



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 2, 2026 – 12:18 AM UTC

PDB ID : 3NIX / pdb_00003nix
Title : Crystal structure of flavoprotein/dehydrogenase from *Cytophaga hutchinsonii*.
Northeast Structural Genomics Consortium Target Chr43.
Authors : Vorobiev, S.; Su, M.; Seetharaman, J.; Sahdev, S.; Xiao, R.; Foote, E.L.;
Ciccosanti, C.; Maglaqui, M.; Everett, J.K.; Nair, R.; Acton, T.B.; Rost,
B.; Montelione, G.T.; Hunt, J.F.; Tong, L.; Northeast Structural Genomics
Consortium (NESG)
Deposited on : 2010-06-16
Resolution : 2.60 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<https://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)

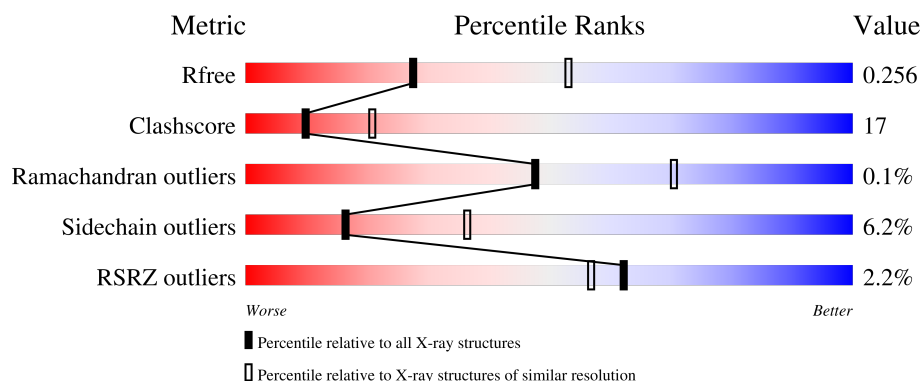
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	421	2% 67% 26% • •
1	B	421	2% 62% 31% • 5%
1	C	421	2% 64% 28% 5% •

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Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
 Validation Pipeline (wwPDB-VP) : 2.49

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Mol	Chain	Length	Quality of chain
1	D	421	
1	E	421	
1	F	421	
1	G	421	
1	H	421	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	FAD	F	506	X	-	-	-
2	FAD	G	507	X	-	-	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 26852 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 Flavoprotein/dehydrogenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	407	Total	C	N	O	S	Se	0	0	0
			3210	2062	544	593	2	9			
1	B	400	Total	C	N	O	S	Se	0	0	0
			3162	2032	537	582	2	9			
1	C	405	Total	C	N	O	S	Se	0	0	0
			3185	2047	542	585	2	9			
1	D	403	Total	C	N	O	S	Se	0	0	0
			3183	2046	541	585	2	9			
1	E	406	Total	C	N	O	S	Se	0	0	0
			3212	2062	547	592	2	9			
1	F	402	Total	C	N	O	S	Se	0	0	0
			3182	2045	540	586	2	9			
1	G	406	Total	C	N	O	S	Se	0	0	0
			3204	2057	546	590	2	9			
1	H	403	Total	C	N	O	S	Se	0	0	0
			3177	2041	540	585	2	9			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	initiating methionine	UNP Q11R94
A	414	LEU	-	expression tag	UNP Q11R94
A	415	GLU	-	expression tag	UNP Q11R94
A	416	HIS	-	expression tag	UNP Q11R94
A	417	HIS	-	expression tag	UNP Q11R94
A	418	HIS	-	expression tag	UNP Q11R94
A	419	HIS	-	expression tag	UNP Q11R94
A	420	HIS	-	expression tag	UNP Q11R94
A	421	HIS	-	expression tag	UNP Q11R94
B	1	MSE	-	initiating methionine	UNP Q11R94
B	414	LEU	-	expression tag	UNP Q11R94
B	415	GLU	-	expression tag	UNP Q11R94
B	416	HIS	-	expression tag	UNP Q11R94

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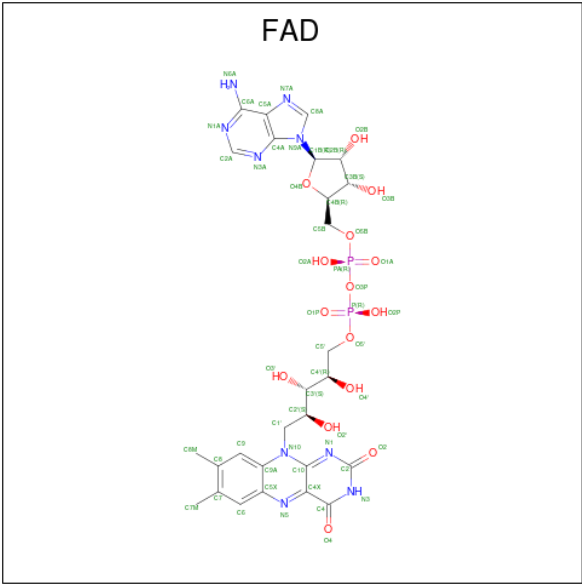
Chain	Residue	Modelled	Actual	Comment	Reference
B	417	HIS	-	expression tag	UNP Q11R94
B	418	HIS	-	expression tag	UNP Q11R94
B	419	HIS	-	expression tag	UNP Q11R94
B	420	HIS	-	expression tag	UNP Q11R94
B	421	HIS	-	expression tag	UNP Q11R94
C	1	MSE	-	initiating methionine	UNP Q11R94
C	414	LEU	-	expression tag	UNP Q11R94
C	415	GLU	-	expression tag	UNP Q11R94
C	416	HIS	-	expression tag	UNP Q11R94
C	417	HIS	-	expression tag	UNP Q11R94
C	418	HIS	-	expression tag	UNP Q11R94
C	419	HIS	-	expression tag	UNP Q11R94
C	420	HIS	-	expression tag	UNP Q11R94
C	421	HIS	-	expression tag	UNP Q11R94
D	1	MSE	-	initiating methionine	UNP Q11R94
D	414	LEU	-	expression tag	UNP Q11R94
D	415	GLU	-	expression tag	UNP Q11R94
D	416	HIS	-	expression tag	UNP Q11R94
D	417	HIS	-	expression tag	UNP Q11R94
D	418	HIS	-	expression tag	UNP Q11R94
D	419	HIS	-	expression tag	UNP Q11R94
D	420	HIS	-	expression tag	UNP Q11R94
D	421	HIS	-	expression tag	UNP Q11R94
E	1	MSE	-	initiating methionine	UNP Q11R94
E	414	LEU	-	expression tag	UNP Q11R94
E	415	GLU	-	expression tag	UNP Q11R94
E	416	HIS	-	expression tag	UNP Q11R94
E	417	HIS	-	expression tag	UNP Q11R94
E	418	HIS	-	expression tag	UNP Q11R94
E	419	HIS	-	expression tag	UNP Q11R94
E	420	HIS	-	expression tag	UNP Q11R94
E	421	HIS	-	expression tag	UNP Q11R94
F	1	MSE	-	initiating methionine	UNP Q11R94
F	414	LEU	-	expression tag	UNP Q11R94
F	415	GLU	-	expression tag	UNP Q11R94
F	416	HIS	-	expression tag	UNP Q11R94
F	417	HIS	-	expression tag	UNP Q11R94
F	418	HIS	-	expression tag	UNP Q11R94
F	419	HIS	-	expression tag	UNP Q11R94
F	420	HIS	-	expression tag	UNP Q11R94
F	421	HIS	-	expression tag	UNP Q11R94
G	1	MSE	-	initiating methionine	UNP Q11R94

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Chain	Residue	Modelled	Actual	Comment	Reference
G	414	LEU	-	expression tag	UNP Q11R94
G	415	GLU	-	expression tag	UNP Q11R94
G	416	HIS	-	expression tag	UNP Q11R94
G	417	HIS	-	expression tag	UNP Q11R94
G	418	HIS	-	expression tag	UNP Q11R94
G	419	HIS	-	expression tag	UNP Q11R94
G	420	HIS	-	expression tag	UNP Q11R94
G	421	HIS	-	expression tag	UNP Q11R94
H	1	MSE	-	initiating methionine	UNP Q11R94
H	414	LEU	-	expression tag	UNP Q11R94
H	415	GLU	-	expression tag	UNP Q11R94
H	416	HIS	-	expression tag	UNP Q11R94
H	417	HIS	-	expression tag	UNP Q11R94
H	418	HIS	-	expression tag	UNP Q11R94
H	419	HIS	-	expression tag	UNP Q11R94
H	420	HIS	-	expression tag	UNP Q11R94
H	421	HIS	-	expression tag	UNP Q11R94

