wwPDB X-ray Structure Validation Summary Report (i)

Nov 22, 2023 – 10:27 PM JST

PDB ID	:	7XQA
Title	:	Orf1-glycine complex
Authors	:	Wang, Y.L.; Li, T.L.
Deposited on	:	2022-05-07
Resolution	:	1.93 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:

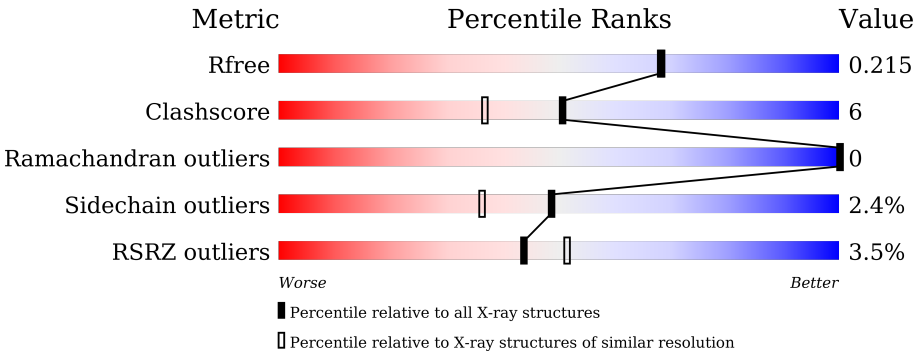
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (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 1.93 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	4310 (1.96-1.92)
Clashscore	141614	1023 (1.94-1.94)
Ramachandran outliers	138981	1007 (1.94-1.94)
Sidechain outliers	138945	1007 (1.94-1.94)
RSRZ outliers	127900	4250 (1.96-1.92)

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	512	85% 8% • 6%
1	B	512	87% 7% • 6%
1	C	512	86% 7% 6%
1	D	512	5% 84% 9% 6%
1	E	512	86% 7% • 6%
1	F	512	86% 7% • 6%

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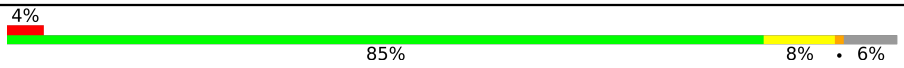
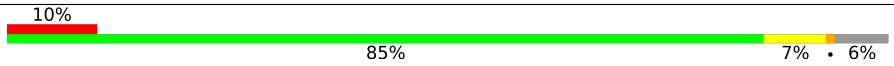
Mol	Chain	Length	Quality of chain
1	G	512	
2	H	512	

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 32835 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 N-formimidoyl fortimicin A synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	482	Total	C	N	O	S	0	1	0
			3662	2298	661	690	13			
1	B	482	Total	C	N	O	S	0	2	0
			3673	2304	665	691	13			
1	C	482	Total	C	N	O	S	0	0	0
			3654	2293	660	689	12			
1	D	479	Total	C	N	O	S	0	0	0
			3630	2280	656	682	12			
1	E	482	Total	C	N	O	S	0	1	0
			3662	2298	661	690	13			
1	F	482	Total	C	N	O	S	0	1	0
			3662	2298	661	690	13			
1	G	482	Total	C	N	O	S	0	0	0
			3654	2293	660	689	12			

There are 147 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	initiating methionine	UNP A0A125SZC1
A	-19	GLY	-	expression tag	UNP A0A125SZC1
A	-18	SER	-	expression tag	UNP A0A125SZC1
A	-17	SER	-	expression tag	UNP A0A125SZC1
A	-16	HIS	-	expression tag	UNP A0A125SZC1
A	-15	HIS	-	expression tag	UNP A0A125SZC1
A	-14	HIS	-	expression tag	UNP A0A125SZC1
A	-13	HIS	-	expression tag	UNP A0A125SZC1
A	-12	HIS	-	expression tag	UNP A0A125SZC1
A	-11	HIS	-	expression tag	UNP A0A125SZC1
A	-10	SER	-	expression tag	UNP A0A125SZC1
A	-9	SER	-	expression tag	UNP A0A125SZC1
A	-8	GLY	-	expression tag	UNP A0A125SZC1
A	-7	LEU	-	expression tag	UNP A0A125SZC1
A	-6	VAL	-	expression tag	UNP A0A125SZC1

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	PRO	-	expression tag	UNP A0A125SZC1
A	-4	ARG	-	expression tag	UNP A0A125SZC1
A	-3	GLY	-	expression tag	UNP A0A125SZC1
A	-2	SER	-	expression tag	UNP A0A125SZC1
A	-1	HIS	-	expression tag	UNP A0A125SZC1
A	0	MET	-	expression tag	UNP A0A125SZC1
B	-20	MET	-	initiating methionine	UNP A0A125SZC1
B	-19	GLY	-	expression tag	UNP A0A125SZC1
B	-18	SER	-	expression tag	UNP A0A125SZC1
B	-17	SER	-	expression tag	UNP A0A125SZC1
B	-16	HIS	-	expression tag	UNP A0A125SZC1
B	-15	HIS	-	expression tag	UNP A0A125SZC1
B	-14	HIS	-	expression tag	UNP A0A125SZC1
B	-13	HIS	-	expression tag	UNP A0A125SZC1
B	-12	HIS	-	expression tag	UNP A0A125SZC1
B	-11	HIS	-	expression tag	UNP A0A125SZC1
B	-10	SER	-	expression tag	UNP A0A125SZC1
B	-9	SER	-	expression tag	UNP A0A125SZC1
B	-8	GLY	-	expression tag	UNP A0A125SZC1
B	-7	LEU	-	expression tag	UNP A0A125SZC1
B	-6	VAL	-	expression tag	UNP A0A125SZC1
B	-5	PRO	-	expression tag	UNP A0A125SZC1
B	-4	ARG	-	expression tag	UNP A0A125SZC1
B	-3	GLY	-	expression tag	UNP A0A125SZC1
B	-2	SER	-	expression tag	UNP A0A125SZC1
B	-1	HIS	-	expression tag	UNP A0A125SZC1
B	0	MET	-	expression tag	UNP A0A125SZC1
C	-20	MET	-	initiating methionine	UNP A0A125SZC1
C	-19	GLY	-	expression tag	UNP A0A125SZC1
C	-18	SER	-	expression tag	UNP A0A125SZC1
C	-17	SER	-	expression tag	UNP A0A125SZC1
C	-16	HIS	-	expression tag	UNP A0A125SZC1
C	-15	HIS	-	expression tag	UNP A0A125SZC1
C	-14	HIS	-	expression tag	UNP A0A125SZC1
C	-13	HIS	-	expression tag	UNP A0A125SZC1
C	-12	HIS	-	expression tag	UNP A0A125SZC1
C	-11	HIS	-	expression tag	UNP A0A125SZC1
C	-10	SER	-	expression tag	UNP A0A125SZC1
C	-9	SER	-	expression tag	UNP A0A125SZC1
C	-8	GLY	-	expression tag	UNP A0A125SZC1
C	-7	LEU	-	expression tag	UNP A0A125SZC1
C	-6	VAL	-	expression tag	UNP A0A125SZC1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-5	PRO	-	expression tag	UNP A0A125SZC1
C	-4	ARG	-	expression tag	UNP A0A125SZC1
C	-3	GLY	-	expression tag	UNP A0A125SZC1
C	-2	SER	-	expression tag	UNP A0A125SZC1
C	-1	HIS	-	expression tag	UNP A0A125SZC1
C	0	MET	-	expression tag	UNP A0A125SZC1
D	-20	MET	-	initiating methionine	UNP A0A125SZC1
D	-19	GLY	-	expression tag	UNP A0A125SZC1
D	-18	SER	-	expression tag	UNP A0A125SZC1
D	-17	SER	-	expression tag	UNP A0A125SZC1
D	-16	HIS	-	expression tag	UNP A0A125SZC1
D	-15	HIS	-	expression tag	UNP A0A125SZC1
D	-14	HIS	-	expression tag	UNP A0A125SZC1
D	-13	HIS	-	expression tag	UNP A0A125SZC1
D	-12	HIS	-	expression tag	UNP A0A125SZC1
D	-11	HIS	-	expression tag	UNP A0A125SZC1
D	-10	SER	-	expression tag	UNP A0A125SZC1
D	-9	SER	-	expression tag	UNP A0A125SZC1
D	-8	GLY	-	expression tag	UNP A0A125SZC1
D	-7	LEU	-	expression tag	UNP A0A125SZC1
D	-6	VAL	-	expression tag	UNP A0A125SZC1
D	-5	PRO	-	expression tag	UNP A0A125SZC1
D	-4	ARG	-	expression tag	UNP A0A125SZC1
D	-3	GLY	-	expression tag	UNP A0A125SZC1
D	-2	SER	-	expression tag	UNP A0A125SZC1
D	-1	HIS	-	expression tag	UNP A0A125SZC1
D	0	MET	-	expression tag	UNP A0A125SZC1
E	-20	MET	-	initiating methionine	UNP A0A125SZC1
E	-19	GLY	-	expression tag	UNP A0A125SZC1
E	-18	SER	-	expression tag	UNP A0A125SZC1
E	-17	SER	-	expression tag	UNP A0A125SZC1
E	-16	HIS	-	expression tag	UNP A0A125SZC1
E	-15	HIS	-	expression tag	UNP A0A125SZC1
E	-14	HIS	-	expression tag	UNP A0A125SZC1
E	-13	HIS	-	expression tag	UNP A0A125SZC1
E	-12	HIS	-	expression tag	UNP A0A125SZC1
E	-11	HIS	-	expression tag	UNP A0A125SZC1
E	-10	SER	-	expression tag	UNP A0A125SZC1
E	-9	SER	-	expression tag	UNP A0A125SZC1
E	-8	GLY	-	expression tag	UNP A0A125SZC1
E	-7	LEU	-	expression tag	UNP A0A125SZC1
E	-6	VAL	-	expression tag	UNP A0A125SZC1

