Full wwPDB X-ray Structure Validation Report (i)

Oct 4, 2023 – 04:52 PM EDT

PDB ID	:	6P00
Title	:	RebH Variant 8F, Tryptamine 6-halogenase
Authors	:	Andorfer, M.C.; Sukumar, N.; Lewis, J.C.
Deposited on	:	2019-05-16
Resolution	:	2.25 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 (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

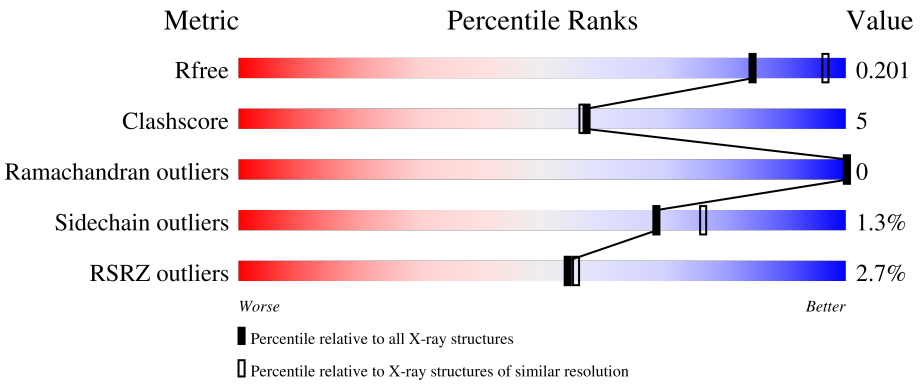
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	2391 (2.26-2.22)
Clashscore	141614	2539 (2.26-2.22)
Ramachandran outliers	138981	2489 (2.26-2.22)
Sidechain outliers	138945	2490 (2.26-2.22)
RSRZ outliers	127900	2353 (2.26-2.22)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	550	2% 85% 9% 6%
1	B	550	3% 84% 10% 6%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 9024 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 Flavin-dependent tryptophan halogenase RebH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	518	4161	2648	725	768	20	0	0	0
1	B	519	4168	2652	726	769	21	0	0	0

There are 62 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP Q8KHZ8
A	-18	GLY	-	expression tag	UNP Q8KHZ8
A	-17	SER	-	expression tag	UNP Q8KHZ8
A	-16	SER	-	expression tag	UNP Q8KHZ8
A	-15	HIS	-	expression tag	UNP Q8KHZ8
A	-14	HIS	-	expression tag	UNP Q8KHZ8
A	-13	HIS	-	expression tag	UNP Q8KHZ8
A	-12	HIS	-	expression tag	UNP Q8KHZ8
A	-11	HIS	-	expression tag	UNP Q8KHZ8
A	-10	HIS	-	expression tag	UNP Q8KHZ8
A	-9	SER	-	expression tag	UNP Q8KHZ8
A	-8	SER	-	expression tag	UNP Q8KHZ8
A	-7	GLY	-	expression tag	UNP Q8KHZ8
A	-6	LEU	-	expression tag	UNP Q8KHZ8
A	-5	VAL	-	expression tag	UNP Q8KHZ8
A	-4	PRO	-	expression tag	UNP Q8KHZ8
A	-3	ARG	-	expression tag	UNP Q8KHZ8
A	-2	GLY	-	expression tag	UNP Q8KHZ8
A	-1	SER	-	expression tag	UNP Q8KHZ8
A	0	HIS	-	expression tag	UNP Q8KHZ8
A	52	MET	ILE	engineered mutation	UNP Q8KHZ8
A	110	PRO	SER	engineered mutation	UNP Q8KHZ8
A	130	LEU	SER	engineered mutation	UNP Q8KHZ8
A	166	SER	ASN	engineered mutation	UNP Q8KHZ8
A	380	PHE	LEU	engineered mutation	UNP Q8KHZ8

