



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:09 AM UTC

PDB ID : 5GXU / pdb_00005gxu
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

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

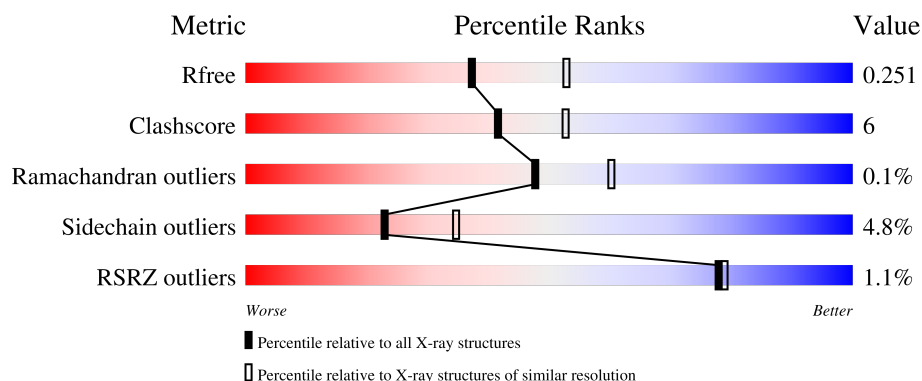
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

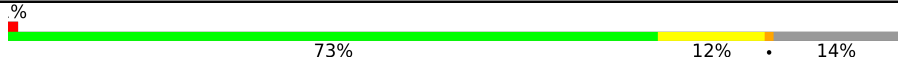
The reported resolution of this entry is 2.30 Å.

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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (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 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	640	
1	B	640	

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.

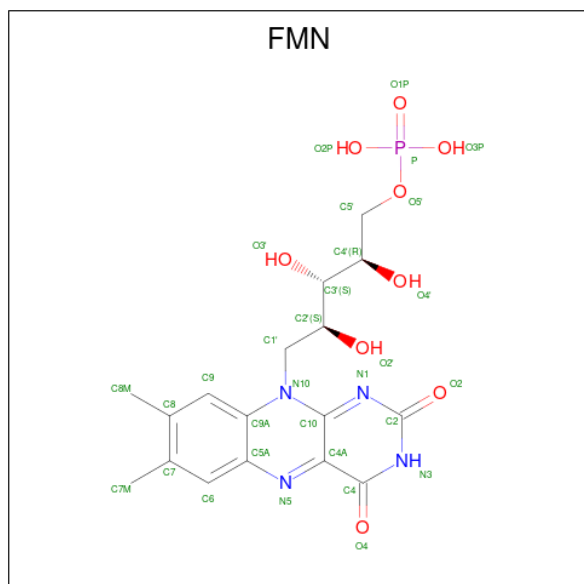
- Molecule 1 is a protein called NADPH-cytochrome P450 reductase 2.

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:

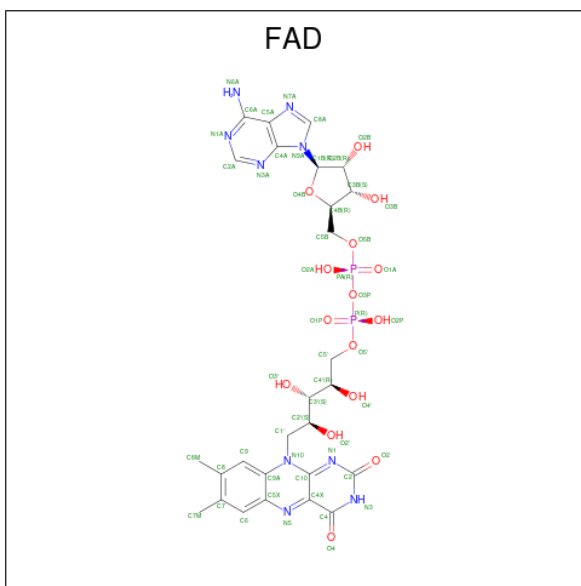
Chain	Residue	Modelled	Actual	Comment	Reference
A	72	GLY	-	expression tag	UNP Q9SUM3
B	72	GLY	-	expression tag	UNP Q9SUM3

- Molecule 2 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: C₁₇H₂₁N₄O₉P).