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

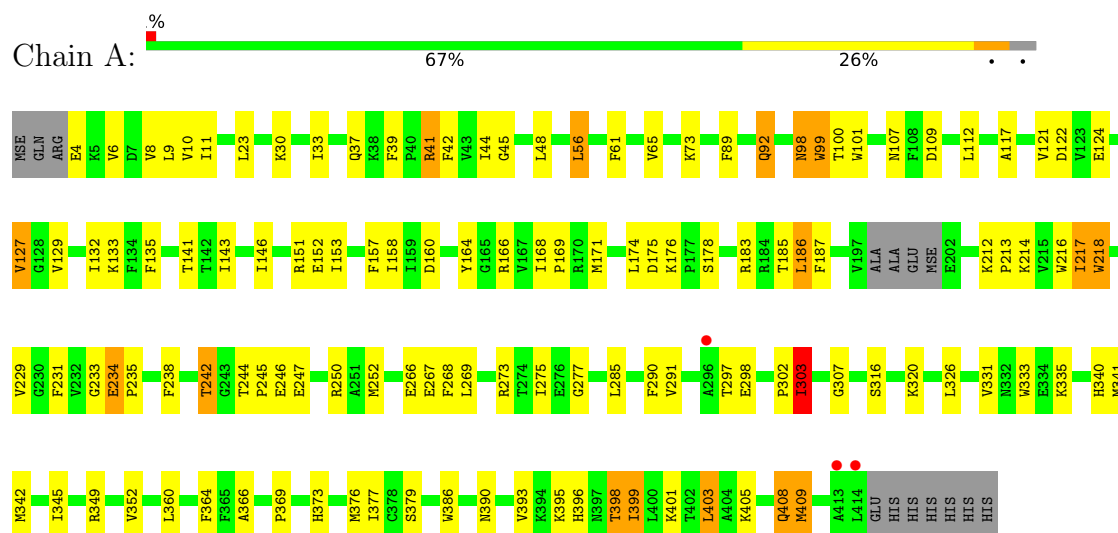
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	125	Total	O	0	0
			125	125		
3	B	139	Total	O	0	0
			139	139		
3	C	119	Total	O	0	0
			119	119		
3	D	81	Total	O	0	0
			81	81		
3	E	148	Total	O	0	0
			148	148		
3	F	99	Total	O	0	0
			99	99		
3	G	89	Total	O	0	0
			89	89		
3	H	113	Total	O	0	0
			113	113		

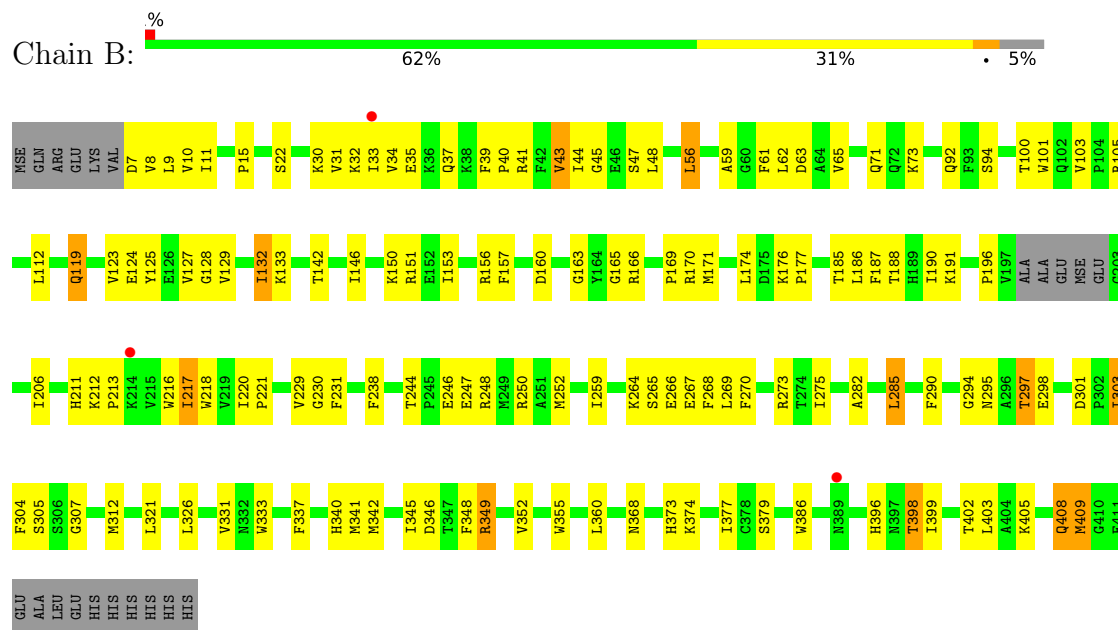
3 Residue-property plots

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: Flavoprotein/dehydrogenase

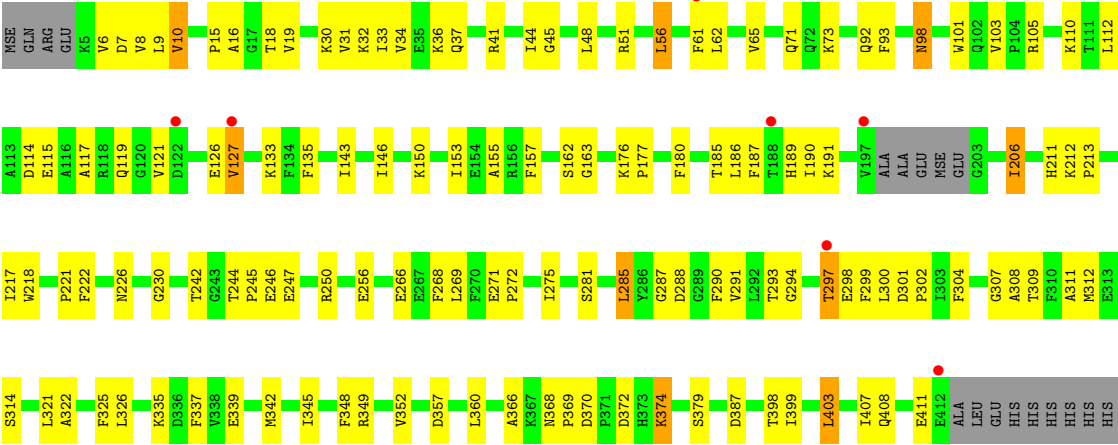


- Molecule 1: Flavoprotein/dehydrogenase



- Molecule 1: Flavoprotein/dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	86.22Å 244.05Å 100.22Å 90.00° 104.10° 90.00°	Depositor
Resolution (Å)	29.33 – 2.60 29.33 – 2.60	Depositor EDS
% Data completeness (in resolution range)	96.8 (29.33-2.60) 97.4 (29.33-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.08 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.6_289, REFMAC	Depositor
R, R_{free}	0.237 , 0.264 0.229 , 0.256	Depositor DCC
R_{free} test set	6026 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	32.0	Xtriage
Anisotropy	0.282	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 40.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	26852	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 61.29 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.3396e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0/3281	0.92	8/4415 (0.2%)
1	B	0.70	2/3233 (0.1%)	0.92	7/4349 (0.2%)
1	C	0.68	4/3256 (0.1%)	0.94	6/4382 (0.1%)
1	D	0.55	0/3254	0.90	8/4377 (0.2%)
1	E	0.62	0/3283	0.93	3/4415 (0.1%)
1	F	0.56	0/3253	0.90	4/4375 (0.1%)
1	G	0.56	0/3275	0.89	3/4406 (0.1%)
1	H	0.61	0/3248	0.90	4/4370 (0.1%)
All	All	0.62	6/26083 (0.0%)	0.91	43/35089 (0.1%)

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	208	ALA	C-N	-10.39	1.20	1.33
1	C	209	VAL	C-N	-6.73	1.26	1.33
1	C	219	VAL	C-O	-6.21	1.17	1.24
1	B	348	PHE	C-O	-5.82	1.17	1.24
1	C	217	ILE	C-O	-5.79	1.18	1.24

The worst 5 of 43 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	96	GLY	N-CA-C	-7.73	99.69	112.83
1	G	167	VAL	N-CA-C	7.37	117.50	110.42
1	C	370	ASP	CA-C-N	6.99	126.81	119.05
1	C	370	ASP	C-N-CA	6.99	126.81	119.05
1	G	96	GLY	N-CA-C	-6.84	101.20	112.83

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	3210	0	3125	109	0
1	B	3162	0	3084	115	0
1	C	3185	0	3097	115	0
1	D	3183	0	3108	107	0
1	E	3212	0	3136	94	0
1	F	3182	0	3110	125	0
1	G	3204	0	3121	129	0
1	H	3177	0	3090	103	0
2	A	53	0	28	4	0
2	B	53	0	29	5	0
2	C	53	0	29	2	0
2	D	53	0	29	8	0
2	E	53	0	29	5	0
2	F	53	0	28	6	0
2	G	53	0	29	14	0
2	H	53	0	29	10	0
3	A	125	0	0	12	0
3	B	139	0	0	7	0
3	C	119	0	0	5	0
3	D	81	0	0	6	0
3	E	148	0	0	15	0
3	F	99	0	0	14	0
3	G	89	0	0	9	0
3	H	113	0	0	5	0
All	All	26852	0	25101	892	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 17.