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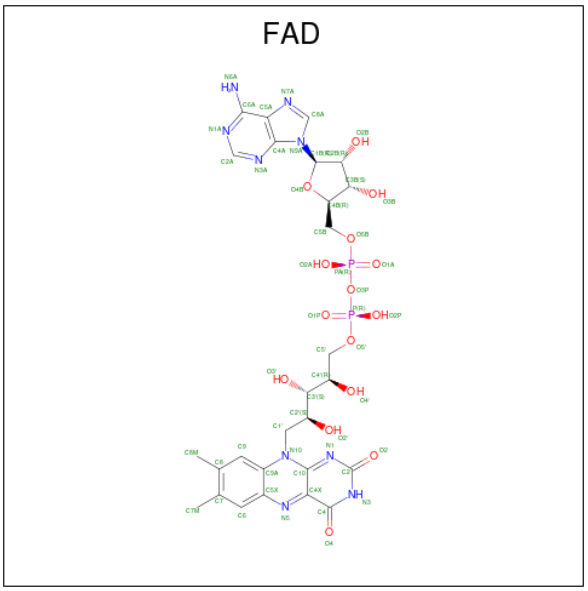
Chain	Residue	Modelled	Actual	Comment	Reference
A	448	PRO	SER	engineered mutation	UNP Q8KHZ8
A	455	TRP	TYR	engineered mutation	UNP Q8KHZ8
A	465	LEU	PHE	engineered mutation	UNP Q8KHZ8
A	470	SER	ASN	engineered mutation	UNP Q8KHZ8
A	494	ARG	GLN	engineered mutation	UNP Q8KHZ8
A	509	GLN	ARG	engineered mutation	UNP Q8KHZ8
B	-19	MET	-	initiating methionine	UNP Q8KHZ8
B	-18	GLY	-	expression tag	UNP Q8KHZ8
B	-17	SER	-	expression tag	UNP Q8KHZ8
B	-16	SER	-	expression tag	UNP Q8KHZ8
B	-15	HIS	-	expression tag	UNP Q8KHZ8
B	-14	HIS	-	expression tag	UNP Q8KHZ8
B	-13	HIS	-	expression tag	UNP Q8KHZ8
B	-12	HIS	-	expression tag	UNP Q8KHZ8
B	-11	HIS	-	expression tag	UNP Q8KHZ8
B	-10	HIS	-	expression tag	UNP Q8KHZ8
B	-9	SER	-	expression tag	UNP Q8KHZ8
B	-8	SER	-	expression tag	UNP Q8KHZ8
B	-7	GLY	-	expression tag	UNP Q8KHZ8
B	-6	LEU	-	expression tag	UNP Q8KHZ8
B	-5	VAL	-	expression tag	UNP Q8KHZ8
B	-4	PRO	-	expression tag	UNP Q8KHZ8
B	-3	ARG	-	expression tag	UNP Q8KHZ8
B	-2	GLY	-	expression tag	UNP Q8KHZ8
B	-1	SER	-	expression tag	UNP Q8KHZ8
B	0	HIS	-	expression tag	UNP Q8KHZ8
B	52	MET	ILE	engineered mutation	UNP Q8KHZ8
B	110	PRO	SER	engineered mutation	UNP Q8KHZ8
B	130	LEU	SER	engineered mutation	UNP Q8KHZ8
B	166	SER	ASN	engineered mutation	UNP Q8KHZ8
B	380	PHE	LEU	engineered mutation	UNP Q8KHZ8
B	448	PRO	SER	engineered mutation	UNP Q8KHZ8
B	455	TRP	TYR	engineered mutation	UNP Q8KHZ8
B	465	LEU	PHE	engineered mutation	UNP Q8KHZ8
B	470	SER	ASN	engineered mutation	UNP Q8KHZ8
B	494	ARG	GLN	engineered mutation	UNP Q8KHZ8
B	509	GLN	ARG	engineered mutation	UNP Q8KHZ8

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is water.