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

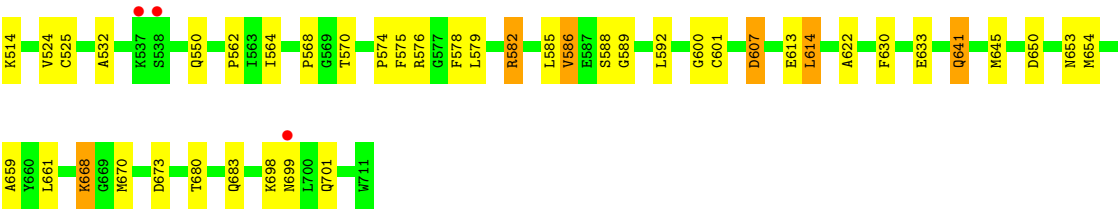
- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

- 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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	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) 89.6 (30.61-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.87 (at 2.31Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, R_{free}	0.200 , 0.248 0.202 , 0.251	Depositor DCC
R_{free} test set	2310 reflections (4.01%)	wwPDB-VP
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 22.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	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 ²)	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.*

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, FMN

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	0.41	0/4358	0.83	7/5909 (0.1%)
1	B	0.43	0/3100	0.81	2/4208 (0.0%)
All	All	0.42	0/7458	0.82	9/10117 (0.1%)

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.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	539	GLU	N-CA-C	-9.63	101.09	113.12
1	A	561	VAL	CA-C-N	6.04	125.98	119.76
1	A	561	VAL	C-N-CA	6.04	125.98	119.76
1	B	468	PHE	CA-C-N	5.78	125.91	119.32
1	B	468	PHE	C-N-CA	5.78	125.91	119.32
1	A	634	GLY	CA-C-N	5.51	125.60	119.32
1	A	634	GLY	C-N-CA	5.51	125.60	119.32
1	A	496	SER	CA-C-N	5.10	125.39	119.47
1	A	496	SER	C-N-CA	5.10	125.39	119.47

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

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	41	0
1	B	3017	0	2934	51	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	93	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 (93) 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:327:GLU:HG2	1:A:606:MET:HE2	1.74	0.69
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:645:MET:HE1	1:B:680:THR:HB	1.89	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:645:MET:HE1	1:A:680:THR:HB	1.90	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:HH11	1:B:582:ARG:CG	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:A:166:ALA:HA	1:A:204:LEU:HB2	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:113:THR:OG1	2:A:801:FMN:O2P	2.30	0.44
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:B:670:MET:HE3	1:B:670:MET:HB3	1.97	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

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	548/640 (86%)	531 (97%)	16 (3%)	1 (0%)	43	55
1	B	393/640 (61%)	379 (96%)	14 (4%)	0	100	100
All	All	941/1280 (74%)	910 (97%)	30 (3%)	1 (0%)	48	60

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%)	414 (95%)	20 (5%)	24	36
1	B	320/538 (60%)	303 (95%)	17 (5%)	20	30
All	All	754/1076 (70%)	717 (95%)	37 (5%)	23	34

All (37) 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

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Mol	Chain	Res	Type
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	433	THR
1	A	489	ARG
1	A	503	ARG
1	A	506	VAL
1	A	524	VAL
1	A	570	THR
1	A	614	LEU
1	A	619	GLU
1	A	661	LEU
1	B	305	THR
1	B	485[A]	ARG
1	B	485[B]	ARG
1	B	489	ARG
1	B	501	GLU
1	B	503	ARG
1	B	506	VAL
1	B	514	LYS
1	B	524	VAL
1	B	570	THR
1	B	582	ARG
1	B	586	VAL
1	B	607	ASP
1	B	614	LEU
1	B	641	GLN
1	B	661	LEU
1	B	668	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	458	GLN
1	A	685	GLN
1	B	381	HIS
1	B	505	HIS
1	B	683	GLN

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 ⓘ

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	$\# Z > 2$
2	FMN	A	801	-	33,33,33	1.36	5 (15%)	48,50,50	1.34	7 (14%)
3	FAD	A	802	-	58,58,58	1.78	15 (25%)	85,89,89	1.61	14 (16%)
3	FAD	B	801	-	58,58,58	1.79	14 (24%)	85,89,89	1.62	14 (16%)

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	FMN	A	801	-	-	5/18/18/18	0/3/3/3
3	FAD	A	802	-	-	4/34/50/50	0/6/6/6
3	FAD	B	801	-	-	5/34/50/50	0/6/6/6

