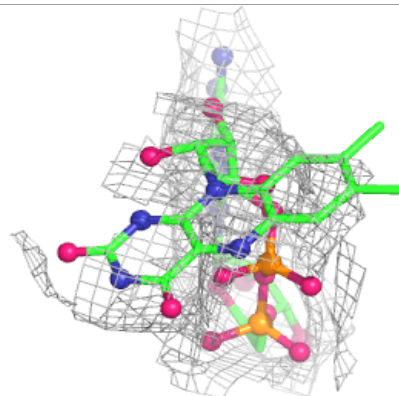
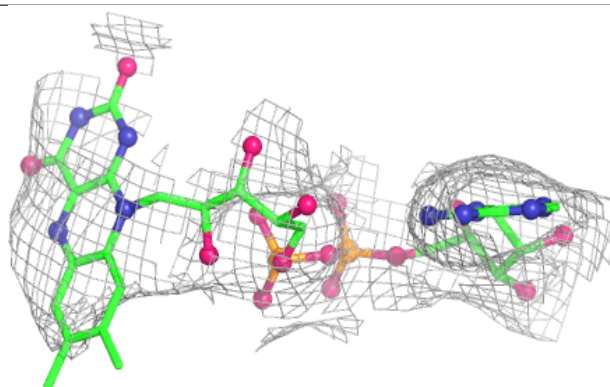
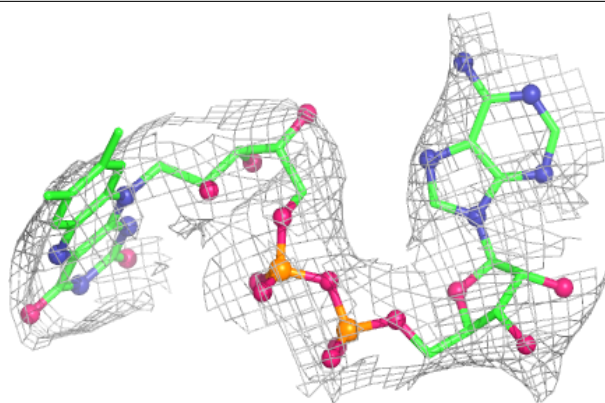
**Electron density around FAD A 401:**

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

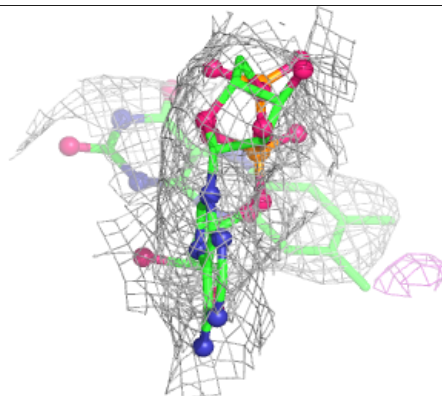
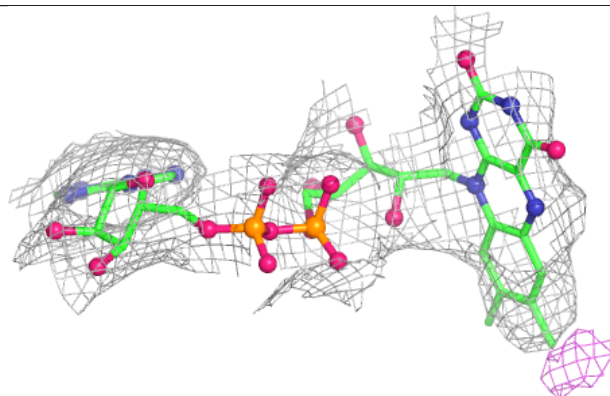
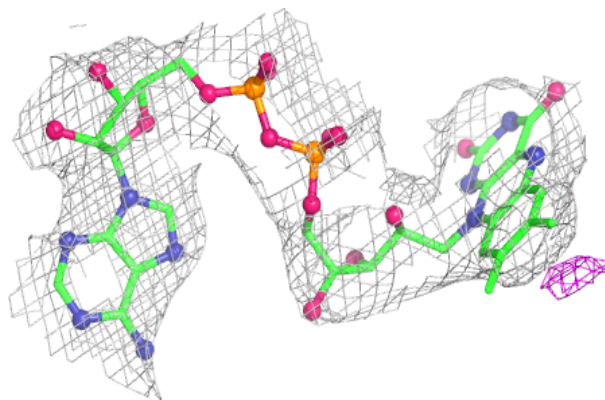