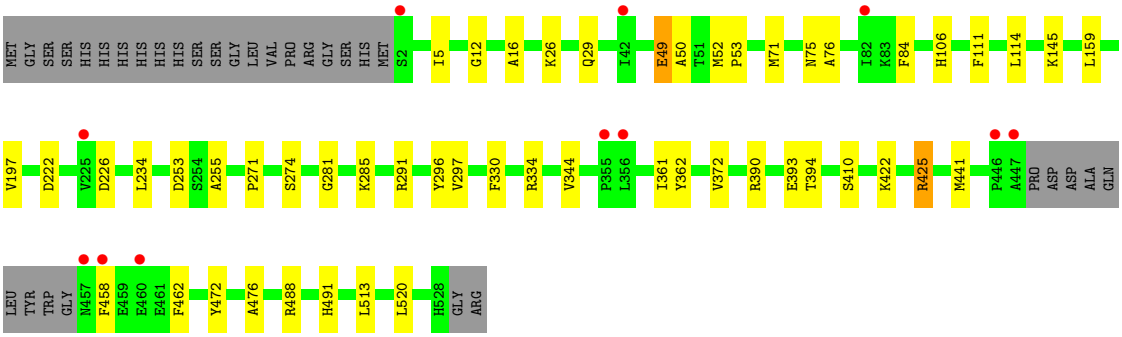
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	302	302	302	0	0
			Total	O		
3	B	287	287	287	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

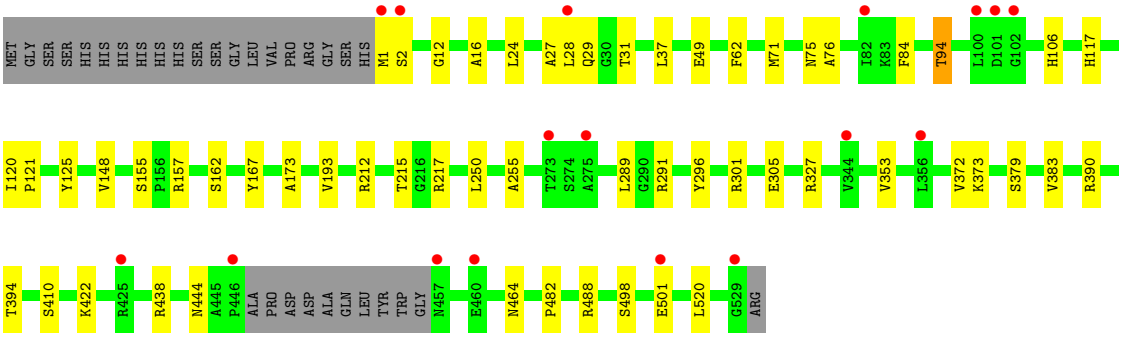
• Molecule 1: Flavin-dependent tryptophan halogenase RebH

Chain A:



• Molecule 1: Flavin-dependent tryptophan halogenase RebH

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 62	Depositor
Cell constants a, b, c, α, β, γ	115.97Å 115.97Å 229.51Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	31.17 – 2.25 31.17 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.1 (31.17-2.25) 90.5 (31.17-2.25)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.61 (at 2.24Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.170 , 0.200 0.177 , 0.201	Depositor DCC
R_{free} test set	4082 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å^2)	34.9	Xtriage
Anisotropy	0.339	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 38.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.049 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	9024	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.91% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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	0.42	0/4276	0.56	0/5801
1	B	0.45	0/4283	0.58	1/5809 (0.0%)
All	All	0.44	0/8559	0.57	1/11610 (0.0%)

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	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	94	THR	CB-CA-C	-5.54	96.65	111.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	379	SER	Mainchain

