



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:17 PM UTC

PDB ID : 7XXD / pdb_00007xxd
Title : Orf1-sarcosine complex
Authors : Wang, Y.L.; Li, T.L.
Deposited on : 2022-05-29
Resolution : 1.98 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 ⓘ 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	:	2022.3.0, CSD as543be (2022)
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	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

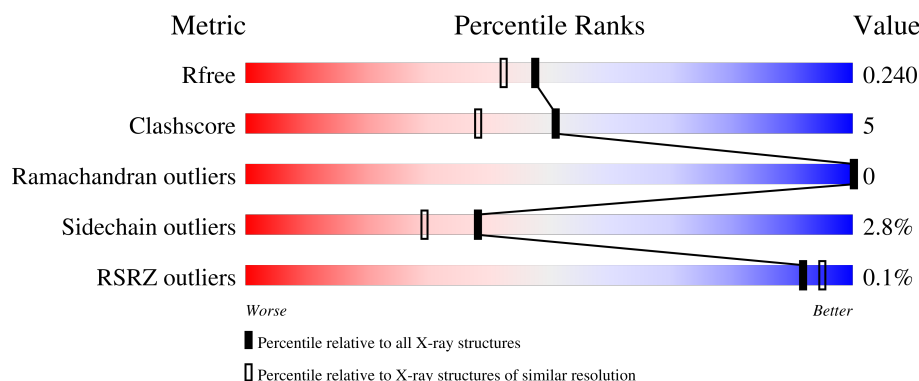
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

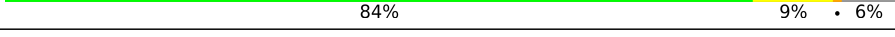
The reported resolution of this entry is 1.98 Å.

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	1506 (1.98-1.98)
Clashscore	190562	1534 (1.98-1.98)
Ramachandran outliers	187476	1518 (1.98-1.98)
Sidechain outliers	187428	1518 (1.98-1.98)
RSRZ outliers	180081	1506 (1.98-1.98)

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	512	 84% 9% • 6%
1	B	512	 86% 8% 6%
1	C	512	 85% 8% 6%
1	D	512	 84% 9% • 6%
1	E	512	 85% 8% • 6%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	512	87% 7% • 6%
1	G	512	85% 8% 6%
1	H	512	85% 8% • 6%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 32439 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	0	0
			3654	2293	660	689	12			
1	B	482	Total	C	N	O	S	0	1	0
			3662	2298	661	690	13			
1	C	482	Total	C	N	O	S	0	1	0
			3662	2298	661	690	13			
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	0	0
			3654	2293	660	689	12			
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			
1	H	479	Total	C	N	O	S	0	0	0
			3630	2280	656	682	12			

There are 168 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

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	LEU	-	expression tag	UNP A0A125SZC1
A	-6	VAL	-	expression tag	UNP A0A125SZC1
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

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	-7	LEU	-	expression tag	UNP A0A125SZC1
C	-6	VAL	-	expression tag	UNP A0A125SZC1
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

Continued on next page...

Continued from previous page...

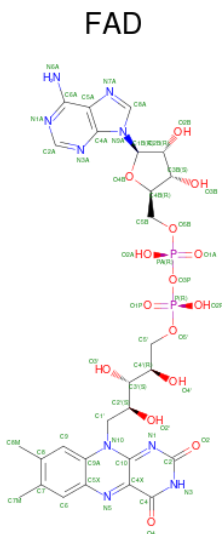
Chain	Residue	Modelled	Actual	Comment	Reference
E	-7	LEU	-	expression tag	UNP A0A125SZC1
E	-6	VAL	-	expression tag	UNP A0A125SZC1
E	-5	PRO	-	expression tag	UNP A0A125SZC1
E	-4	ARG	-	expression tag	UNP A0A125SZC1
E	-3	GLY	-	expression tag	UNP A0A125SZC1
E	-2	SER	-	expression tag	UNP A0A125SZC1
E	-1	HIS	-	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

Continued on next page...

Continued from previous page...

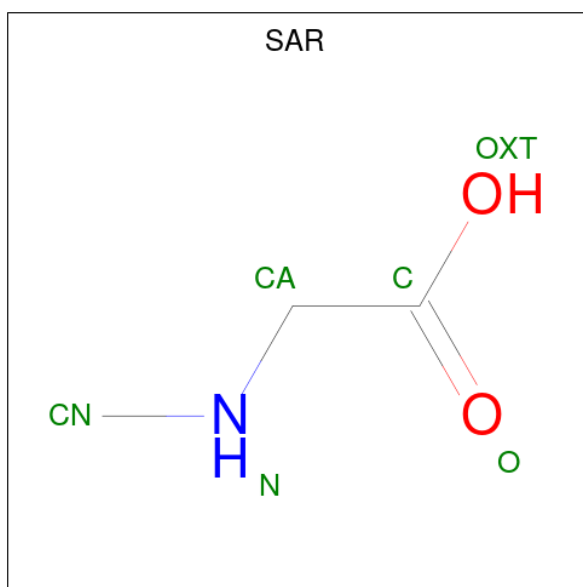
Chain	Residue	Modelled	Actual	Comment	Reference
G	-7	LEU	-	expression tag	UNP A0A125SZC1
G	-6	VAL	-	expression tag	UNP A0A125SZC1
G	-5	PRO	-	expression tag	UNP A0A125SZC1
G	-4	ARG	-	expression tag	UNP A0A125SZC1
G	-3	GLY	-	expression tag	UNP A0A125SZC1
G	-2	SER	-	expression tag	UNP A0A125SZC1
G	-1	HIS	-	expression tag	UNP A0A125SZC1
G	0	MET	-	expression tag	UNP A0A125SZC1
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 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is SARCOSINE (CCD ID: SAR) (formula: $\text{C}_3\text{H}_7\text{NO}_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			6	3	1	2		
3	B	1	Total	C	N	O	0	0
			6	3	1	2		
3	C	1	Total	C	N	O	0	0
			6	3	1	2		
3	D	1	Total	C	N	O	0	0
			6	3	1	2		
3	E	1	Total	C	N	O	0	0
			6	3	1	2		
3	F	1	Total	C	N	O	0	0
			6	3	1	2		
3	G	1	Total	C	N	O	0	0
			6	3	1	2		
3	H	1	Total	C	N	O	0	0
			6	3	1	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	399	Total	O	0	0
			399	399		
4	B	409	Total	O	0	0
			409	409		
4	C	344	Total	O	0	0
			344	344		
4	D	250	Total	O	0	0
			250	250		

Continued on next page...

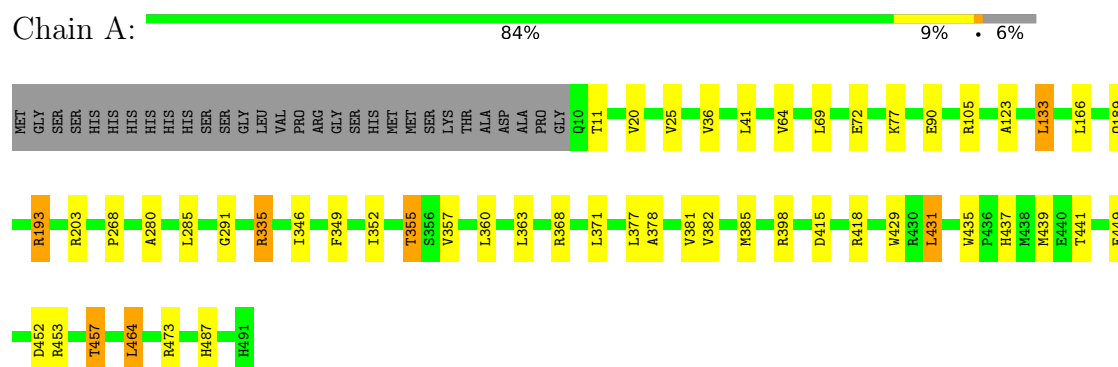
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	385	Total 385	O 385	0	0
4	F	421	Total 421	O 421	0	0
4	G	322	Total 322	O 322	0	0
4	H	229	Total 229	O 229	0	0

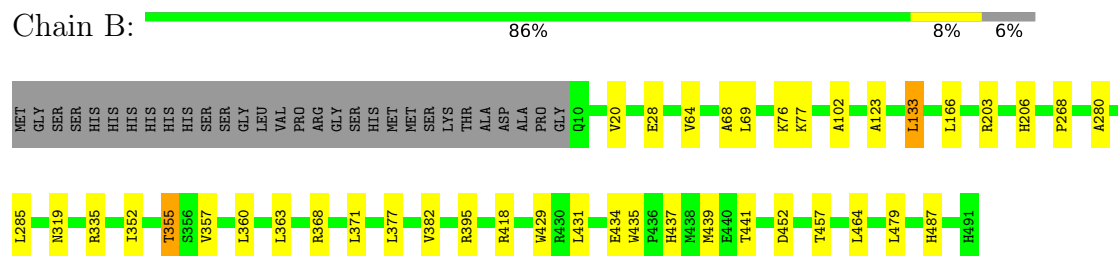
3 Residue-property plots

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.

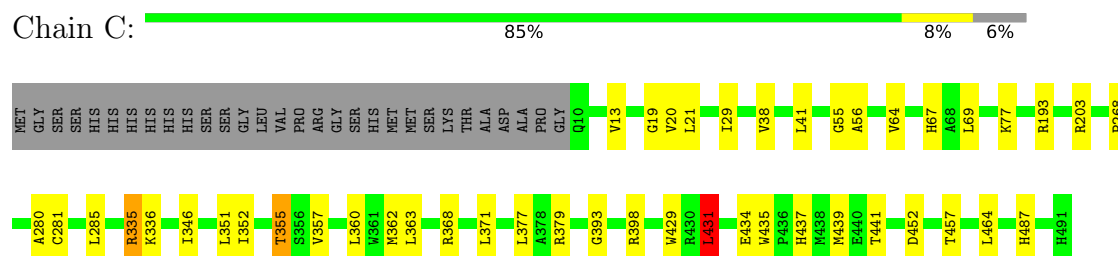
- Molecule 1: N-formimidoyl fortimicin A synthase



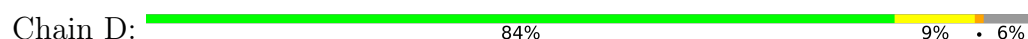
- Molecule 1: N-formimidoyl fortimicin A synthase

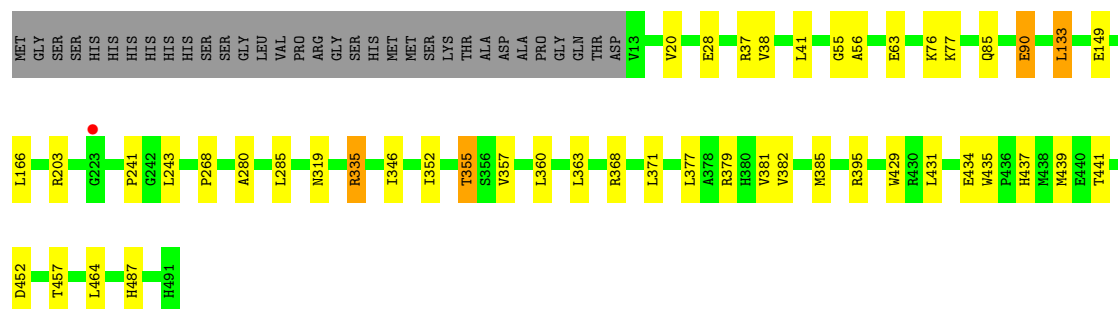


- Molecule 1: N-formimidoyl fortimicin A synthase



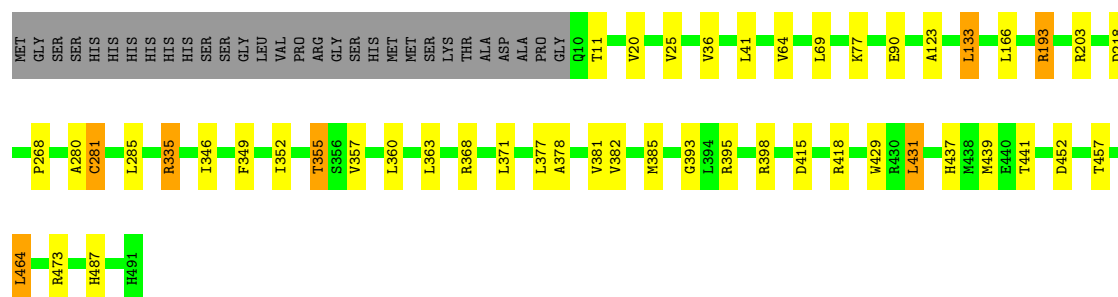
- Molecule 1: N-formimidoyl fortimicin A synthase





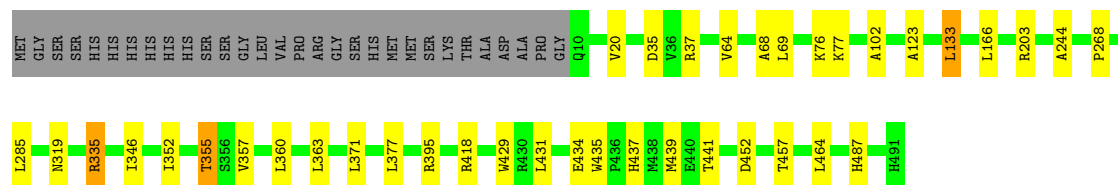
- Molecule 1: N-formimidoyl fortimicin A synthase

Chain E: 85% 8% 6%



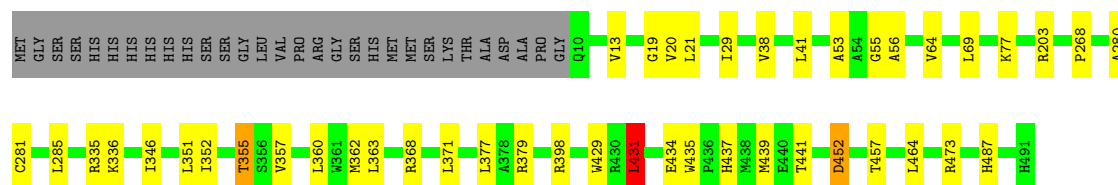
- Molecule 1: N-formimidoyl fortimicin A synthase

Chain F: 87% 7% 6%



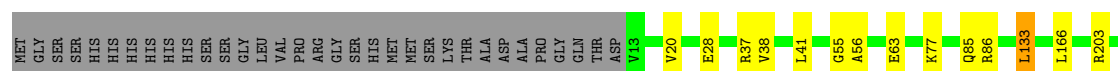
- Molecule 1: N-formimidoyl fortimicin A synthase

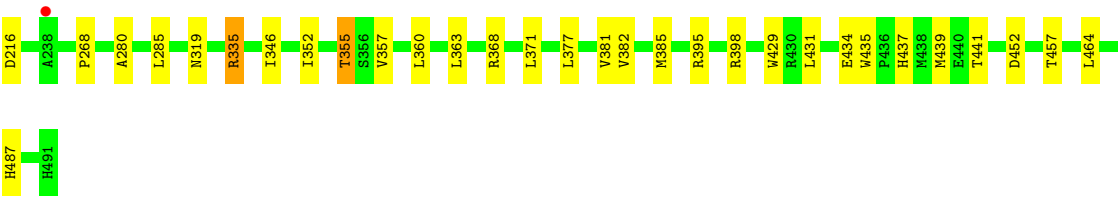
Chain G: 85% 8% 6%



- Molecule 1: N-formimidoyl fortimicin A synthase

Chain H: 85% 8% 6%





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	103.94Å 108.87Å 134.73Å 90.02° 90.00° 83.30°	Depositor
Resolution (Å)	28.73 – 1.98 28.73 – 1.98	Depositor EDS
% Data completeness (in resolution range)	93.2 (28.73-1.98) 93.2 (28.73-1.98)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.62 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.211 , 0.235 0.217 , 0.240	Depositor DCC
R_{free} test set	18447 reflections (4.53%)	wwPDB-VP
Wilson B-factor (Å ²)	24.4	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 27.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.478 for -h,-k,l 0.012 for -k,-h,-l 0.012 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	32439	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.15% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FAD, SAR

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	1.02	1/3737 (0.0%)	1.32	3/5092 (0.1%)
1	B	1.03	0/3745	1.31	2/5102 (0.0%)
1	C	1.04	2/3745 (0.1%)	1.32	2/5102 (0.0%)
1	D	0.98	0/3713	1.30	2/5059 (0.0%)
1	E	1.02	1/3737 (0.0%)	1.32	3/5092 (0.1%)
1	F	1.02	0/3745	1.31	2/5102 (0.0%)
1	G	1.04	2/3737 (0.1%)	1.32	2/5092 (0.0%)
1	H	0.98	0/3713	1.30	2/5059 (0.0%)
All	All	1.02	6/29872 (0.0%)	1.31	18/40700 (0.0%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	431	LEU	N-CA	6.38	1.51	1.45
1	C	19	GLY	C-O	6.28	1.29	1.23
1	G	431	LEU	N-CA	6.26	1.50	1.45
1	G	19	GLY	C-O	5.48	1.28	1.23
1	E	431	LEU	N-CA	5.37	1.50	1.45
1	A	431	LEU	N-CA	5.35	1.50	1.45