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	801	FAD	C5X-N5	5.75	1.50	1.39
3	A	802	FAD	C5X-N5	5.63	1.49	1.39
3	A	802	FAD	C2-N1	4.59	1.47	1.36
3	B	801	FAD	C6A-N6A	4.38	1.45	1.34
3	B	801	FAD	C2-N1	4.37	1.46	1.36
3	A	802	FAD	C6A-N6A	4.22	1.45	1.34
3	B	801	FAD	C2B-C3B	-3.58	1.43	1.53
3	A	802	FAD	C2B-C3B	-3.49	1.43	1.53
3	B	801	FAD	C5A-N7A	-2.81	1.34	1.39
3	A	802	FAD	C5A-N7A	-2.78	1.34	1.39
2	A	801	FMN	C5A-N5	-2.71	1.34	1.39
2	A	801	FMN	O2'-C2'	-2.68	1.37	1.43
3	B	801	FAD	C3B-C4B	-2.64	1.46	1.53
3	A	802	FAD	O4'-C4'	-2.62	1.37	1.43
3	A	802	FAD	C3B-C4B	-2.61	1.46	1.53
3	A	802	FAD	C2B-C1B	-2.56	1.45	1.53
3	B	801	FAD	C8A-N9A	-2.50	1.33	1.37
3	B	801	FAD	C2B-C1B	-2.48	1.45	1.53
3	A	802	FAD	O3'-C3'	-2.47	1.36	1.43
2	A	801	FMN	O4'-C4'	-2.46	1.38	1.43
3	A	802	FAD	C8A-N9A	-2.43	1.33	1.37
3	B	801	FAD	O3'-C3'	-2.42	1.37	1.43
3	B	801	FAD	C10-N10	2.39	1.42	1.37
3	A	802	FAD	C4'-C3'	-2.37	1.49	1.53
3	B	801	FAD	C4X-C10	2.37	1.51	1.44
3	B	801	FAD	O4'-C4'	-2.35	1.38	1.43
3	B	801	FAD	O4B-C1B	2.34	1.47	1.42
2	A	801	FMN	O3'-C3'	-2.32	1.37	1.43
2	A	801	FMN	C4-N3	-2.27	1.34	1.38
3	A	802	FAD	C10-N10	2.21	1.42	1.37
3	A	802	FAD	O4B-C1B	2.19	1.47	1.42
3	A	802	FAD	C4X-C10	2.18	1.50	1.44
3	B	801	FAD	O2'-C2'	-2.16	1.38	1.43
3	A	802	FAD	C2A-N3A	2.07	1.37	1.33

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	801	FAD	N3A-C2A-N1A	-5.41	120.39	128.58
3	B	801	FAD	C5A-C4A-N3A	-5.35	119.34	126.72
3	A	802	FAD	C5A-C4A-N3A	-5.35	119.36	126.72
3	A	802	FAD	N3A-C2A-N1A	-5.29	120.57	128.58
3	A	802	FAD	N3A-C4A-N9A	4.47	134.77	127.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	801	FAD	N3A-C4A-N9A	4.24	134.39	127.17
3	A	802	FAD	C5A-N7A-C8A	3.70	109.26	103.45
3	B	801	FAD	C2A-N3A-C4A	3.66	120.76	111.83
3	B	801	FAD	C5A-N7A-C8A	3.57	109.06	103.45
3	A	802	FAD	C2A-N3A-C4A	3.40	120.14	111.83
2	A	801	FMN	C4-N3-C2	-3.08	120.18	125.64
3	B	801	FAD	C4A-C5A-N7A	-2.91	107.26	110.58
2	A	801	FMN	C4A-C4-N3	2.85	120.51	113.25
3	B	801	FAD	C4-C4X-N5	2.83	122.11	118.21
3	A	802	FAD	C4A-C5A-N7A	-2.82	107.35	110.58
2	A	801	FMN	C4A-C10-N1	-2.73	117.90	124.59
3	A	802	FAD	C4-N3-C2	-2.73	120.80	125.64
3	B	801	FAD	C4-N3-C2	-2.72	120.82	125.64
2	A	801	FMN	O4-C4-C4A	-2.71	119.38	126.53
3	B	801	FAD	C4X-C4-N3	2.65	119.99	113.25
3	A	802	FAD	O4-C4-C4X	-2.62	119.61	126.53
3	A	802	FAD	C4X-C4-N3	2.56	119.76	113.25
3	A	802	FAD	N9A-C8A-N7A	-2.54	110.33	113.94
3	B	801	FAD	C4'-C3'-C2'	-2.50	109.41	113.57
2	A	801	FMN	C10-N1-C2	2.43	122.11	116.85
3	A	802	FAD	C3B-C2B-C1B	2.38	105.97	101.46
3	B	801	FAD	N9A-C8A-N7A	-2.31	110.66	113.94
3	A	802	FAD	C4'-C3'-C2'	-2.28	109.77	113.57
2	A	801	FMN	C10-C4A-N5	-2.28	120.15	124.81
3	B	801	FAD	C10-C4X-N5	-2.24	120.24	124.81
3	B	801	FAD	O4-C4-C4X	-2.20	120.72	126.53
3	B	801	FAD	C3B-C2B-C1B	2.08	105.40	101.46
2	A	801	FMN	C4-C4A-N5	2.06	121.06	118.21
3	A	802	FAD	O2A-PA-O3P	2.02	112.73	107.27
3	A	802	FAD	C10-C4X-N5	-2.02	120.69	124.81