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	4161	0	3983	45	0
1	B	4168	0	3993	40	0
2	A	53	0	30	12	0
2	B	53	0	29	7	0
3	A	302	0	0	1	0
3	B	287	0	0	6	0
All	All	9024	0	8035	81	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 (81) 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:49:GLU:OE1	2:B:601:FAD:C2	2.01	1.07
1:B:49:GLU:OE1	2:B:601:FAD:N1	1.93	1.02
1:A:12:GLY:HA2	2:A:601:FAD:H1B	1.53	0.90
1:A:49:GLU:OE2	2:A:601:FAD:C2'	2.23	0.87
1:B:12:GLY:HA2	2:B:601:FAD:H1B	1.57	0.85
1:A:458:PHE:CE2	1:A:462:PHE:HB2	2.11	0.85
1:B:24:LEU:O	1:B:28:LEU:HD13	1.77	0.84
1:A:49:GLU:OE2	2:A:601:FAD:H2'	1.77	0.83
1:B:49:GLU:OE1	2:B:601:FAD:C10	2.35	0.74
1:B:49:GLU:HG3	1:B:173:ALA:HB2	1.70	0.71
1:B:410:SER:O	1:B:422:LYS:NZ	2.23	0.71
1:A:50:ALA:O	2:A:601:FAD:N3	2.21	0.71
1:A:410:SER:O	1:A:422:LYS:NZ	2.24	0.70
1:B:49:GLU:OE2	2:B:601:FAD:H2'	1.94	0.67
1:A:197:VAL:H	2:A:601:FAD:H62A	1.41	0.67
1:A:491:HIS:HA	1:B:31:THR:HG21	1.80	0.64
1:B:27:ALA:HB2	1:B:62:PHE:HZ	1.64	0.62
1:B:212:ARG:NH1	3:B:702:HOH:O	2.28	0.62
1:A:49:GLU:HG3	2:A:601:FAD:C2	2.30	0.61
1:B:75:ASN:HB3	1:B:289:LEU:HD12	1.83	0.60
1:B:390:ARG:O	1:B:394:THR:HG23	2.00	0.60
1:A:49:GLU:OE2	2:A:601:FAD:C1'	2.49	0.59
1:B:27:ALA:HB2	1:B:62:PHE:CZ	2.39	0.57
1:A:49:GLU:HG2	1:A:361:ILE:CD1	2.35	0.57
1:A:253:ASP:HA	1:A:330:PHE:CE1	2.39	0.56
1:A:49:GLU:HG2	1:A:361:ILE:HD12	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:425:ARG:HH11	1:A:425:ARG:HG3	1.72	0.54
1:B:327:ARG:NE	3:B:706:HOH:O	2.40	0.53
1:A:26:LYS:O	1:A:29:GLN:HG2	2.09	0.53
1:B:28:LEU:HD12	1:B:28:LEU:N	2.26	0.51
1:A:372:VAL:O	1:B:488:ARG:HD2	2.12	0.50
1:A:71:MET:HG2	1:A:76:ALA:HB3	1.94	0.50
1:B:444:ASN:HB3	1:B:464:ASN:OD1	2.12	0.49
1:A:49:GLU:CD	2:A:601:FAD:C10	2.80	0.49
1:A:5:ILE:HG23	1:A:222:ASP:HB2	1.95	0.49
1:A:49:GLU:OE2	2:A:601:FAD:N1	2.47	0.48
1:B:37:LEU:HD12	1:B:193:VAL:HB	1.94	0.48
1:A:488:ARG:HD2	1:B:372:VAL:O	2.13	0.48
1:A:197:VAL:N	2:A:601:FAD:H62A	2.09	0.48
1:B:305:GLU:OE2	1:B:327:ARG:NH1	2.45	0.47
1:A:390:ARG:O	1:A:394:THR:HG23	2.15	0.47