The worst 5 of 892 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:C:297:THR:HG22	1:C:298:GLU:H	1.12	1.11
1:D:41:ARG:HG2	1:D:41:ARG:HH11	0.88	1.02
1:B:297:THR:HG22	1:B:298:GLU:H	1.22	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:23:LEU:HG	3:D:1156:HOH:O	1.58	1.01
1:E:41:ARG:HG2	1:E:41:ARG:HH11	0.87	1.00

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	403/421 (96%)	380 (94%)	22 (6%)	1 (0%)	43	66
1	B	396/421 (94%)	377 (95%)	19 (5%)	0	100	100
1	C	401/421 (95%)	380 (95%)	20 (5%)	1 (0%)	43	66
1	D	399/421 (95%)	378 (95%)	21 (5%)	0	100	100
1	E	402/421 (96%)	387 (96%)	14 (4%)	1 (0%)	43	66
1	F	398/421 (94%)	380 (96%)	18 (4%)	0	100	100
1	G	402/421 (96%)	377 (94%)	24 (6%)	1 (0%)	43	66
1	H	399/421 (95%)	381 (96%)	18 (4%)	0	100	100
All	All	3200/3368 (95%)	3040 (95%)	156 (5%)	4 (0%)	48	70

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	294	GLY
1	G	210	VAL
1	A	303	ILE
1	E	294	GLY

5.3.2 Protein sidechains ⓘ

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	336/340 (99%)	315 (94%)	21 (6%)	16	36
1	B	332/340 (98%)	310 (93%)	22 (7%)	15	34
1	C	332/340 (98%)	312 (94%)	20 (6%)	17	37
1	D	334/340 (98%)	308 (92%)	26 (8%)	11	26
1	E	337/340 (99%)	313 (93%)	24 (7%)	13	31
1	F	335/340 (98%)	316 (94%)	19 (6%)	18	40
1	G	335/340 (98%)	316 (94%)	19 (6%)	18	40
1	H	332/340 (98%)	316 (95%)	16 (5%)	23	47
All	All	2673/2720 (98%)	2506 (94%)	167 (6%)	16	36

5 of 167 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	48	LEU
1	G	255	ASN
1	F	122	ASP
1	F	398	THR
1	G	403	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 67 such sidechains are listed below:

Mol	Chain	Res	Type
1	G	368	ASN
1	H	37	GLN
1	H	295	ASN
1	D	92	GLN
1	D	68	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	502	-	58,58,58	2.52	15 (25%)	85,89,89	1.96	26 (30%)
2	FAD	D	504	-	58,58,58	2.55	18 (31%)	85,89,89	1.92	26 (30%)
2	FAD	C	503	-	58,58,58	2.62	17 (29%)	85,89,89	1.98	25 (29%)
2	FAD	F	506	-	58,58,58	2.46	15 (25%)	85,89,89	1.88	23 (27%)
2	FAD	E	505	-	58,58,58	2.52	19 (32%)	85,89,89	1.90	24 (28%)
2	FAD	A	501	-	58,58,58	2.49	19 (32%)	85,89,89	2.01	28 (32%)
2	FAD	G	507	-	58,58,58	2.48	16 (27%)	85,89,89	2.16	27 (31%)
2	FAD	H	508	-	58,58,58	2.50	17 (29%)	85,89,89	1.73	20 (23%)

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	502	-	-	6/34/50/50	0/6/6/6
2	FAD	D	504	-	-	5/34/50/50	0/6/6/6
2	FAD	C	503	-	-	6/34/50/50	0/6/6/6
2	FAD	F	506	-	1/1/9/9	4/34/50/50	0/6/6/6

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FAD	E	505	-	-	8/34/50/50	0/6/6/6
2	FAD	A	501	-	-	5/34/50/50	0/6/6/6
2	FAD	G	507	-	1/1/9/9	10/34/50/50	0/6/6/6
2	FAD	H	508	-	-	4/34/50/50	0/6/6/6

The worst 5 of 136 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	503	FAD	O4-C4	12.18	1.46	1.23
2	B	502	FAD	O4-C4	11.92	1.46	1.23
2	E	505	FAD	O4-C4	11.84	1.46	1.23
2	D	504	FAD	O4-C4	11.78	1.45	1.23
2	G	507	FAD	O4-C4	11.65	1.45	1.23

The worst 5 of 199 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	507	FAD	C5A-C4A-N3A	-6.86	117.28	126.72
2	D	504	FAD	C5A-C4A-N3A	-6.10	118.32	126.72
2	E	505	FAD	C5A-C4A-N3A	-5.26	119.47	126.72
2	F	506	FAD	C5A-C4A-N3A	-5.22	119.53	126.72
2	C	503	FAD	N3A-C2A-N1A	-5.14	120.80	128.58

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	F	506	FAD	C2'
2	G	507	FAD	C2'

5 of 48 torsion outliers are listed below:

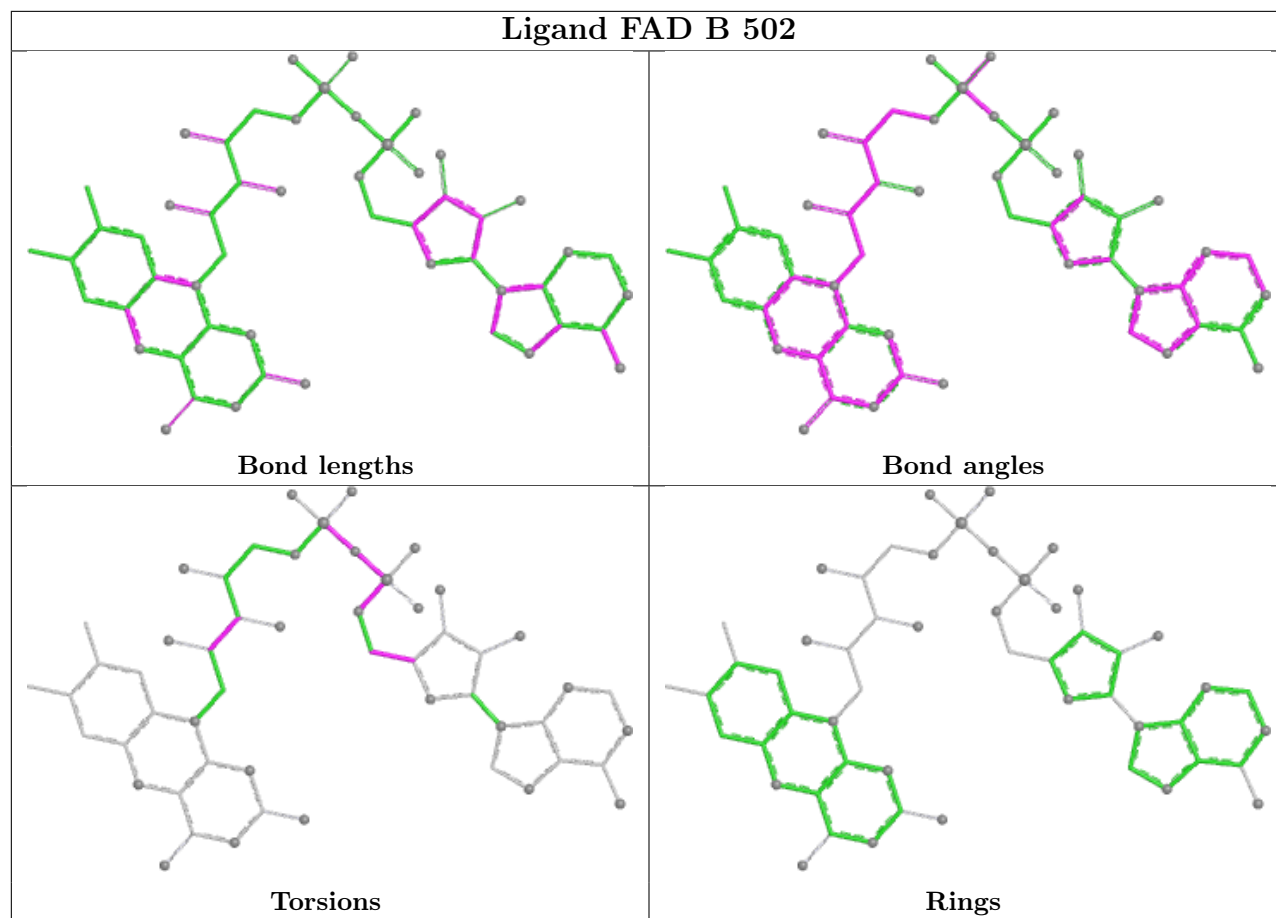
Mol	Chain	Res	Type	Atoms
2	A	501	FAD	C1'-C2'-C3'-C4'
2	A	501	FAD	PA-O3P-P-O5'
2	B	502	FAD	C1'-C2'-C3'-O3'
2	B	502	FAD	C1'-C2'-C3'-C4'
2	B	502	FAD	PA-O3P-P-O5'

