



# Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 09:17 am BST

PDB ID : 5GXU  
Title : Crystal structure of Arabidopsis ATR2  
Authors : Niu, G.; Liu, L.  
Deposited on : 2016-09-20  
Resolution : 2.30 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

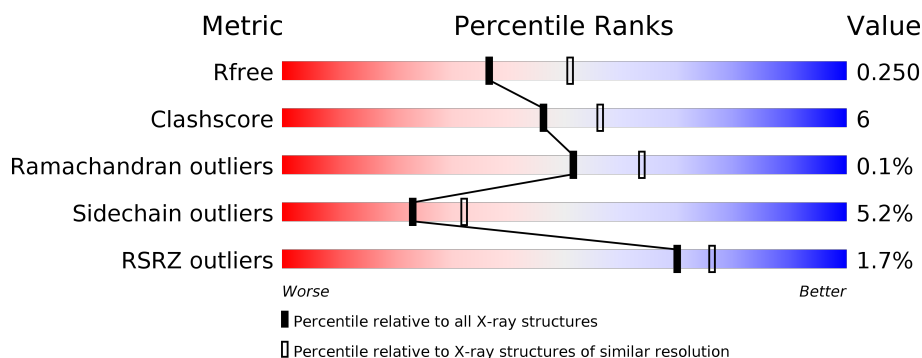
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	640	<div> <div>2%</div> <div> <div></div> <div>74%</div> <div>12%</div> <div>•</div> <div>14%</div> </div> </div>
1	B	640	<div> <div>%</div> <div> <div></div> <div>50%</div> <div>10%</div> <div>•</div> <div>38%</div> </div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7780 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 NADPH–cytochrome P450 reductase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	552	Total	C	N	O	S	0	2	0
			4251	2715	725	790	21			
1	B	394	Total	C	N	O	S	0	3	0
			3017	1918	519	560	20			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	72	GLY	-	expression tag	UNP Q9SUM3
B	72	GLY	-	expression tag	UNP Q9SUM3

- Molecule 2 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: C<sub>17</sub>H<sub>21</sub>N<sub>4</sub>O<sub>9</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	17	4	9	1		


- # FAD

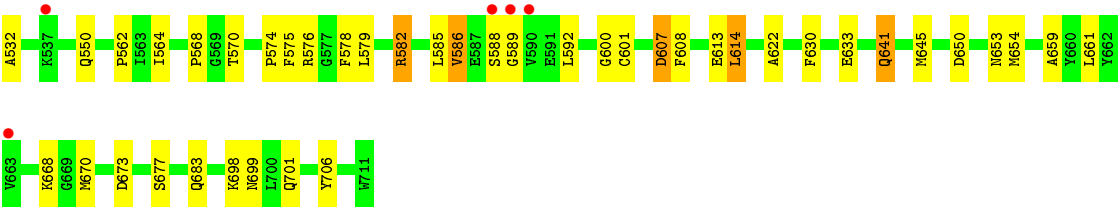
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	205	Total O 205 205	0	0
4	B	170	Total O 170 170	0	0

These plots are drawn for all protein, RNA and DNA 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.

- [illegible]

- Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	55.33Å 61.82Å 88.33Å 97.84° 100.51° 90.27°	Depositor
Resolution (Å)	30.61 – 2.30 30.61 – 2.30	Depositor EDS
% Data completeness (in resolution range)	90.8 (30.61-2.30) 89.7 (30.61-2.30)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.87 (at 2.31Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, $R_{free}$	0.200 , 0.248 0.201 , 0.250	Depositor DCC
$R_{free}$ test set	2310 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	31.0	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 31.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	7780	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.14% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: FMN, 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	0.31	0/4358	0.49	0/5909
1	B	0.33	0/3100	0.52	0/4208
All	All	0.32	0/7458	0.50	0/10117

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	538	SER	Peptide
1	A	606	MET	Peptide