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	487	HIS	CA-C-N	5.57	127.69	120.44
1	C	487	HIS	C-N-CA	5.57	127.69	120.44
1	E	487	HIS	CA-C-N	5.57	127.68	120.44
1	E	487	HIS	C-N-CA	5.57	127.68	120.44
1	A	487	HIS	CA-C-N	5.56	127.67	120.44
1	A	487	HIS	C-N-CA	5.56	127.67	120.44
1	H	487	HIS	CA-C-N	5.52	127.62	120.44

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	487	HIS	C-N-CA	5.52	127.62	120.44
1	G	487	HIS	CA-C-N	5.52	127.61	120.44
1	G	487	HIS	C-N-CA	5.52	127.61	120.44
1	D	487	HIS	CA-C-N	5.50	127.59	120.44
1	D	487	HIS	C-N-CA	5.50	127.59	120.44
1	B	487	HIS	CA-C-N	5.38	127.44	120.44
1	B	487	HIS	C-N-CA	5.38	127.44	120.44
1	F	487	HIS	CA-C-N	5.38	127.43	120.44
1	F	487	HIS	C-N-CA	5.38	127.43	120.44
1	A	349	PHE	CB-CA-C	5.14	116.84	108.87
1	E	349	PHE	CB-CA-C	5.05	116.69	108.87

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	3654	0	3584	50	0
1	B	3662	0	3592	38	0
1	C	3662	0	3592	40	0
1	D	3630	0	3565	43	0
1	E	3654	0	3584	47	0
1	F	3662	0	3592	38	0
1	G	3654	0	3584	36	0
1	H	3630	0	3565	37	0
2	A	53	0	31	2	0
2	B	53	0	31	1	0
2	C	53	0	31	0	0
2	D	53	0	31	3	0
2	E	53	0	31	1	0
2	F	53	0	31	0	0
2	G	53	0	31	1	0
2	H	53	0	31	4	0
3	A	6	0	6	1	0
3	B	6	0	6	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	6	0	6	0	0
3	D	6	0	6	1	0
3	E	6	0	6	1	0
3	F	6	0	6	0	0
3	G	6	0	6	0	0
3	H	6	0	6	1	0
4	A	399	0	0	18	0
4	B	409	0	0	8	0
4	C	344	0	0	11	0
4	D	250	0	0	9	0
4	E	385	0	0	16	0
4	F	421	0	0	11	0
4	G	322	0	0	8	0
4	H	229	0	0	4	0
All	All	32439	0	28954	306	0

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

All (306) 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:439:MET:HE3	1:G:439:MET:CE	1.85	1.06
1:E:439:MET:CE	1:G:439:MET:HE3	1.85	1.06
1:A:439:MET:HE3	1:C:439[A]:MET:CE	1.86	1.04
1:A:439:MET:CE	1:C:439[A]:MET:HE3	1.85	1.04
1:F:439[A]:MET:CE	1:H:439:MET:HE3	1.90	1.01
1:B:439[A]:MET:CE	1:D:439:MET:HE3	1.91	0.99
1:F:439[A]:MET:HE3	1:H:439:MET:CE	1.91	0.99
1:B:102:ALA:HB1	4:B:972:HOH:O	1.61	0.99
1:A:398:ARG:HG2	4:A:903:HOH:O	1.62	0.99
1:E:398:ARG:HG2	4:E:897:HOH:O	1.62	0.98
1:B:439[A]:MET:HE3	1:D:439:MET:CE	1.92	0.97
1:E:415:ASP:HB2	4:E:886:HOH:O	1.62	0.97
1:E:281:CYS:SG	4:E:1103:HOH:O	2.22	0.96
1:F:441:THR:HB	4:F:807:HOH:O	1.71	0.91
1:D:319:ASN:HB3	4:D:984:HOH:O	1.73	0.88
1:E:335:ARG:HG2	4:E:1116:HOH:O	1.74	0.87
1:E:439:MET:CE	1:G:439:MET:CE	2.50	0.87
1:A:335:ARG:HG2	4:A:1116:HOH:O	1.75	0.86
1:A:439:MET:HE3	1:C:439[A]:MET:HE2	1.58	0.85

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:355:THR:HG22	1:A:357:VAL:H	1.42	0.85
1:A:439:MET:CE	1:C:439[A]:MET:CE	2.50	0.84
1:B:355:THR:HG22	1:B:357:VAL:H	1.42	0.84
1:C:355:THR:HG22	1:C:357:VAL:H	1.42	0.84
1:D:241:PRO:HA	4:D:1012:HOH:O	1.77	0.84
1:D:355:THR:HG22	1:D:357:VAL:H	1.42	0.84
1:E:439:MET:HE3	1:G:439:MET:HE2	1.57	0.83
1:A:439:MET:HE2	1:C:439[A]:MET:HE3	1.60	0.83
1:E:355:THR:HG22	1:E:357:VAL:H	1.41	0.83
1:G:355:THR:HG22	1:G:357:VAL:H	1.42	0.83
1:B:441:THR:HB	4:B:835:HOH:O	1.76	0.82
1:H:355:THR:HG22	1:H:357:VAL:H	1.42	0.82
1:F:355:THR:HG22	1:F:357:VAL:H	1.42	0.82
4:A:819:HOH:O	1:C:434:GLU:HG2	1.78	0.82
1:F:439[A]:MET:HE2	1:H:439:MET:HE3	1.61	0.82
1:A:382:VAL:HA	1:A:385:MET:HE3	1.62	0.81
1:E:439:MET:HE2	1:G:439:MET:HE3	1.61	0.81
1:B:439[A]:MET:CE	1:D:439:MET:CE	2.55	0.81
1:B:439[A]:MET:HE2	1:D:439:MET:HE3	1.61	0.81
1:D:382:VAL:HA	1:D:385:MET:HE3	1.63	0.81
1:H:382:VAL:HA	1:H:385:MET:HE3	1.63	0.81
1:E:382:VAL:HA	1:E:385:MET:HE3	1.62	0.80
1:C:398:ARG:HD2	4:C:1092:HOH:O	1.80	0.80
1:C:335:ARG:HG2	4:C:1096:HOH:O	1.82	0.79
1:B:439[A]:MET:HE3	1:D:439:MET:HE2	1.65	0.79
4:F:870:HOH:O	1:H:434:GLU:HG2	1.83	0.79
1:F:439[A]:MET:CE	1:H:439:MET:CE	2.55	0.78
4:B:858:HOH:O	1:D:434:GLU:HG2	1.84	0.77
1:F:439[A]:MET:HE3	1:H:439:MET:HE2	1.65	0.77
1:A:415:ASP:HB2	4:A:936:HOH:O	1.82	0.77
1:F:434:GLU:HG2	4:H:807:HOH:O	1.86	0.75
4:E:809:HOH:O	1:G:434:GLU:HG2	1.88	0.74
1:C:363:LEU:HD11	1:C:377:LEU:HB3	1.70	0.74
1:F:441:THR:CB	4:F:807:HOH:O	2.34	0.72
1:F:102:ALA:HB1	4:F:941:HOH:O	1.90	0.71
1:E:382:VAL:HA	1:E:385:MET:CE	2.21	0.71
1:G:363:LEU:HD11	1:G:377:LEU:HB3	1.70	0.71
1:D:382:VAL:HA	1:D:385:MET:CE	2.21	0.70
1:H:382:VAL:HA	1:H:385:MET:CE	2.21	0.70
1:G:452:ASP:HB3	4:G:820:HOH:O	1.91	0.70
1:F:335:ARG:HG2	4:F:1167:HOH:O	1.91	0.69