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	802	FAD	C2B-C1B-N9A-C8A
3	B	801	FAD	C2B-C1B-N9A-C8A
2	A	801	FMN	C2'-C3'-C4'-O4'
3	B	801	FAD	O2'-C2'-C3'-C4'
2	A	801	FMN	C2'-C3'-C4'-C5'
2	A	801	FMN	O3'-C3'-C4'-O4'
3	A	802	FAD	C2B-C1B-N9A-C4A

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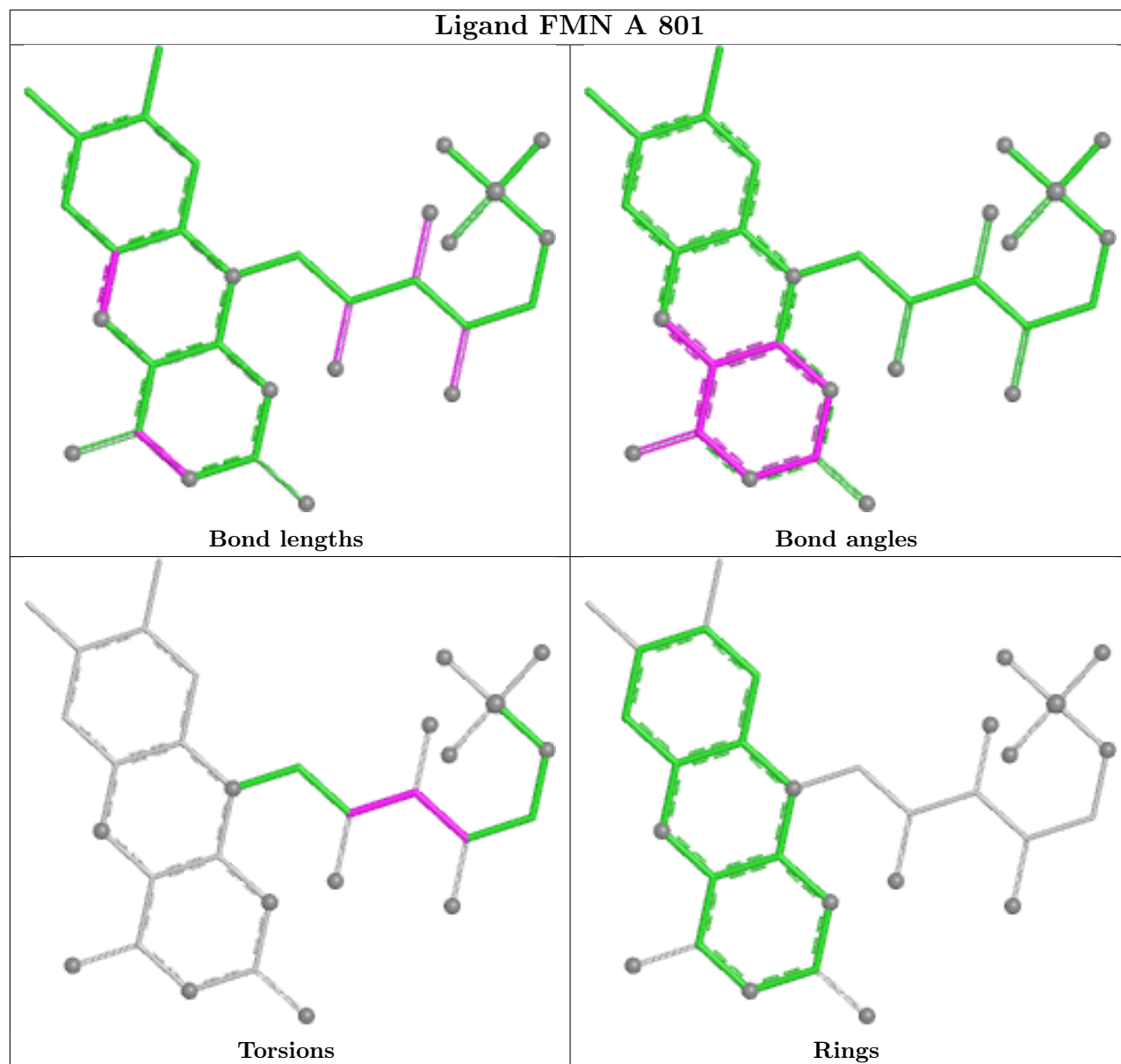
Mol	Chain	Res	Type	Atoms
2	A	801	FMN	O3'-C3'-C4'-C5'
3	B	801	FAD	C2B-C1B-N9A-C4A
3	B	801	FAD	O2'-C2'-C3'-O3'
3	A	802	FAD	O4B-C1B-N9A-C8A
3	B	801	FAD	O4B-C1B-N9A-C8A
3	A	802	FAD	O2'-C2'-C3'-C4'
2	A	801	FMN	O2'-C2'-C3'-C4'