### 5.2 Too-close contacts [i](#)

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



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4251	0	4088	40	0
1	B	3017	0	2934	50	0
2	A	31	0	19	1	0
3	A	53	0	31	4	0
3	B	53	0	31	3	0
4	A	205	0	0	6	0
4	B	170	0	0	10	0
All	All	7780	0	7103	91	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (91) 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:B:505:HIS:HD2	1:B:576:ARG:HH22	1.19	0.91
1:B:588:SER:H	1:B:589:GLY:HA2	1.35	0.90
1:B:505:HIS:CD2	1:B:576:ARG:HH22	2.03	0.76
1:B:582:ARG:HH11	1:B:582:ARG:HG3	1.54	0.71
1:A:121:LYS:HD3	1:A:138:ILE:HD13	1.75	0.67
1:B:683:GLN:OE1	4:B:901:HOH:O	2.14	0.66
1:B:588:SER:N	1:B:589:GLY:HA2	2.09	0.66
1:A:604:ARG:NH1	1:A:633:GLU:OE2	2.29	0.66
1:B:601:CYS:HB2	1:B:607:ASP:HB3	1.78	0.64
1:B:582:ARG:NH1	1:B:622:ALA:O	2.31	0.64
1:B:607:ASP:OD2	4:B:902:HOH:O	2.16	0.63
1:A:145:ALA:HB1	1:A:150:GLU:HB3	1.81	0.63
1:A:595:SER:HB3	1:A:624:ALA:H	1.64	0.62
1:B:650:ASP:HA	1:B:653:ASN:HB2	1.81	0.62
1:A:605:ARG:NH2	4:A:908:HOH:O	2.33	0.62
1:A:489:ARG:HG2	3:A:802:FAD:H3'	1.83	0.61
1:A:560:LYS:HE2	1:A:591:GLU:HB3	1.82	0.60
1:A:186:GLU:OE2	1:A:708[B]:ARG:NH2	2.35	0.60
1:A:576:ARG:NH1	1:A:613:GLU:OE1	2.36	0.58
1:B:485[A]:ARG:NH1	4:B:919:HOH:O	2.36	0.58
1:B:489:ARG:HD2	1:B:525:CYS:HB2	1.84	0.58
1:A:641:GLN:HG2	1:A:673:ASP:HB3	1.86	0.57
1:B:630:PHE:HB3	1:B:633:GLU:HG3	1.88	0.56
1:B:653:ASN:OD1	4:B:904:HOH:O	2.17	0.56
1:A:489:ARG:HD2	1:A:525:CYS:HB2	1.86	0.56
1:B:320:LYS:HD3	1:B:532:ALA:O	2.05	0.56
1:B:425:LEU:HD23	1:B:473:PRO:HG3	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:578:PHE:O	1:B:582:ARG:HG2	2.07	0.54
1:B:641:GLN:HG2	1:B:673:ASP:HB3	1.88	0.54
1:A:582:ARG:NH2	4:A:917:HOH:O	2.40	0.54
1:A:130:ARG:NH1	4:A:914:HOH:O	2.38	0.53
1:B:505:HIS:HD2	1:B:576:ARG:NH2	1.99	0.53
1:B:378:PHE:O	4:B:905:HOH:O	2.19	0.52
1:A:615:GLN:NE2	4:A:902:HOH:O	2.20	0.52
1:A:409:ARG:NH2	1:A:410:TYR:OH	2.43	0.51
1:B:489:ARG:HD3	3:B:801:FAD:O1P	2.11	0.51
1:B:564:ILE:HD11	1:B:654:MET:HG3	1.91	0.51
1:B:588:SER:H	1:B:589:GLY:CA	2.16	0.51
1:A:198:LYS:HB3	1:A:231:LEU:HD21	1.93	0.51
1:A:355:LEU:HD23	1:A:488:PRO:HB3	1.93	0.51
1:B:582:ARG:CG	1:B:582:ARG:HH11	2.23	0.51
1:A:378:PHE:O	4:A:901:HOH:O	2.17	0.50
1:A:188:ASN:O	1:A:190:ARG:HG2	2.10	0.50
1:A:662:TYR:CE1	1:A:707:LEU:HD12	2.47	0.49
1:B:434:GLU:OE2	1:B:469:PRO:HD2	2.13	0.49
1:B:505:HIS:CD2	1:B:576:ARG:NH2	2.76	0.49
1:B:585:LEU:O	1:B:588:SER:HB3	2.12	0.49
1:A:568:PRO:HG3	1:A:670:MET:HG3	1.94	0.48
1:B:698:LYS:HA	1:B:701:GLN:HE21	1.78	0.48
1:B:492:SER:HB3	1:B:574:PRO:HG3	1.96	0.48
1:B:562:PRO:HB2	1:B:659:ALA:HB2	1.94	0.48
1:A:511:VAL:HG21	3:A:802:FAD:H5'2	1.94	0.48
1:B:645:MET:HE1	1:B:677:SER:HA	1.95	0.48
1:A:166:ALA:HA	1:A:204:LEU:HB2	1.95	0.47
1:B:489:ARG:HG2	3:B:801:FAD:H3'	1.96	0.47
1:A:505:HIS:ND1	1:A:576:ARG:NH2	2.64	0.46
1:B:586:VAL:HG13	1:B:592:LEU:HG	1.98	0.46
1:A:595:SER:O	1:A:624:ALA:HB3	2.16	0.46
1:A:528:TRP:CH2	1:A:544:ALA:HB2	2.51	0.45
1:B:668:LYS:HB2	1:B:668:LYS:NZ	2.32	0.45
1:B:613:GLU:OE2	4:B:906:HOH:O	2.21	0.45
1:A:489:ARG:HD3	3:A:802:FAD:O1P	2.17	0.45
1:A:414:LEU:HA	1:A:460:SER:HB2	1.99	0.45
1:B:479:PHE:O	1:B:485[A]:ARG:NH1	2.49	0.45
1:A:511:VAL:HG11	3:A:802:FAD:H3B	1.99	0.44
1:A:564:ILE:HB	1:A:661:LEU:HD22	1.98	0.44
1:A:645:MET:HE1	1:A:680:THR:HB	1.99	0.44
1:A:113:THR:OG1	2:A:801:FMN:O2P	2.30	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:651:ILE:O	1:A:655:ILE:HG12	2.19	0.43
1:B:511:VAL:HG11	3:B:801:FAD:H3B	2.01	0.43
1:B:365:GLU:OE1	1:B:368:ARG:NH1	2.52	0.43
1:B:568:PRO:HG2	1:B:670:MET:HG3	2.01	0.43
1:A:160:VAL:HG22	1:A:198:LYS:HB2	2.01	0.42
1:B:503:ARG:HD3	4:B:980:HOH:O	2.18	0.42
1:B:399:PRO:HA	1:B:400:PRO:HD3	1.86	0.42
1:A:434:GLU:OE2	1:A:469:PRO:HD2	2.20	0.42
1:B:575:PHE:O	1:B:579:LEU:HG	2.20	0.42
1:B:568:PRO:HA	1:B:600:GLY:H	1.85	0.42
1:A:473:PRO:HA	1:A:474:PRO:HD3	1.95	0.42
1:B:396:PRO:HA	1:B:397:PRO:HD3	1.89	0.42
1:B:431:ASP:HA	1:B:432:PRO:HD3	1.92	0.42
1:B:550:GLN:HG2	4:B:910:HOH:O	2.19	0.42
1:B:416:SER:HA	1:B:417:PRO:HD3	1.90	0.41
1:A:104:LYS:HB2	1:A:104:LYS:HE3	1.79	0.41
1:A:416:SER:HA	1:A:417:PRO:HD3	1.88	0.41
1:B:417:PRO:HG2	1:B:464:VAL:HG21	2.02	0.41
1:B:614:LEU:HD12	1:B:614:LEU:HA	1.86	0.40
1:A:380:LEU:HD21	1:A:476:GLY:HA2	2.03	0.40
1:B:699:ASN:ND2	4:B:911:HOH:O	2.25	0.40
1:A:489:ARG:HG3	4:A:1072:HOH:O	2.22	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	548/640 (86%)	531 (97%)	16 (3%)	1 (0%)	47	58
1	B	393/640 (61%)	379 (96%)	14 (4%)	0	100	100
All	All	941/1280 (74%)	910 (97%)	30 (3%)	1 (0%)	51	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	169	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	434/538 (81%)	412 (95%)	22 (5%)	24	33
1	B	320/538 (60%)	302 (94%)	18 (6%)	21	29
All	All	754/1076 (70%)	714 (95%)	40 (5%)	23	31