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-5	PRO	-	expression tag	UNP A0A125SZC1
E	-4	ARG	-	expression tag	UNP A0A125SZC1
E	-3	GLY	-	expression tag	UNP A0A125SZC1
E	-2	SER	-	expression tag	UNP A0A125SZC1
E	-1	HIS	-	expression tag	UNP A0A125SZC1
E	0	MET	-	expression tag	UNP A0A125SZC1
F	-20	MET	-	initiating methionine	UNP A0A125SZC1
F	-19	GLY	-	expression tag	UNP A0A125SZC1
F	-18	SER	-	expression tag	UNP A0A125SZC1
F	-17	SER	-	expression tag	UNP A0A125SZC1
F	-16	HIS	-	expression tag	UNP A0A125SZC1
F	-15	HIS	-	expression tag	UNP A0A125SZC1
F	-14	HIS	-	expression tag	UNP A0A125SZC1
F	-13	HIS	-	expression tag	UNP A0A125SZC1
F	-12	HIS	-	expression tag	UNP A0A125SZC1
F	-11	HIS	-	expression tag	UNP A0A125SZC1
F	-10	SER	-	expression tag	UNP A0A125SZC1
F	-9	SER	-	expression tag	UNP A0A125SZC1
F	-8	GLY	-	expression tag	UNP A0A125SZC1
F	-7	LEU	-	expression tag	UNP A0A125SZC1
F	-6	VAL	-	expression tag	UNP A0A125SZC1
F	-5	PRO	-	expression tag	UNP A0A125SZC1
F	-4	ARG	-	expression tag	UNP A0A125SZC1
F	-3	GLY	-	expression tag	UNP A0A125SZC1
F	-2	SER	-	expression tag	UNP A0A125SZC1
F	-1	HIS	-	expression tag	UNP A0A125SZC1
F	0	MET	-	expression tag	UNP A0A125SZC1
G	-20	MET	-	initiating methionine	UNP A0A125SZC1
G	-19	GLY	-	expression tag	UNP A0A125SZC1
G	-18	SER	-	expression tag	UNP A0A125SZC1
G	-17	SER	-	expression tag	UNP A0A125SZC1
G	-16	HIS	-	expression tag	UNP A0A125SZC1
G	-15	HIS	-	expression tag	UNP A0A125SZC1
G	-14	HIS	-	expression tag	UNP A0A125SZC1
G	-13	HIS	-	expression tag	UNP A0A125SZC1
G	-12	HIS	-	expression tag	UNP A0A125SZC1
G	-11	HIS	-	expression tag	UNP A0A125SZC1
G	-10	SER	-	expression tag	UNP A0A125SZC1
G	-9	SER	-	expression tag	UNP A0A125SZC1
G	-8	GLY	-	expression tag	UNP A0A125SZC1
G	-7	LEU	-	expression tag	UNP A0A125SZC1
G	-6	VAL	-	expression tag	UNP A0A125SZC1

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-5	PRO	-	expression tag	UNP A0A125SZC1
G	-4	ARG	-	expression tag	UNP A0A125SZC1
G	-3	GLY	-	expression tag	UNP A0A125SZC1
G	-2	SER	-	expression tag	UNP A0A125SZC1
G	-1	HIS	-	expression tag	UNP A0A125SZC1
G	0	MET	-	expression tag	UNP A0A125SZC1

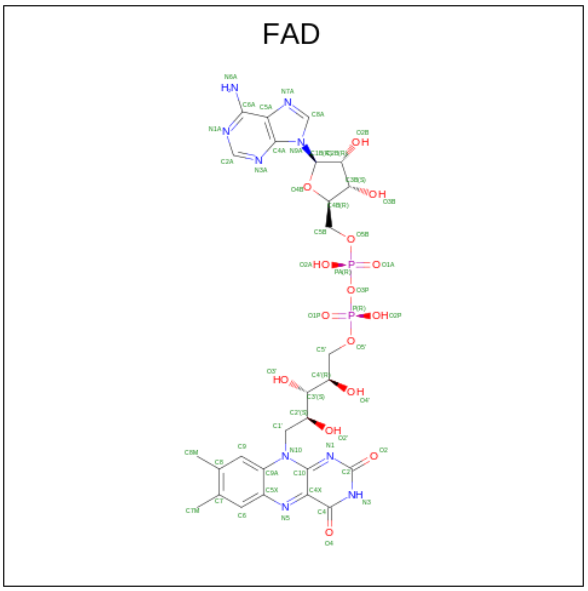
- Molecule 2 is a protein called N-formimidoyl fortimicin A synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	479	Total	C	N	O	S	0	1	0
			3642	2285	658	686	13			

There are 21 discrepancies between the modelled and reference sequences:

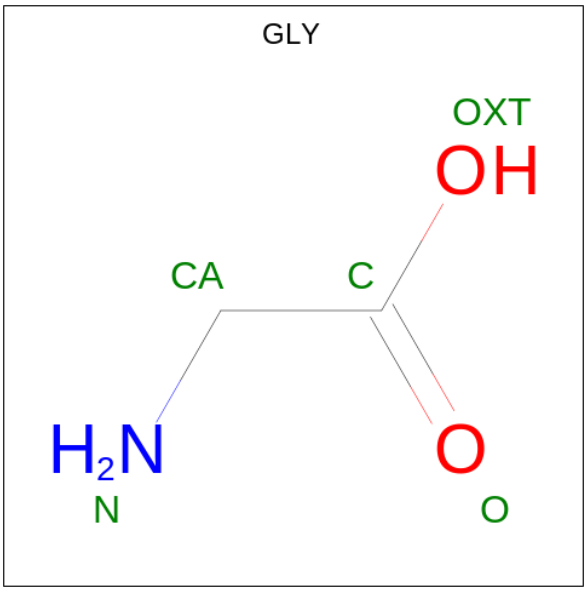
Chain	Residue	Modelled	Actual	Comment	Reference
H	-20	MET	-	initiating methionine	UNP A0A125SZC1
H	-19	GLY	-	expression tag	UNP A0A125SZC1
H	-18	SER	-	expression tag	UNP A0A125SZC1
H	-17	SER	-	expression tag	UNP A0A125SZC1
H	-16	HIS	-	expression tag	UNP A0A125SZC1
H	-15	HIS	-	expression tag	UNP A0A125SZC1
H	-14	HIS	-	expression tag	UNP A0A125SZC1
H	-13	HIS	-	expression tag	UNP A0A125SZC1
H	-12	HIS	-	expression tag	UNP A0A125SZC1
H	-11	HIS	-	expression tag	UNP A0A125SZC1
H	-10	SER	-	expression tag	UNP A0A125SZC1
H	-9	SER	-	expression tag	UNP A0A125SZC1
H	-8	GLY	-	expression tag	UNP A0A125SZC1
H	-7	LEU	-	expression tag	UNP A0A125SZC1
H	-6	VAL	-	expression tag	UNP A0A125SZC1
H	-5	PRO	-	expression tag	UNP A0A125SZC1
H	-4	ARG	-	expression tag	UNP A0A125SZC1
H	-3	GLY	-	expression tag	UNP A0A125SZC1
H	-2	SER	-	expression tag	UNP A0A125SZC1
H	-1	HIS	-	expression tag	UNP A0A125SZC1
H	0	MET	-	expression tag	UNP A0A125SZC1

- 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
3	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
3	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
3	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
3	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
3	E	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
3	F	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
3	G	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
3	H	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 4 is GLYCINE (three-letter code: GLY) (formula: $C_2H_5NO_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O 5 2 1 2	0	0
4	A	1	Total C N O 5 2 1 2	0	0
4	B	1	Total C N O 5 2 1 2	0	0
4	C	1	Total C N O 5 2 1 2	0	0
4	D	1	Total C N O 5 2 1 2	0	0
4	E	1	Total C N O 5 2 1 2	0	0
4	F	1	Total C N O 5 2 1 2	0	0
4	F	1	Total C N O 5 2 1 2	0	0
4	G	1	Total C N O 5 2 1 2	0	0
4	H	1	Total C N O 5 2 1 2	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	414	Total O 414 414	0	0
5	B	423	Total O 423 423	0	0

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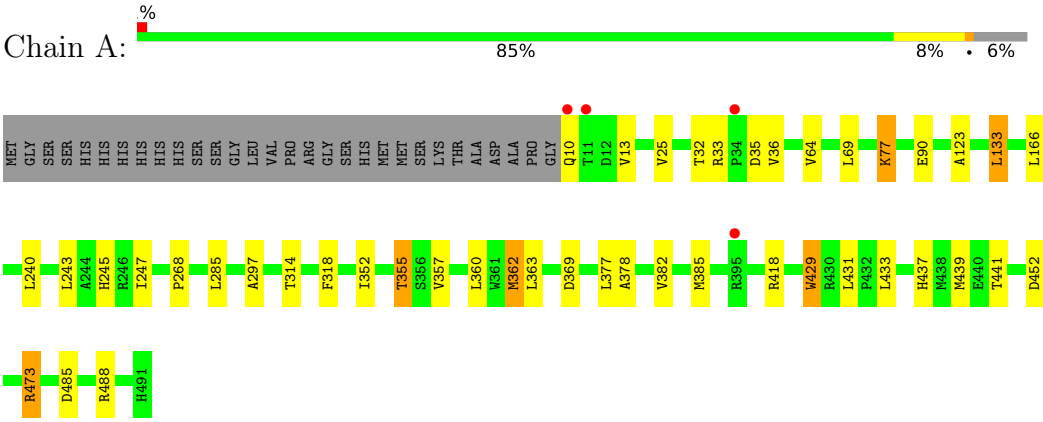
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	403	Total O 403 403	0	0
5	D	316	Total O 316 316	0	0
5	E	404	Total O 404 404	0	0
5	F	448	Total O 448 448	0	0
5	G	409	Total O 409 409	0	0
5	H	305	Total O 305 305	0	0

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.

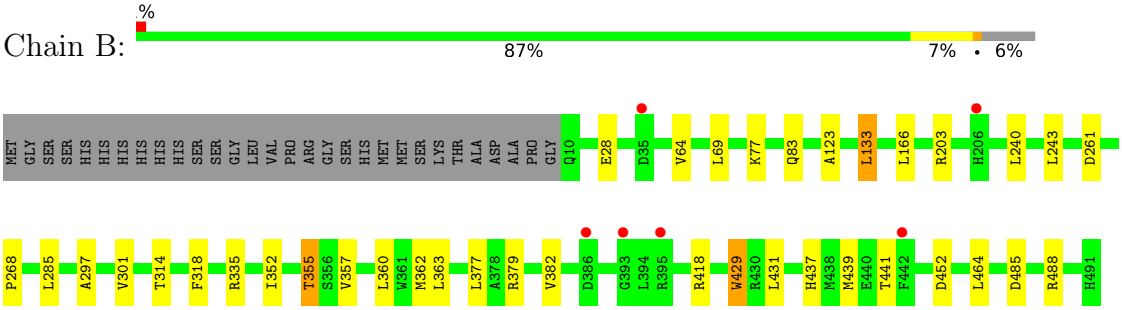
• Molecule 1: N-formimidoyl fortimicin A synthase

Chain A:



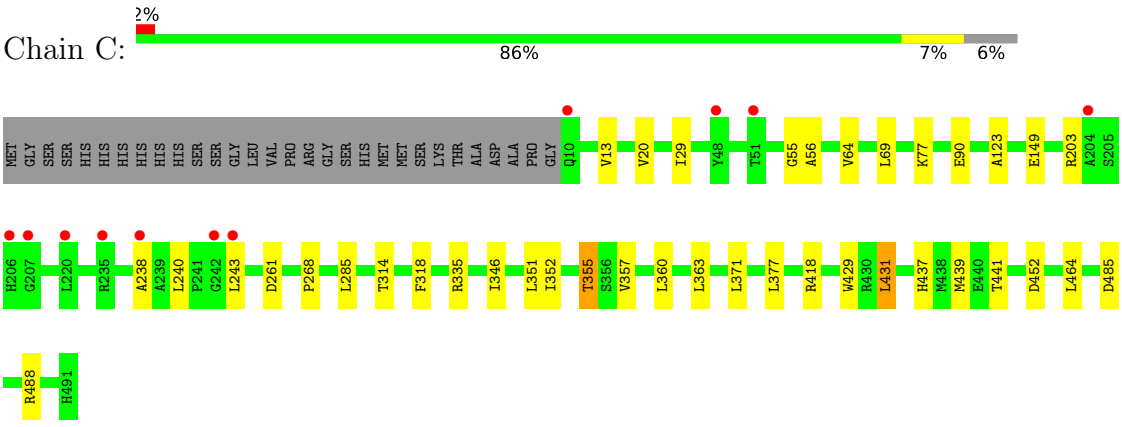
• Molecule 1: N-formimidoyl fortimicin A synthase