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:382:VAL:HA	1:A:385:MET:CE	2.21	0.69
1:A:449:GLU:HB3	4:A:1131:HOH:O	1.92	0.69
1:F:352:ILE:HD13	1:F:377:LEU:HD22	1.75	0.68
1:B:431:LEU:HD21	1:B:439[A]:MET:HE1	1.76	0.68
1:C:429:TRP:CD1	1:C:431:LEU:HD11	2.29	0.68
1:G:429:TRP:CD1	1:G:431:LEU:HD11	2.28	0.68
1:D:431:LEU:HD21	1:D:439:MET:HE1	1.76	0.67
1:F:431:LEU:HD21	1:F:439[A]:MET:HE1	1.76	0.67
1:A:431:LEU:HD21	1:A:439:MET:HE1	1.77	0.67
1:B:319:ASN:HB3	4:B:980:HOH:O	1.94	0.67
1:F:35:ASP:HB3	4:F:1155:HOH:O	1.95	0.67
1:E:431:LEU:HD21	1:E:439:MET:HE1	1.77	0.66
1:H:352:ILE:HD13	1:H:377:LEU:HD22	1.77	0.66
1:D:352:ILE:HD13	1:D:377:LEU:HD22	1.77	0.66
1:B:441:THR:CB	4:B:835:HOH:O	2.41	0.66
1:E:415:ASP:CB	4:E:886:HOH:O	2.32	0.65
1:G:355:THR:HG21	1:G:360:LEU:HD23	1.78	0.65
1:H:431:LEU:HD21	1:H:439:MET:HE1	1.77	0.65
1:B:355:THR:HG21	1:B:360:LEU:HD23	1.78	0.65
1:B:352:ILE:HD13	1:B:377:LEU:HD22	1.80	0.64
1:G:53:ALA:HB1	4:G:850:HOH:O	1.97	0.63
1:B:434:GLU:HG2	4:D:836:HOH:O	1.97	0.63
1:C:355:THR:HG21	1:C:360:LEU:HD23	1.79	0.63
1:E:355:THR:HG21	1:E:360:LEU:HD23	1.80	0.63
1:F:355:THR:HG21	1:F:360:LEU:HD23	1.79	0.63
2:G:701:FAD:HM73	4:G:850:HOH:O	1.99	0.63
1:D:355:THR:HG21	1:D:360:LEU:HD23	1.80	0.62
1:A:355:THR:HG21	1:A:360:LEU:HD23	1.80	0.62
1:H:355:THR:HG21	1:H:360:LEU:HD23	1.81	0.62
1:A:381:VAL:HG12	1:A:385:MET:HE2	1.82	0.62
1:A:352:ILE:HD13	1:A:377:LEU:HD22	1.82	0.61
1:F:346:ILE:HG13	1:F:464:LEU:CD2	2.30	0.61
1:G:429:TRP:NE1	1:G:431:LEU:HD11	2.15	0.61
1:C:429:TRP:NE1	1:C:431:LEU:HD11	2.15	0.61
1:E:381:VAL:HG12	1:E:385:MET:HE2	1.83	0.61
1:A:335:ARG:CD	4:A:1113:HOH:O	2.49	0.60
1:D:346:ILE:HG13	1:D:464:LEU:CD2	2.32	0.60
1:E:346:ILE:HG13	1:E:464:LEU:CD2	2.32	0.60
1:H:346:ILE:HG13	1:H:464:LEU:CD2	2.32	0.60
1:G:429:TRP:CE2	1:G:431:LEU:HD11	2.37	0.60
1:C:429:TRP:CE2	1:C:431:LEU:HD11	2.37	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:381:VAL:HG12	1:H:385:MET:HE2	1.83	0.59
1:D:381:VAL:HG12	1:D:385:MET:HE2	1.83	0.59
1:D:363:LEU:HD11	1:D:377:LEU:HB3	1.84	0.59
1:E:395:ARG:N	4:E:803:HOH:O	2.35	0.59
1:H:363:LEU:HD11	1:H:377:LEU:HB3	1.85	0.59
1:E:352:ILE:HD13	1:E:377:LEU:HD22	1.84	0.58
1:G:346:ILE:HG13	1:G:464:LEU:CD2	2.34	0.58
1:E:363:LEU:HD11	1:E:377:LEU:HB3	1.86	0.58
1:C:346:ILE:HG13	1:C:464:LEU:CD2	2.34	0.58
1:E:335:ARG:CD	4:E:1114:HOH:O	2.51	0.57
1:B:206:HIS:O	1:H:86:ARG:HD2	2.03	0.57
1:E:441:THR:HG22	4:E:1084:HOH:O	2.05	0.57
1:A:189:GLN:HG2	4:A:1119:HOH:O	2.03	0.57
1:F:363:LEU:HD11	1:F:377:LEU:HB3	1.87	0.56
1:H:355:THR:CG2	1:H:357:VAL:H	2.17	0.56
1:B:363:LEU:HD11	1:B:377:LEU:HB3	1.87	0.56
1:E:335:ARG:HD2	4:E:1114:HOH:O	2.05	0.56
1:D:335:ARG:HG2	4:D:925:HOH:O	2.06	0.56
1:E:346:ILE:HG13	1:E:464:LEU:HD21	1.86	0.56
1:G:473:ARG:HB2	4:G:1047:HOH:O	2.04	0.56
1:C:398:ARG:HG2	4:C:802:HOH:O	2.05	0.55
1:C:398:ARG:CG	4:C:802:HOH:O	2.54	0.55
1:A:335:ARG:HD3	4:A:1113:HOH:O	2.07	0.55
1:A:363:LEU:HD11	1:A:377:LEU:HB3	1.88	0.55
1:E:355:THR:CG2	1:E:357:VAL:H	2.17	0.55
1:A:90:GLU:HG3	4:A:1118:HOH:O	2.07	0.54
1:G:352:ILE:HD12	1:G:363:LEU:HD12	1.89	0.54
1:C:352:ILE:HD12	1:C:363:LEU:HD12	1.90	0.54
1:G:355:THR:CG2	1:G:357:VAL:H	2.18	0.54
1:A:346:ILE:HG13	1:A:464:LEU:CD1	2.37	0.54
1:A:352:ILE:HG23	1:A:363:LEU:HD13	1.90	0.54
1:A:335:ARG:HD2	4:A:1113:HOH:O	2.07	0.54
1:E:352:ILE:HG23	1:E:363:LEU:HD13	1.89	0.54
1:A:355:THR:CG2	1:A:357:VAL:H	2.17	0.54
1:C:393:GLY:N	4:C:804:HOH:O	2.42	0.53
1:H:335:ARG:HG2	4:H:904:HOH:O	2.08	0.53
1:F:68:ALA:HA	4:F:990:HOH:O	2.09	0.53
1:F:355:THR:CG2	1:F:357:VAL:H	2.17	0.52
1:A:457:THR:HG23	4:A:1159:HOH:O	2.08	0.52
1:E:90:GLU:HG3	4:E:1115:HOH:O	2.09	0.52
1:B:68:ALA:HA	4:B:987:HOH:O	2.10	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:355:THR:CG2	1:B:357:VAL:H	2.17	0.51
1:D:355:THR:CG2	1:D:357:VAL:H	2.17	0.51
1:B:352:ILE:HG23	1:B:363:LEU:HD13	1.92	0.51
1:C:355:THR:CG2	1:C:357:VAL:H	2.18	0.51
1:E:268:PRO:HG3	1:E:285:LEU:HD22	1.93	0.51
1:F:268:PRO:HG3	1:F:285:LEU:HD22	1.93	0.50
1:C:268:PRO:HG3	1:C:285:LEU:HD22	1.94	0.50
1:E:193:ARG:HG2	1:E:193:ARG:HH11	1.77	0.50
1:G:398:ARG:HG2	4:G:802:HOH:O	2.11	0.50
1:B:133:LEU:HD13	1:B:166:LEU:HD22	1.93	0.50
1:A:133:LEU:HD13	1:A:166:LEU:HD22	1.93	0.50
1:D:268:PRO:HG3	1:D:285:LEU:HD22	1.94	0.49
1:F:133:LEU:HD13	1:F:166:LEU:HD22	1.93	0.49
1:A:268:PRO:HG3	1:A:285:LEU:HD22	1.94	0.49
1:B:268:PRO:HG3	1:B:285:LEU:HD22	1.94	0.49
1:G:398:ARG:CG	4:G:802:HOH:O	2.60	0.49
1:F:352:ILE:HG23	1:F:363:LEU:HD13	1.93	0.49
1:C:398:ARG:HB2	4:C:1092:HOH:O	2.13	0.49
1:H:63:GLU:OE1	3:H:702:SAR:HN2	2.13	0.49
1:E:335:ARG:HD3	4:E:1114:HOH:O	2.13	0.48
1:A:193:ARG:HH11	1:A:193:ARG:HG2	1.79	0.48
1:F:319:ASN:HB3	4:F:1074:HOH:O	2.13	0.48
1:H:268:PRO:HG3	1:H:285:LEU:HD22	1.94	0.48
1:G:55:GLY:N	1:G:56:ALA:HA	2.29	0.48
1:A:123:ALA:HB1	1:A:418:ARG:HD2	1.96	0.48
1:B:431:LEU:HD11	1:D:431:LEU:CD1	2.44	0.48
1:E:25:VAL:HG22	1:E:378:ALA:HA	1.96	0.48
1:C:55:GLY:N	1:C:56:ALA:HA	2.28	0.48
1:E:123:ALA:HB1	1:E:418:ARG:HD2	1.96	0.48
1:H:133:LEU:HD13	1:H:166:LEU:HD22	1.95	0.48
2:H:701:FAD:H1'1	2:H:701:FAD:O4'	2.14	0.47
1:D:133:LEU:HD13	1:D:166:LEU:HD22	1.96	0.47
1:F:431:LEU:HD11	1:H:431:LEU:CD1	2.44	0.47
1:A:25:VAL:HG22	1:A:378:ALA:HA	1.95	0.47
1:G:268:PRO:HG3	1:G:285:LEU:HD22	1.95	0.47
1:B:355:THR:HG22	1:B:357:VAL:N	2.22	0.47
1:A:346:ILE:HG13	1:A:464:LEU:HD12	1.96	0.47
1:D:335:ARG:HD2	4:D:1035:HOH:O	2.13	0.47
1:E:133:LEU:HD13	1:E:166:LEU:HD22	1.95	0.47
1:A:64:VAL:HG13	1:A:69:LEU:HD22	1.96	0.47
1:C:379:ARG:HD3	4:C:915:HOH:O	2.15	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:291:GLY:N	4:A:802:HOH:O	2.35	0.47
2:D:701:FAD:H9	2:D:701:FAD:H1'2	1.74	0.47
1:A:355:THR:HG22	1:A:357:VAL:N	2.22	0.46
1:C:67:HIS:HD2	4:C:1112:HOH:O	1.98	0.46
1:C:362:MET:HE2	1:C:362:MET:HB3	1.81	0.46
1:D:63:GLU:OE1	3:D:702:SAR:HN2	2.16	0.46
1:A:72:GLU:HG3	4:A:1037:HOH:O	2.14	0.46
1:B:431:LEU:CD1	1:D:431:LEU:HD11	2.46	0.46
1:E:64:VAL:HG13	1:E:69:LEU:HD22	1.97	0.46
1:B:352:ILE:HG23	1:B:363:LEU:CD1	2.46	0.46
2:D:701:FAD:H1'1	2:D:701:FAD:O4'	2.15	0.46
1:A:415:ASP:CB	4:A:936:HOH:O	2.51	0.46
1:B:64:VAL:HG13	1:B:69:LEU:HD22	1.98	0.46
1:D:355:THR:HG22	1:D:357:VAL:N	2.22	0.46
1:G:355:THR:HG22	1:G:357:VAL:N	2.22	0.45
1:F:431:LEU:CD1	1:H:431:LEU:HD11	2.46	0.45
1:G:352:ILE:HG23	1:G:363:LEU:CD1	2.46	0.45
1:H:352:ILE:HG23	1:H:363:LEU:CD1	2.47	0.45
1:C:352:ILE:HG23	1:C:363:LEU:CD1	2.46	0.45
1:C:20:VAL:HG11	1:C:371:LEU:HA	1.99	0.45
1:D:352:ILE:HG23	1:D:363:LEU:CD1	2.47	0.45
1:D:76:LYS:HE2	4:D:966:HOH:O	2.17	0.45
1:F:64:VAL:HG13	1:F:69:LEU:HD22	1.98	0.45
1:E:437:HIS:O	1:E:441:THR:HG23	2.17	0.45
1:E:368:ARG:NH2	3:E:702:SAR:HA3	2.32	0.44
1:F:352:ILE:HG23	1:F:363:LEU:CD1	2.47	0.44
1:H:352:ILE:HG23	1:H:363:LEU:HD13	2.00	0.44
1:A:437:HIS:O	1:A:441:THR:HG23	2.17	0.44
1:A:453:ARG:CG	4:A:1078:HOH:O	2.65	0.44
1:E:393:GLY:C	4:E:803:HOH:O	2.61	0.44
1:E:473:ARG:HD3	4:E:1091:HOH:O	2.17	0.44
1:G:336:LYS:HG3	4:G:1079:HOH:O	2.17	0.44
1:D:352:ILE:HG23	1:D:363:LEU:HD13	2.00	0.44
1:E:218:ASP:CB	4:E:1087:HOH:O	2.66	0.44
1:E:218:ASP:HB2	4:E:1087:HOH:O	2.17	0.44
1:G:437:HIS:O	1:G:441:THR:HG23	2.18	0.44
2:B:701:FAD:H9	2:B:701:FAD:H1'2	1.75	0.44
1:C:437:HIS:O	1:C:441:THR:HG23	2.18	0.44
1:F:244:ALA:CB	4:F:1039:HOH:O	2.66	0.44
1:D:149:GLU:HG3	4:D:863:HOH:O	2.17	0.43
1:G:379:ARG:HD3	4:G:951:HOH:O	2.18	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:701:FAD:H1'2	2:H:701:FAD:H9	1.74	0.43
1:G:351:LEU:O	1:G:352:ILE:HD13	2.19	0.43
1:C:351:LEU:O	1:C:352:ILE:HD13	2.18	0.43
1:G:363:LEU:HD12	1:G:363:LEU:HA	1.84	0.43
1:C:13:VAL:HG11	1:C:29:ILE:HD13	2.01	0.43
1:D:90:GLU:HG3	4:D:825:HOH:O	2.17	0.43
1:F:437:HIS:O	1:F:441:THR:HG23	2.19	0.43
1:E:352:ILE:HG23	1:E:363:LEU:CD1	2.48	0.43
1:G:64:VAL:HG13	1:G:69:LEU:HD22	2.01	0.43
1:H:355:THR:HG22	1:H:357:VAL:N	2.22	0.43
1:B:76:LYS:HE3	4:B:1118:HOH:O	2.18	0.42
1:D:280:ALA:HB1	1:D:368:ARG:HD2	2.00	0.42
2:D:701:FAD:O4'	2:D:701:FAD:C1'	2.66	0.42
1:E:193:ARG:HH11	1:E:193:ARG:CG	2.32	0.42
1:G:20:VAL:HG11	1:G:371:LEU:HA	2.00	0.42
1:A:473:ARG:HD3	4:A:1093:HOH:O	2.19	0.42
1:B:437:HIS:O	1:B:441:THR:HG23	2.19	0.42
1:B:123:ALA:HB1	1:B:418:ARG:HD2	2.01	0.42
2:E:701:FAD:H1'2	2:E:701:FAD:H9	1.73	0.42
1:G:13:VAL:HG11	1:G:29:ILE:HD13	2.02	0.42
1:F:435:TRP:CE3	1:H:439:MET:HE2	2.54	0.42
1:H:216:ASP:C	4:H:907:HOH:O	2.62	0.42
1:A:193:ARG:HH11	1:A:193:ARG:CG	2.32	0.42
1:A:352:ILE:HG23	1:A:363:LEU:CD1	2.48	0.42
1:F:123:ALA:HB1	1:F:418:ARG:HD2	2.02	0.42
1:C:336:LYS:HG3	4:C:1114:HOH:O	2.19	0.42
1:D:379:ARG:HD3	4:D:993:HOH:O	2.19	0.42
1:F:439[A]:MET:HE2	1:H:435:TRP:CE3	2.55	0.42
1:H:280:ALA:HB1	1:H:368:ARG:HD2	2.00	0.42
1:D:437:HIS:O	1:D:441:THR:HG23	2.19	0.42
1:B:20:VAL:HG11	1:B:371:LEU:HA	2.01	0.42
1:B:435:TRP:CE3	1:D:439:MET:HE2	2.55	0.42
1:C:64:VAL:HG13	1:C:69:LEU:HD22	2.01	0.42
1:A:280:ALA:HB1	1:A:368:ARG:HD2	2.02	0.41
1:B:431:LEU:CD1	1:D:431:LEU:CD1	2.98	0.41
1:E:11:THR:O	1:E:36:VAL:HG22	2.20	0.41
1:D:28:GLU:HB3	1:D:382:VAL:HG21	2.02	0.41
1:E:439:MET:HE2	1:G:435:TRP:CE3	2.55	0.41
1:H:437:HIS:O	1:H:441:THR:HG23	2.19	0.41
2:A:701:FAD:C4	3:A:702:SAR:HN3	2.50	0.41
1:B:280:ALA:HB1	1:B:368:ARG:HD2	2.02	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:20:VAL:HG11	1:F:371:LEU:HA	2.01	0.41
1:G:362:MET:HE2	1:G:362:MET:HB3	1.82	0.41
2:H:701:FAD:O4'	2:H:701:FAD:C1'	2.69	0.41
2:A:701:FAD:H1'2	2:A:701:FAD:H9	1.72	0.41
1:F:76:LYS:HE3	4:F:1146:HOH:O	2.20	0.41
1:F:431:LEU:CD1	1:H:431:LEU:CD1	2.98	0.41
1:G:280:ALA:HB1	1:G:368:ARG:HD2	2.02	0.41
1:A:105:ARG:HD3	4:A:842:HOH:O	2.19	0.41
1:B:429:TRP:CG	1:B:431:LEU:HD13	2.55	0.41
1:C:280:ALA:HB1	1:C:368:ARG:HD2	2.02	0.41
1:D:55:GLY:N	1:D:56:ALA:HA	2.35	0.41
1:D:363:LEU:HD12	1:D:363:LEU:HA	1.89	0.41
1:E:20:VAL:HG11	1:E:371:LEU:HA	2.01	0.41
1:F:37:ARG:NH1	4:F:836:HOH:O	2.54	0.41
1:H:28:GLU:HB3	1:H:382:VAL:HG21	2.03	0.41
1:H:55:GLY:N	1:H:56:ALA:HA	2.34	0.41
1:A:435:TRP:CE3	1:C:439[A]:MET:HE2	2.56	0.41
1:H:429:TRP:CG	1:H:431:LEU:HD13	2.56	0.41
1:A:453:ARG:HD2	4:A:1078:HOH:O	2.20	0.41
1:C:21:LEU:HD23	1:C:21:LEU:HA	1.90	0.41
1:E:280:ALA:HB1	1:E:368:ARG:HD2	2.02	0.41
1:F:363:LEU:HD12	1:F:363:LEU:HA	1.93	0.41
1:H:20:VAL:HG11	1:H:371:LEU:HA	2.03	0.41
1:D:429:TRP:CG	1:D:431:LEU:HD13	2.56	0.41
1:B:439[A]:MET:HE2	1:D:435:TRP:CE3	2.56	0.40
1:C:69:LEU:HD23	4:C:920:HOH:O	2.21	0.40
1:D:20:VAL:HG11	1:D:371:LEU:HA	2.03	0.40
1:F:429:TRP:CG	1:F:431:LEU:HD13	2.56	0.40
2:H:701:FAD:N1	2:H:701:FAD:O2'	2.44	0.40
1:A:11:THR:O	1:A:36:VAL:HG22	2.21	0.40
1:A:20:VAL:HG11	1:A:371:LEU:HA	2.02	0.40
1:B:28:GLU:HB3	1:B:382:VAL:HG21	2.04	0.40
1:E:429:TRP:CG	1:E:431:LEU:HD13	2.57	0.40
1:C:193:ARG:HA	4:C:965:HOH:O	2.20	0.40
1:C:285:LEU:HD23	1:C:285:LEU:C	2.46	0.40
1:G:21:LEU:HD23	1:G:21:LEU:HA	1.91	0.40
1:A:429:TRP:CG	1:A:431:LEU:HD13	2.56	0.40
1:A:439:MET:HE2	1:C:435:TRP:CE3	2.56	0.40
1:B:431:LEU:HD11	1:D:431:LEU:HD11	2.03	0.40
1:B:479:LEU:HD22	4:B:872:HOH:O	2.22	0.40
1:H:398:ARG:HB2	4:H:1001:HOH:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	480/512 (94%)	470 (98%)	10 (2%)	0	100	100
1	B	481/512 (94%)	470 (98%)	11 (2%)	0	100	100
1	C	481/512 (94%)	470 (98%)	11 (2%)	0	100	100
1	D	477/512 (93%)	466 (98%)	11 (2%)	0	100	100
1	E	480/512 (94%)	469 (98%)	11 (2%)	0	100	100
1	F	481/512 (94%)	471 (98%)	10 (2%)	0	100	100
1	G	480/512 (94%)	469 (98%)	11 (2%)	0	100	100
1	H	477/512 (93%)	466 (98%)	11 (2%)	0	100	100
All	All	3837/4096 (94%)	3751 (98%)	86 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	376/400 (94%)	366 (97%)	10 (3%)	39	31
1	B	377/400 (94%)	368 (98%)	9 (2%)	43	35
1	C	377/400 (94%)	367 (97%)	10 (3%)	39	31
1	D	373/400 (93%)	359 (96%)	14 (4%)	29	19

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	376/400 (94%)	365 (97%)	11 (3%)	37	28
1	F	377/400 (94%)	369 (98%)	8 (2%)	47	40
1	G	376/400 (94%)	366 (97%)	10 (3%)	39	31
1	H	373/400 (93%)	360 (96%)	13 (4%)	32	21
All	All	3005/3200 (94%)	2920 (97%)	85 (3%)	38	29

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	LEU
1	A	77	LYS
1	A	133	LEU
1	A	193	ARG
1	A	203	ARG
1	A	335	ARG
1	A	355	THR
1	A	452	ASP
1	A	457	THR
1	A	464	LEU
1	B	77	LYS
1	B	133	LEU
1	B	203	ARG
1	B	335	ARG
1	B	355	THR
1	B	395	ARG
1	B	452	ASP
1	B	457	THR
1	B	464	LEU
1	C	38	VAL
1	C	41	LEU
1	C	77	LYS
1	C	203	ARG
1	C	281	CYS
1	C	335	ARG
1	C	355	THR
1	C	431	LEU
1	C	452	ASP
1	C	457	THR
1	D	37	ARG
1	D	38	VAL
1	D	41	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	77	LYS
1	D	85	GLN
1	D	90	GLU
1	D	133	LEU
1	D	203	ARG
1	D	243	LEU
1	D	335	ARG
1	D	355	THR
1	D	395	ARG
1	D	452	ASP
1	D	457	THR
1	E	41	LEU
1	E	77	LYS
1	E	133	LEU
1	E	193	ARG
1	E	203	ARG
1	E	281	CYS
1	E	335	ARG
1	E	355	THR
1	E	452	ASP
1	E	457	THR
1	E	464	LEU
1	F	77	LYS
1	F	133	LEU
1	F	203	ARG
1	F	335	ARG
1	F	355	THR
1	F	395	ARG
1	F	452	ASP
1	F	457	THR
1	G	38	VAL
1	G	41	LEU
1	G	77	LYS
1	G	203	ARG
1	G	281	CYS
1	G	335	ARG
1	G	355	THR
1	G	431	LEU
1	G	452	ASP
1	G	457	THR
1	H	37	ARG
1	H	38	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	H	41	LEU
1	H	77	LYS
1	H	85	GLN
1	H	133	LEU
1	H	203	ARG
1	H	319	ASN
1	H	335	ARG
1	H	355	THR
1	H	395	ARG
1	H	452	ASP
1	H	457	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	HIS
1	A	437	HIS
1	B	10	GLN
1	B	67	HIS
1	B	159	ASN
1	B	437	HIS
1	C	10	GLN
1	C	67	HIS
1	C	159	ASN
1	C	338	GLN
1	C	437	HIS
1	D	159	ASN
1	D	319	ASN
1	D	437	HIS
1	D	491	HIS
1	E	159	ASN
1	E	437	HIS
1	F	10	GLN
1	F	159	ASN
1	F	437	HIS
1	G	10	GLN
1	G	67	HIS
1	G	159	ASN
1	G	319	ASN
1	G	437	HIS
1	H	159	ASN
1	H	319	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	H	437	HIS
1	H	491	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

16 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	FAD	B	701	-	58,58,58	0.80	2 (3%)	85,89,89	0.85	3 (3%)
3	SAR	F	702	-	5,5,5	1.21	0	3,5,5	0.38	0
3	SAR	H	702	-	5,5,5	0.87	0	3,5,5	1.75	1 (33%)
3	SAR	D	702	-	5,5,5	0.87	0	3,5,5	2.02	1 (33%)
3	SAR	B	702	-	5,5,5	1.22	0	3,5,5	0.53	0
2	FAD	G	701	-	58,58,58	0.67	1 (1%)	85,89,89	0.76	1 (1%)
2	FAD	E	701	-	58,58,58	0.81	2 (3%)	85,89,89	0.95	3 (3%)
2	FAD	D	701	-	58,58,58	0.70	2 (3%)	85,89,89	0.70	2 (2%)
3	SAR	C	702	-	5,5,5	1.04	0	3,5,5	1.74	2 (66%)
3	SAR	E	702	-	5,5,5	0.82	0	3,5,5	1.19	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FAD	F	701	-	58,58,58	0.74	1 (1%)	85,89,89	0.86	3 (3%)
3	SAR	A	702	-	5,5,5	0.77	0	3,5,5	1.31	0
2	FAD	H	701	-	58,58,58	0.72	1 (1%)	85,89,89	0.68	2 (2%)
2	FAD	A	701	-	58,58,58	0.79	2 (3%)	85,89,89	0.95	4 (4%)
3	SAR	G	702	-	5,5,5	1.08	0	3,5,5	1.60	0
2	FAD	C	701	-	58,58,58	0.65	1 (1%)	85,89,89	0.76	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
2	FAD	B	701	-	-	3/34/50/50	0/6/6/6
3	SAR	F	702	-	-	0/2/3/3	-
3	SAR	H	702	-	-	0/2/3/3	-
3	SAR	D	702	-	-	0/2/3/3	-
3	SAR	B	702	-	-	0/2/3/3	-
2	FAD	G	701	-	-	3/34/50/50	0/6/6/6
2	FAD	E	701	-	-	2/34/50/50	0/6/6/6
2	FAD	D	701	-	-	3/34/50/50	0/6/6/6
3	SAR	C	702	-	-	0/2/3/3	-
3	SAR	E	702	-	-	0/2/3/3	-
2	FAD	F	701	-	-	3/34/50/50	0/6/6/6
3	SAR	A	702	-	-	0/2/3/3	-
2	FAD	H	701	-	-	3/34/50/50	0/6/6/6
2	FAD	A	701	-	-	3/34/50/50	0/6/6/6
3	SAR	G	702	-	-	0/2/3/3	-
2	FAD	C	701	-	-	1/34/50/50	0/6/6/6