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	104	LYS
1	A	134	THR
1	A	142	ASP
1	A	150	GLU
1	A	152	GLU
1	A	172	GLU
1	A	197	LEU
1	A	201	VAL
1	A	248	THR
1	A	305	THR
1	A	320	LYS
1	A	330	ARG
1	A	433	THR
1	A	489	ARG
1	A	503	ARG
1	A	524	VAL
1	A	570	THR
1	A	607	ASP
1	A	608	PHE
1	A	614	LEU
1	A	619	GLU
1	A	661	LEU

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Mol	Chain	Res	Type
1	B	321	ARG
1	B	485[A]	ARG
1	B	485[B]	ARG
1	B	489	ARG
1	B	501	GLU
1	B	503	ARG
1	B	505	HIS
1	B	506	VAL
1	B	524	VAL
1	B	570	THR
1	B	582	ARG
1	B	586	VAL
1	B	607	ASP
1	B	608	PHE
1	B	614	LEU
1	B	641	GLN
1	B	661	LEU
1	B	706	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	505	HIS
1	B	683	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

3 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	FAD	B	801	-	51,58,58	2.90	16 (31%)	60,89,89	1.75	10 (16%)
2	FMN	A	801	-	31,33,33	1.30	4 (12%)	40,50,50	1.88	6 (15%)
3	FAD	A	802	-	51,58,58	2.84	16 (31%)	60,89,89	1.67	8 (13%)

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
3	FAD	B	801	-	-	2/30/50/50	0/6/6/6
2	FMN	A	801	-	-	4/18/18/18	0/3/3/3
3	FAD	A	802	-	-	1/30/50/50	0/6/6/6