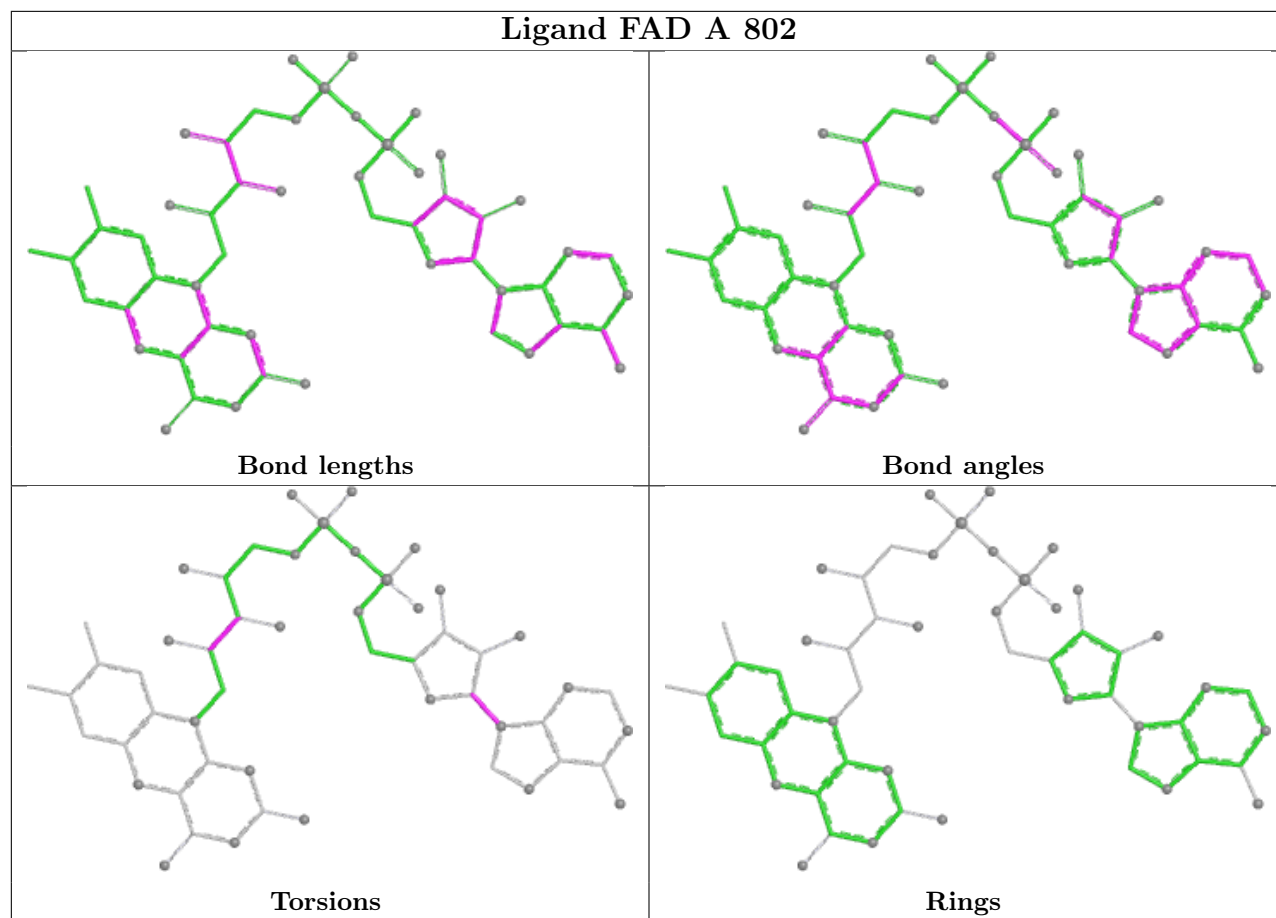
There are no ring outliers.