Chain B:



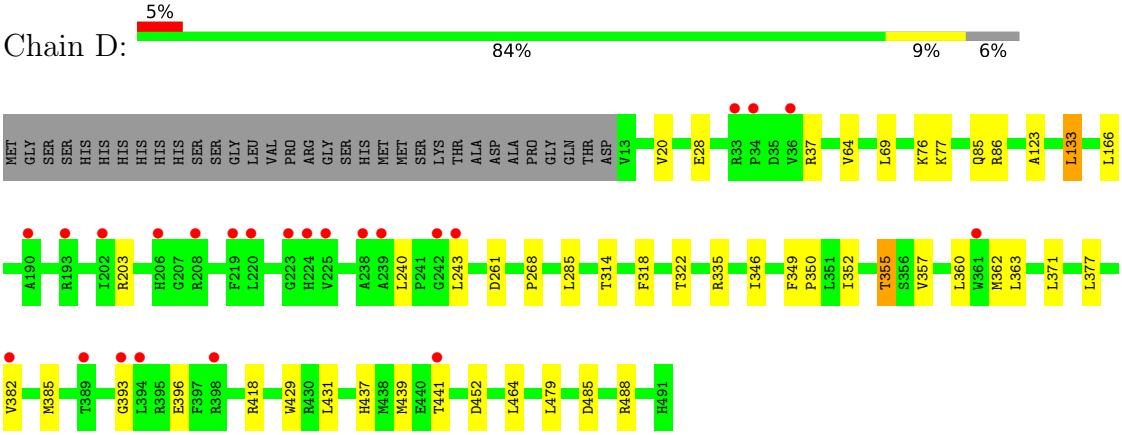
• Molecule 1: N-formimidoyl fortimicin A synthase

Chain C:



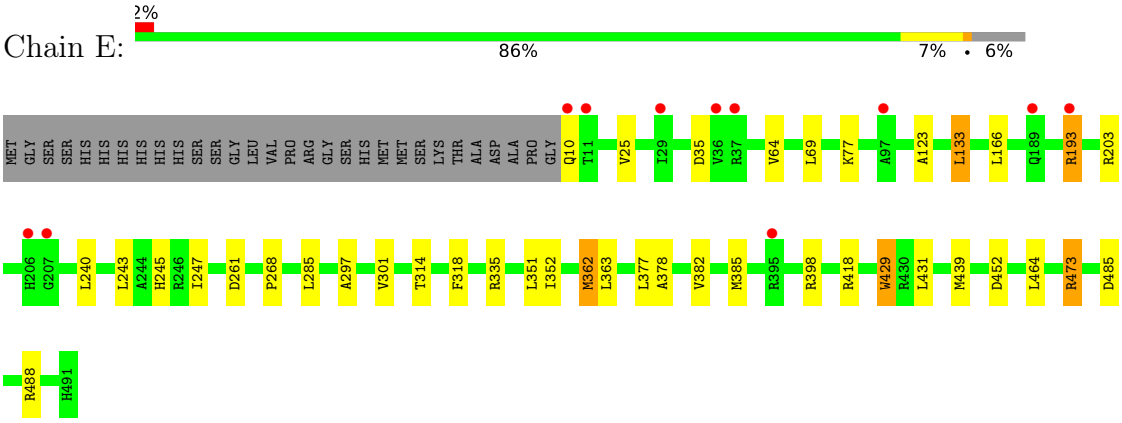
• Molecule 1: N-formimidoyl fortimicin A synthase

Chain D:



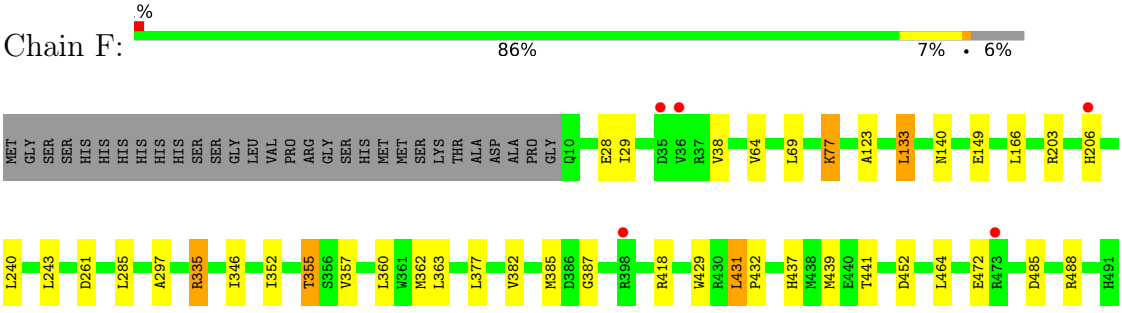
• Molecule 1: N-formimidoyl fortimicin A synthase

Chain E:



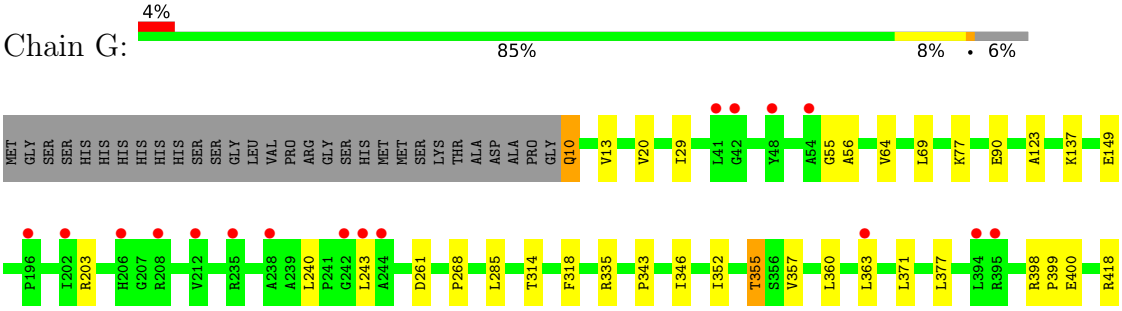
• Molecule 1: N-formimidoyl fortimicin A synthase

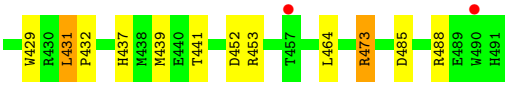
Chain F:



• Molecule 1: N-formimidoyl fortimicin A synthase

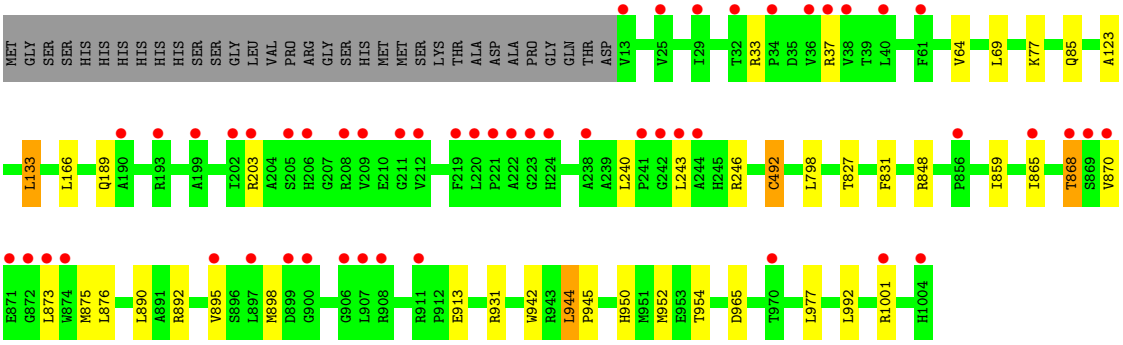
Chain G:





• Molecule 2: N-formimidoyl fortimicin A synthase

Chain H:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	103.39Å 107.80Å 134.75Å 89.85° 89.99° 83.58°	Depositor
Resolution (Å)	29.72 – 1.93 29.70 – 1.93	Depositor EDS
% Data completeness (in resolution range)	97.7 (29.72-1.93) 97.2 (29.70-1.93)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.73 (at 1.93Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.192 , 0.208 0.199 , 0.215	Depositor DCC
R_{free} test set	21320 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	17.9	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 28.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.467 for -h,-k,l 0.009 for k,h,-l 0.009 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	32835	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.79% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GQI, 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	0.66	0/3745	0.78	0/5102
1	B	0.66	0/3756	0.76	0/5116
1	C	0.69	0/3737	0.77	0/5092
1	D	0.64	0/3713	0.76	0/5059
1	E	0.66	0/3745	0.78	0/5102
1	F	0.66	0/3745	0.76	0/5102
1	G	0.68	0/3737	0.77	0/5092
2	H	0.64	0/3706	0.76	0/5048
All	All	0.66	0/29884	0.77	0/40713

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

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	3662	0	3592	54	0
1	B	3673	0	3604	41	0
1	C	3654	0	3584	41	0
1	D	3630	0	3565	51	0
1	E	3662	0	3592	45	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	3662	0	3592	47	0
1	G	3654	0	3584	51	0
2	H	3642	0	3559	39	0
3	A	53	0	31	2	0
3	B	53	0	31	2	0
3	C	53	0	31	1	0
3	D	53	0	31	2	0
3	E	53	0	31	3	0
3	F	53	0	31	2	0
3	G	53	0	31	0	0
3	H	53	0	31	2	0
4	A	10	0	4	0	0
4	B	5	0	2	0	0
4	C	5	0	2	0	0
4	D	5	0	2	0	0
4	E	5	0	2	0	0
4	F	10	0	4	1	0
4	G	5	0	2	0	0
4	H	5	0	2	0	0
5	A	414	0	0	9	0
5	B	423	0	0	2	0
5	C	403	0	0	5	0
5	D	316	0	0	4	0
5	E	404	0	0	5	0
5	F	448	0	0	4	0
5	G	409	0	0	7	0
5	H	305	0	0	5	0
All	All	32835	0	28940	335	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 6.