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	FAD	PA-O3P	2.61	1.62	1.59
2	B	701	FAD	PA-O3P	2.52	1.62	1.59
2	B	701	FAD	C9-C8	-2.46	1.36	1.39
2	F	701	FAD	C9-C8	-2.45	1.36	1.39
2	E	701	FAD	C9-C8	-2.38	1.36	1.39
2	H	701	FAD	PA-O3P	2.37	1.62	1.59
2	C	701	FAD	PA-O3P	2.25	1.61	1.59

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	701	FAD	PA-O3P	2.16	1.61	1.59
2	D	701	FAD	P-O3P	2.06	1.61	1.59
2	E	701	FAD	PA-O3P	2.05	1.61	1.59
2	D	701	FAD	PA-O3P	2.04	1.61	1.59
2	A	701	FAD	C9-C8	-2.04	1.36	1.39

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	701	FAD	C4'-C3'-C2'	3.69	119.71	113.57
2	A	701	FAD	O2A-PA-O1A	3.30	127.82	112.44
2	E	701	FAD	O2A-PA-O1A	3.28	127.71	112.44
2	A	701	FAD	C4'-C3'-C2'	3.14	118.80	113.57
2	F	701	FAD	O2A-PA-O1A	3.02	126.50	112.44
2	B	701	FAD	O2A-PA-O1A	2.98	126.32	112.44
3	D	702	SAR	OXT-C-CA	2.75	123.27	112.81
2	A	701	FAD	O3P-PA-O1A	-2.61	102.84	110.70
2	D	701	FAD	O2P-P-O1P	2.45	123.85	112.44
2	F	701	FAD	C4'-C3'-C2'	2.45	117.64	113.57
2	E	701	FAD	O3P-PA-O1A	-2.40	103.49	110.70
3	H	702	SAR	OXT-C-CA	2.40	121.92	112.81
2	H	701	FAD	O2P-P-O1P	2.29	123.12	112.44
2	B	701	FAD	O3P-PA-O1A	-2.22	104.03	110.70
3	C	702	SAR	OXT-C-CA	2.13	120.90	112.81
2	G	701	FAD	C4-N3-C2	-2.12	121.87	125.64
3	C	702	SAR	OXT-C-O	-2.11	117.92	123.33
2	C	701	FAD	C4-N3-C2	-2.09	121.94	125.64
2	C	701	FAD	O2P-P-O1P	2.08	122.13	112.44
2	H	701	FAD	C4-N3-C2	-2.04	122.02	125.64
2	D	701	FAD	C4-N3-C2	-2.02	122.05	125.64
2	B	701	FAD	C4-N3-C2	-2.02	122.06	125.64
2	A	701	FAD	O4'-C4'-C3'	2.02	113.96	109.25
2	F	701	FAD	O3P-PA-O1A	-2.00	104.69	110.70

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	701	FAD	P-O3P-PA-O5B
2	A	701	FAD	C2B-C1B-N9A-C8A
2	B	701	FAD	C2B-C1B-N9A-C8A
2	D	701	FAD	C2B-C1B-N9A-C8A

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
2	F	701	FAD	C2B-C1B-N9A-C8A
2	G	701	FAD	C2B-C1B-N9A-C8A
2	H	701	FAD	C2B-C1B-N9A-C8A
2	D	701	FAD	C2'-C1'-N10-C10
2	H	701	FAD	C2'-C1'-N10-C10
2	B	701	FAD	O4B-C4B-C5B-O5B
2	E	701	FAD	O4B-C4B-C5B-O5B
2	C	701	FAD	C2B-C1B-N9A-C8A
2	E	701	FAD	C2B-C1B-N9A-C8A
2	A	701	FAD	O4B-C4B-C5B-O5B
2	F	701	FAD	O4B-C4B-C5B-O5B
2	F	701	FAD	O4B-C1B-N9A-C8A
2	H	701	FAD	O4B-C1B-N9A-C8A
2	D	701	FAD	O4B-C4B-C5B-O5B
2	B	701	FAD	O4B-C1B-N9A-C8A
2	G	701	FAD	N10-C1'-C2'-O2'
2	A	701	FAD	O4B-C1B-N9A-C8A

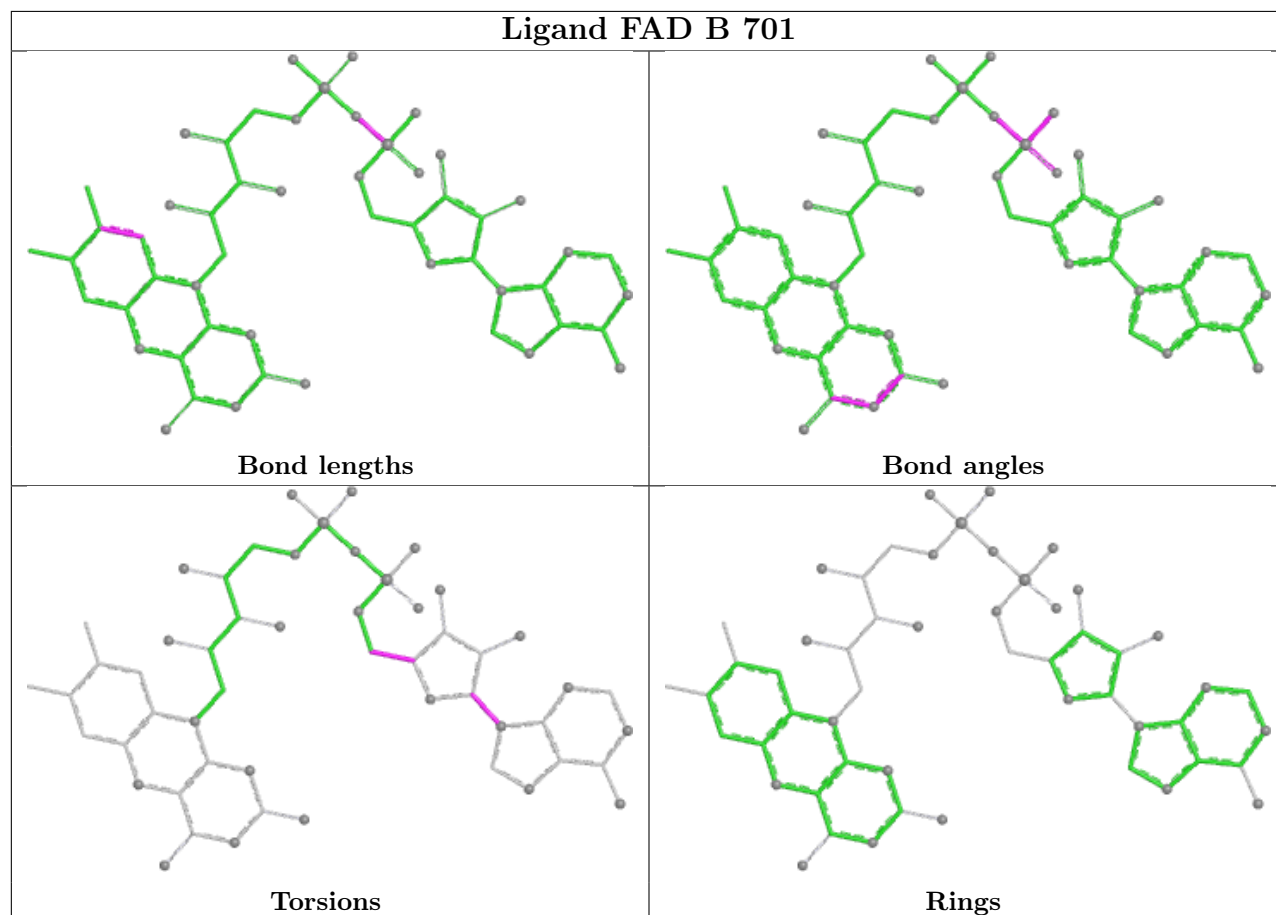
There are no ring outliers.

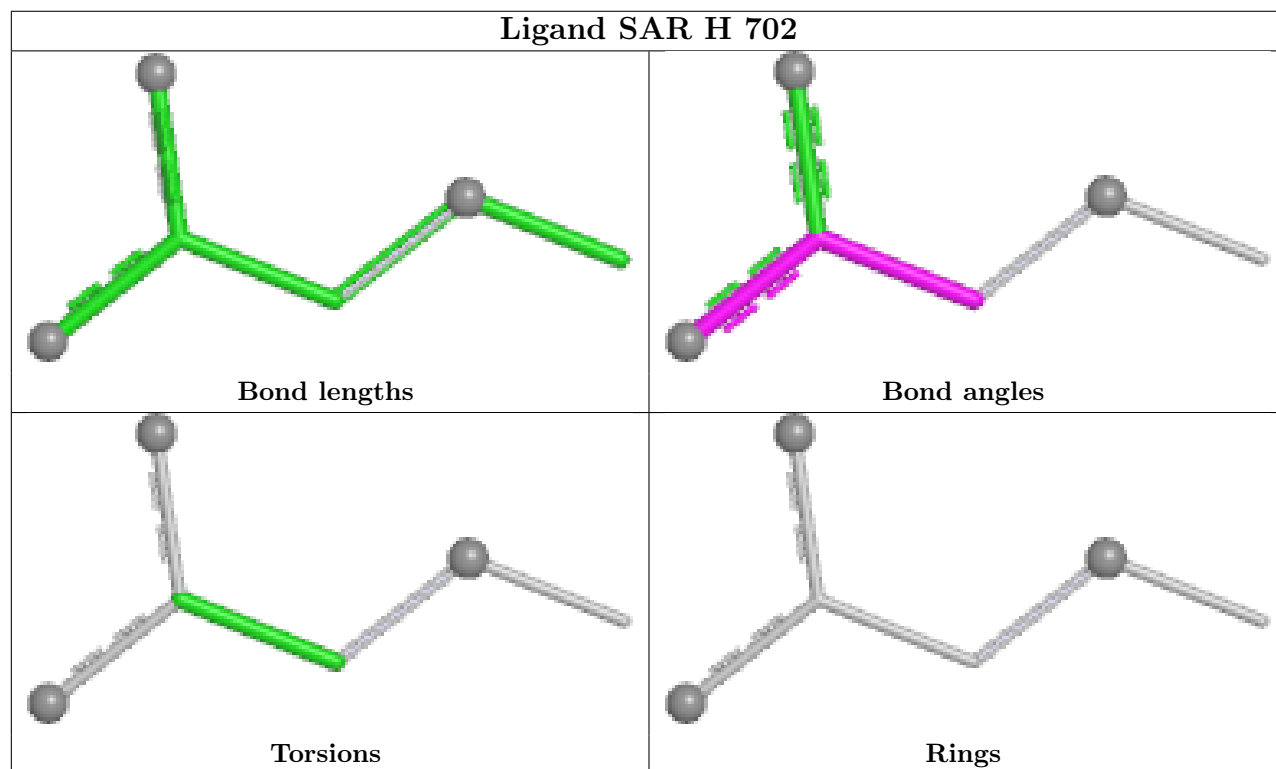
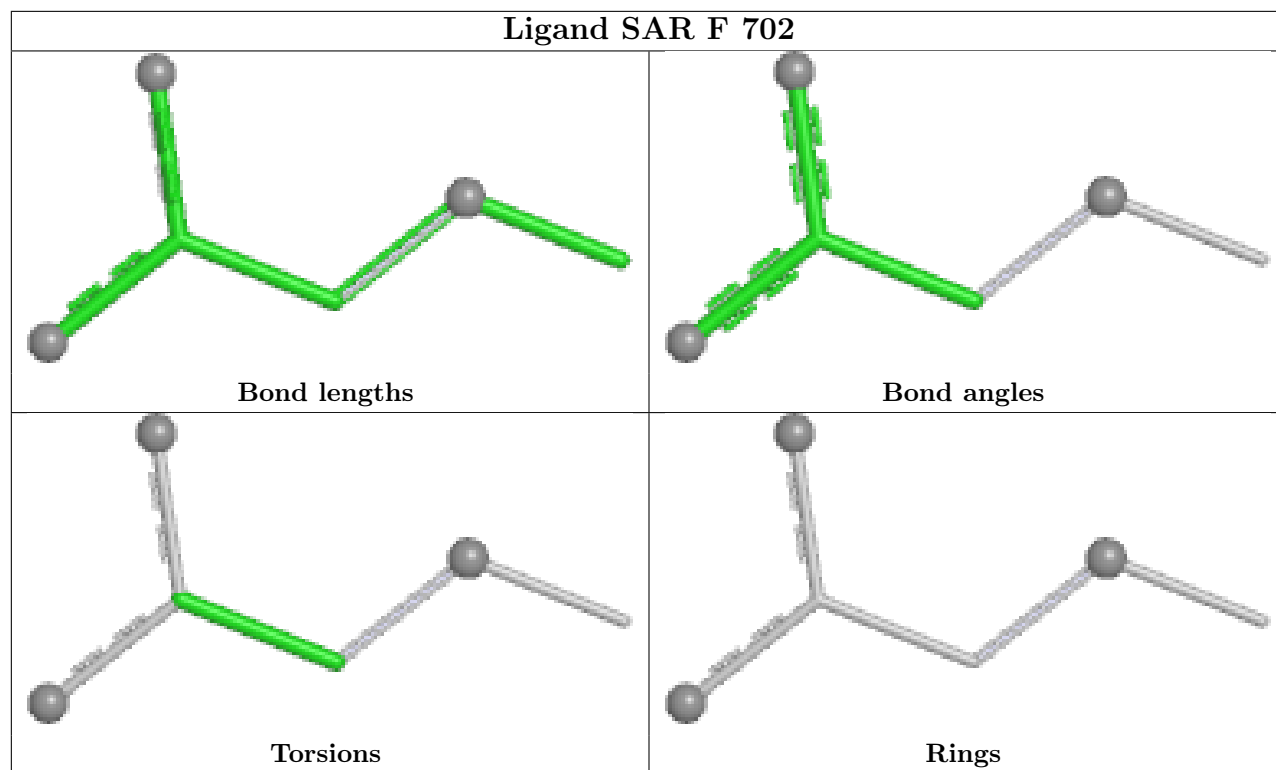
10 monomers are involved in 15 short contacts:

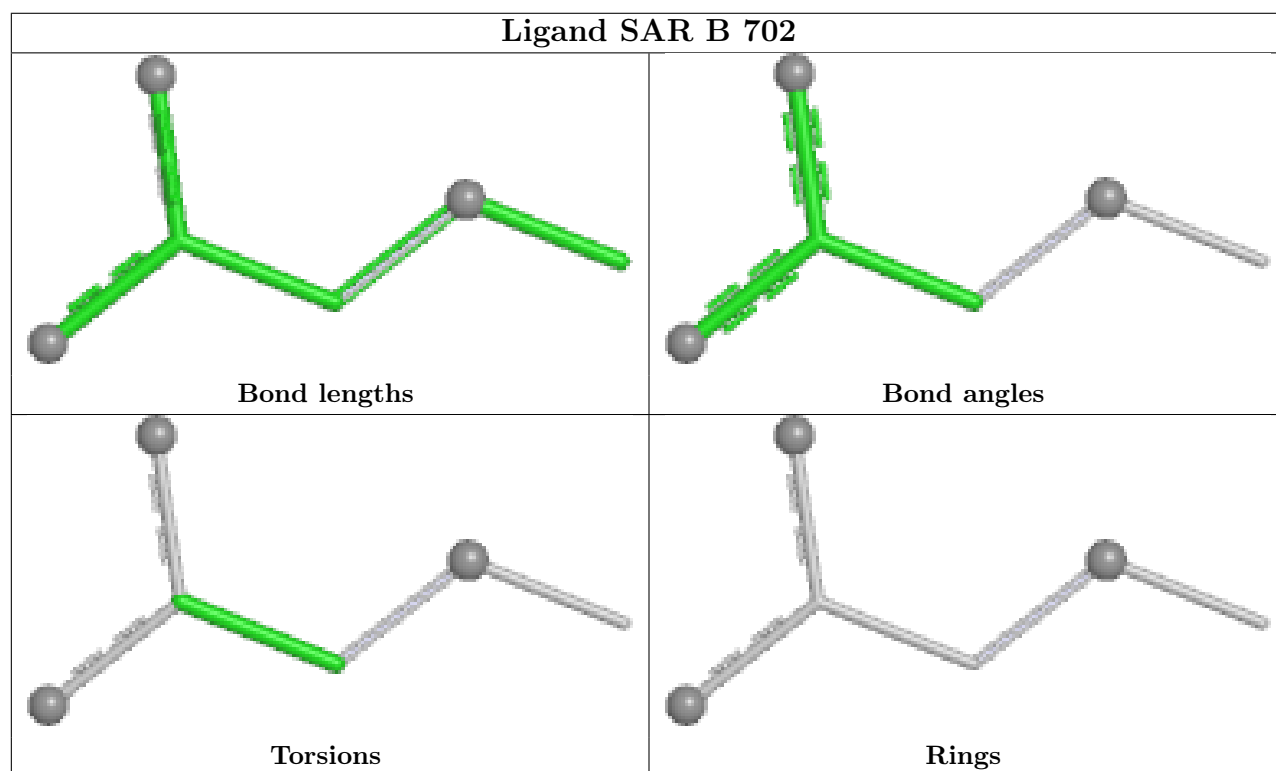
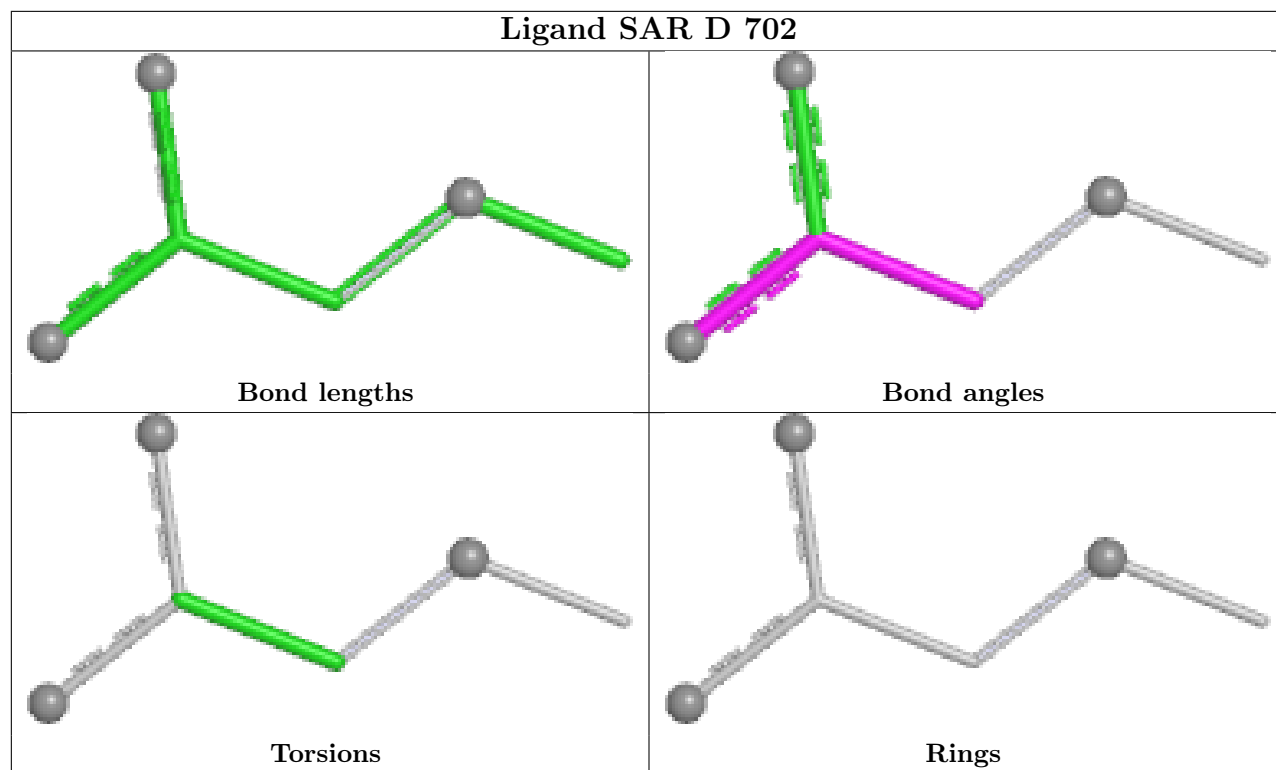
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	701	FAD	1	0
3	H	702	SAR	1	0
3	D	702	SAR	1	0
2	G	701	FAD	1	0
2	E	701	FAD	1	0
2	D	701	FAD	3	0
3	E	702	SAR	1	0
3	A	702	SAR	1	0
2	H	701	FAD	4	0
2	A	701	FAD	2	0