Electron density around FAD F 301:

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

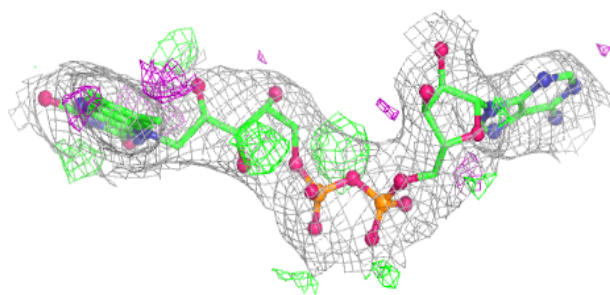
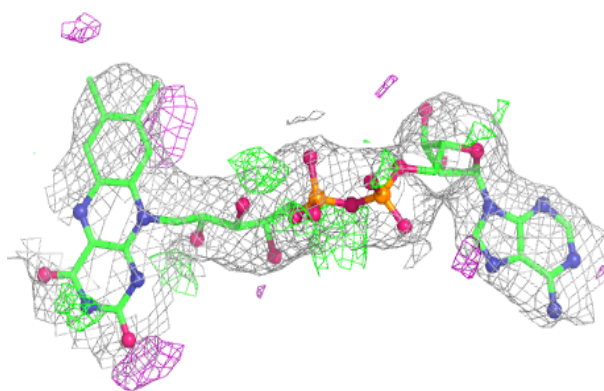
**Electron density around FAD B 301:**

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

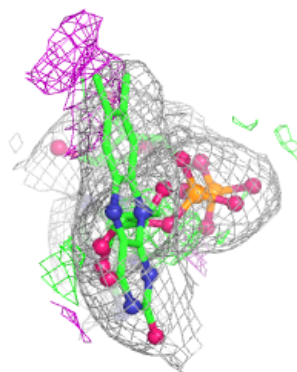
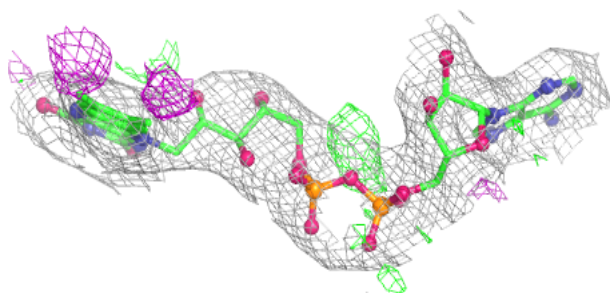
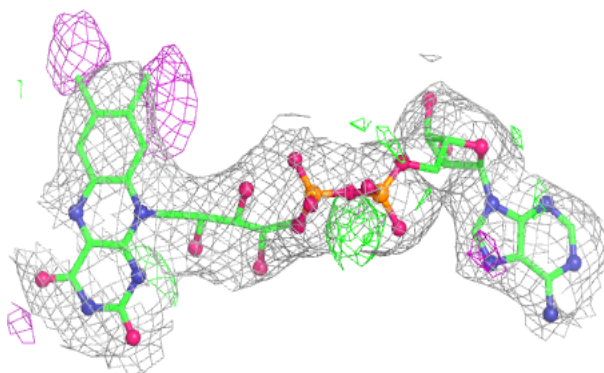


Electron density around FAD D 401:

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

**Electron density around FAD C 401:**

$2mF_o-DF_c$ (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.