The worst 5 of 335 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:439[A]:MET:CE	1:C:439:MET:CE	2.38	1.02
1:E:352:ILE:HD12	1:E:363:LEU:HD12	1.42	1.00
1:E:439[A]:MET:CE	1:G:439:MET:CE	2.38	1.00
1:B:439[A]:MET:CE	1:D:439:MET:CE	2.41	0.98
1:F:439[A]:MET:CE	2:H:952:MET:CE	2.43	0.96

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	481/512 (94%)	469 (98%)	12 (2%)	0	100	100
1	B	482/512 (94%)	470 (98%)	12 (2%)	0	100	100
1	C	480/512 (94%)	468 (98%)	12 (2%)	0	100	100
1	D	477/512 (93%)	464 (97%)	13 (3%)	0	100	100
1	E	481/512 (94%)	469 (98%)	12 (2%)	0	100	100
1	F	481/512 (94%)	468 (97%)	13 (3%)	0	100	100
1	G	480/512 (94%)	468 (98%)	12 (2%)	0	100	100
2	H	476/512 (93%)	463 (97%)	13 (3%)	0	100	100
All	All	3838/4096 (94%)	3739 (97%)	99 (3%)	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	377/400 (94%)	367 (97%)	10 (3%)	44	31
1	B	378/400 (94%)	371 (98%)	7 (2%)	57	45
1	C	376/400 (94%)	370 (98%)	6 (2%)	62	52
1	D	373/400 (93%)	364 (98%)	9 (2%)	49	36

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	377/400 (94%)	368 (98%)	9 (2%)	49	36
1	F	377/400 (94%)	368 (98%)	9 (2%)	49	36
1	G	376/400 (94%)	367 (98%)	9 (2%)	49	36
2	H	372/399 (93%)	359 (96%)	13 (4%)	36	21
All	All	3006/3199 (94%)	2934 (98%)	72 (2%)	49	36

5 of 72 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	452	ASP
2	H	1001	ARG
2	H	33	ARG
2	H	848	ARG
1	D	77	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
1	G	245	HIS
1	G	319	ASN
1	C	245	HIS
1	E	245	HIS
1	F	245	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	GQI	H	492[B]	2	4,5,12	1.08	0	1,5,14	0.29	0
2	GQI	H	492[A]	2	8,11,12	1.59	3 (37%)	6,12,14	1.64	2 (33%)

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	GQI	H	492[B]	2	-	0/1/4/12	-
2	GQI	H	492[A]	2	-	1/2/10/12	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	492[A]	GQI	O1-N2	-2.61	1.40	1.45
2	H	492[A]	GQI	O3-C4	-2.36	1.24	1.30
2	H	492[A]	GQI	C3-C4	2.12	1.51	1.47

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	492[A]	GQI	O2-C4-C3	-2.95	116.24	124.16
2	H	492[A]	GQI	O3-C4-O2	2.05	126.93	122.67

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	H	492[A]	GQI	N-CA-CB-S1

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	492[B]	GQI	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

18 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	$\#\|Z\| > 2$
4	GLY	A	703	-	4,4,4	1.32	0	3,4,4	0.52	0
3	FAD	C	701	-	53,58,58	0.68	0	68,89,89	0.87	2 (2%)
3	FAD	A	701	-	53,58,58	0.75	1 (1%)	68,89,89	0.85	3 (4%)
4	GLY	D	702	-	4,4,4	0.80	0	3,4,4	1.87	1 (33%)
4	GLY	E	702	-	4,4,4	1.00	0	3,4,4	1.59	1 (33%)
3	FAD	B	701	-	53,58,58	0.69	0	68,89,89	0.86	1 (1%)
4	GLY	F	702	-	4,4,4	1.25	1 (25%)	3,4,4	0.95	0
4	GLY	H	1102	-	4,4,4	0.84	0	3,4,4	1.69	1 (33%)
4	GLY	G	702	-	4,4,4	1.05	0	3,4,4	1.12	0
3	FAD	G	701	-	53,58,58	0.64	0	68,89,89	0.86	2 (2%)
4	GLY	A	702	-	4,4,4	1.03	0	3,4,4	1.62	1 (33%)
3	FAD	D	701	-	53,58,58	0.63	0	68,89,89	0.77	2 (2%)
3	FAD	E	701	-	53,58,58	0.73	1 (1%)	68,89,89	0.83	3 (4%)
3	FAD	H	1101	-	53,58,58	0.65	0	68,89,89	0.77	2 (2%)
4	GLY	F	703	-	4,4,4	1.16	0	3,4,4	0.74	0
4	GLY	B	702	-	4,4,4	1.22	1 (25%)	3,4,4	1.13	0
4	GLY	C	702	-	4,4,4	1.22	1 (25%)	3,4,4	1.06	0
3	FAD	F	701	-	53,58,58	0.66	0	68,89,89	0.83	2 (2%)

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
4	GLY	A	703	-	-	0/2/2/2	-
3	FAD	C	701	-	-	2/30/50/50	0/6/6/6
3	FAD	A	701	-	-	2/30/50/50	0/6/6/6

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GLY	D	702	-	-	0/2/2/2	-
4	GLY	E	702	-	-	0/2/2/2	-
3	FAD	B	701	-	-	2/30/50/50	0/6/6/6
4	GLY	F	702	-	-	0/2/2/2	-
4	GLY	H	1102	-	-	0/2/2/2	-
4	GLY	G	702	-	-	0/2/2/2	-
3	FAD	G	701	-	-	2/30/50/50	0/6/6/6
4	GLY	A	702	-	-	0/2/2/2	-
3	FAD	D	701	-	-	2/30/50/50	0/6/6/6
3	FAD	E	701	-	-	2/30/50/50	0/6/6/6
3	FAD	H	1101	-	-	3/30/50/50	0/6/6/6
4	GLY	F	703	-	-	0/2/2/2	-
4	GLY	B	702	-	-	0/2/2/2	-
4	GLY	C	702	-	-	0/2/2/2	-
3	FAD	F	701	-	-	1/30/50/50	0/6/6/6

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	702	GLY	OXT-C	-2.29	1.23	1.30
3	A	701	FAD	C1'-C2'	-2.28	1.49	1.52
4	F	702	GLY	OXT-C	-2.07	1.23	1.30
4	B	702	GLY	OXT-C	-2.07	1.23	1.30
3	E	701	FAD	C1'-C2'	-2.04	1.49	1.52

The worst 5 of 21 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	701	FAD	O2A-PA-O1A	2.56	124.90	112.24
3	C	701	FAD	O2P-P-O1P	2.53	124.76	112.24
3	F	701	FAD	O2A-PA-O1A	2.53	124.75	112.24
4	D	702	GLY	OXT-C-CA	2.52	123.48	113.45
3	G	701	FAD	O2P-P-O1P	2.51	124.64	112.24

There are no chirality outliers.

5 of 16 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	701	FAD	O4B-C4B-C5B-O5B
3	G	701	FAD	O4B-C4B-C5B-O5B

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Mol	Chain	Res	Type	Atoms
3	D	701	FAD	C2'-C1'-N10-C10
3	E	701	FAD	C2'-C1'-N10-C10
3	H	1101	FAD	C2'-C1'-N10-C10

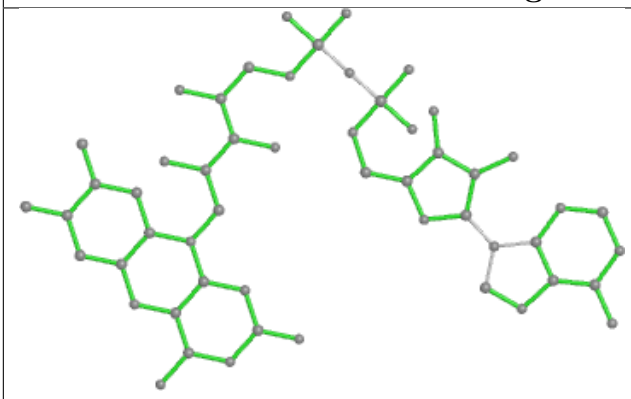
There are no ring outliers.

8 monomers are involved in 15 short contacts:

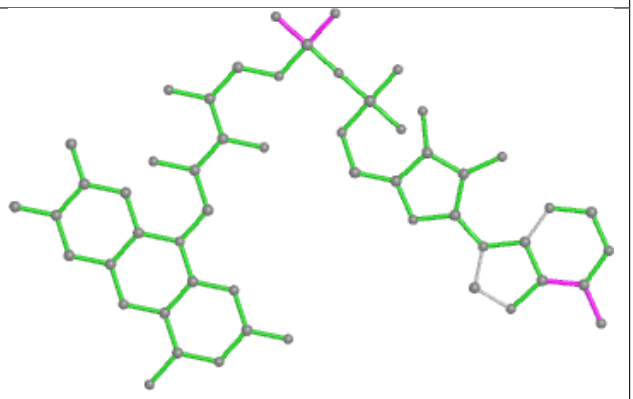
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	701	FAD	1	0
3	A	701	FAD	2	0
3	B	701	FAD	2	0
3	D	701	FAD	2	0
3	E	701	FAD	3	0
3	H	1101	FAD	2	0
4	F	703	GLY	1	0
3	F	701	FAD	2	0

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.