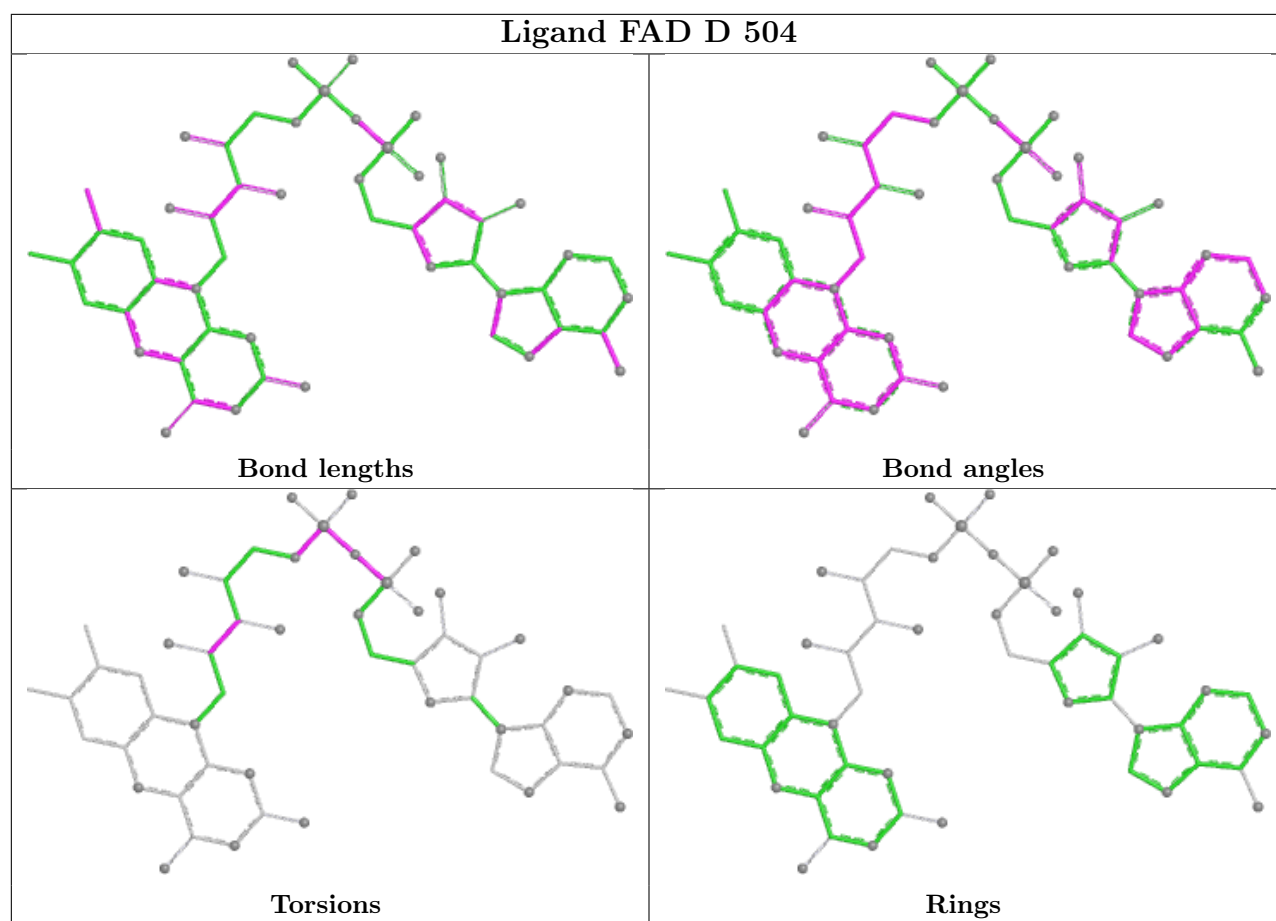
There are no ring outliers.

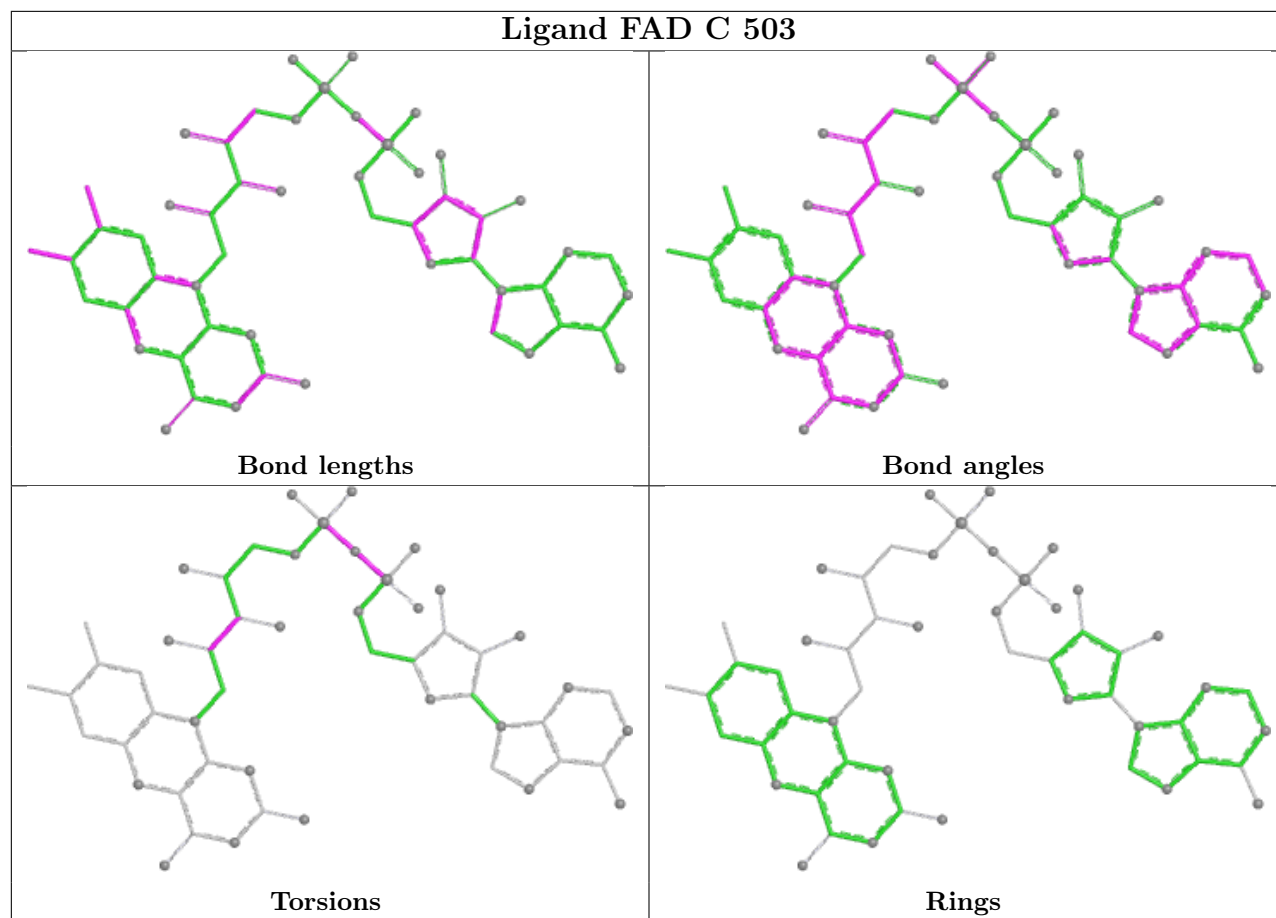
8 monomers are involved in 54 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	502	FAD	5	0
2	D	504	FAD	8	0
2	C	503	FAD	2	0
2	F	506	FAD	6	0
2	E	505	FAD	5	0
2	A	501	FAD	4	0
2	G	507	FAD	14	0
2	H	508	FAD	10	0

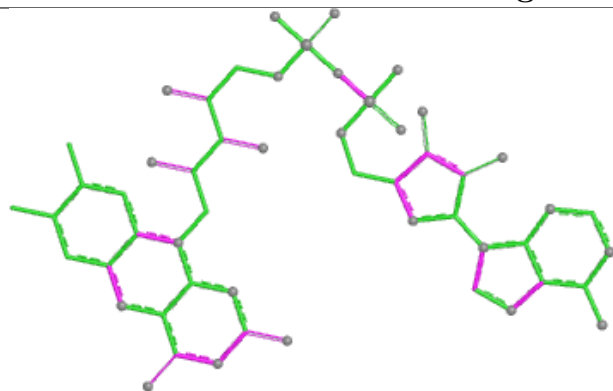
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