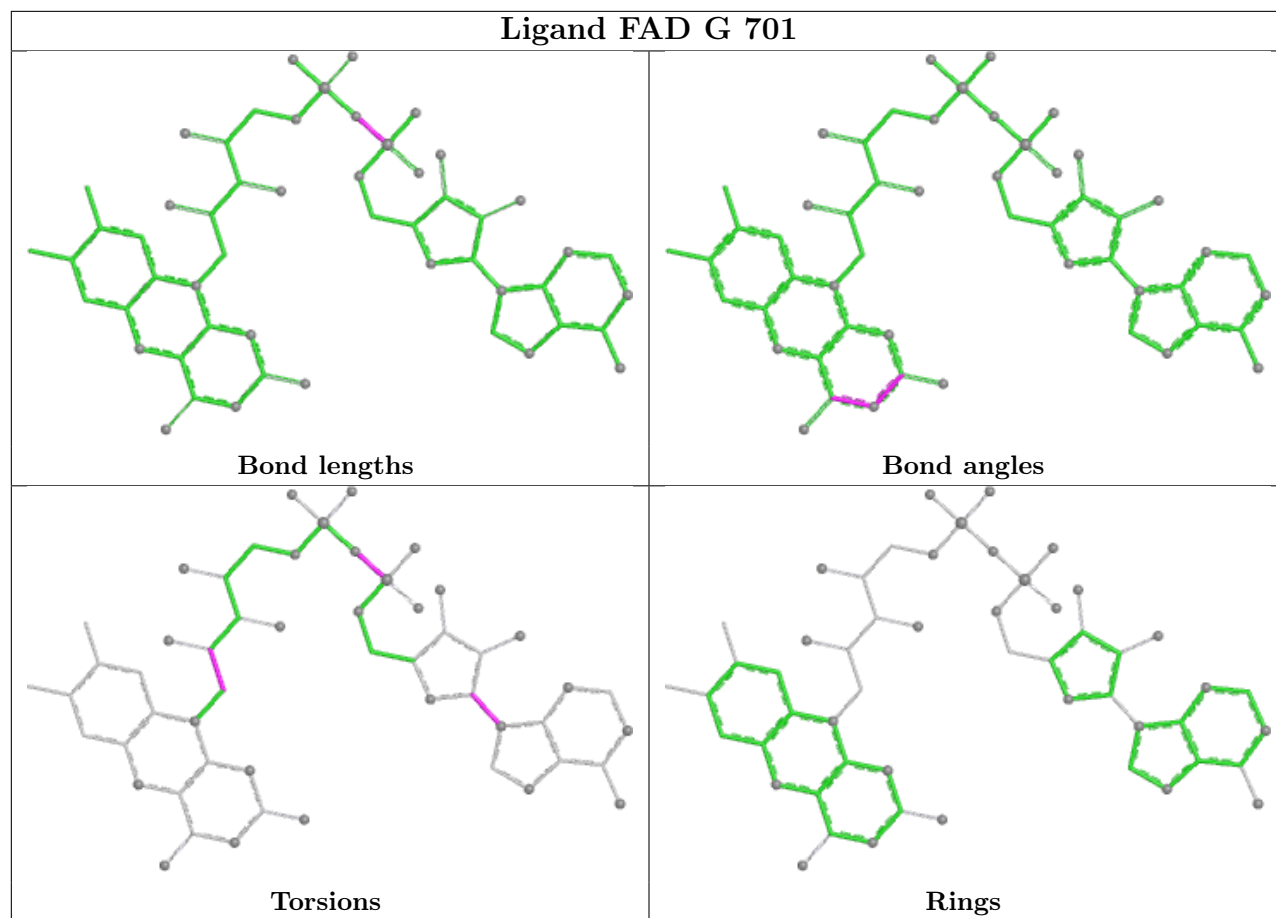
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

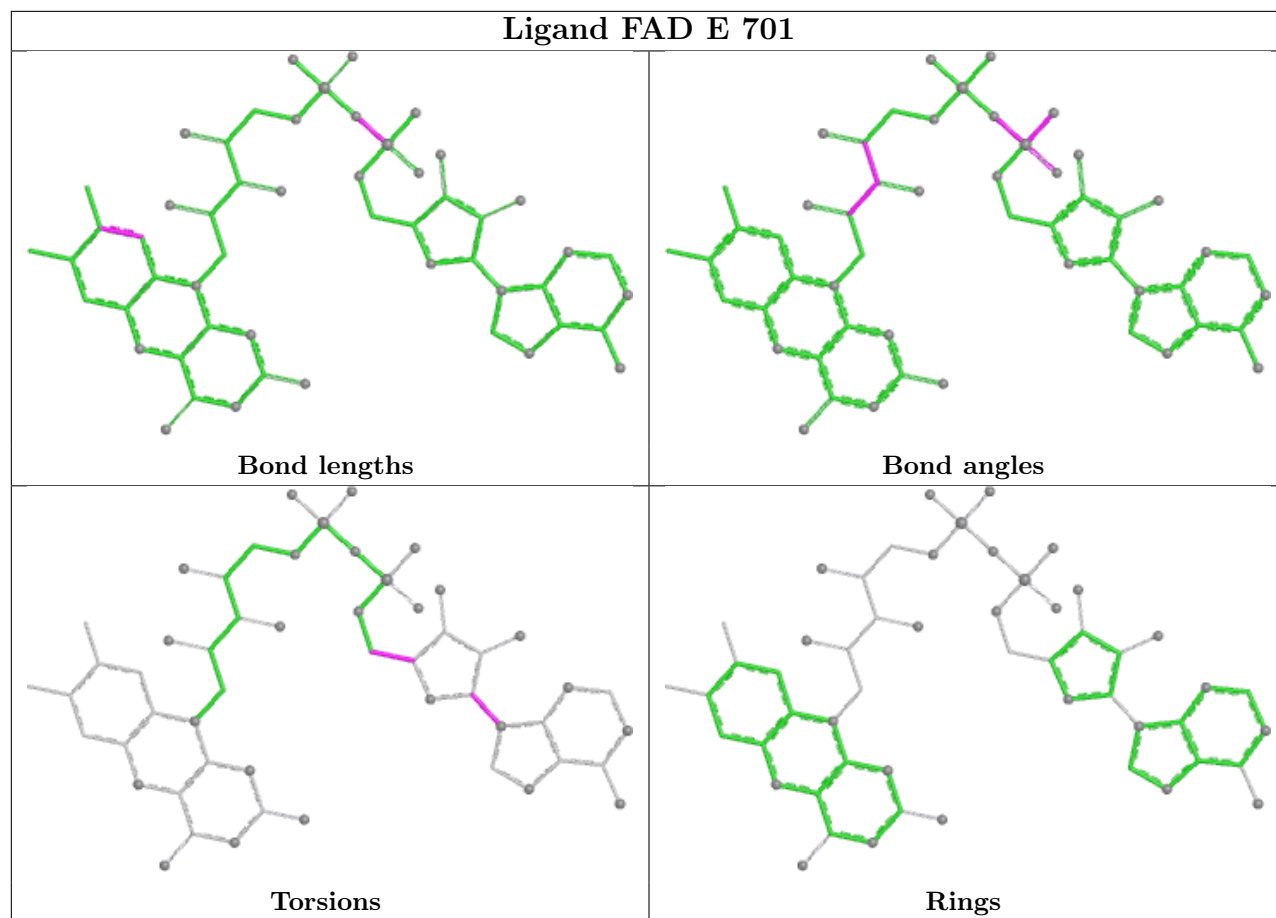
any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