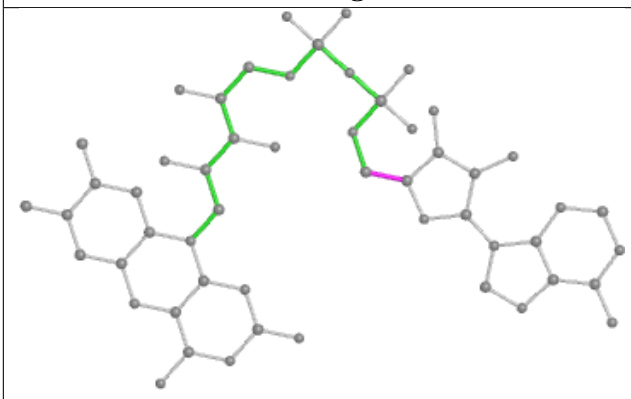
Ligand FAD C 701



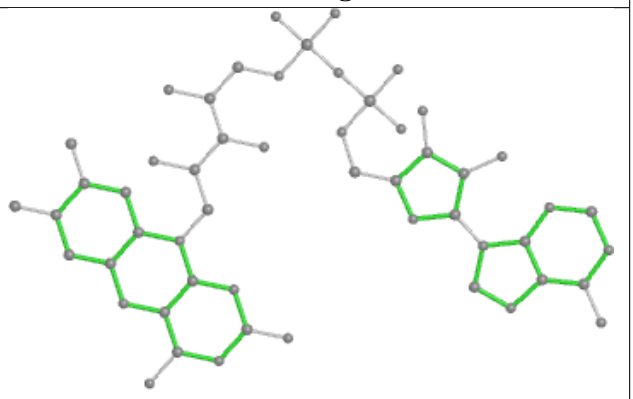
Bond lengths



Bond angles

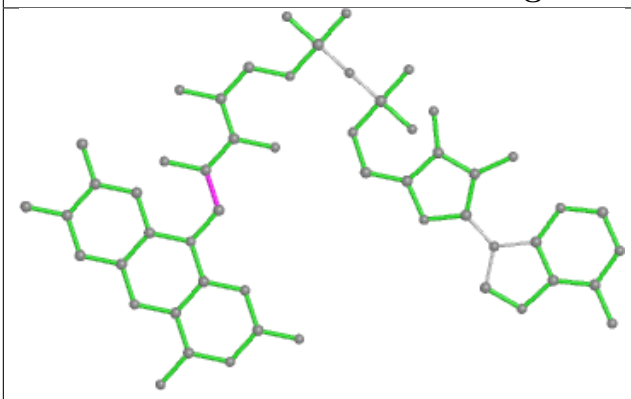


Torsions

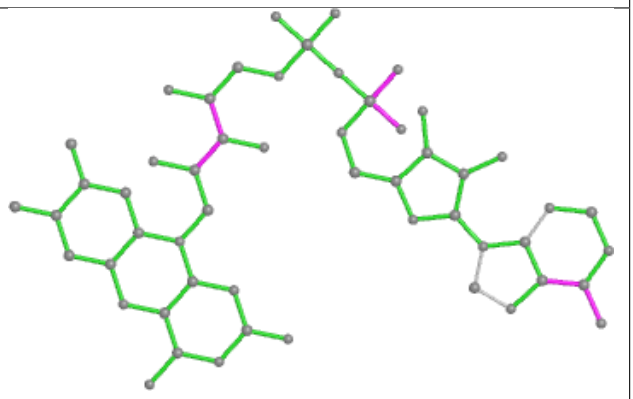


Rings

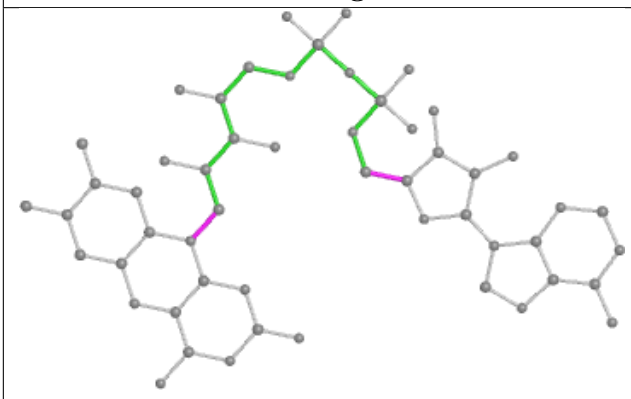
Ligand FAD A 701



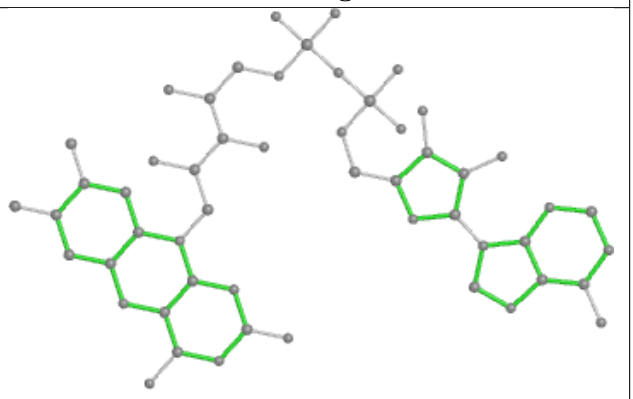
Bond lengths



Bond angles

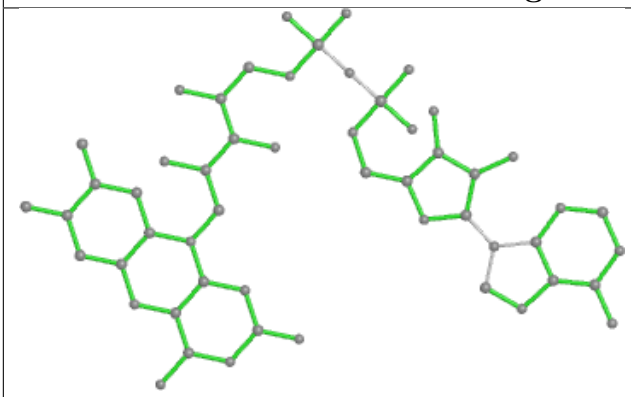


Torsions

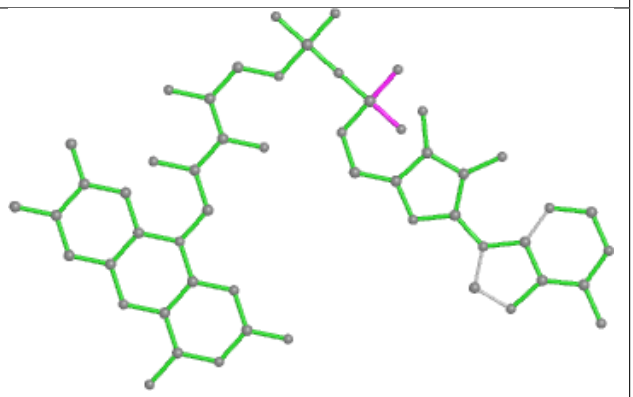


Rings

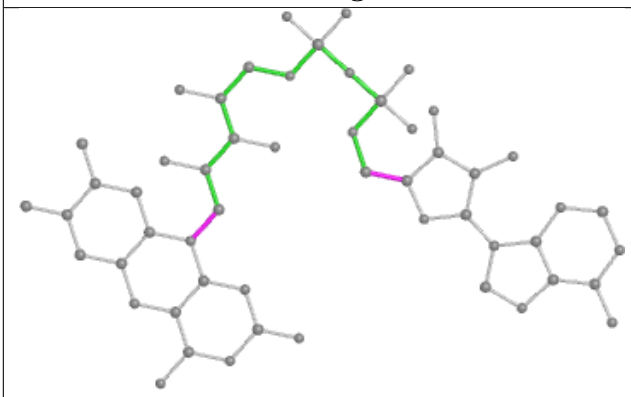
Ligand FAD B 701



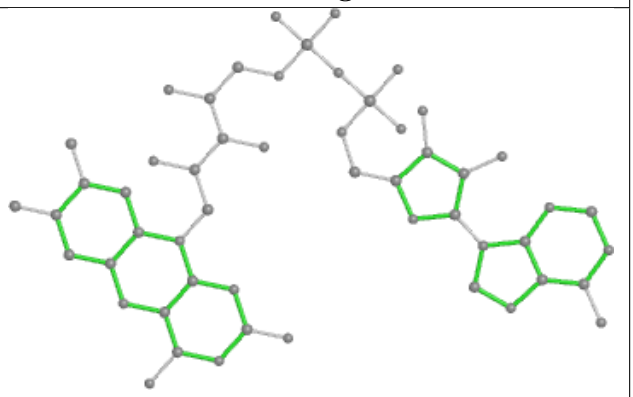
Bond lengths



Bond angles

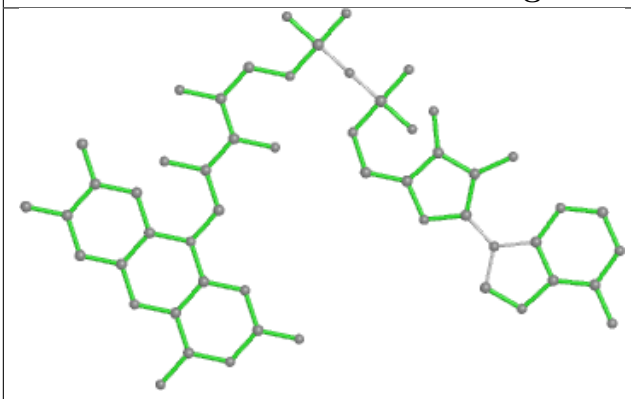


Torsions

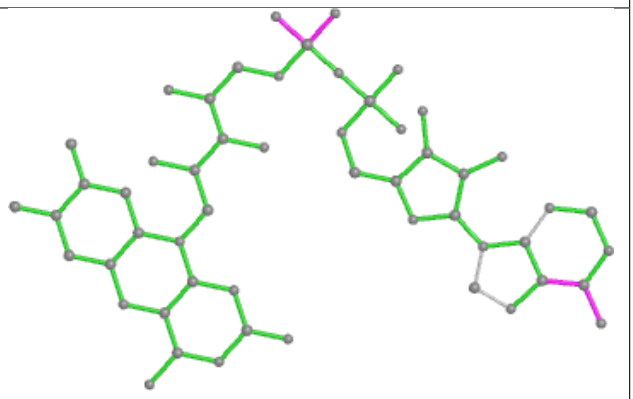


Rings

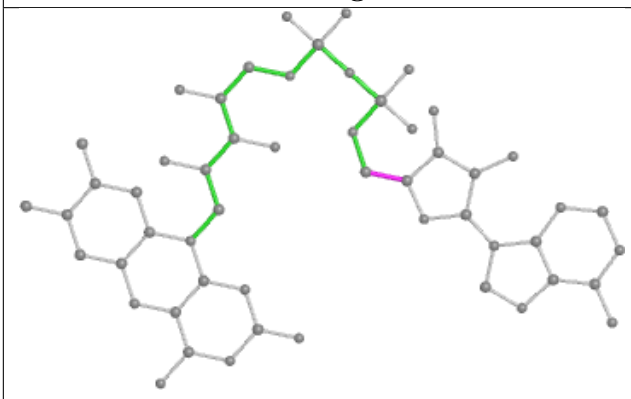
Ligand FAD G 701



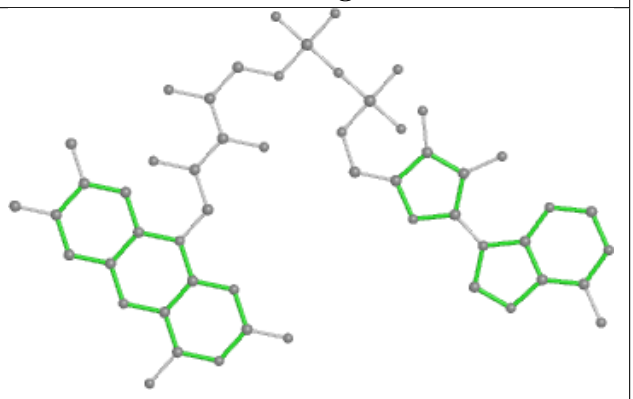
Bond lengths



Bond angles

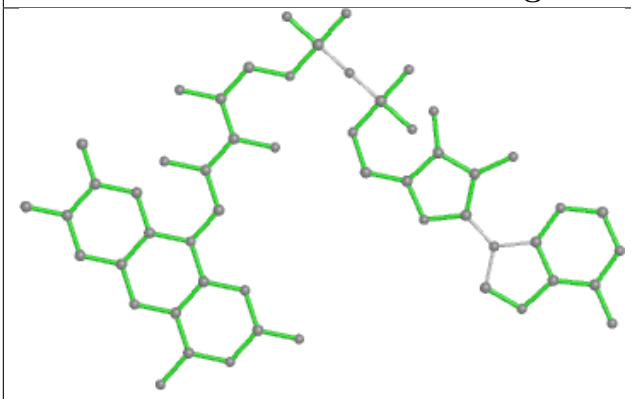


Torsions

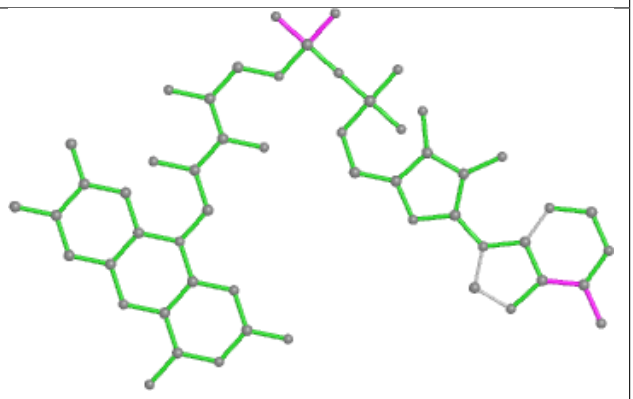


Rings

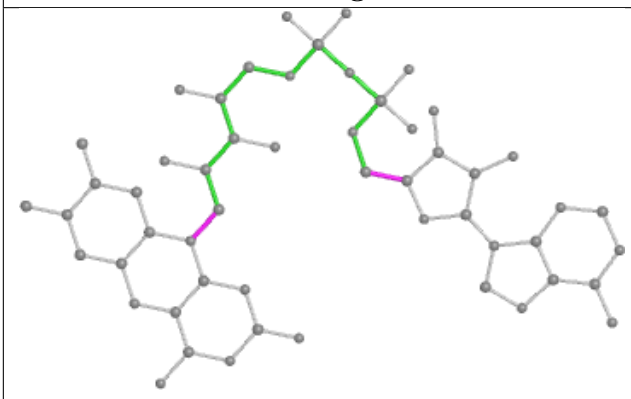
Ligand FAD D 701



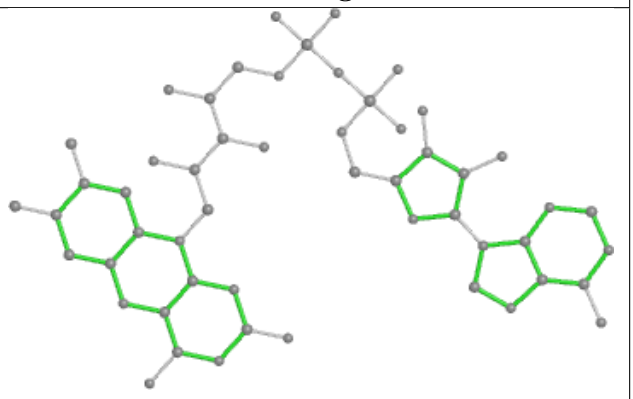
Bond lengths



Bond angles

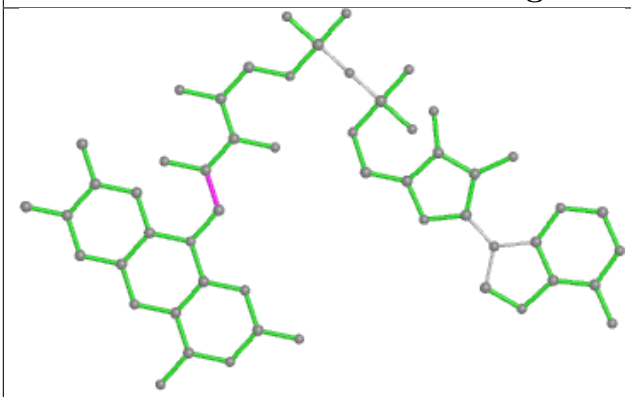


Torsions

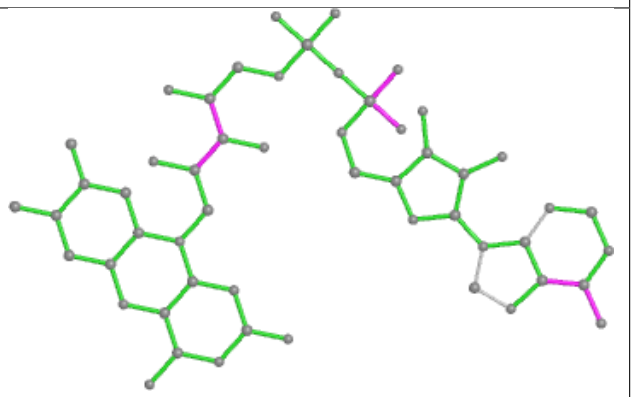


Rings

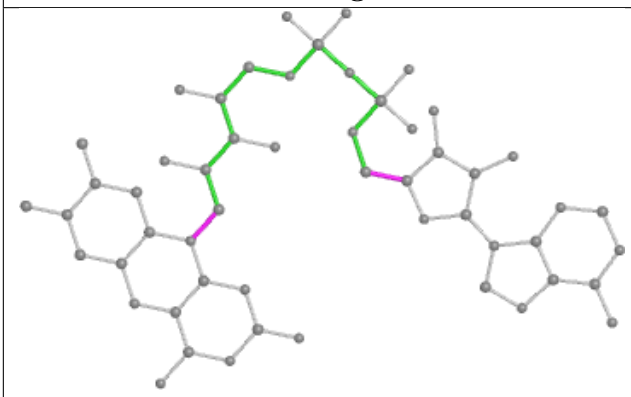