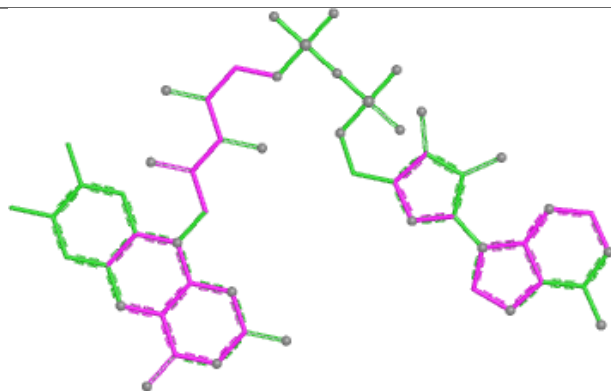




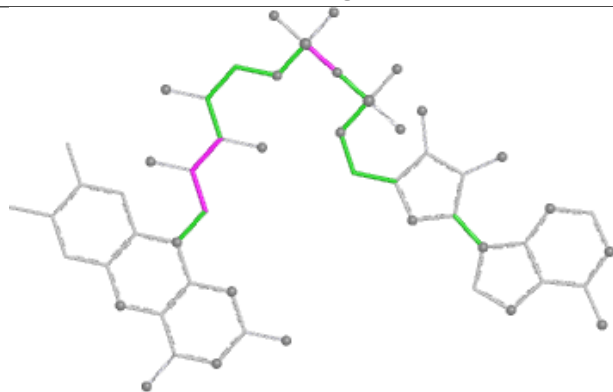
Ligand FAD F 506



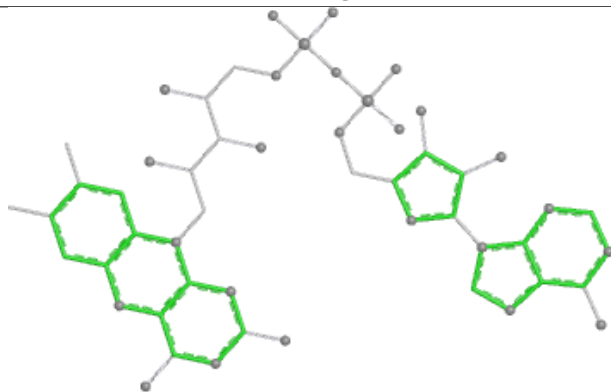
Bond lengths



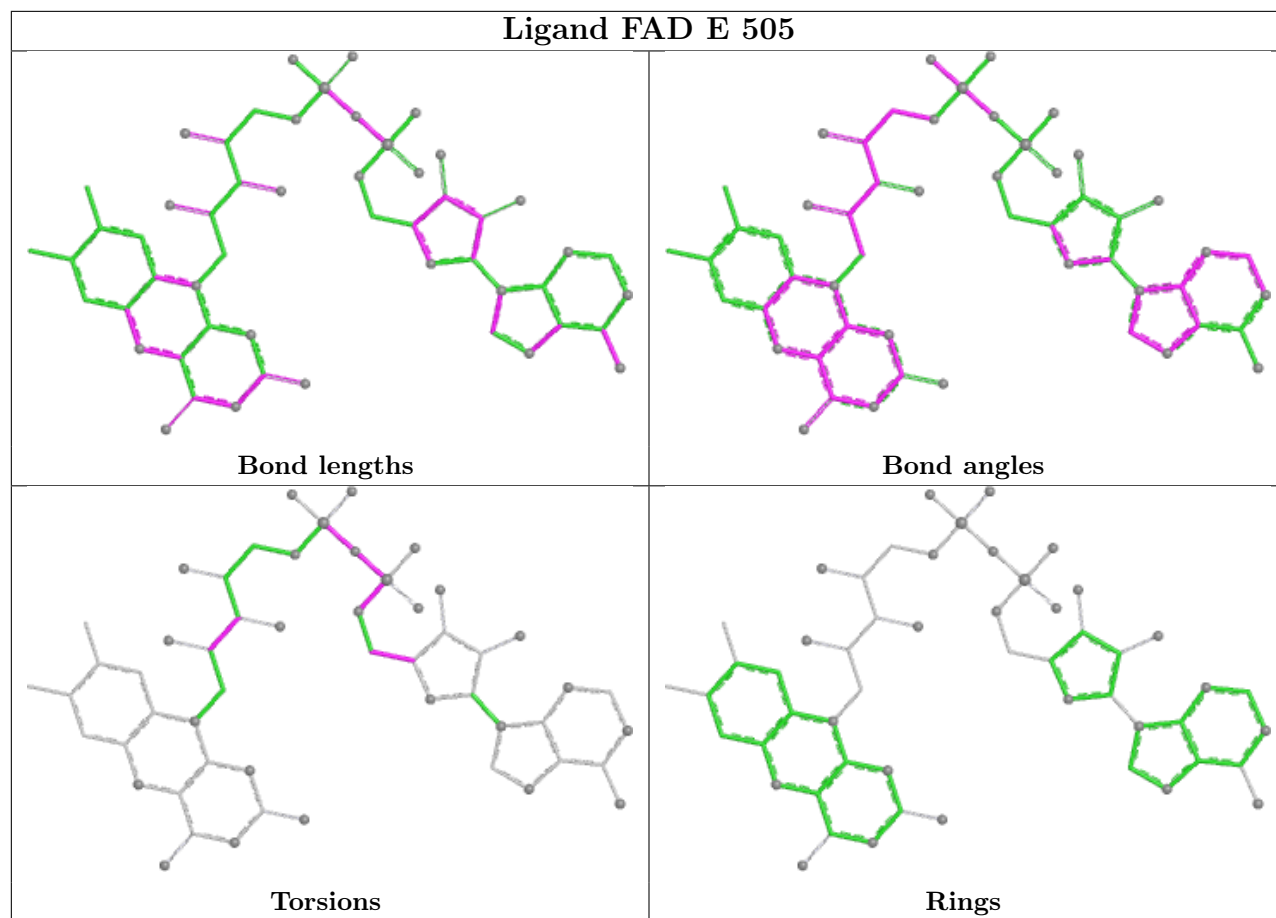
Bond angles

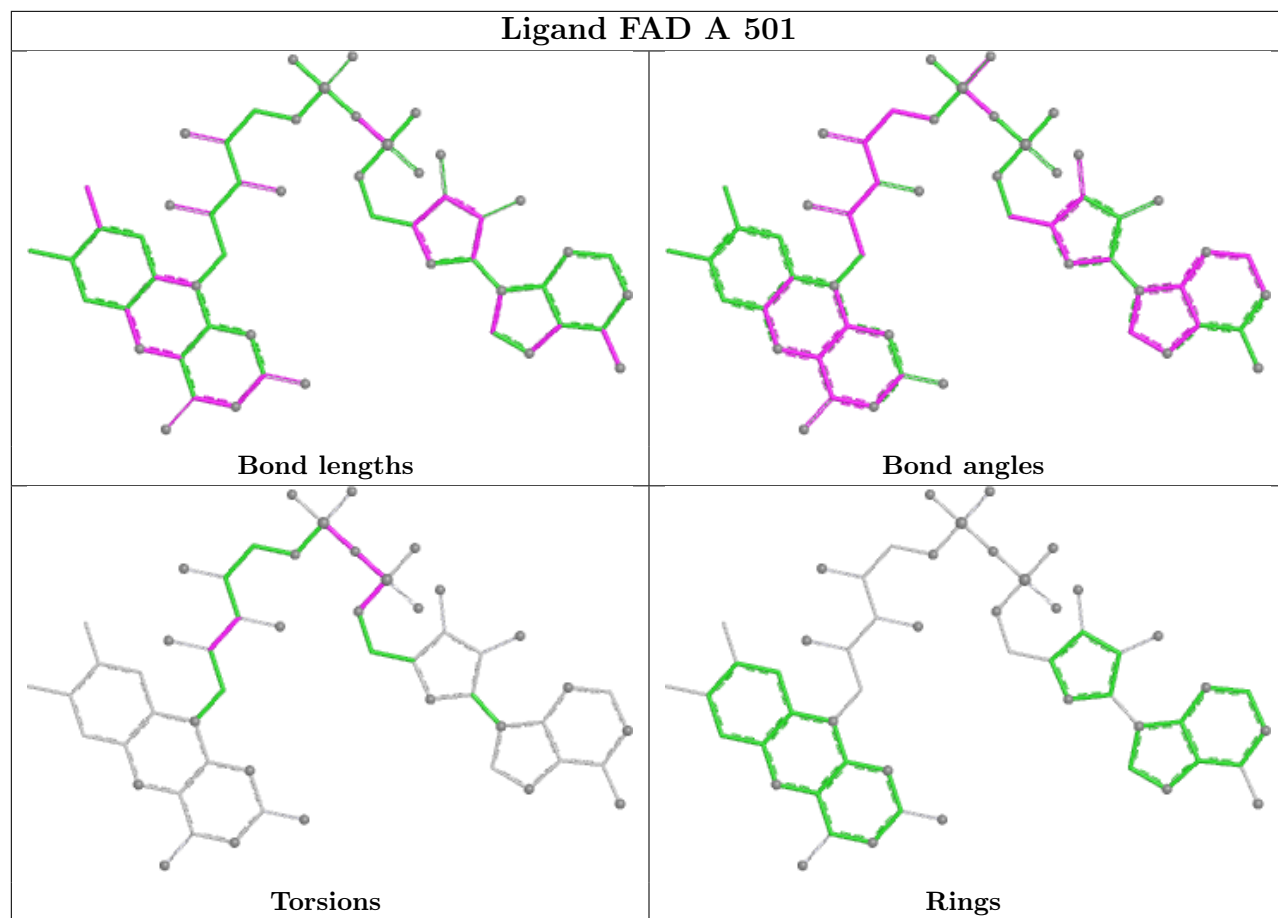


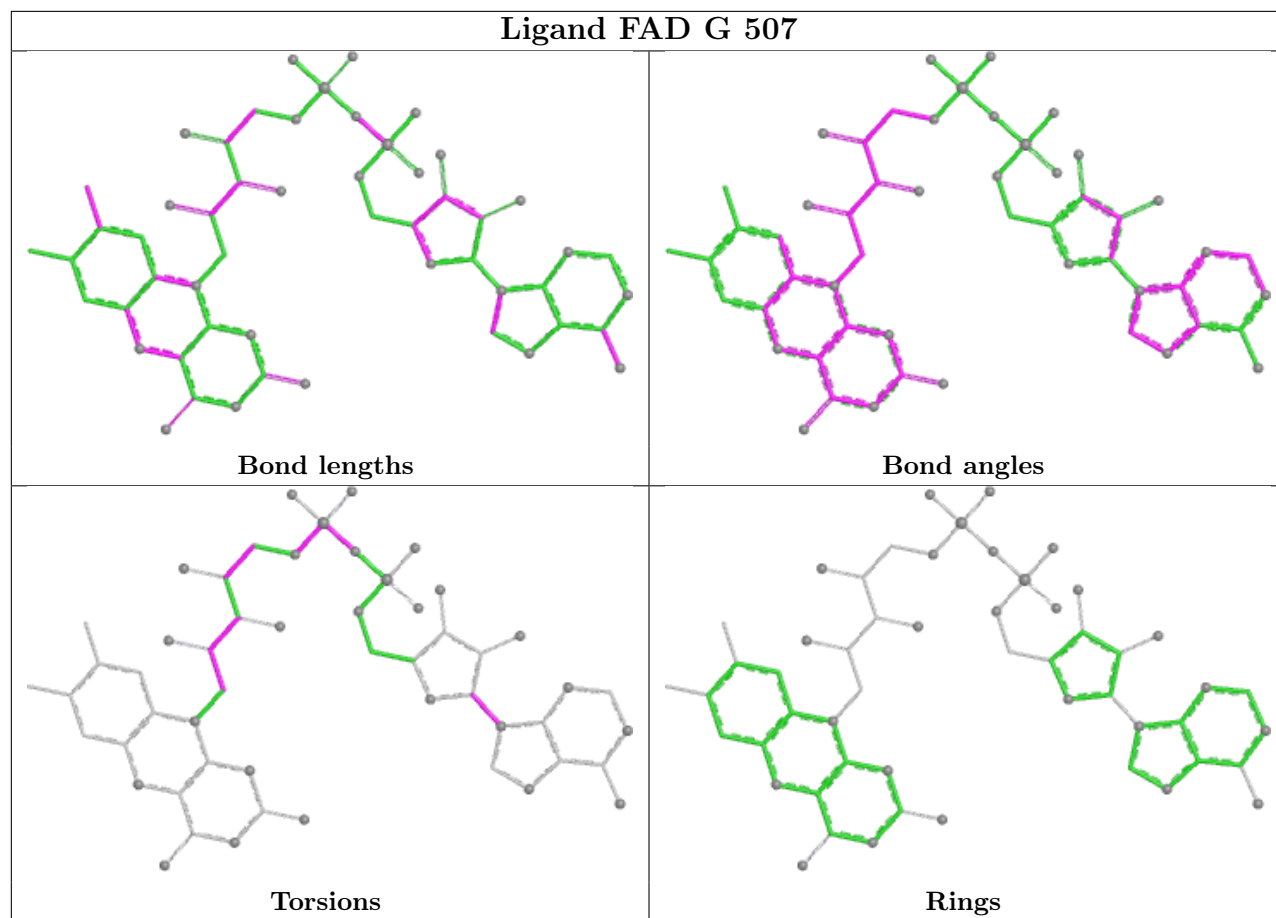
Torsions

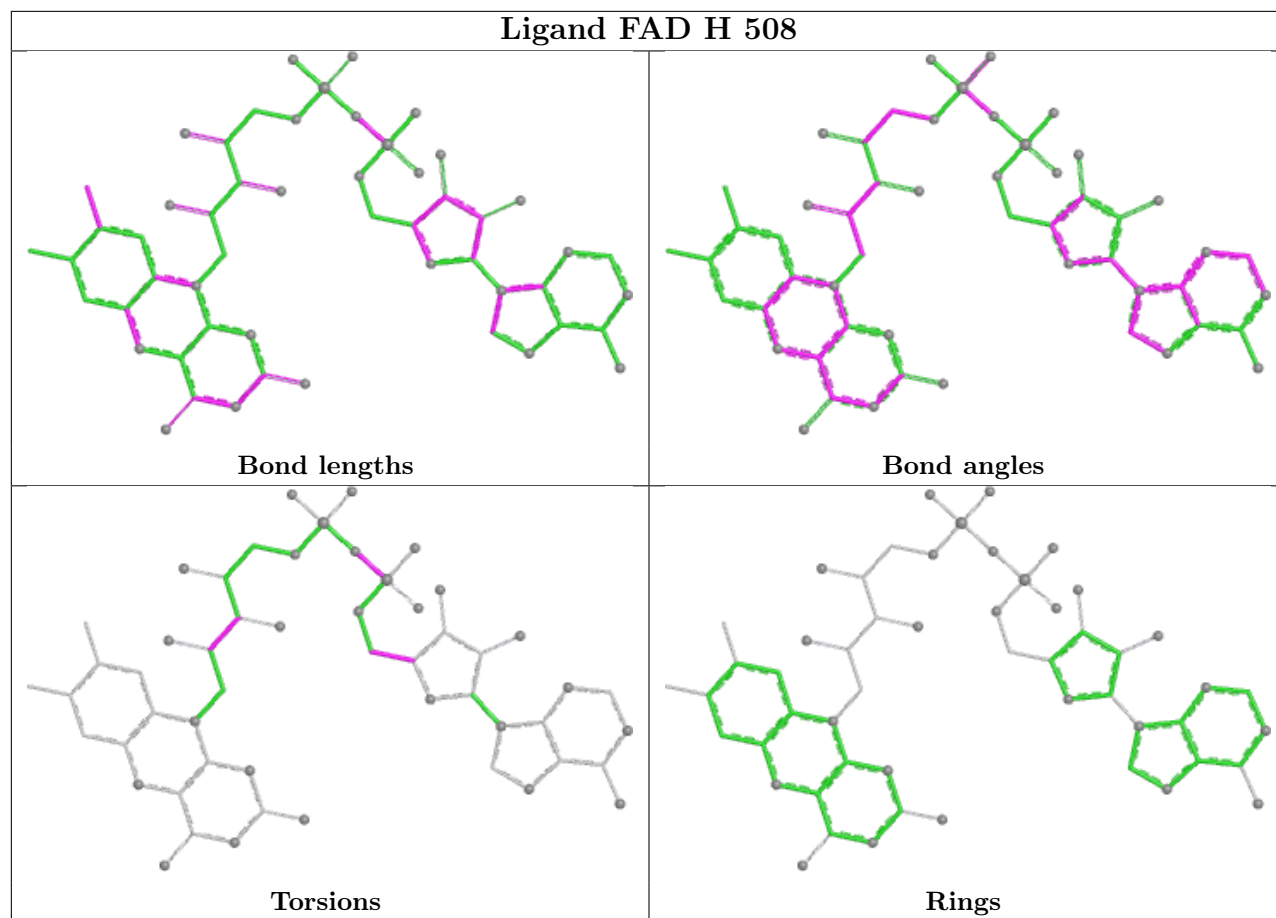


Rings









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	398/421 (94%)	0.10	3 (0%) 82 80	10, 26, 48, 62	0
1	B	391/421 (92%)	0.23	3 (0%) 82 80	14, 29, 50, 65	0
1	C	396/421 (94%)	0.35	8 (2%) 65 60	13, 29, 50, 66	0
1	D	394/421 (93%)	0.27	15 (3%) 44 38	12, 31, 51, 71	0
1	E	397/421 (94%)	0.21	1 (0%) 90 88	11, 26, 46, 64	0
1	F	393/421 (93%)	0.31	11 (2%) 55 49	16, 31, 52, 71	0
1	G	397/421 (94%)	0.42	21 (5%) 32 26	15, 33, 63, 78	0
1	H	394/421 (93%)	0.33	7 (1%) 67 63	18, 33, 51, 65	0
All	All	3160/3368 (93%)	0.28	69 (2%) 62 57	10, 30, 52, 78	0