1:B:250:LEU:HD13	1:B:353:VAL:HG23	1.97	0.46
1:B:162:SER:HA	3:B:728:HOH:O	2.15	0.46
1:B:438:ARG:HA	1:B:482:PRO:HA	1.96	0.46
1:B:71:MET:HG3	1:B:76:ALA:HB3	1.97	0.46
1:A:472:TYR:O	1:A:476:ALA:HB3	2.16	0.45
1:A:52:MET:HE1	1:A:362:TYR:CG	2.51	0.45
1:A:281:GLY:HA3	1:A:297:VAL:O	2.17	0.45
1:A:75:ASN:HB3	3:A:875:HOH:O	2.17	0.45
1:B:49:GLU:OE1	2:B:601:FAD:N3	2.46	0.45
1:B:84:PHE:O	1:B:106:HIS:HA	2.17	0.44
1:B:120:ILE:HG23	1:B:121:PRO:HD2	2.00	0.44
1:B:498:SER:O	1:B:501:GLU:HG2	2.17	0.44
1:A:344:VAL:HG13	1:A:344:VAL:O	2.18	0.44
1:B:488:ARG:HD3	3:B:777:HOH:O	2.17	0.44
1:B:16:ALA:HB2	2:B:601:FAD:H5'2	1.98	0.44
1:A:255:ALA:HA	1:A:296:TYR:O	2.18	0.44
1:A:271:PRO:HB3	1:A:520:LEU:HD22	2.00	0.44
1:A:16:ALA:HB2	2:A:601:FAD:H5'2	2.00	0.43
1:A:226:ASP:HB2	1:A:234:LEU:HB2	2.00	0.43
1:A:49:GLU:OE2	2:A:601:FAD:C10	2.66	0.43
1:A:53:PRO:HD3	1:A:111:PHE:CE1	2.54	0.43
1:A:84:PHE:O	1:A:106:HIS:HA	2.18	0.43
1:A:425:ARG:HH11	1:A:425:ARG:CG	2.32	0.43
1:B:155:SER:HB2	1:B:520:LEU:HA	2.01	0.42
1:B:28:LEU:N	1:B:28:LEU:CD1	2.82	0.42
1:A:52:MET:HE2	1:A:362:TYR:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:117:HIS:HB3	1:B:125:TYR:CE2	2.54	0.42
1:A:274:SER:HB2	1:A:285:LYS:HB3	2.00	0.42
1:B:301:ARG:NH1	3:B:704:HOH:O	2.36	0.42
1:B:215:THR:OG1	1:B:217:ARG:HD2	2.19	0.42
1:B:373:LYS:NZ	3:B:705:HOH:O	2.39	0.41
1:A:334:ARG:NH2	1:A:393:GLU:HG3	2.35	0.41
1:A:145:LYS:HD2	1:A:513:LEU:HD12	2.03	0.41
1:A:52:MET:CE	1:A:362:TYR:CD1	3.04	0.41
1:A:145:LYS:HD2	1:A:513:LEU:CD1	2.51	0.41
1:A:491:HIS:CA	1:B:31:THR:HG21	2.49	0.41
1:A:159:LEU:HD23	1:A:159:LEU:HA	1.88	0.40
1:B:255:ALA:HA	1:B:296:TYR:O	2.21	0.40
1:A:441:MET:HE2	1:A:441:MET:HB3	1.93	0.40
1:B:157:ARG:NH1	1:B:167:TYR:OH	2.53	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	514/550 (94%)	501 (98%)	13 (2%)	0	100	100
1	B	515/550 (94%)	503 (98%)	12 (2%)	0	100	100
All	All	1029/1100 (94%)	1004 (98%)	25 (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	437/463 (94%)	433 (99%)	4 (1%)	78	84
1	B	438/463 (95%)	431 (98%)	7 (2%)	62	70
All	All	875/926 (94%)	864 (99%)	11 (1%)	69	76