Ligand FAD E 701



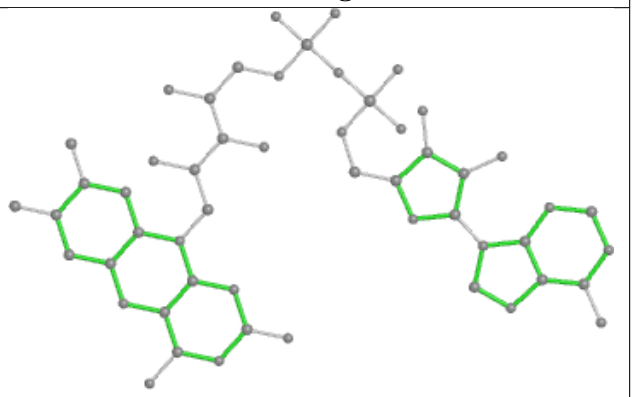
Bond lengths



Bond angles

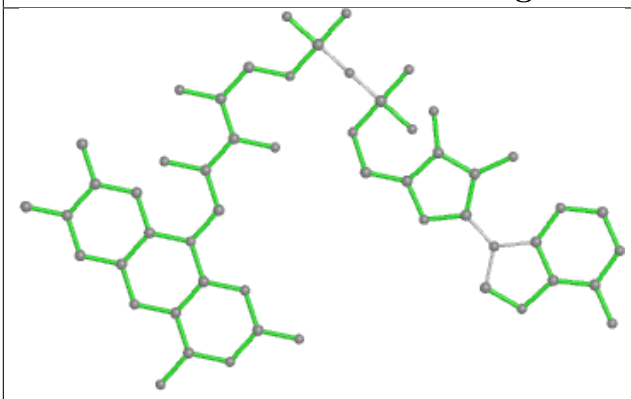


Torsions

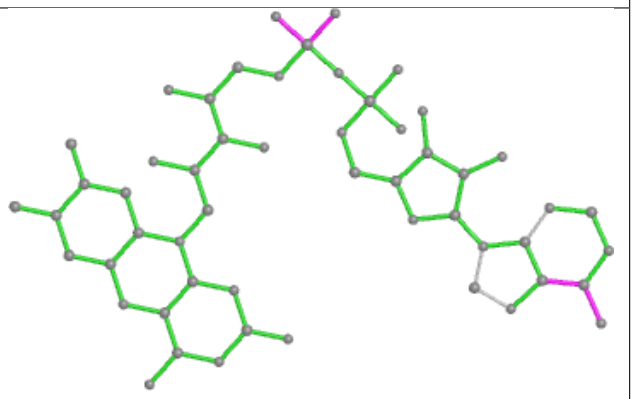


Rings

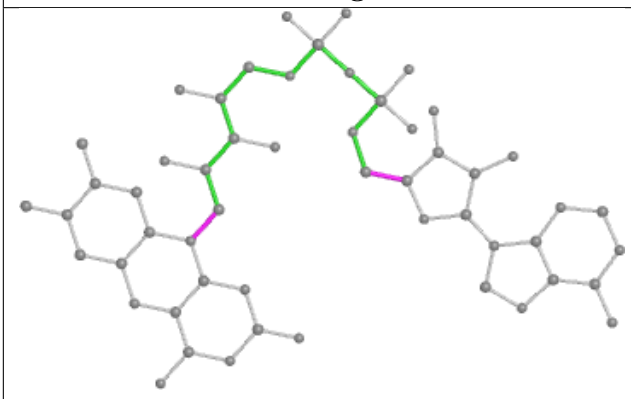
Ligand FAD H 1101



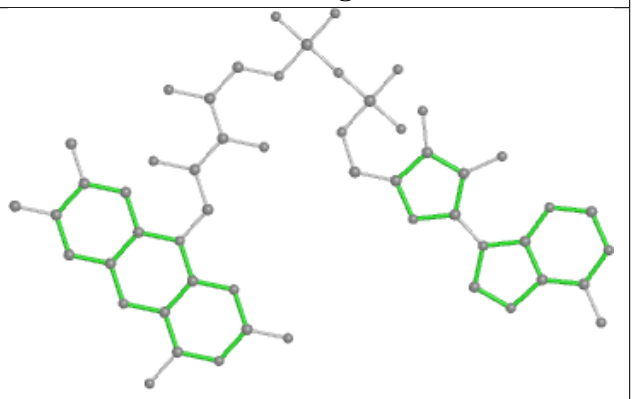
Bond lengths



Bond angles

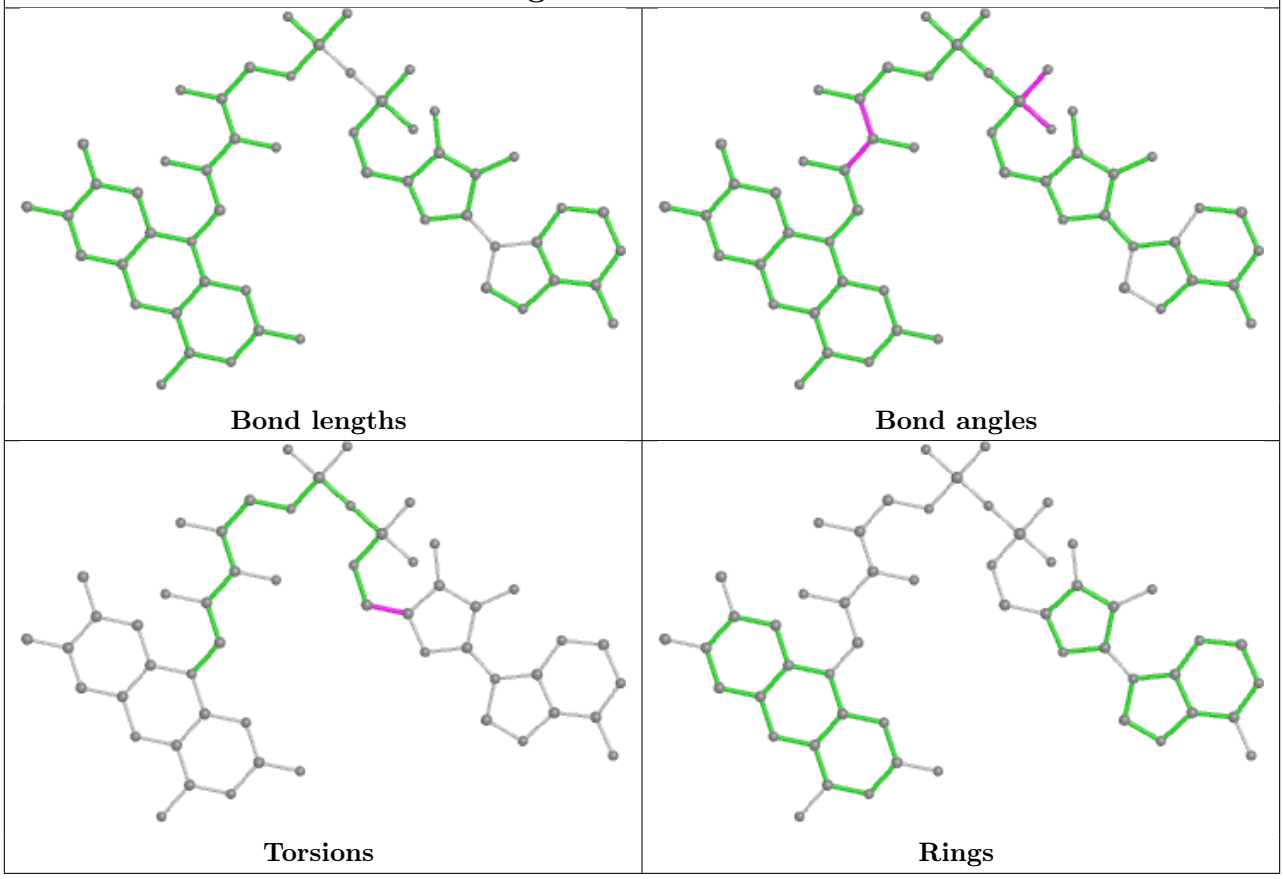


Torsions



Rings

Ligand FAD F 701



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	482/512 (94%)	0.23	5 (1%) 82 86	9, 20, 39, 72	1 (0%)
1	B	482/512 (94%)	0.15	6 (1%) 79 83	7, 18, 37, 64	1 (0%)
1	C	482/512 (94%)	0.32	11 (2%) 60 67	6, 22, 49, 72	0
1	D	479/512 (93%)	0.47	24 (5%) 28 36	10, 29, 67, 86	0
1	E	482/512 (94%)	0.26	11 (2%) 60 67	9, 20, 41, 70	1 (0%)
1	F	482/512 (94%)	0.13	5 (1%) 82 86	7, 17, 36, 67	1 (0%)
1	G	482/512 (94%)	0.40	19 (3%) 39 47	7, 22, 49, 72	0
2	H	478/512 (93%)	0.74	52 (10%) 5 8	10, 30, 68, 88	0
All	All	3849/4096 (93%)	0.34	133 (3%) 44 51	6, 22, 52, 88	4 (0%)

The worst 5 of 133 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	906	GLY	8.2
1	D	393	GLY	7.0
2	H	206	HIS	6.8
2	H	211	GLY	6.1
2	H	36	VAL	5.7

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

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