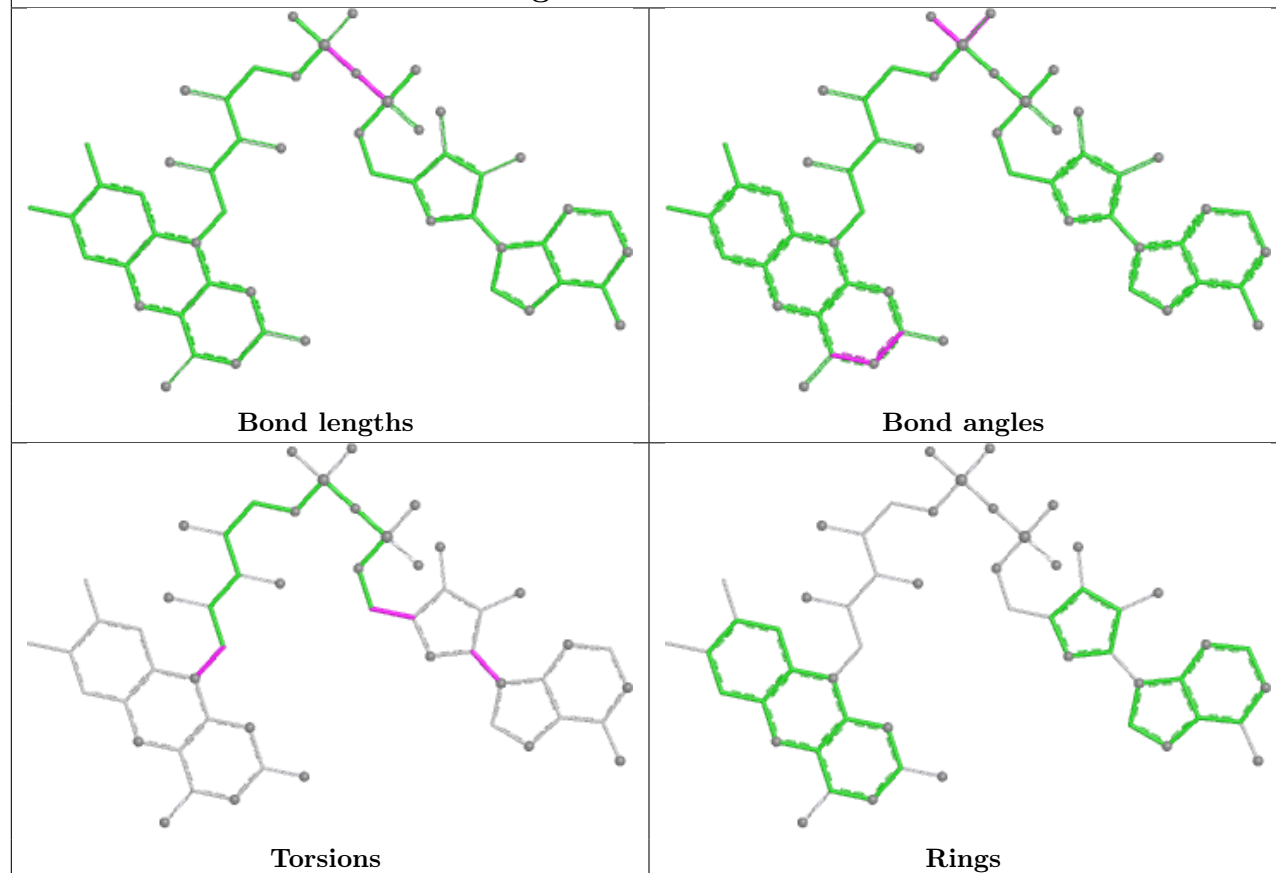




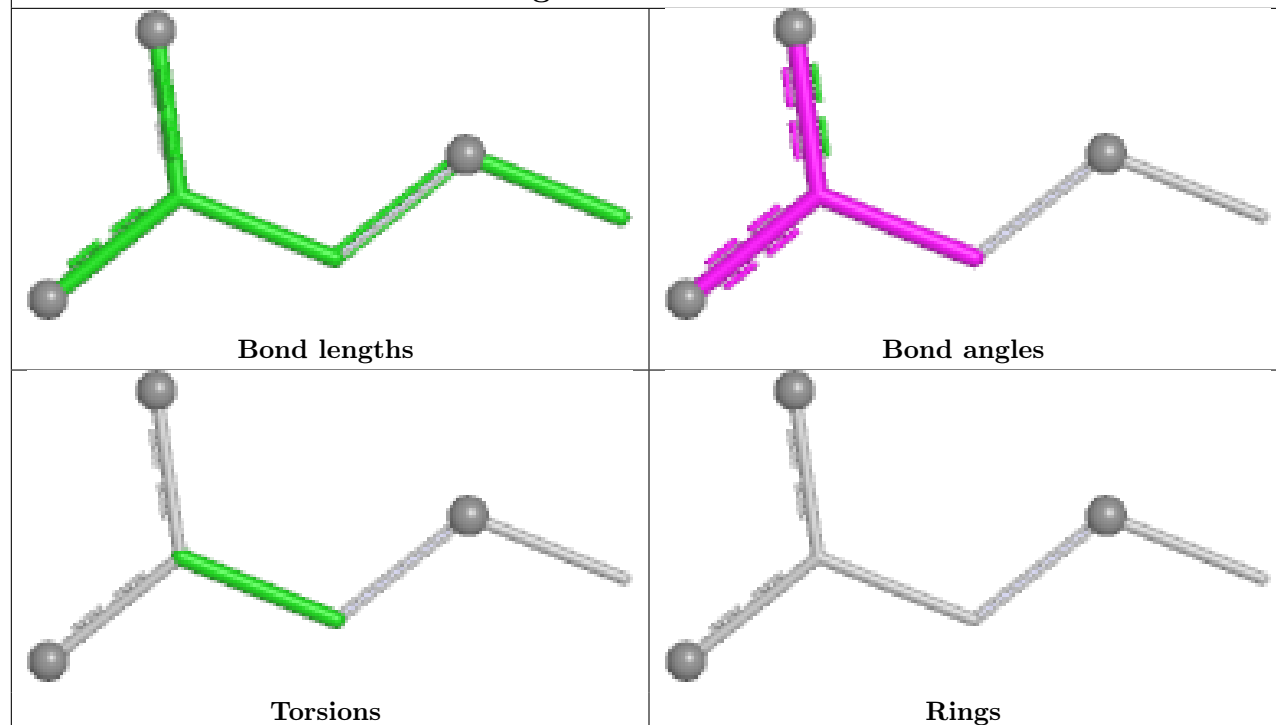




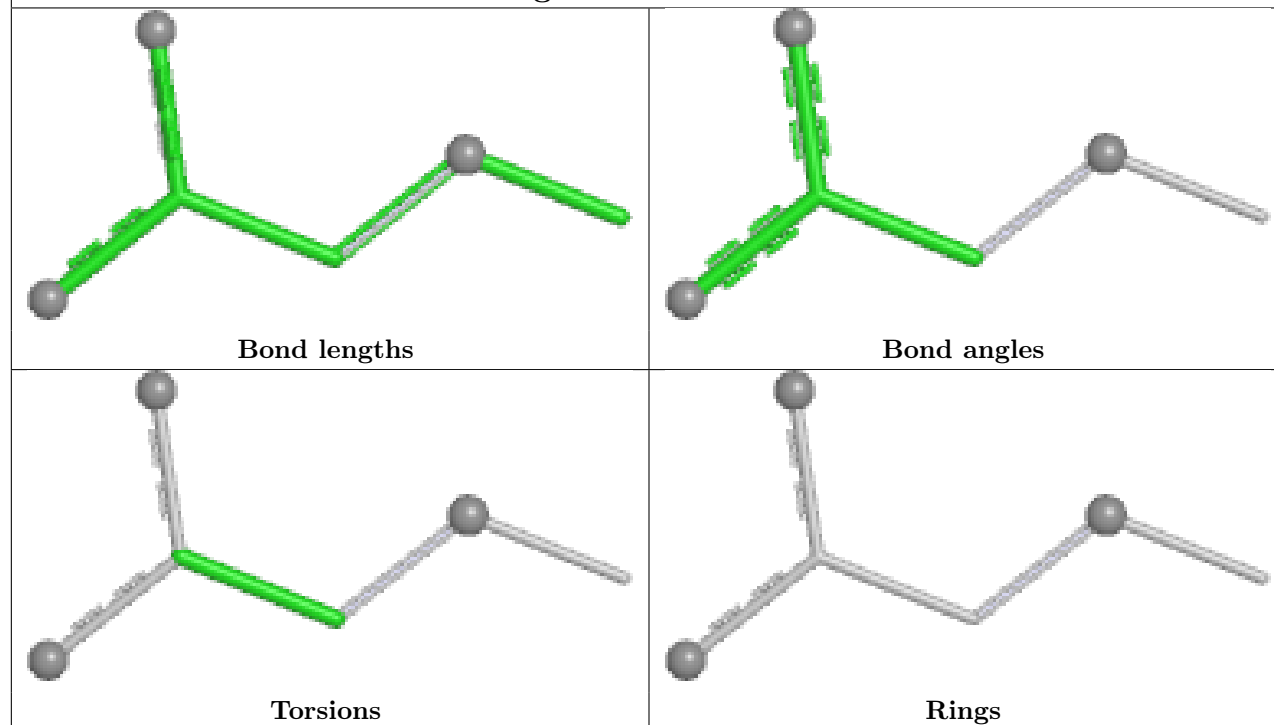
Ligand FAD D 701



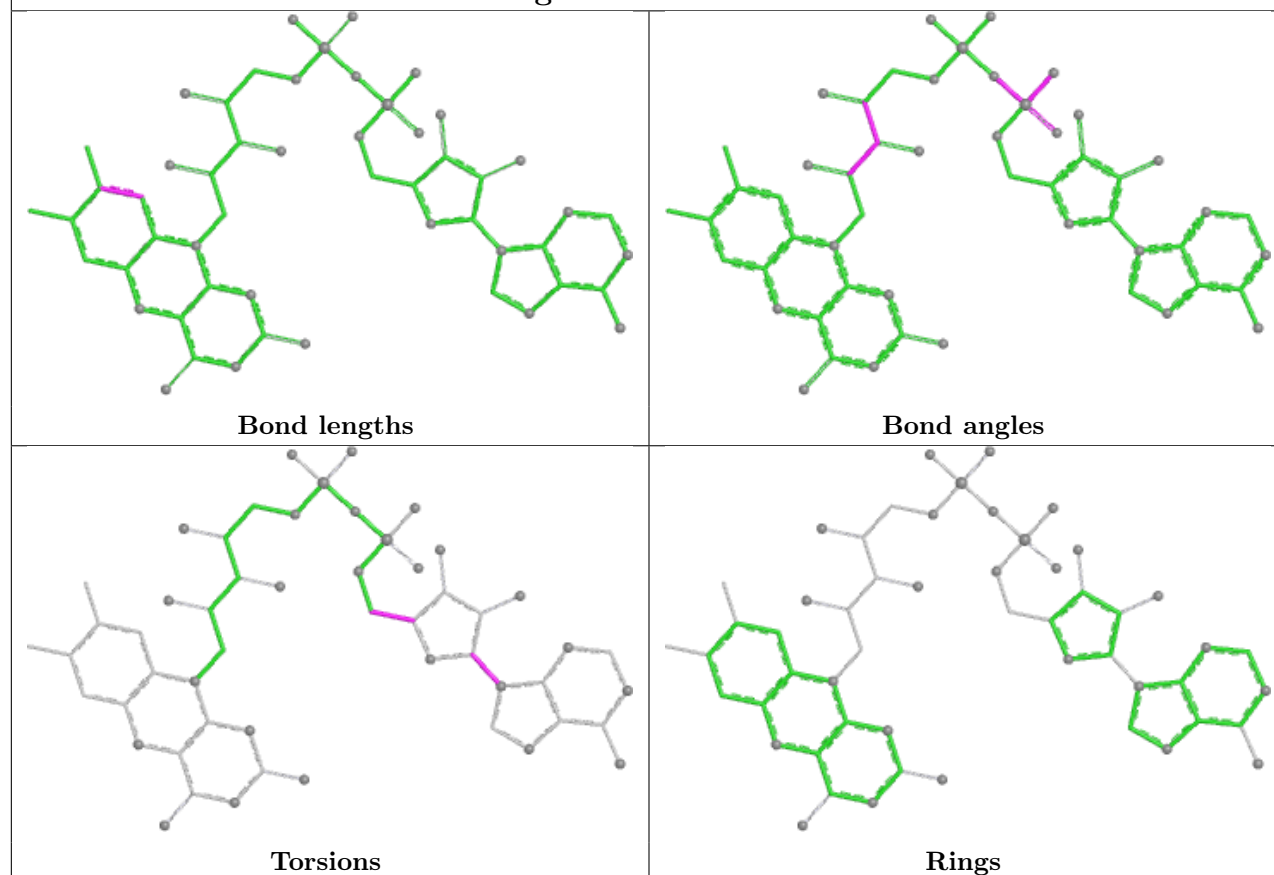
Ligand SAR C 702



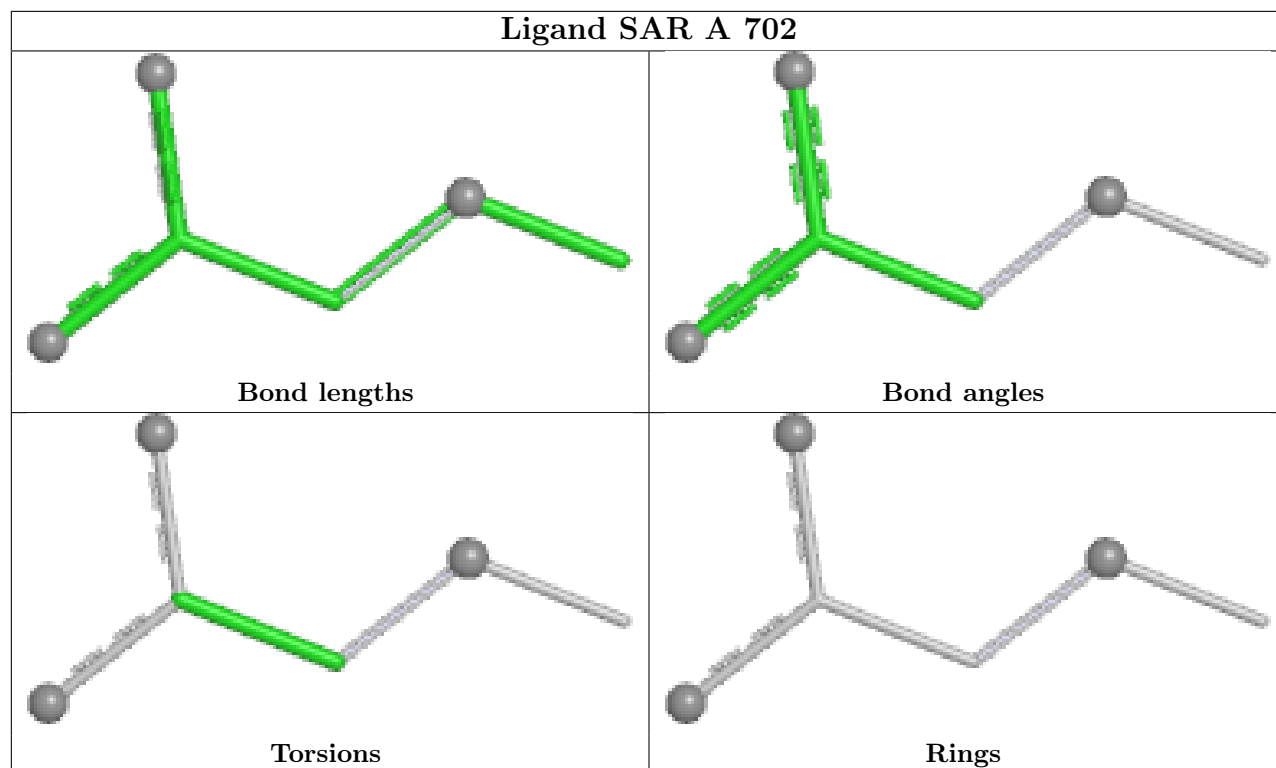
Ligand SAR E 702



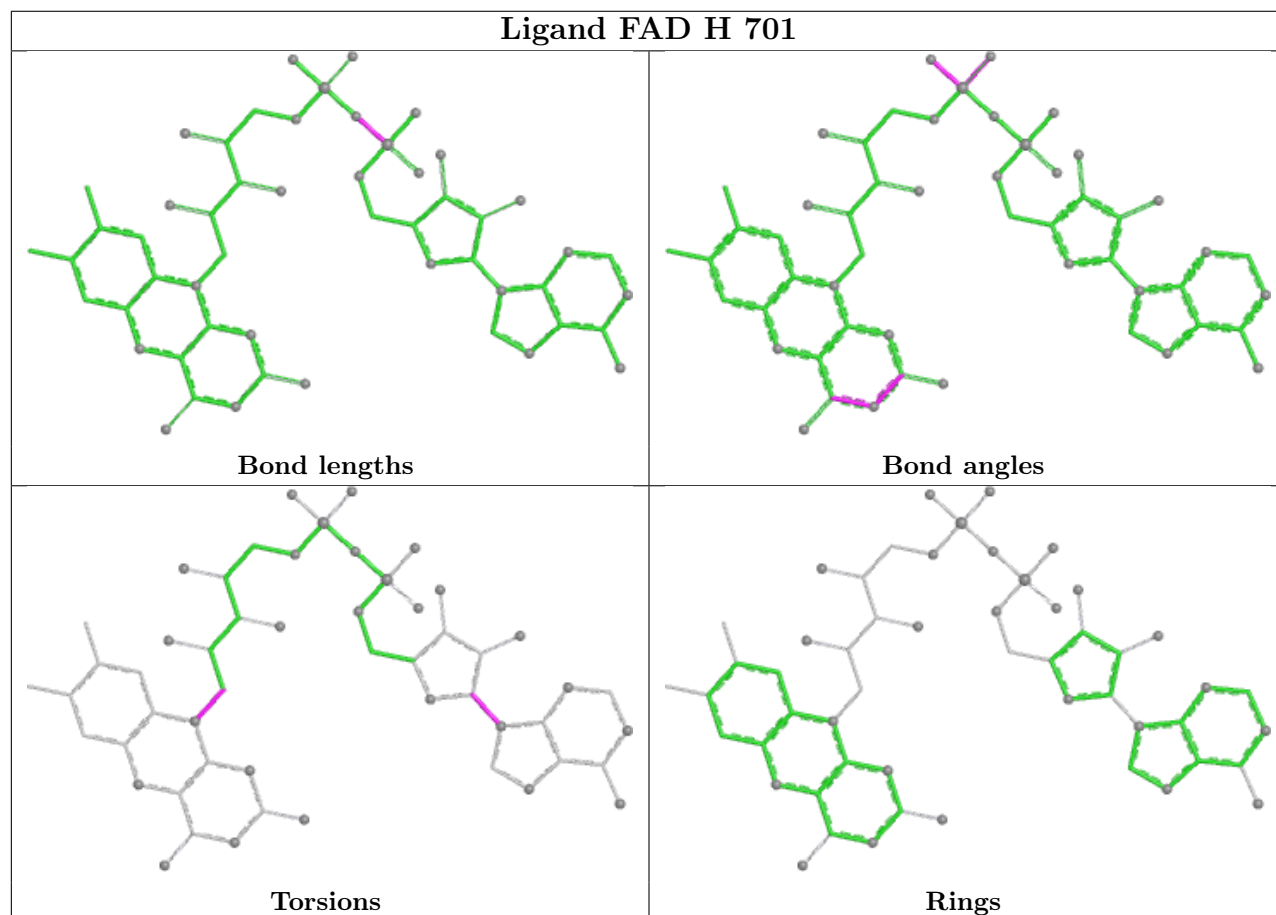
Ligand FAD F 701



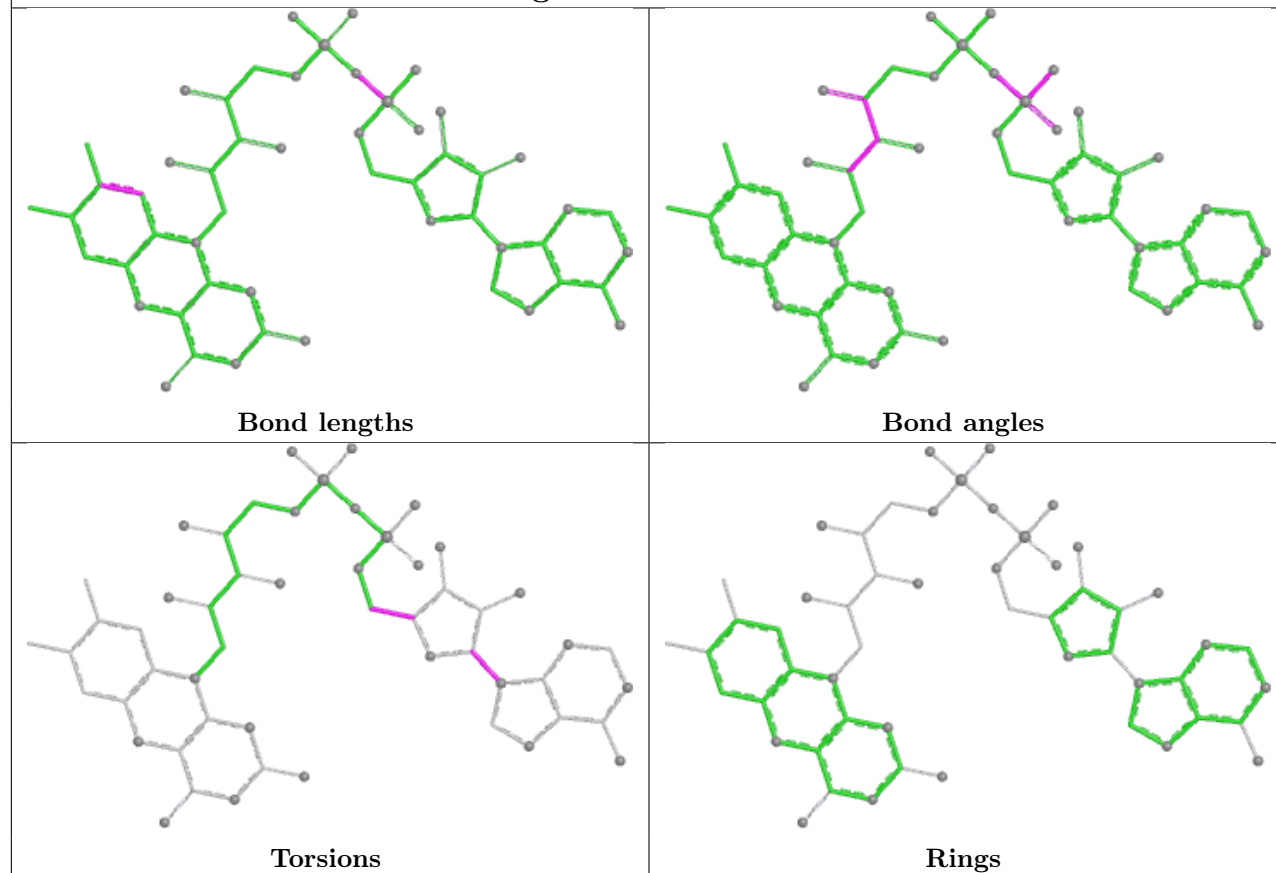
Ligand SAR A 702



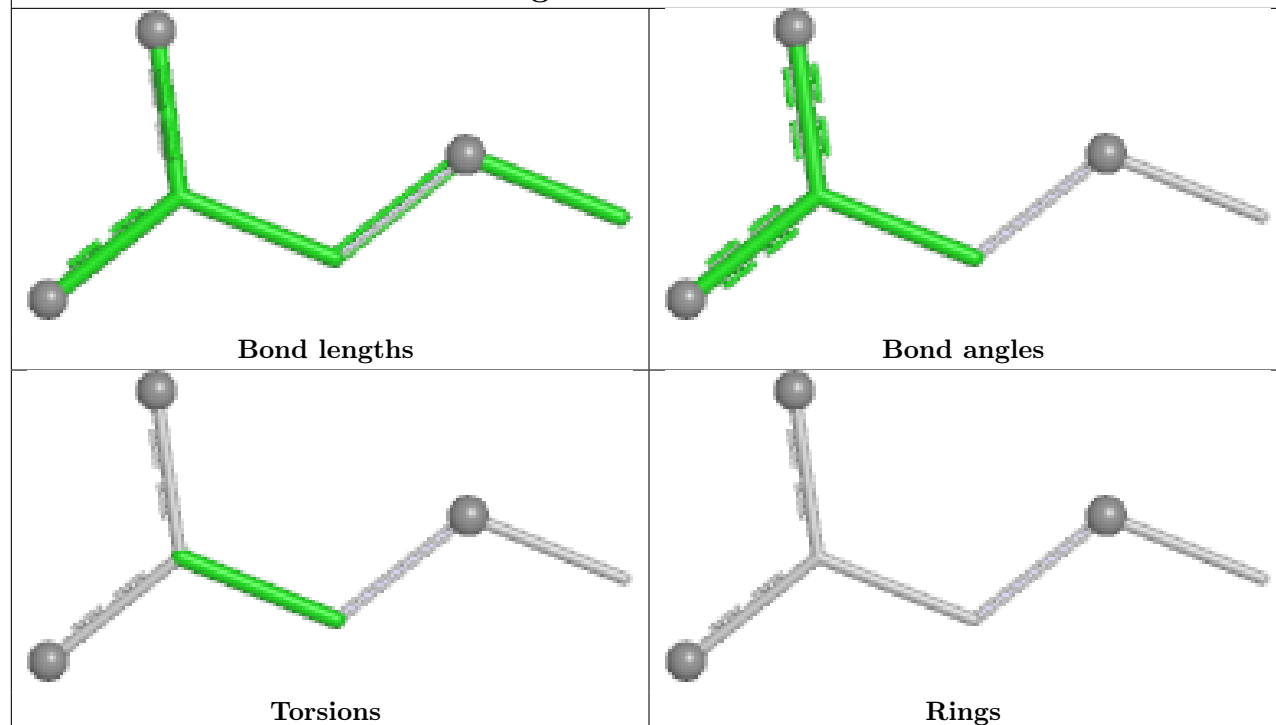
Ligand FAD H 701

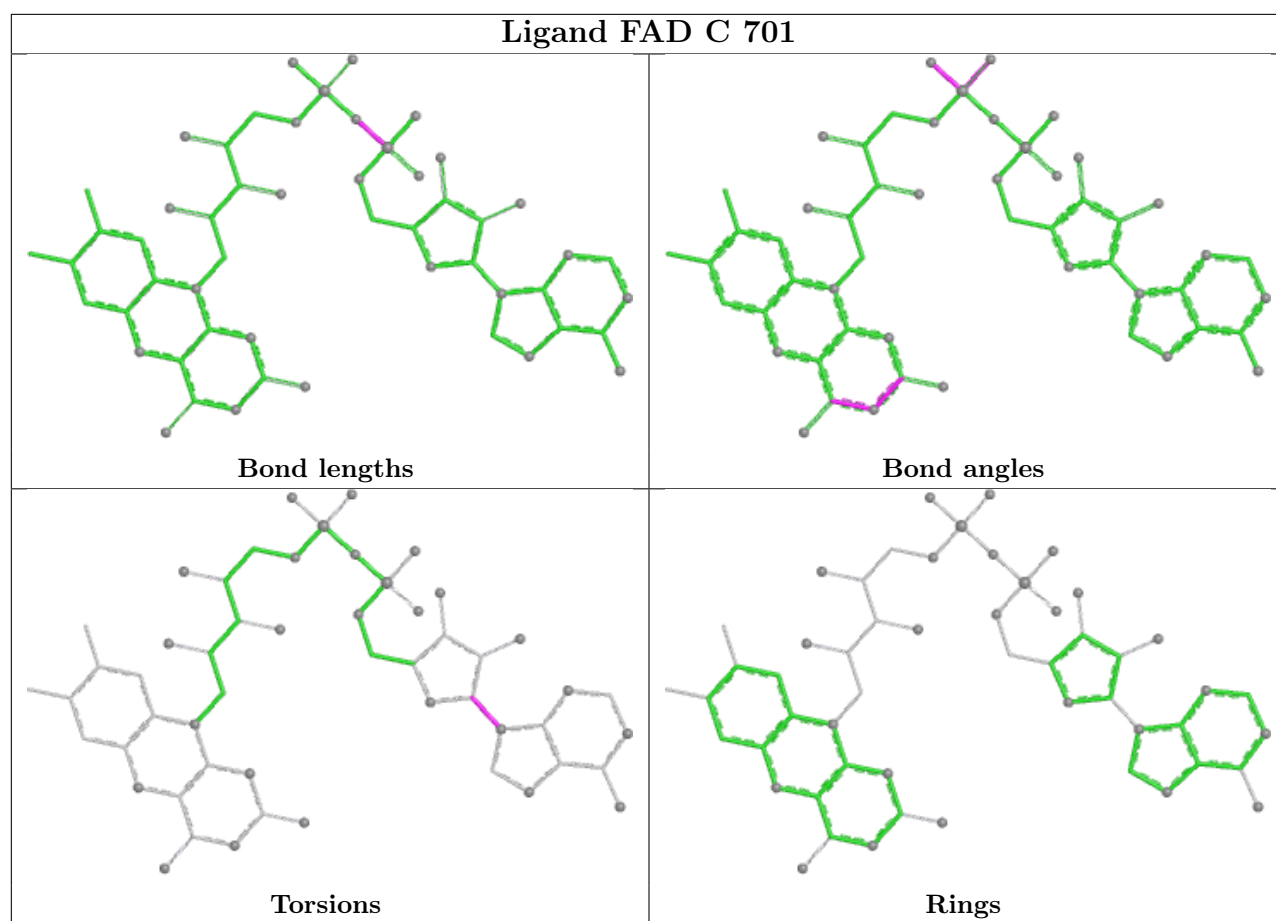


Ligand FAD A 701



Ligand SAR G 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	A	482/512 (94%)	-1.10	0	100 100	14, 27, 49, 80	0
1	B	482/512 (94%)	-1.16	0	100 100	11, 24, 46, 67	1 (0%)
1	C	482/512 (94%)	-1.03	0	100 100	11, 29, 65, 87	1 (0%)
1	D	479/512 (93%)	-0.72	1 (0%)	91 94	16, 43, 96, 116	0
1	E	482/512 (94%)	-1.09	0	100 100	14, 27, 50, 75	0
1	F	482/512 (94%)	-1.17	0	100 100	11, 24, 47, 69	1 (0%)
1	G	482/512 (94%)	-1.03	0	100 100	11, 29, 65, 88	0
1	H	479/512 (93%)	-0.74	1 (0%)	91 94	16, 43, 97, 118	0
All	All	3850/4096 (93%)	-1.01	2 (0%)	92 95	11, 29, 73, 118	3 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	223	GLY	2.3
1	H	238	ALA	2.2

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 ⓘ

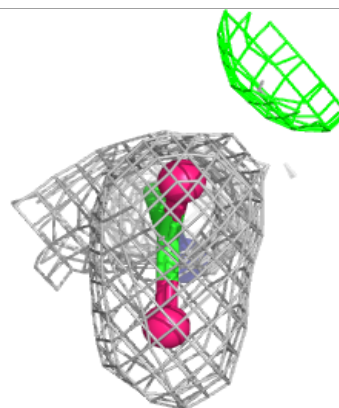
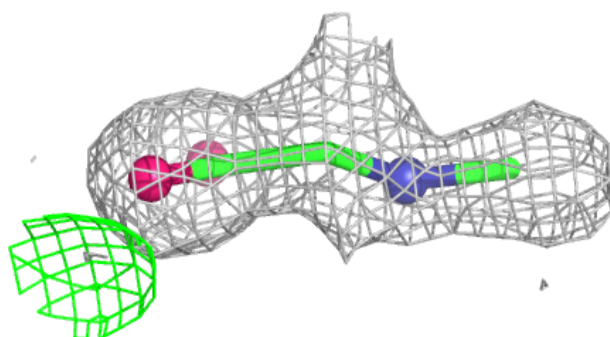
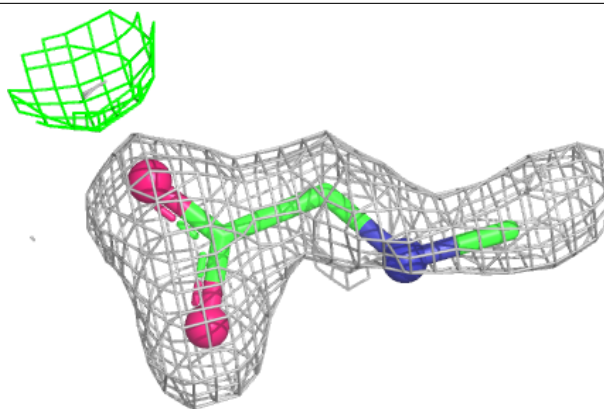
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	SAR	G	702	6/6	0.97	0.07	48,50,52,54	0