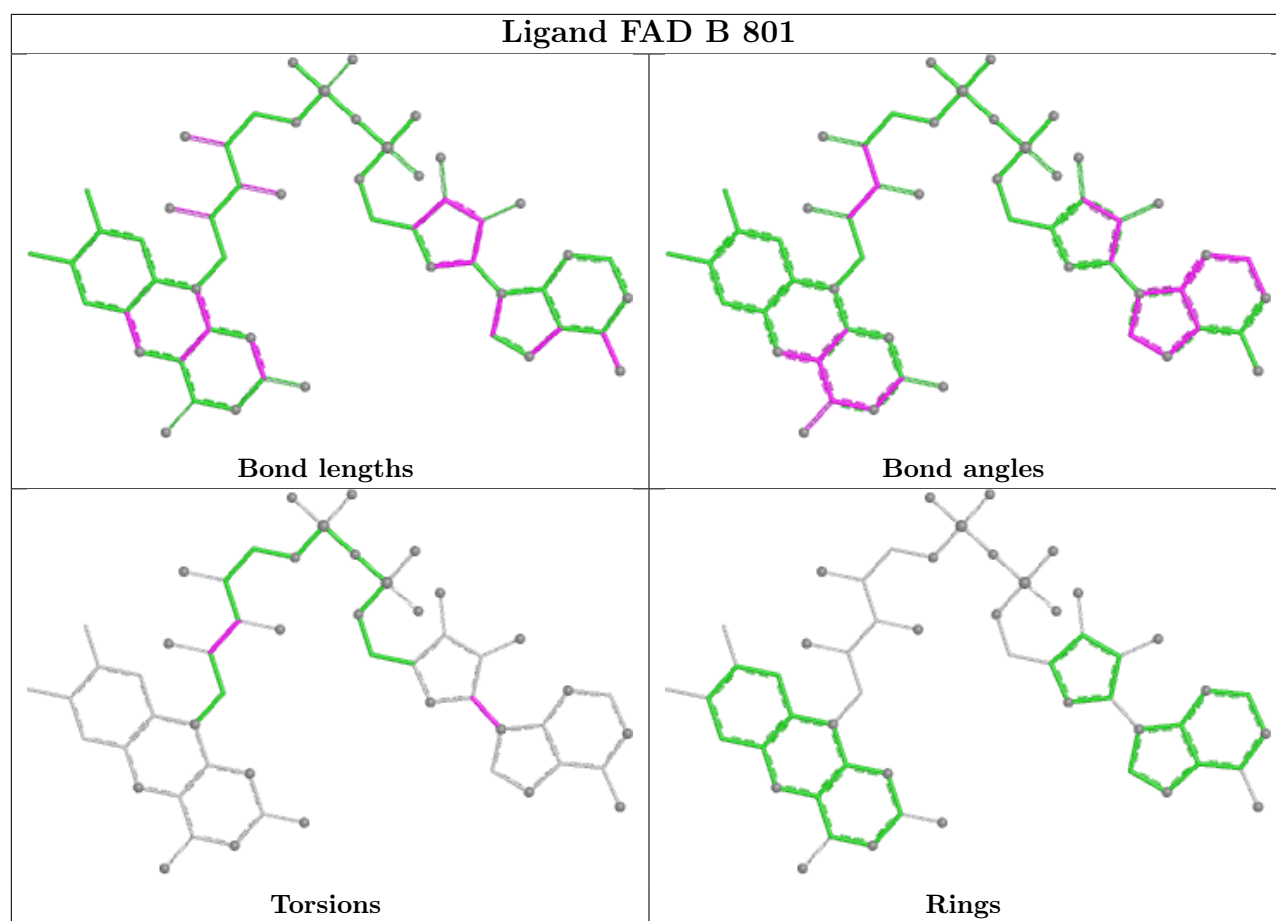
3 monomers are involved in 8 short contacts:

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

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







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	552/640 (86%)	0.10	6 (1%) 78 79	8, 22, 43, 51	7 (1%)
1	B	394/640 (61%)	-0.07	4 (1%) 79 80	8, 18, 32, 39	5 (1%)
All	All	946/1280 (73%)	0.03	10 (1%) 78 79	8, 20, 40, 51	12 (1%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	538	SER	3.6
1	A	129	ALA	3.5
1	B	537	LYS	3.1
1	A	397	PRO	2.8
1	B	538	SER	2.6
1	B	356	CYS	2.4
1	B	699	ASN	2.3
1	A	143	ASP	2.1
1	A	539	GLU	2.0
1	A	238	ASP	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](#)

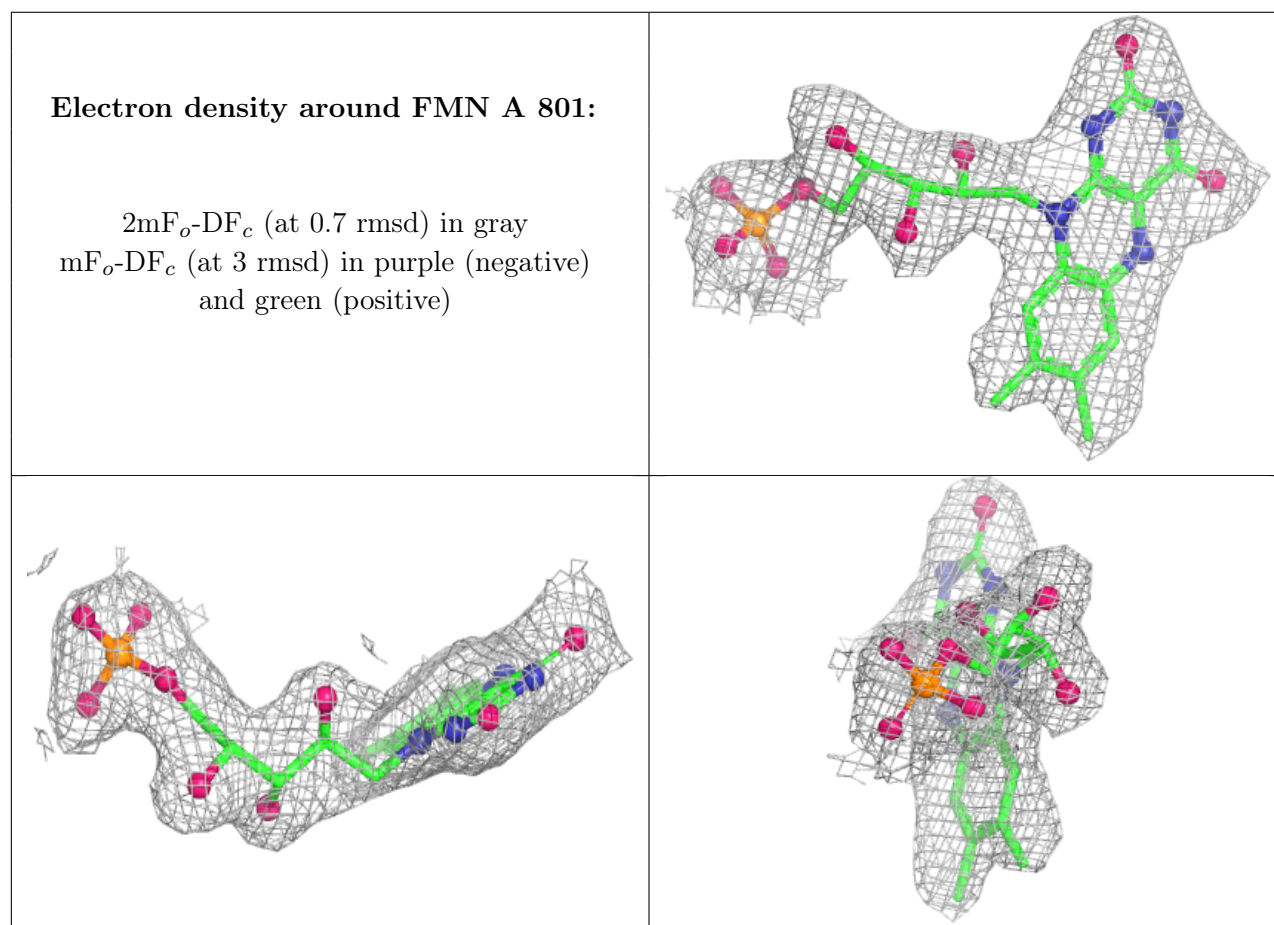
There are no oligosaccharides in this entry.