2	GQI	H	492[A]	12/13	0.88	0.14	24,29,29,34	12

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	GQI	H	492[B]	6/13	0.88	0.14	17,19,20,21	6

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

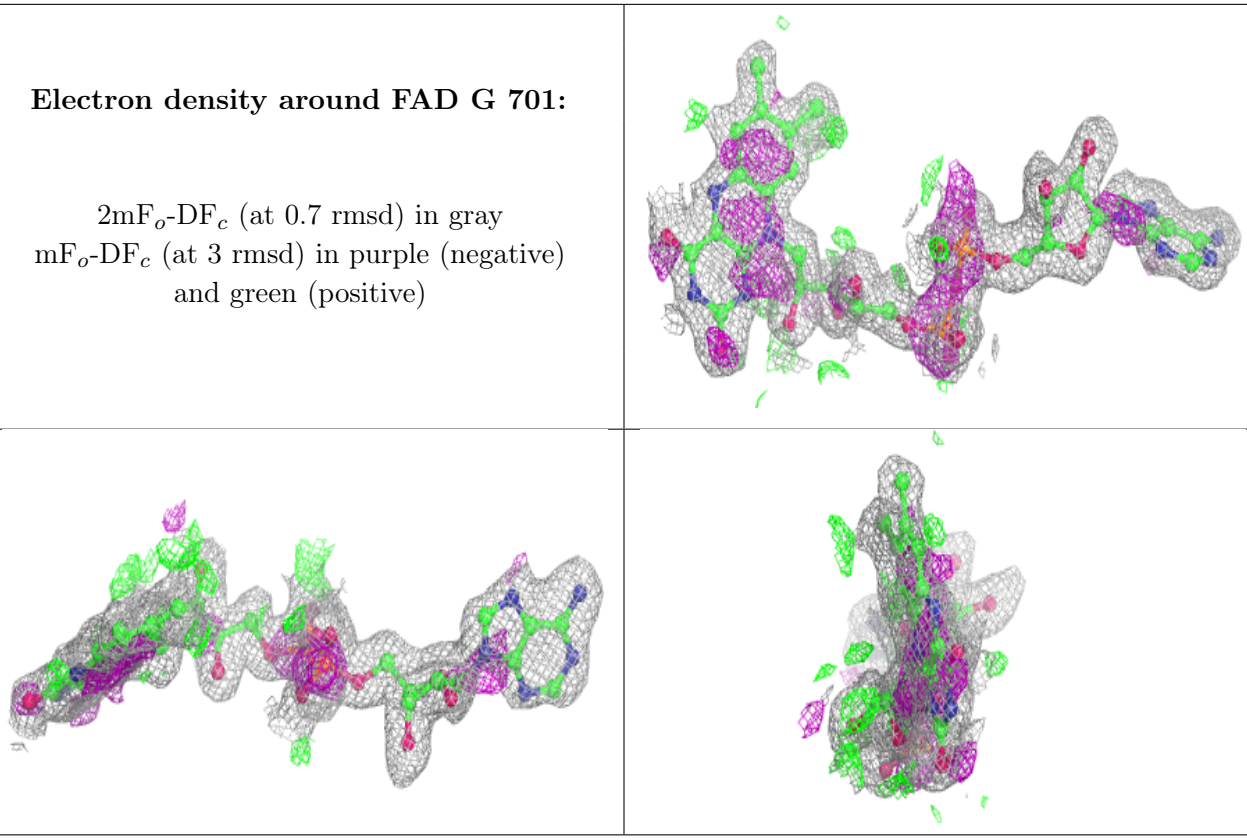
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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
4	GLY	F	703	5/5	0.75	0.17	30,30,34,36	0
4	GLY	A	703	5/5	0.84	0.19	21,25,28,30	0
3	FAD	G	701	53/53	0.88	0.19	23,28,40,40	0
3	FAD	H	1101	53/53	0.90	0.13	21,31,44,48	0
3	FAD	C	701	53/53	0.91	0.14	21,25,40,40	0
4	GLY	G	702	5/5	0.91	0.17	23,24,26,27	0
4	GLY	D	702	5/5	0.93	0.10	24,24,26,26	0
4	GLY	C	702	5/5	0.94	0.12	24,25,27,28	0
4	GLY	H	1102	5/5	0.94	0.09	24,25,26,26	0
3	FAD	D	701	53/53	0.95	0.11	21,29,44,48	0
3	FAD	E	701	53/53	0.95	0.11	12,15,19,19	0
4	GLY	E	702	5/5	0.96	0.07	15,16,17,17	0
3	FAD	A	701	53/53	0.96	0.10	13,16,18,19	0
3	FAD	F	701	53/53	0.97	0.09	9,11,14,15	0
3	FAD	B	701	53/53	0.97	0.09	9,11,13,15	0
4	GLY	B	702	5/5	0.98	0.09	11,11,12,12	0
4	GLY	A	702	5/5	0.98	0.10	15,15,16,16	0
4	GLY	F	702	5/5	0.98	0.10	11,11,11,12	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.

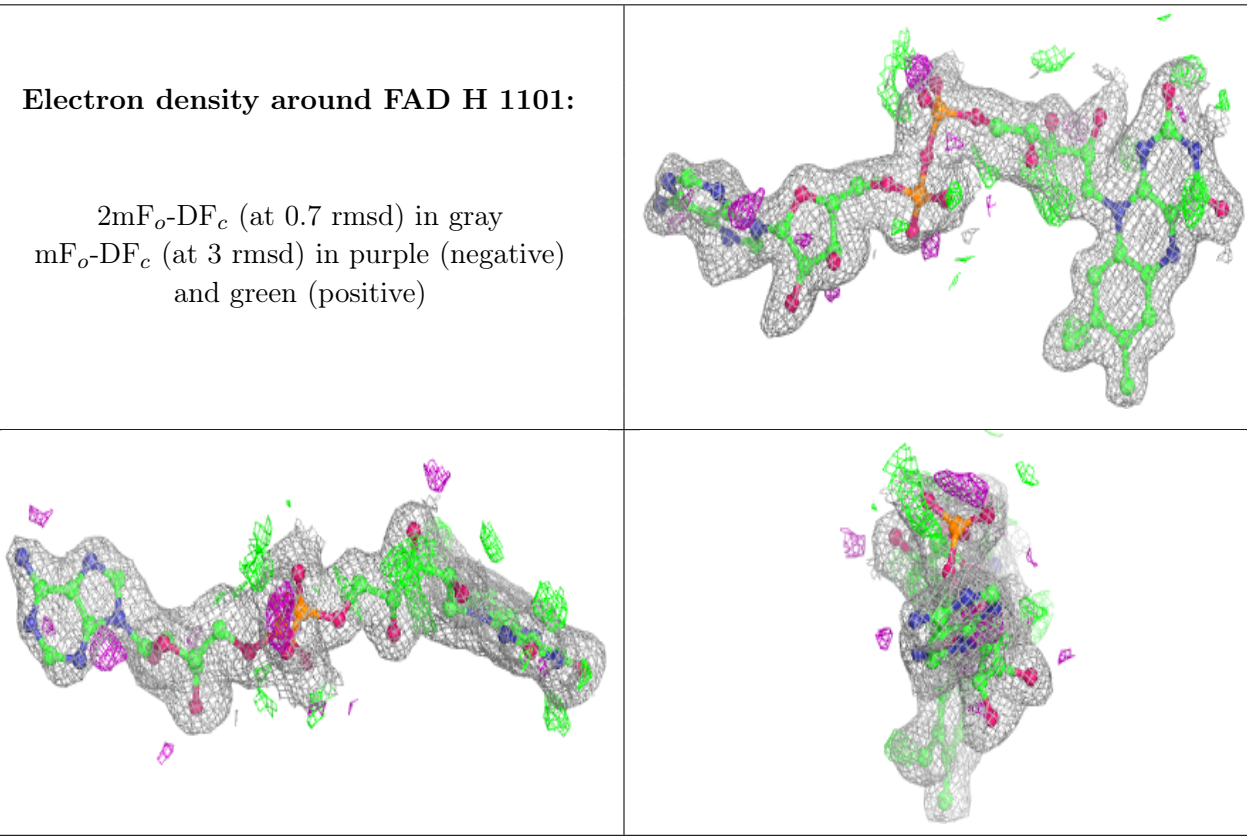
Electron density around FAD G 701:

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



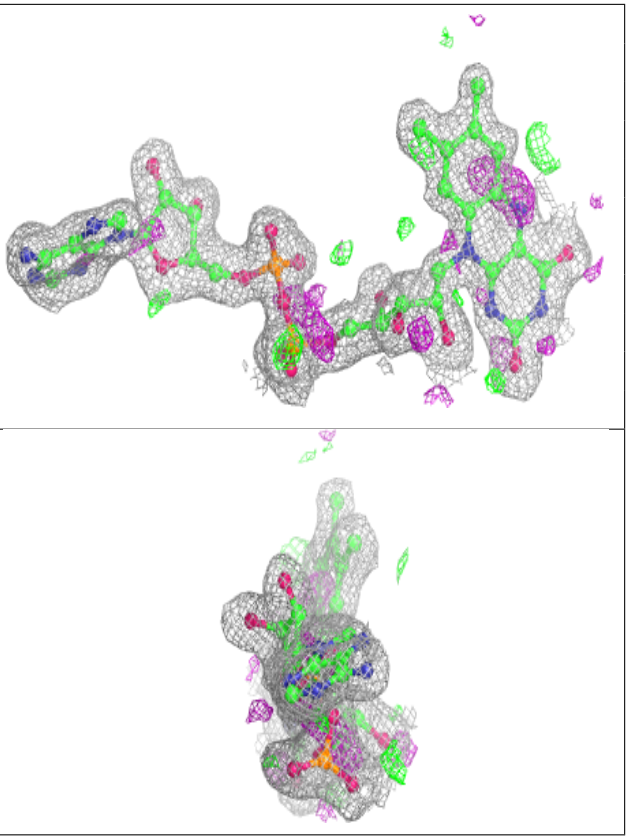
Electron density around FAD H 1101:

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



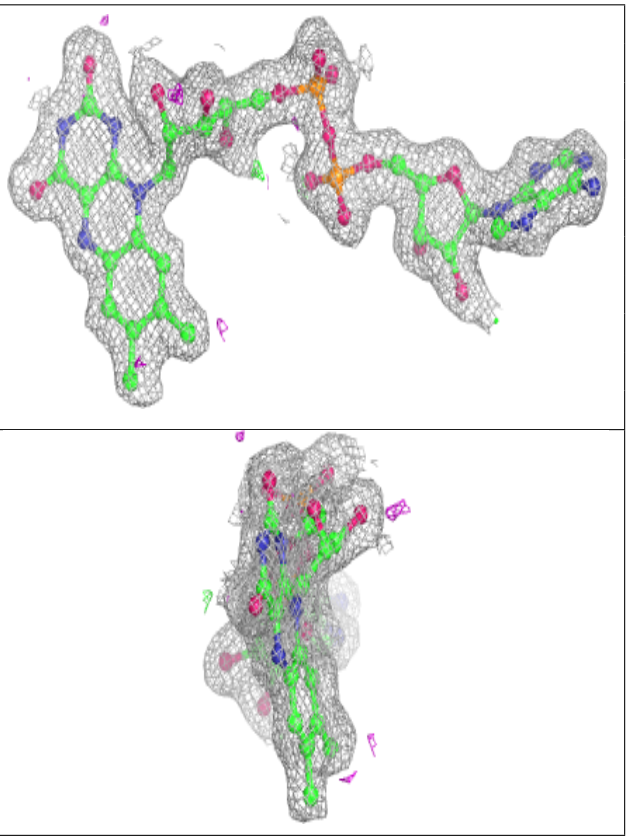
Electron density around FAD C 701:

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



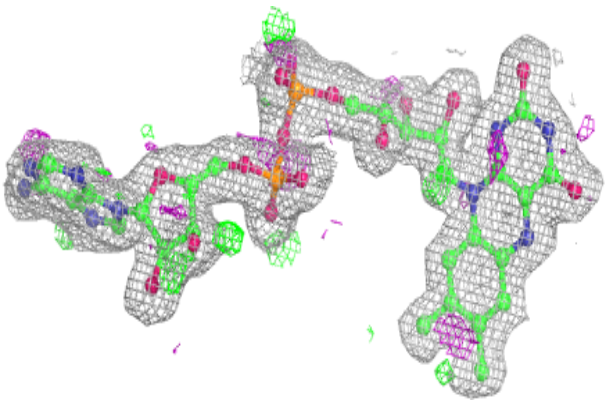
Electron density around FAD D 701:

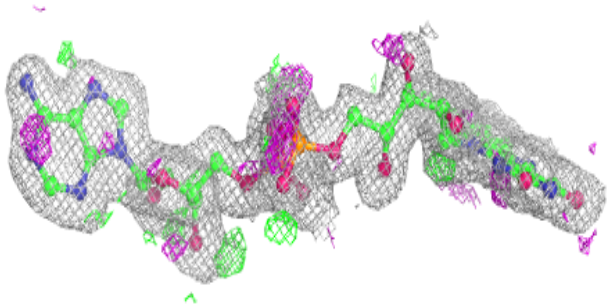
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

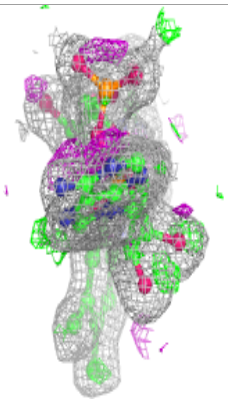


Electron density around FAD E 701:

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

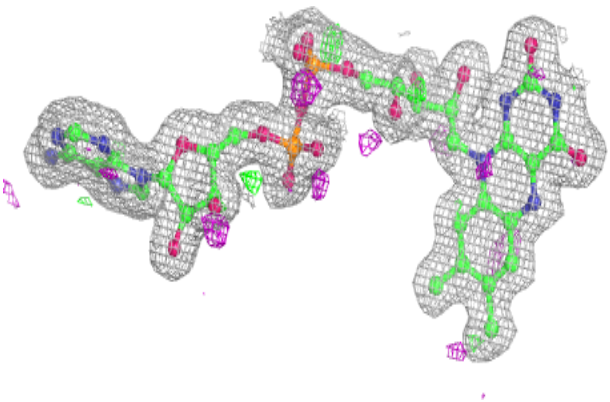


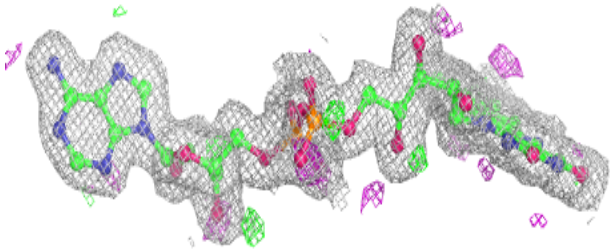


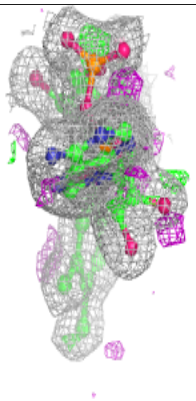


Electron density around FAD A 701:

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

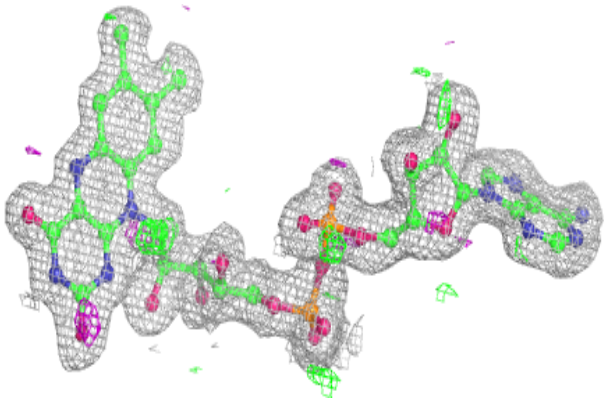


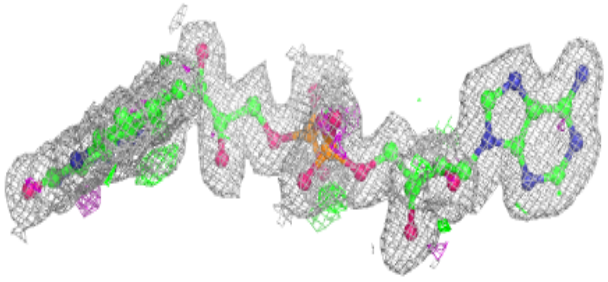


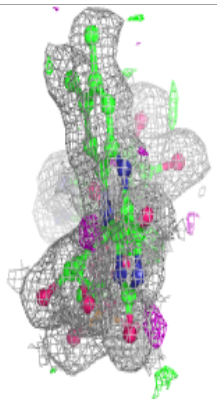


Electron density around FAD F 701:

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

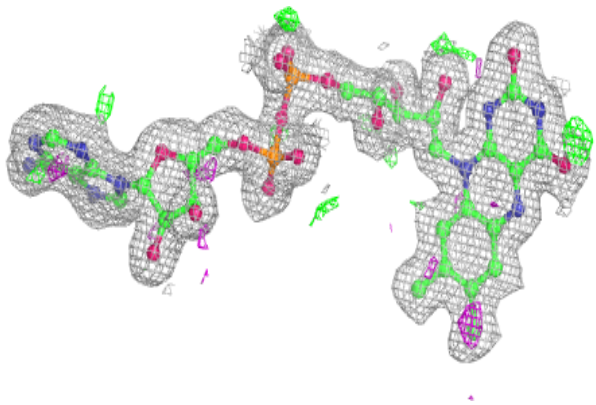


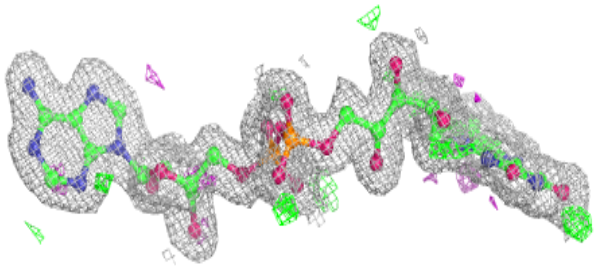


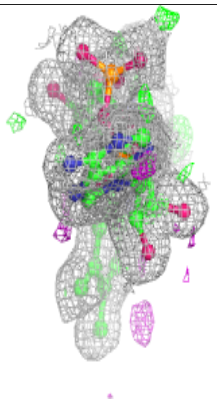


Electron density around FAD B 701:

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







6.5 Other polymers (i)

There are no such residues in this entry.