3	SAR	H	702	6/6	0.97	0.10	42,43,48,49	0
3	SAR	C	702	6/6	0.98	0.06	48,51,57,57	0
3	SAR	D	702	6/6	0.98	0.05	41,45,46,47	0
2	FAD	C	701	53/53	0.98	0.06	32,49,66,67	0
2	FAD	G	701	53/53	0.98	0.06	35,51,67,68	0
2	FAD	B	701	53/53	0.99	0.02	13,16,18,20	0
2	FAD	H	701	53/53	0.99	0.05	38,47,77,81	0
3	SAR	A	702	6/6	0.99	0.04	24,28,31,34	0
3	SAR	B	702	6/6	0.99	0.03	19,22,24,28	0
2	FAD	A	701	53/53	0.99	0.03	15,21,23,24	0
2	FAD	D	701	53/53	0.99	0.05	35,46,74,79	0
3	SAR	F	702	6/6	0.99	0.04	19,22,25,27	0
2	FAD	E	701	53/53	0.99	0.03	15,20,23,24	0
2	FAD	F	701	53/53	0.99	0.03	12,16,18,20	0
3	SAR	E	702	6/6	1.00	0.02	23,29,30,32	0

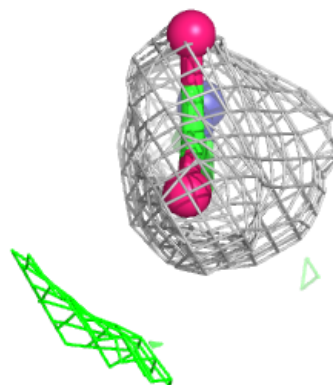
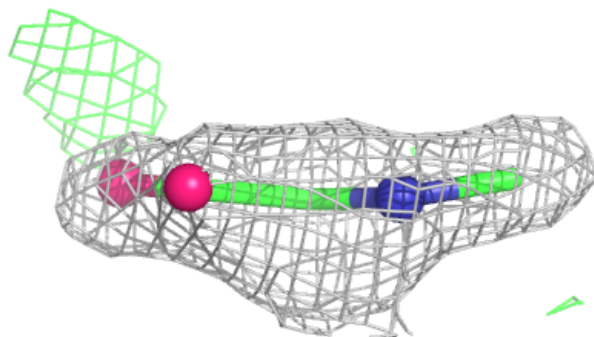
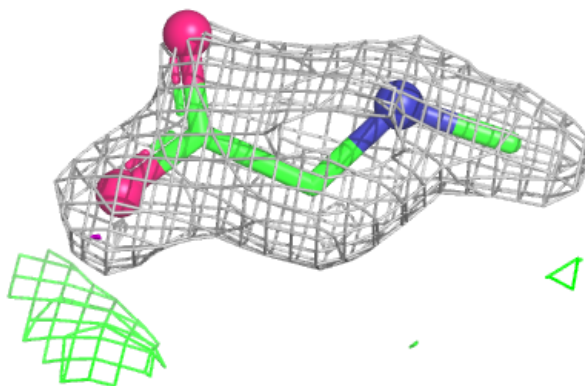
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around SAR G 702:

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

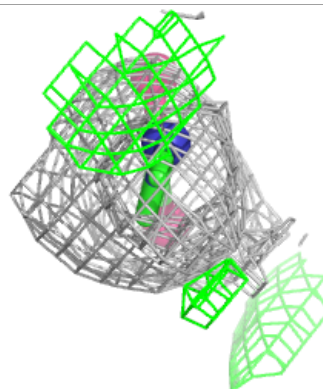
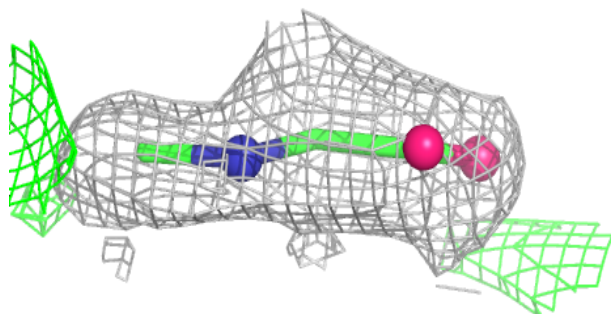
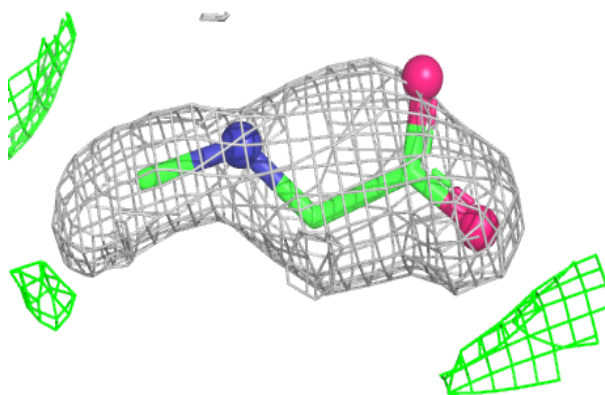
**Electron density around SAR H 702:**

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

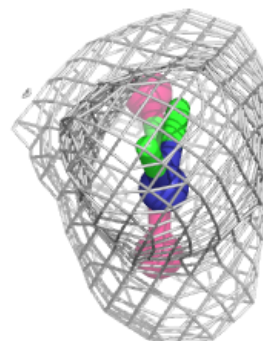
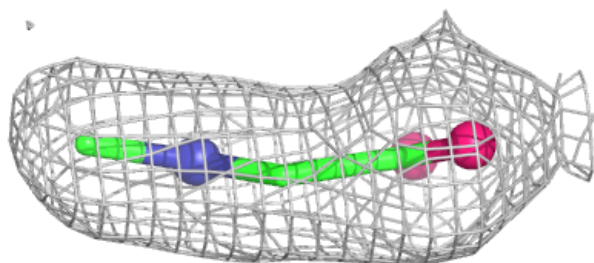
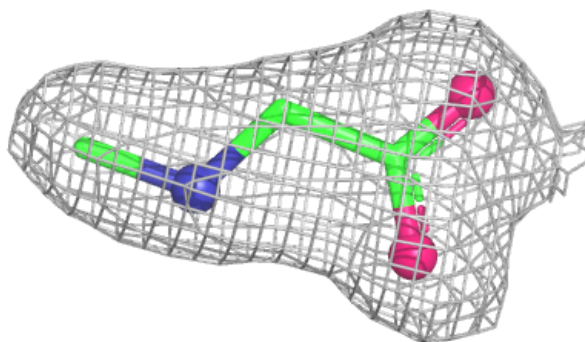


Electron density around SAR C 702:

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

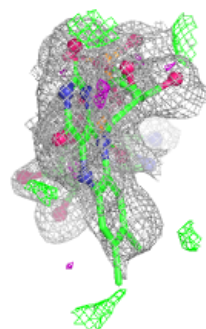
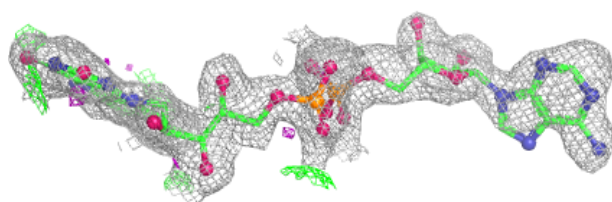
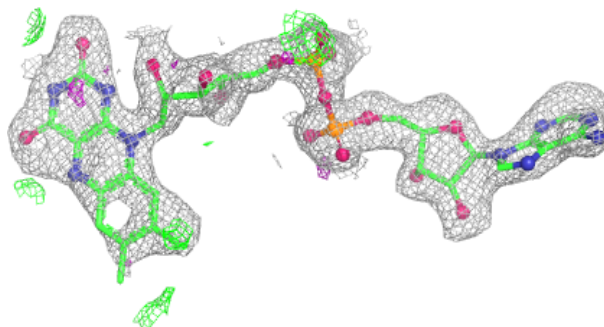
**Electron density around SAR D 702:**