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	801	FAD	C4X-C10	12.33	1.51	1.38
3	A	802	FAD	C4X-C10	11.78	1.50	1.38
3	B	801	FAD	C5X-N5	8.85	1.50	1.35
3	A	802	FAD	C5X-N5	8.72	1.49	1.35
3	A	802	FAD	C2B-C1B	-5.49	1.45	1.53
3	B	801	FAD	C2B-C1B	-5.33	1.45	1.53
3	B	801	FAD	O4B-C1B	4.54	1.47	1.41
3	A	802	FAD	C2-N1	4.47	1.47	1.38
3	A	802	FAD	O4B-C1B	4.29	1.47	1.41
3	B	801	FAD	C2-N1	4.22	1.46	1.38
3	B	801	FAD	C9A-N10	3.78	1.43	1.38
3	B	801	FAD	C4-N3	3.70	1.39	1.33
3	A	802	FAD	C4-N3	3.59	1.39	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	802	FAD	C9A-N10	3.58	1.43	1.38
3	B	801	FAD	C2B-C3B	-3.53	1.43	1.53
3	A	802	FAD	C2B-C3B	-3.45	1.43	1.53
3	B	801	FAD	C4-C4X	3.43	1.47	1.41
3	A	802	FAD	C2A-N3A	3.42	1.37	1.32
3	B	801	FAD	C6A-N6A	3.11	1.45	1.34
3	B	801	FAD	C2A-N3A	3.06	1.37	1.32
3	A	802	FAD	C6A-N6A	3.00	1.45	1.34
3	A	802	FAD	C4-C4X	2.93	1.46	1.41
2	A	801	FMN	C4A-C10	2.81	1.41	1.38
2	A	801	FMN	O2'-C2'	-2.66	1.37	1.43
3	B	801	FAD	C3B-C4B	-2.62	1.46	1.53
3	A	802	FAD	O3'-C3'	-2.60	1.36	1.43
3	A	802	FAD	O4'-C4'	-2.60	1.37	1.43
3	A	802	FAD	C3B-C4B	-2.59	1.46	1.53
3	B	801	FAD	O3'-C3'	-2.55	1.37	1.43
3	B	801	FAD	C2-N3	2.50	1.43	1.38
2	A	801	FMN	O3'-C3'	-2.45	1.37	1.43
2	A	801	FMN	O4'-C4'	-2.44	1.38	1.43
3	B	801	FAD	O4'-C4'	-2.33	1.38	1.43
3	A	802	FAD	C2-N3	2.33	1.42	1.38
3	A	802	FAD	C4'-C3'	-2.22	1.49	1.53
3	B	801	FAD	O2'-C2'	-2.15	1.38	1.43

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	801	FAD	C4-N3-C2	6.72	120.82	115.14
3	A	802	FAD	C4-N3-C2	6.70	120.80	115.14
2	A	801	FMN	C1'-N10-C9A	6.56	123.45	118.29
2	A	801	FMN	C4-N3-C2	5.97	120.18	115.14
3	B	801	FAD	N3A-C2A-N1A	-5.30	120.39	128.68
3	A	802	FAD	N3A-C2A-N1A	-5.19	120.57	128.68
2	A	801	FMN	C4A-N5-C5A	4.41	121.18	116.77
3	B	801	FAD	C4X-N5-C5X	3.69	120.46	116.77
3	A	802	FAD	C4X-N5-C5X	3.57	120.33	116.77
3	B	801	FAD	C4-C4X-C10	-3.51	117.63	119.95
3	A	802	FAD	C3B-C2B-C1B	3.31	105.97	100.98
3	B	801	FAD	C4-C4X-N5	3.07	122.11	118.60
3	B	801	FAD	C3B-C2B-C1B	2.94	105.40	100.98
3	B	801	FAD	P-O3P-PA	-2.85	123.04	132.83
3	A	802	FAD	C4X-C10-N10	-2.81	117.42	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	802	FAD	C4X-C4-N3	-2.69	119.76	123.43
3	B	801	FAD	C4X-C4-N3	-2.51	119.99	123.43
3	B	801	FAD	C4X-C10-N10	-2.42	117.81	120.30
3	A	802	FAD	C4-C4X-C10	-2.38	118.37	119.95
2	A	801	FMN	C4-C4A-N5	2.15	121.06	118.60
2	A	801	FMN	C4A-C4-N3	-2.13	120.51	123.43
2	A	801	FMN	C9A-N10-C10	-2.12	119.14	121.91
3	B	801	FAD	C4A-C5A-N7A	-2.06	107.26	109.40
3	A	802	FAD	C4-C4X-N5	2.04	120.93	118.60