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	GLU
1	A	114	LEU
1	A	291	ARG
1	A	425	ARG
1	B	1	MET
1	B	2	SER
1	B	29	GLN
1	B	94	THR
1	B	148	VAL
1	B	291	ARG
1	B	383	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

2 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	A	601	-	53,58,58	4.03	20 (37%)	68,89,89	2.46	19 (27%)
2	FAD	B	601	-	53,58,58	3.98	23 (43%)	68,89,89	2.38	13 (19%)

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	A	601	-	-	12/30/50/50	0/6/6/6
2	FAD	B	601	-	-	15/30/50/50	0/6/6/6

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	FAD	C2B-C1B	-16.34	1.29	1.53
2	A	601	FAD	C2B-C1B	-15.90	1.29	1.53
2	A	601	FAD	O4B-C1B	15.13	1.62	1.41
2	B	601	FAD	O4B-C1B	13.85	1.60	1.41
2	B	601	FAD	C4X-N5	7.71	1.45	1.30
2	A	601	FAD	C4X-N5	7.60	1.45	1.30
2	B	601	FAD	O4B-C4B	-7.55	1.28	1.45
2	A	601	FAD	O4B-C4B	-7.23	1.28	1.45
2	B	601	FAD	C10-N1	5.93	1.45	1.33
2	A	601	FAD	C9A-N10	5.76	1.51	1.41
2	A	601	FAD	C10-N1	5.75	1.45	1.33
2	B	601	FAD	C5X-N5	5.38	1.49	1.39
2	A	601	FAD	O2-C2	-4.72	1.15	1.24
2	A	601	FAD	C5X-N5	4.22	1.47	1.39
2	A	601	FAD	C2-N3	4.21	1.48	1.39
2	B	601	FAD	C9A-N10	4.20	1.48	1.41
2	B	601	FAD	C2-N1	4.19	1.46	1.36
2	B	601	FAD	C2-N3	4.09	1.48	1.39
2	A	601	FAD	O4-C4	-4.01	1.16	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	FAD	O4-C4	-3.99	1.16	1.23
2	B	601	FAD	O2-C2	-3.97	1.16	1.24
2	A	601	FAD	C2-N1	3.97	1.46	1.36
2	A	601	FAD	C4-N3	3.61	1.45	1.38
2	B	601	FAD	O3B-C3B	-3.57	1.34	1.43
2	B	601	FAD	C5A-C4A	-3.47	1.31	1.40
2	A	601	FAD	C5'-C4'	3.41	1.56	1.51
2	A	601	FAD	C5A-C4A	-3.34	1.32	1.40
2	B	601	FAD	C10-N10	3.25	1.44	1.37
2	A	601	FAD	O3B-C3B	-3.02	1.35	1.43
2	A	601	FAD	C10-N10	2.91	1.43	1.37
2	A	601	FAD	O2B-C2B	2.71	1.49	1.43
2	B	601	FAD	O3'-C3'	-2.64	1.36	1.43
2	A	601	FAD	C6A-N6A	2.60	1.43	1.34
2	B	601	FAD	O4'-C4'	-2.58	1.37	1.43
2	B	601	FAD	O2B-C2B	2.51	1.48	1.43
2	B	601	FAD	C4-N3	2.46	1.43	1.38
2	B	601	FAD	C6A-N6A	2.43	1.42	1.34
2	B	601	FAD	C4X-C4	2.40	1.53	1.44
2	B	601	FAD	P-O2P	-2.38	1.44	1.55
2	A	601	FAD	O4'-C4'	-2.33	1.38	1.43
2	B	601	FAD	O2'-C2'	-2.33	1.38	1.43
2	A	601	FAD	O2'-C2'	-2.19	1.38	1.43
2	B	601	FAD	C5'-C4'	2.18	1.54	1.51

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	601	FAD	C5A-C6A-N6A	10.52	136.34	120.35
2	B	601	FAD	C5A-C6A-N6A	10.08	135.68	120.35
2	A	601	FAD	N6A-C6A-N1A	-7.59	102.83	118.57
2	B	601	FAD	N6A-C6A-N1A	-7.23	103.56	118.57
2	A	601	FAD	C7M-C7-C6	-6.56	107.35	119.49
2	B	601	FAD	C7M-C7-C6	-6.38	107.69	119.49
2	B	601	FAD	N3A-C2A-N1A	-5.76	119.67	128.68
2	A	601	FAD	C7M-C7-C8	5.37	131.75	120.74
2	A	601	FAD	N3A-C2A-N1A	-5.22	120.53	128.68
2	B	601	FAD	C7M-C7-C8	5.03	131.05	120.74
2	B	601	FAD	C4-C4X-N5	3.47	123.17	118.23
2	A	601	FAD	C1'-C2'-C3'	-3.32	100.49	109.79
2	B	601	FAD	C1'-C2'-C3'	-3.30	100.57	109.79
2	B	601	FAD	C5'-C4'-C3'	-3.29	105.84	112.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	601	FAD	C1'-N10-C9A	3.29	125.99	120.51
2	A	601	FAD	C4-N3-C2	-3.26	119.62	125.64
2	A	601	FAD	C4X-C10-N10	3.25	121.24	116.48
2	B	601	FAD	C4X-C10-N10	3.22	121.19	116.48
2	B	601	FAD	C10-C4X-N5	-3.07	118.35	124.86
2	A	601	FAD	C3B-C2B-C1B	2.90	105.34	100.98
2	A	601	FAD	C4X-C4-N3	2.81	120.32	113.19
2	A	601	FAD	C10-C4X-N5	-2.74	119.04	124.86
2	B	601	FAD	O4B-C1B-C2B	-2.67	103.03	106.93
2	A	601	FAD	C4-C4X-N5	2.65	122.00	118.23
2	A	601	FAD	C10-N1-C2	2.46	121.83	116.90
2	A	601	FAD	O4-C4-C4X	-2.42	120.17	126.60
2	B	601	FAD	C4X-C4-N3	2.34	119.14	113.19
2	B	601	FAD	C4-N3-C2	-2.22	121.53	125.64
2	A	601	FAD	C4A-C5A-N7A	-2.20	107.10	109.40
2	A	601	FAD	O4B-C4B-C3B	-2.20	100.76	105.11
2	A	601	FAD	O2'-C2'-C3'	2.20	114.44	109.10
2	A	601	FAD	P-O3P-PA	-2.01	125.92	132.83