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

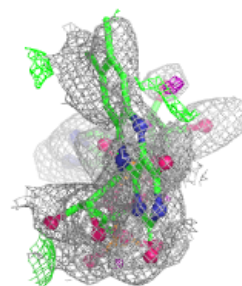
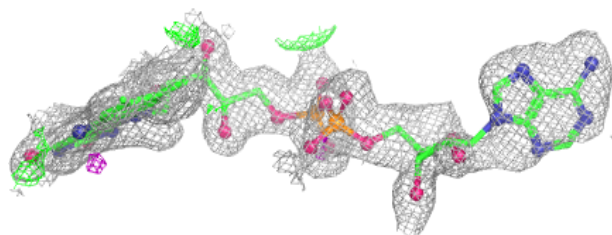
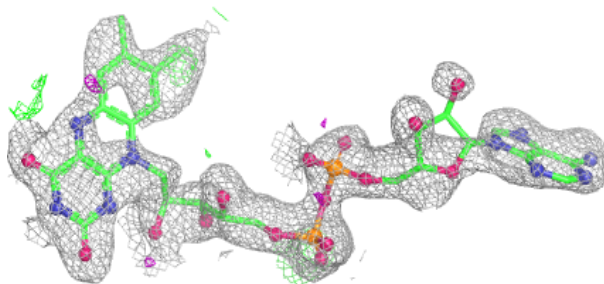


Electron density around FAD C 701:

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

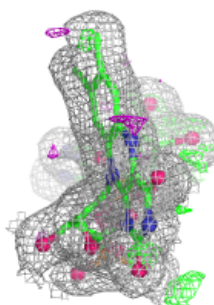
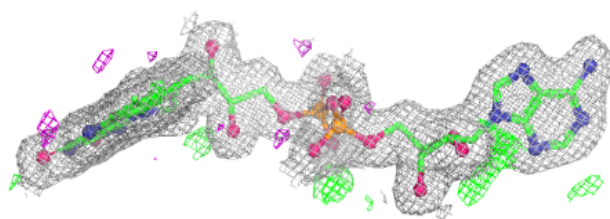
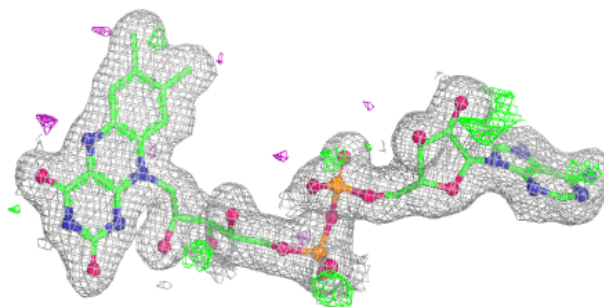
**Electron density around FAD G 701:**

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

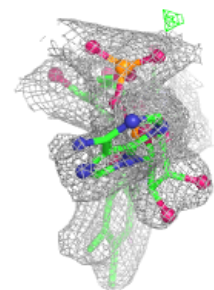
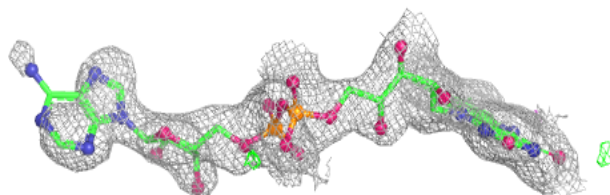
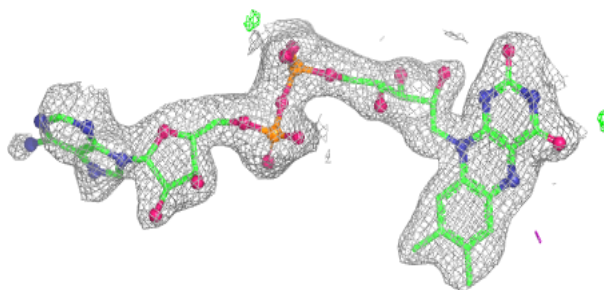


Electron density around FAD B 701:

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

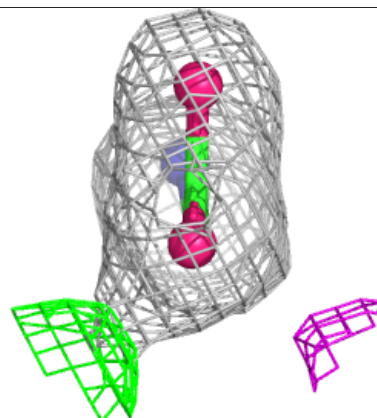
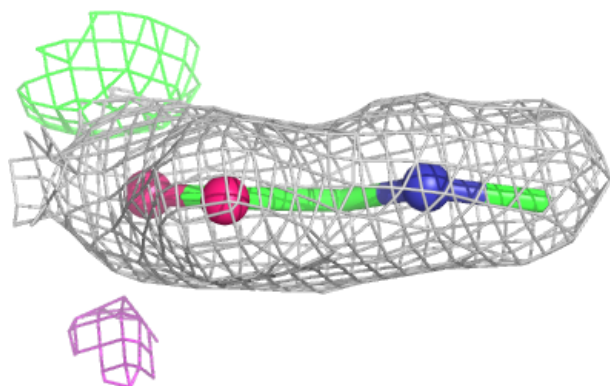
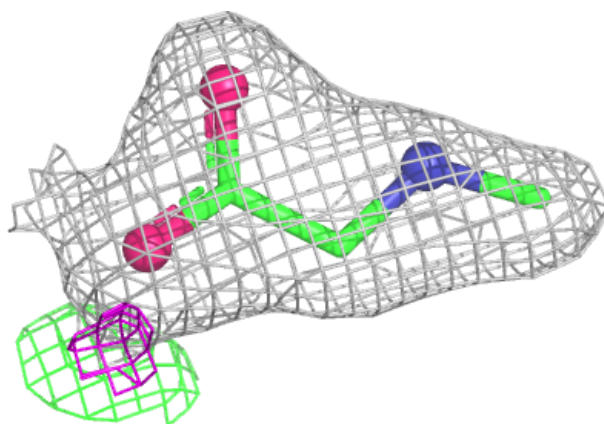
**Electron density around FAD H 701:**

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

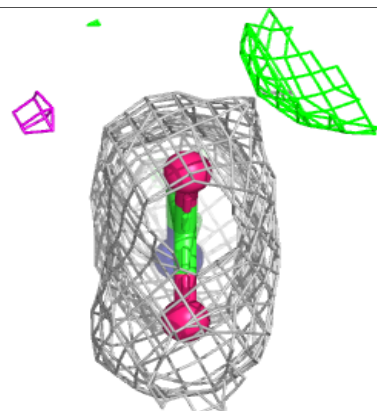
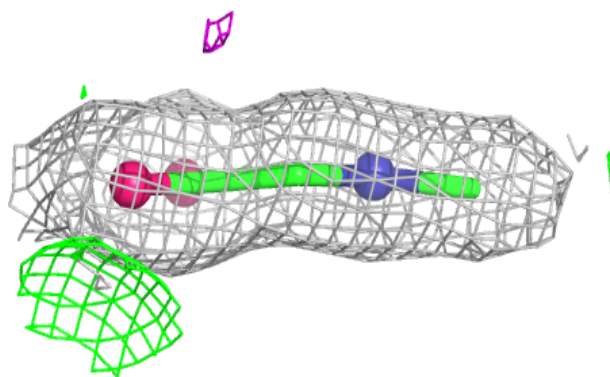
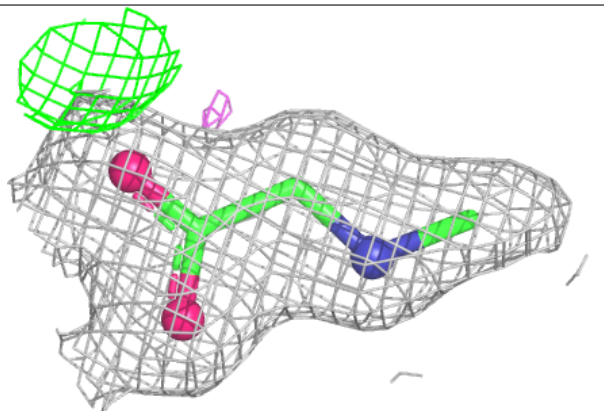


Electron density around SAR A 702:

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

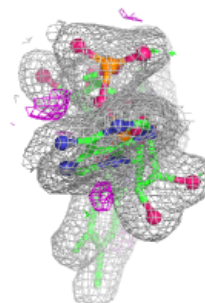
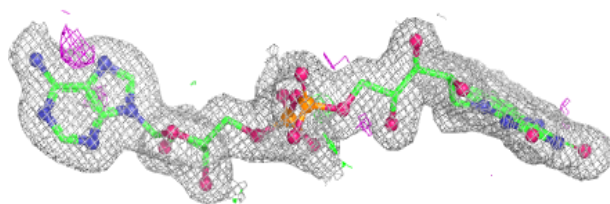
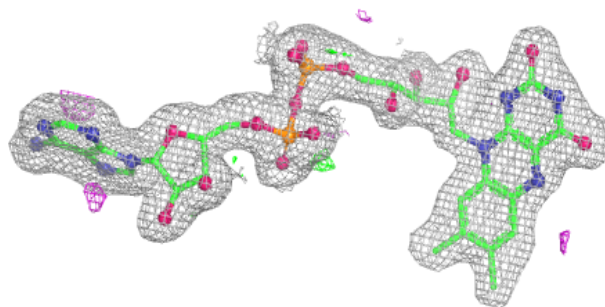
**Electron density around SAR B 702:**

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

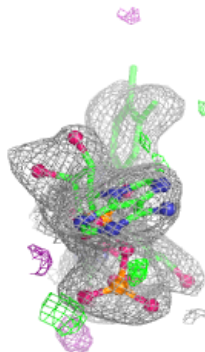
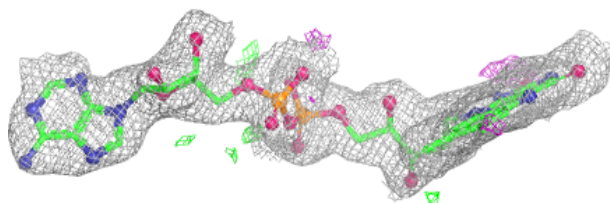
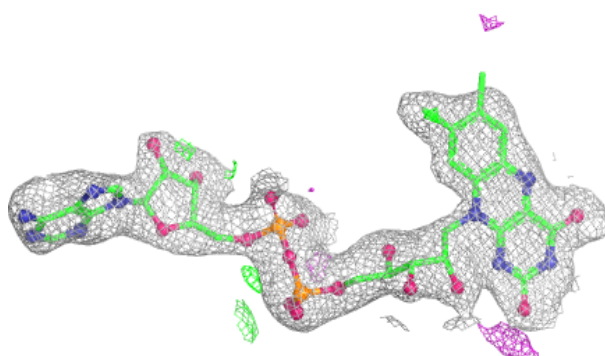


Electron density around FAD A 701:

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

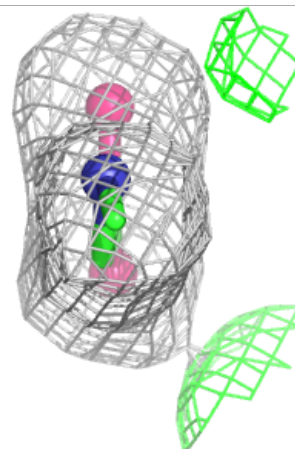
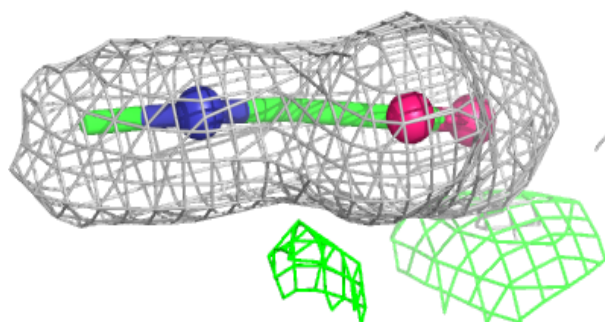
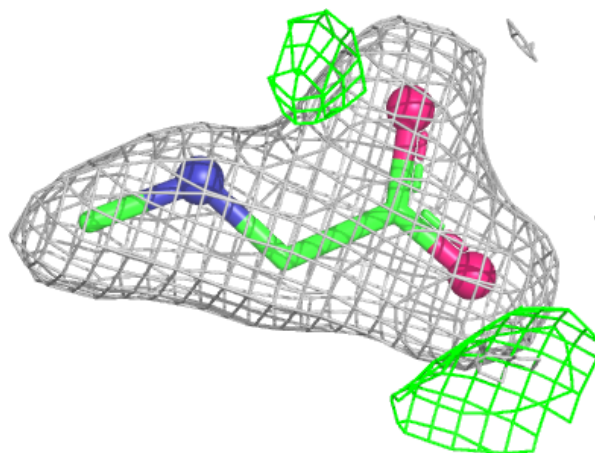
**Electron density around FAD D 701:**

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



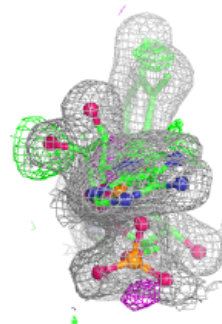
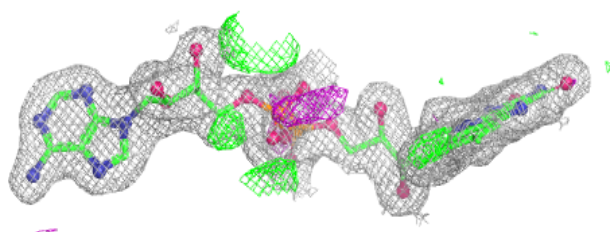
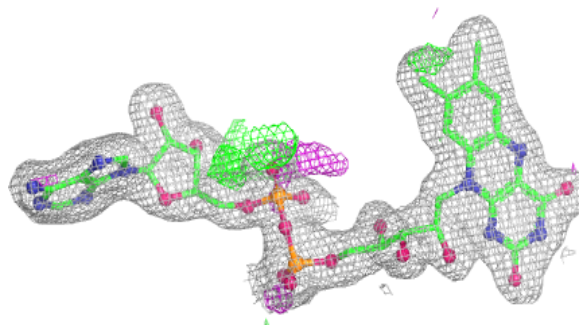
Electron density around SAR F 702:

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

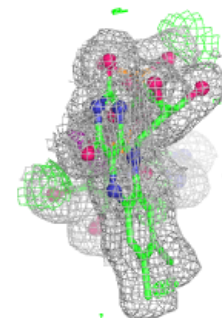
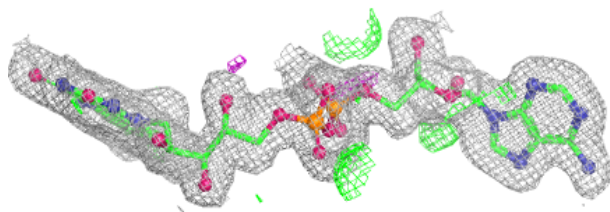
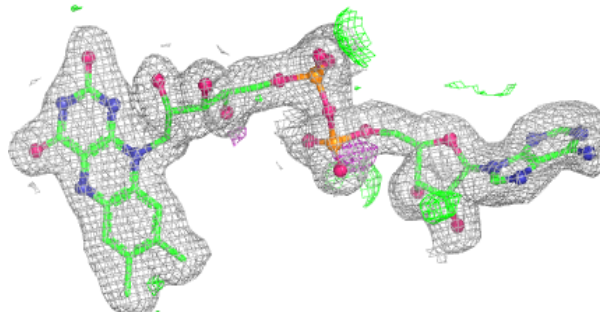


Electron density around FAD E 701:

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

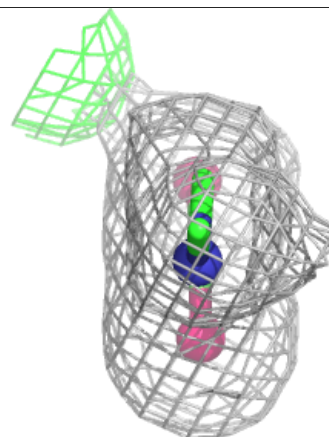
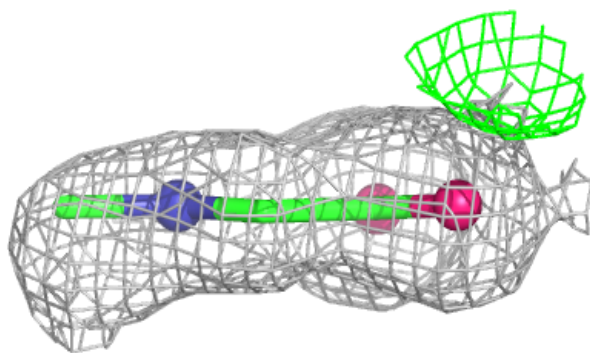
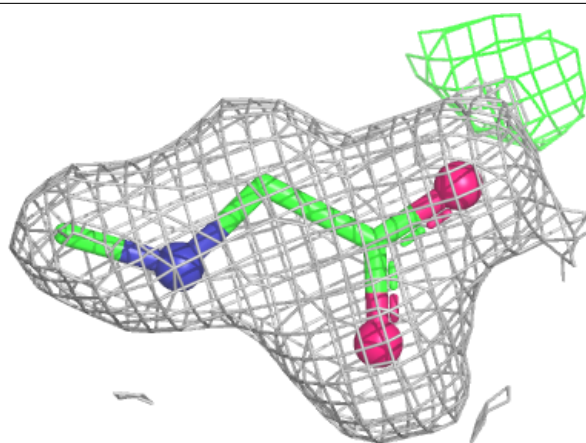
**Electron density around FAD F 701:**

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



Electron density around SAR E 702:

$2mF_o-DF_c$ (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.