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	801	FAD	O2'-C2'-C3'-C4'
2	A	801	FMN	C2'-C3'-C4'-O4'
2	A	801	FMN	O3'-C3'-C4'-O4'
2	A	801	FMN	C2'-C3'-C4'-C5'
3	B	801	FAD	O2'-C2'-C3'-O3'
2	A	801	FMN	O3'-C3'-C4'-C5'
3	A	802	FAD	O2'-C2'-C3'-C4'

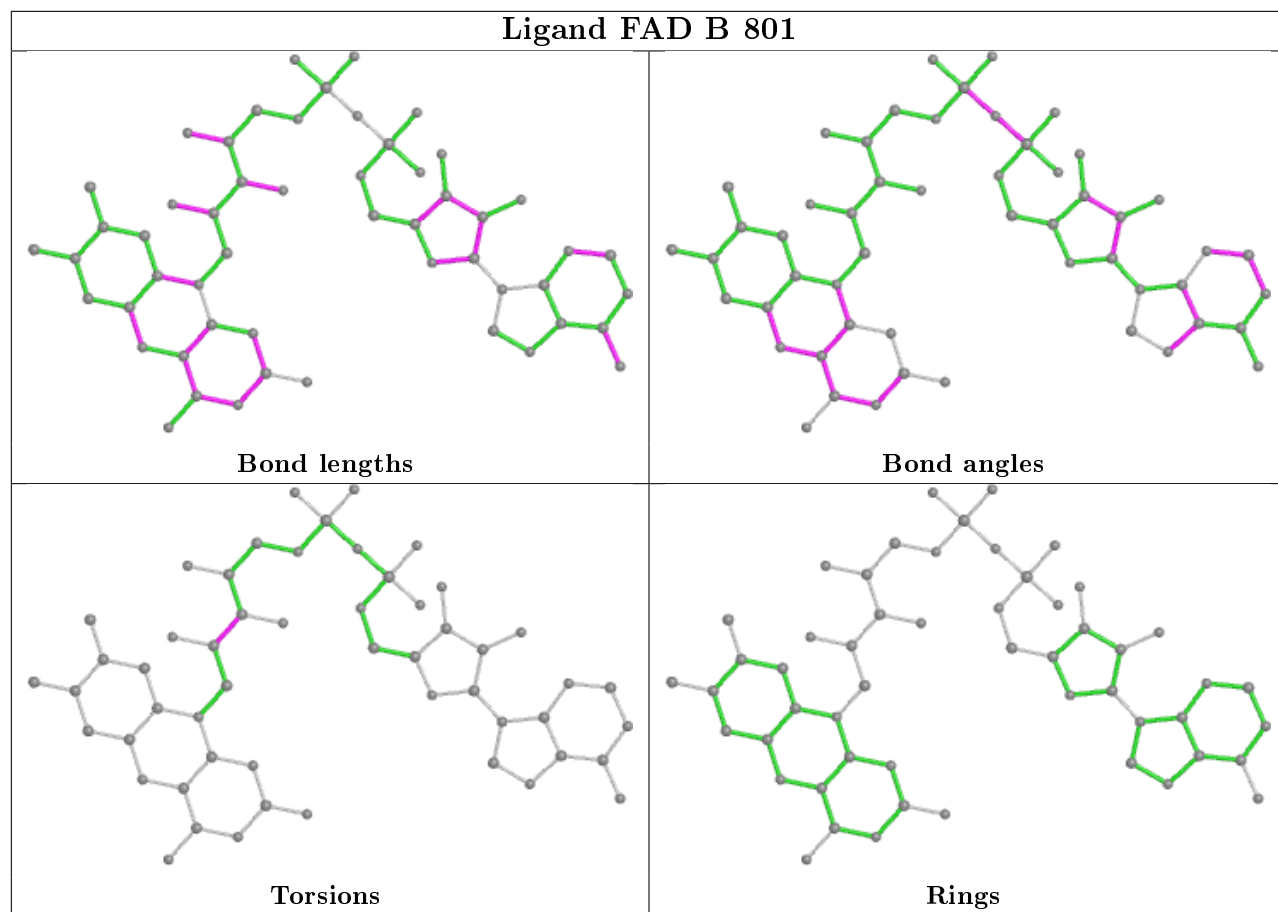
There are no ring outliers.