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, 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.

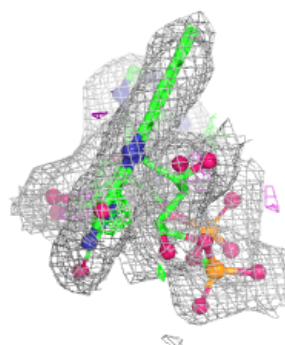
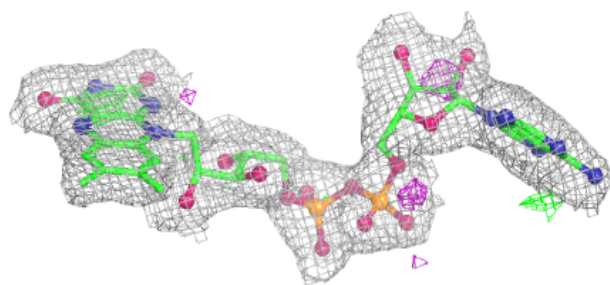
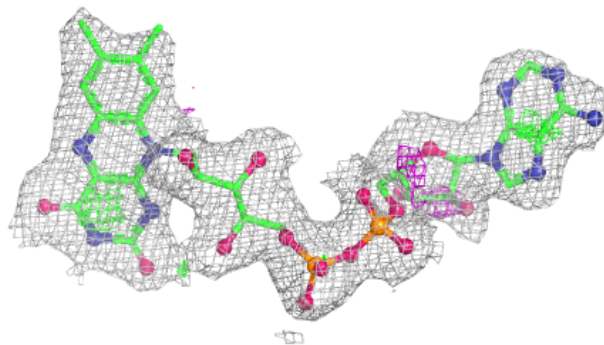
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	FMN	A	801	31/31	0.94	0.08	27,34,36,36	0
3	FAD	A	802	53/53	0.97	0.07	8,11,21,21	0
3	FAD	B	801	53/53	0.97	0.07	7,10,18,20	0

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.

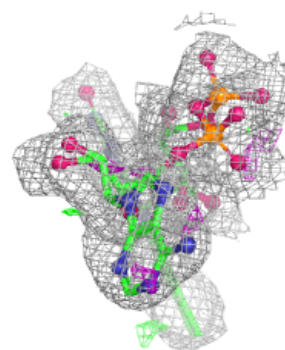
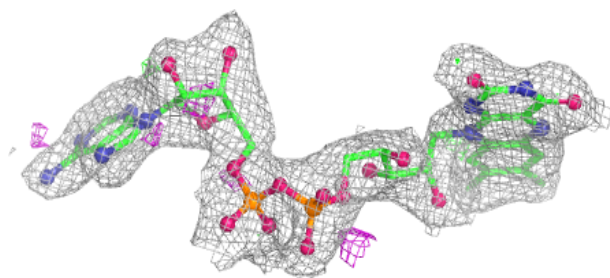
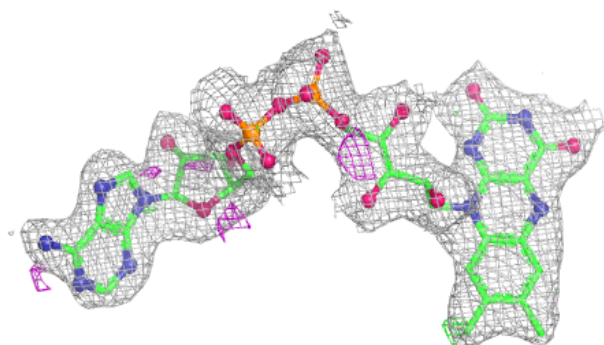


Electron density around FAD A 802:

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

**Electron density around FAD B 801:**

$2mF_o-DF_c$ (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.