The worst 5 of 69 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	278	TYR	5.5
1	C	412	GLU	4.9
1	G	280	ILE	4.7
1	C	48	LEU	4.6
1	D	412	GLU	4.0

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](#)

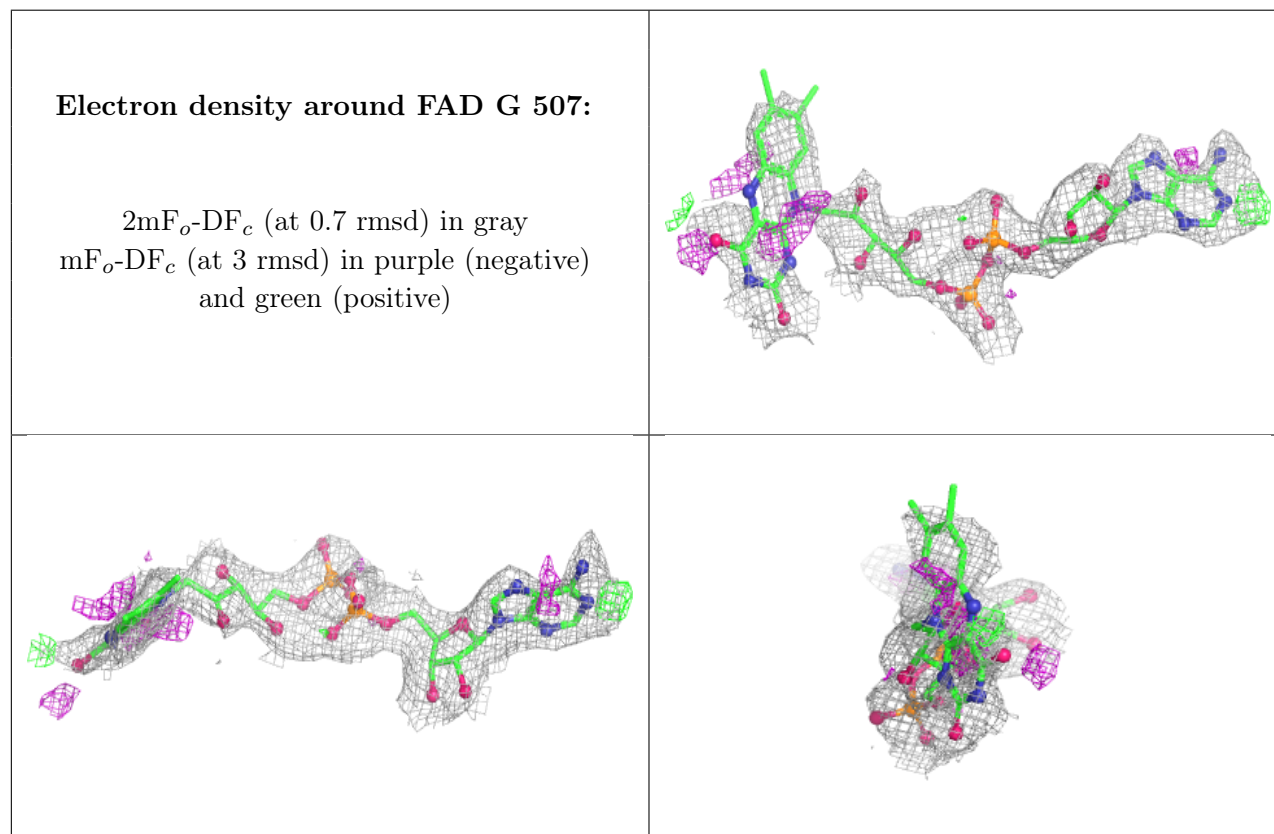
There are no oligosaccharides in this entry.

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.

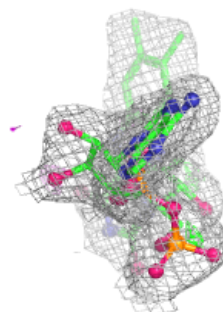
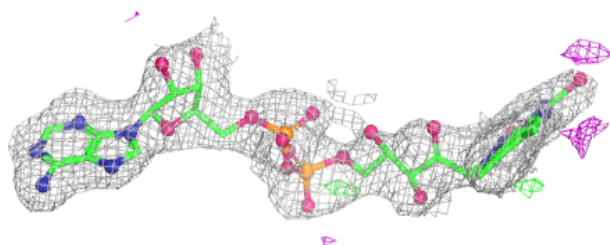
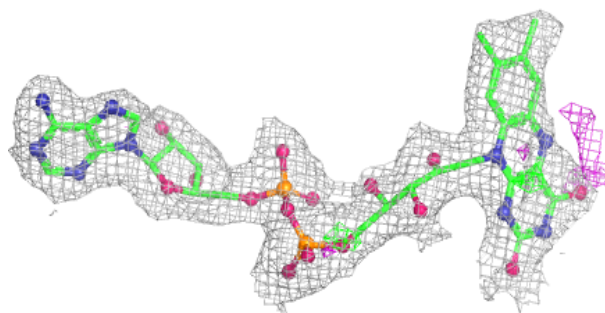
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	FAD	G	507	53/53	0.82	0.17	31,50,64,82	0
2	FAD	F	506	53/53	0.92	0.10	16,33,51,85	0
2	FAD	B	502	53/53	0.93	0.10	3,18,35,63	0
2	FAD	D	504	53/53	0.93	0.10	14,29,42,54	0
2	FAD	C	503	53/53	0.94	0.09	6,20,36,48	0
2	FAD	E	505	53/53	0.94	0.09	7,23,34,39	0
2	FAD	H	508	53/53	0.94	0.09	12,28,38,48	0
2	FAD	A	501	53/53	0.96	0.07	5,19,30,48	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.

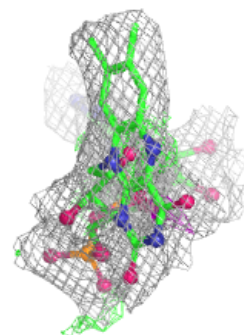
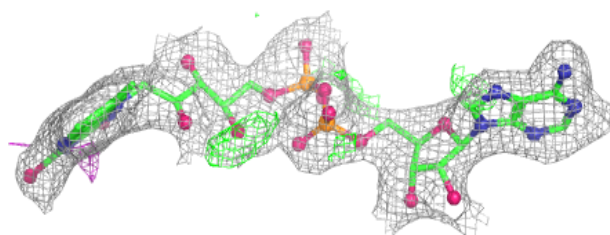
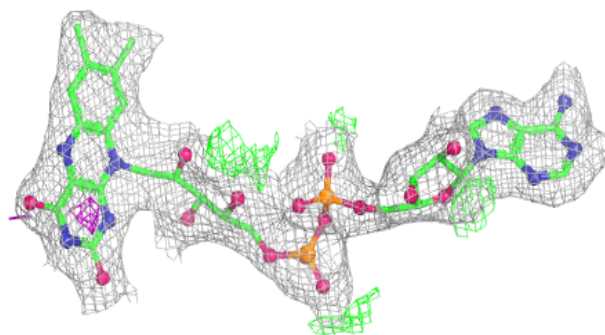


Electron density around FAD F 506:

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

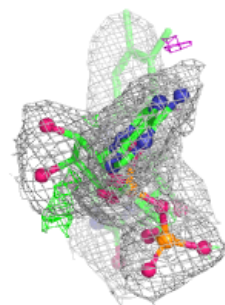
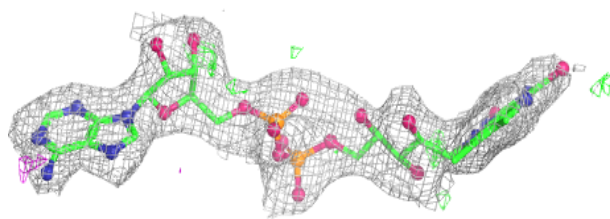
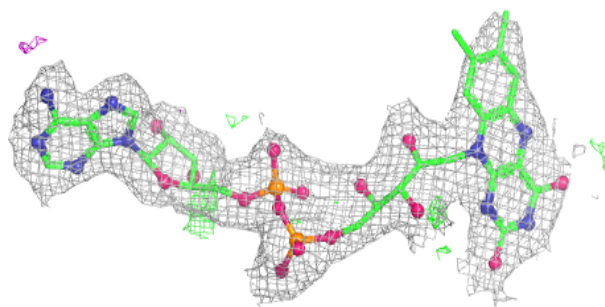
**Electron density around FAD B 502:**

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

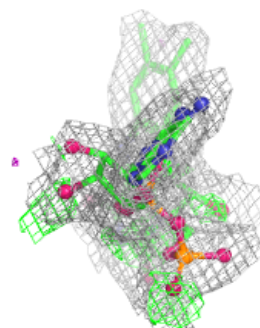
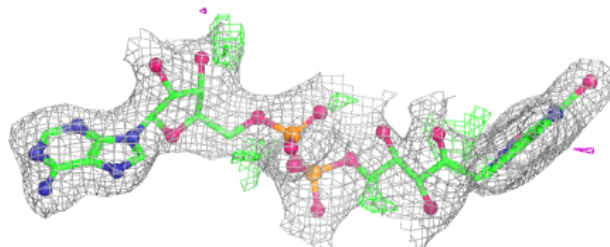
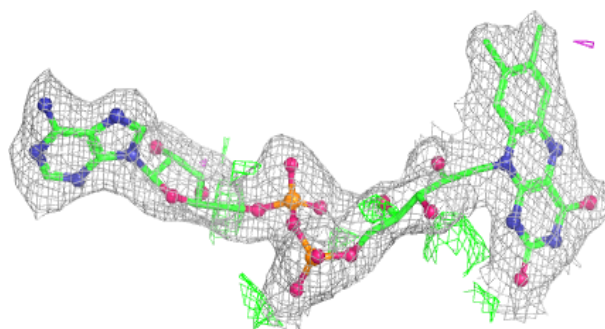


Electron density around FAD D 504:

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

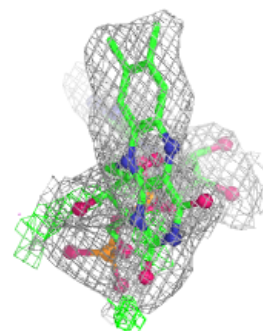
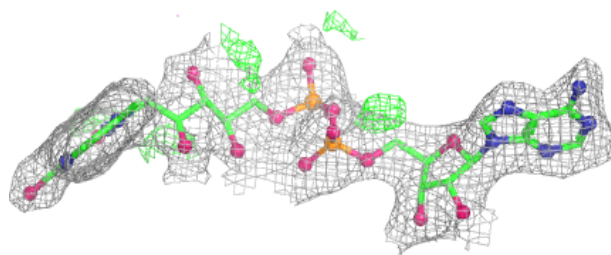
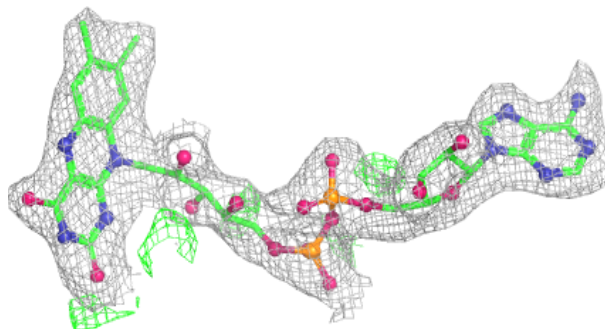
**Electron density around FAD C 503:**

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

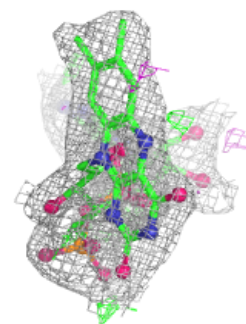
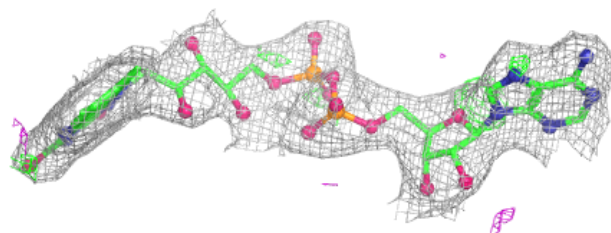
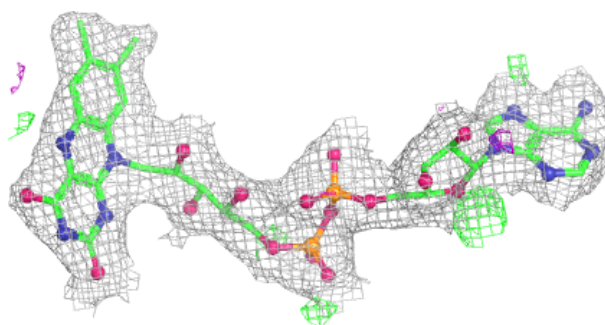


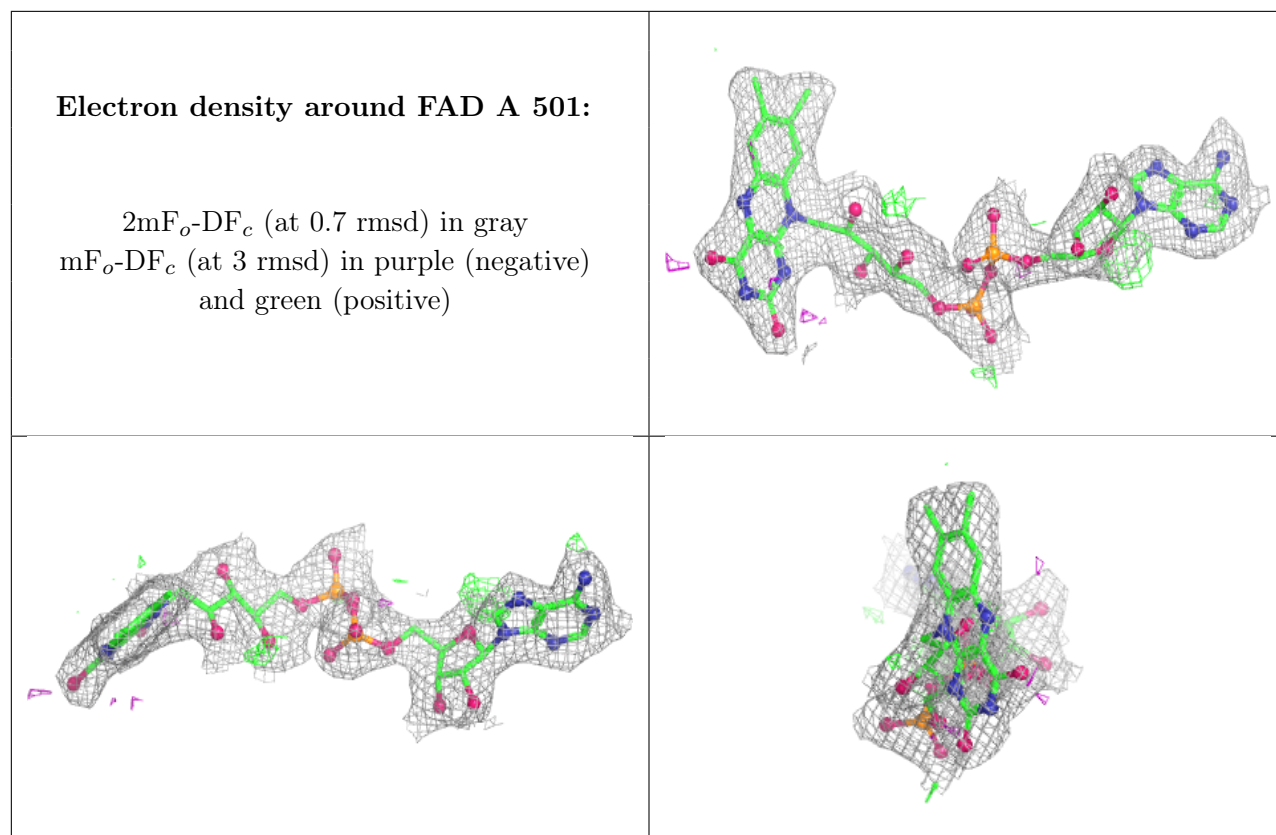
Electron density around FAD E 505:

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

**Electron density around FAD H 508:**

$2mF_o-DF_c$ (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.