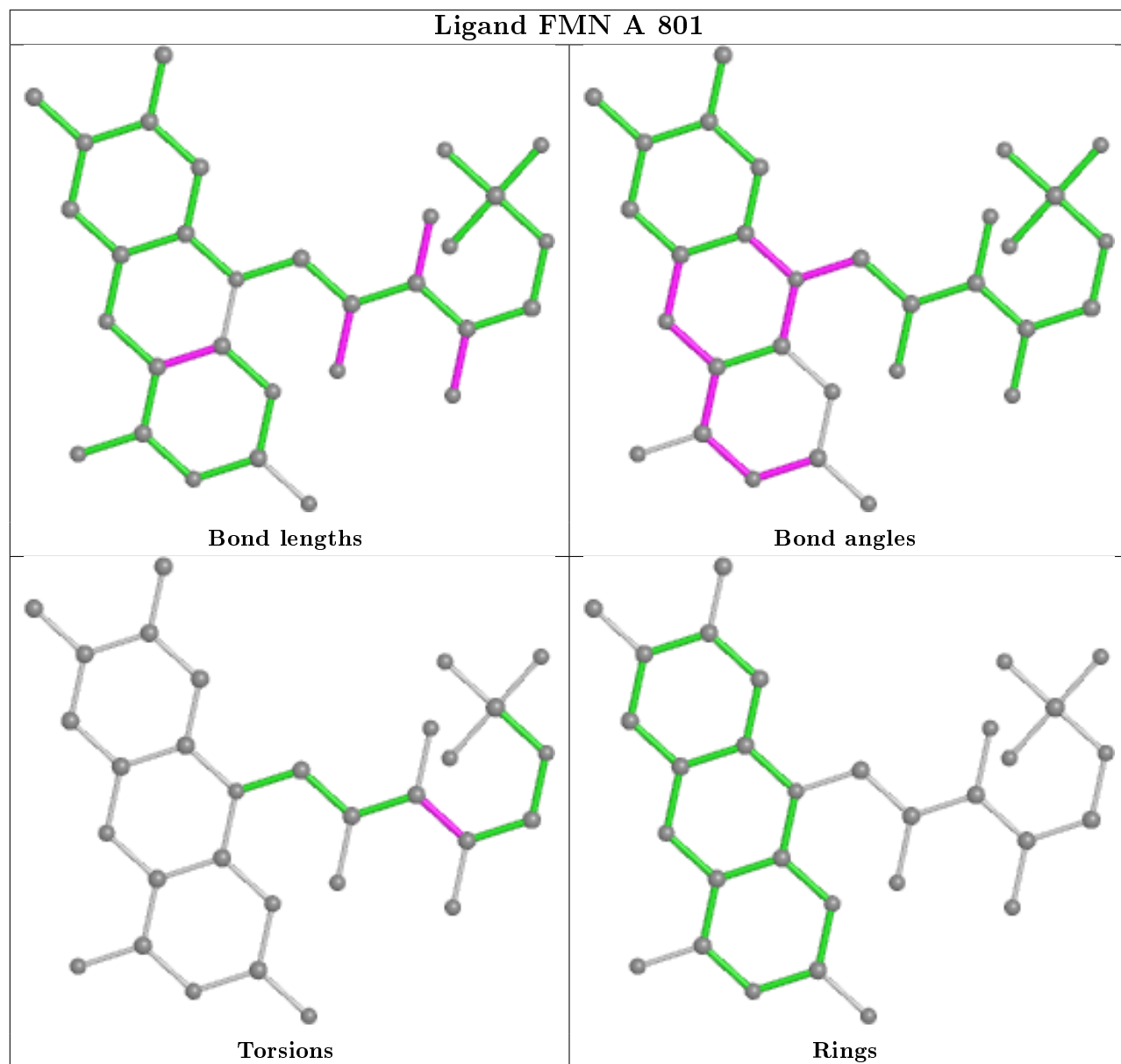
3 monomers are involved in 8 short contacts:

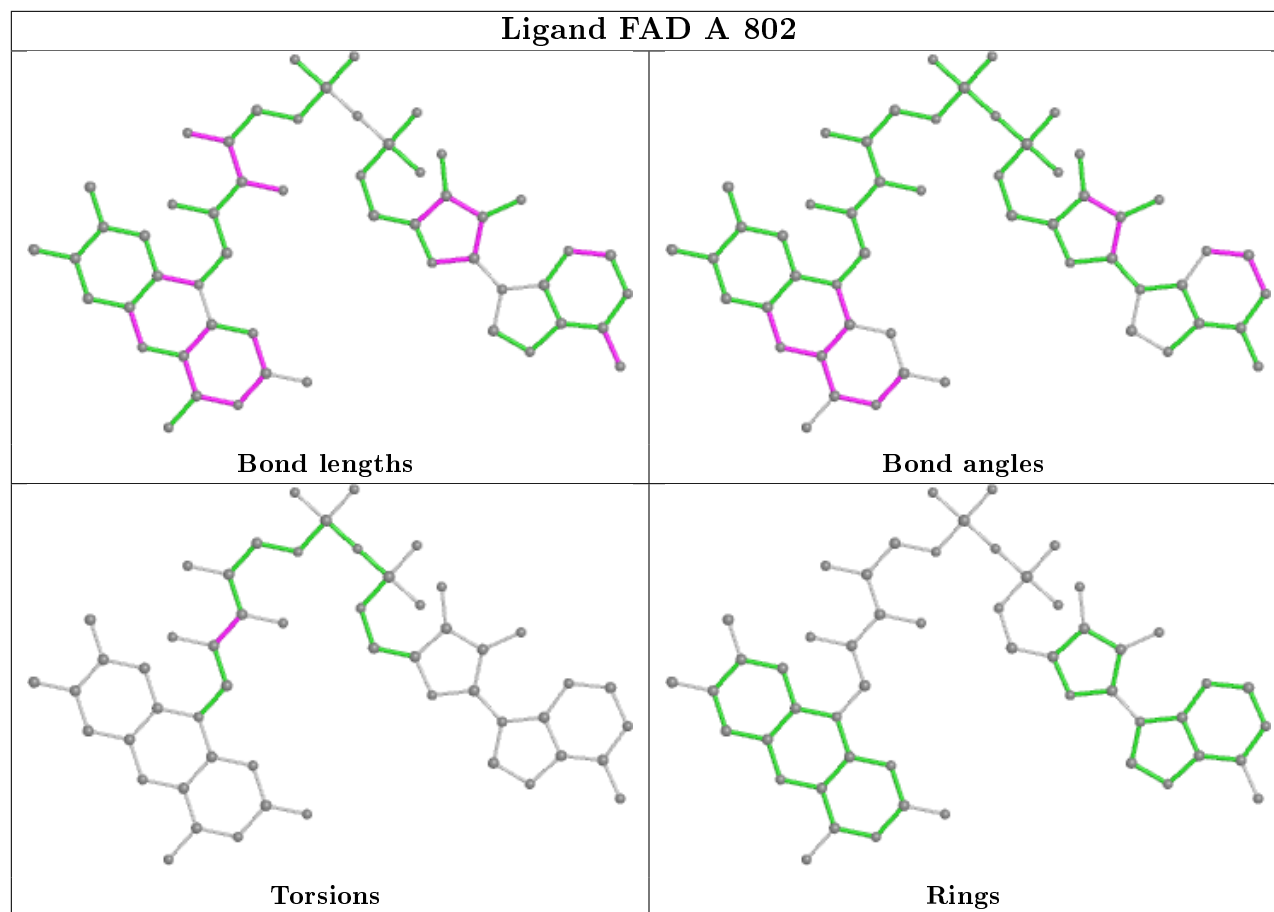
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	801	FAD	3	0
2	A	801	FMN	1	0
3	A	802	FAD	4	0

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.









## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q < 0.9
1	A	552/640 (86%)	0.04	11 (1%) 65 71	8, 23, 43, 51	5 (0%)
1	B	394/640 (61%)	-0.05	5 (1%) 77 81	8, 18, 32, 39	2 (0%)
All	All	946/1280 (73%)	0.00	16 (1%) 70 76	8, 20, 40, 51	7 (0%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	589	GLY	3.4
1	A	156	LYS	3.4
1	A	129	ALA	3.0
1	A	195	LYS	2.7
1	A	262	LEU	2.7
1	A	133	LYS	2.6
1	A	240	ASP	2.5
1	B	590	VAL	2.4
1	A	538	SER	2.3
1	A	379	SER	2.2
1	A	211	HIS	2.1
1	B	588	SER	2.1
1	B	537	LYS	2.1
1	A	132	GLU	2.1
1	B	663	VAL	2.0
1	A	663	VAL	2.0

### 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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

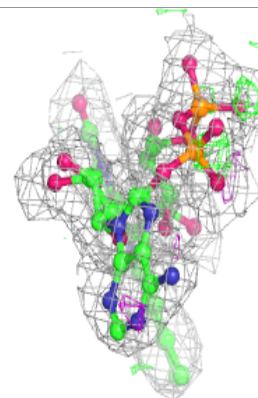
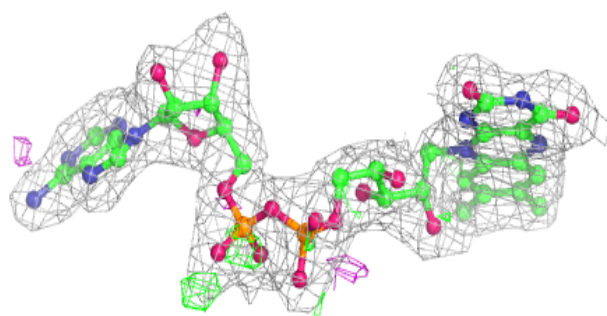
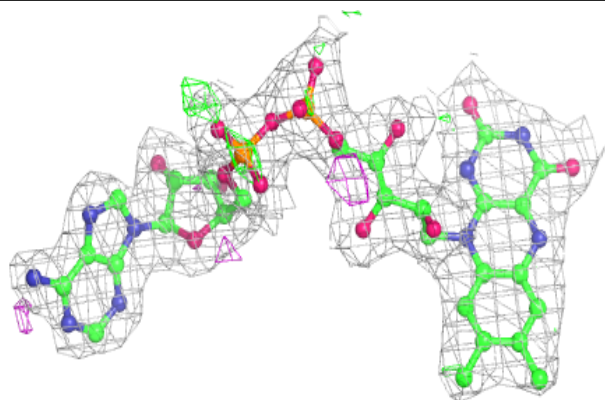
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	FAD	B	801	53/53	0.95	0.13	7,10,18,20	0
2	FMN	A	801	31/31	0.95	0.11	27,34,36,36	0
3	FAD	A	802	53/53	0.95	0.15	8,11,21,21	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.

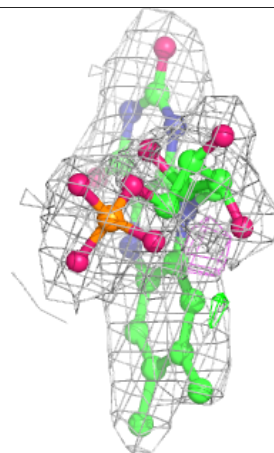
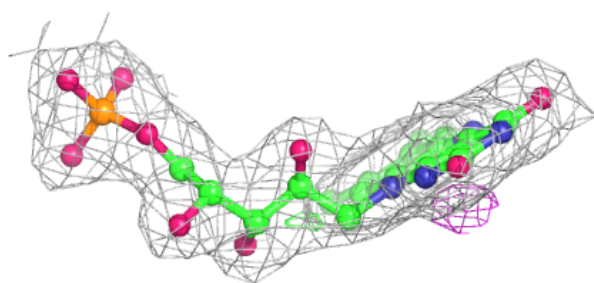
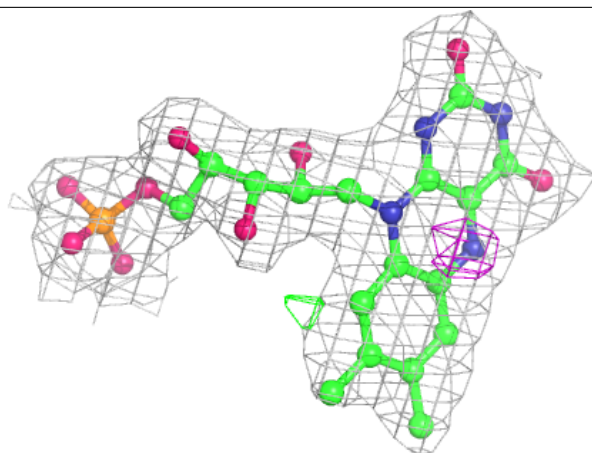
### Electron density around FAD B 801:

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)

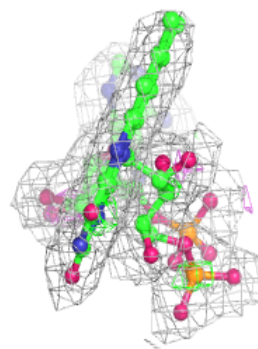
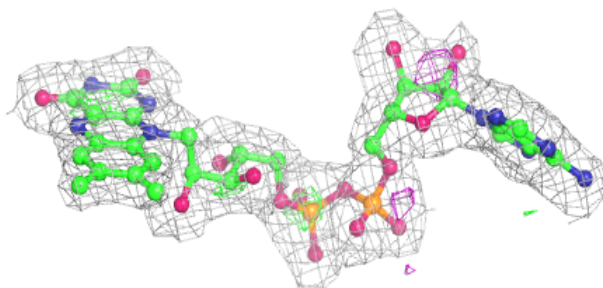
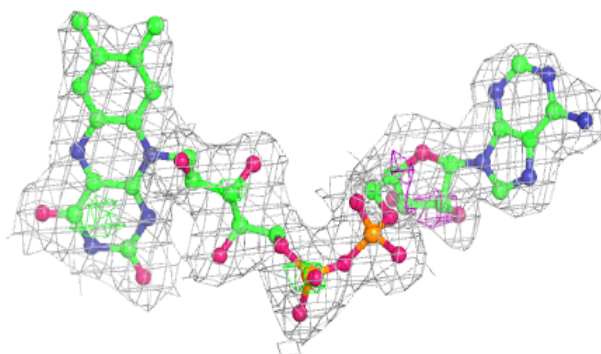


**Electron density around FMN A 801:**

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

**Electron density around FAD A 802:**

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



## 6.5 Other polymers

There are no such residues in this entry.