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	FAD	N10-C1'-C2'-O2'
2	A	601	FAD	N10-C1'-C2'-C3'
2	A	601	FAD	C3'-C4'-C5'-O5'
2	A	601	FAD	O4'-C4'-C5'-O5'
2	A	601	FAD	C5'-O5'-P-O1P
2	A	601	FAD	C5'-O5'-P-O2P
2	B	601	FAD	C5B-O5B-PA-O1A
2	B	601	FAD	P-O3P-PA-O5B
2	B	601	FAD	C3'-C4'-C5'-O5'
2	B	601	FAD	O4'-C4'-C5'-O5'
2	B	601	FAD	C5'-O5'-P-O1P
2	B	601	FAD	C5'-O5'-P-O2P
2	A	601	FAD	O4B-C4B-C5B-O5B
2	A	601	FAD	C3B-C4B-C5B-O5B
2	B	601	FAD	O4B-C4B-C5B-O5B
2	B	601	FAD	C3B-C4B-C5B-O5B
2	B	601	FAD	O3'-C3'-C4'-C5'
2	B	601	FAD	O3'-C3'-C4'-O4'
2	B	601	FAD	C4'-C5'-O5'-P

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Mol	Chain	Res	Type	Atoms
2	A	601	FAD	C4'-C5'-O5'-P
2	B	601	FAD	C5B-O5B-PA-O3P
2	B	601	FAD	C5B-O5B-PA-O2A
2	B	601	FAD	C2'-C3'-C4'-O4'
2	A	601	FAD	C2'-C3'-C4'-O4'
2	A	601	FAD	C5'-O5'-P-O3P
2	B	601	FAD	C5'-O5'-P-O3P
2	A	601	FAD	C5B-O5B-PA-O1A

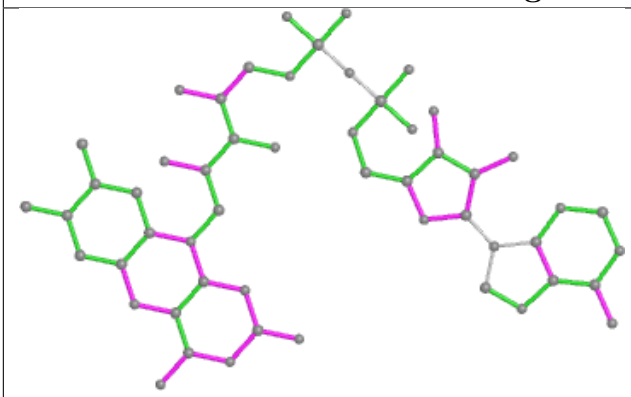
There are no ring outliers.

2 monomers are involved in 19 short contacts:

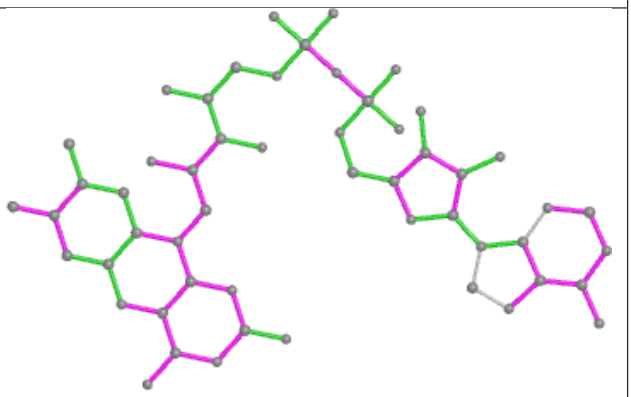
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	FAD	12	0
2	B	601	FAD	7	0

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

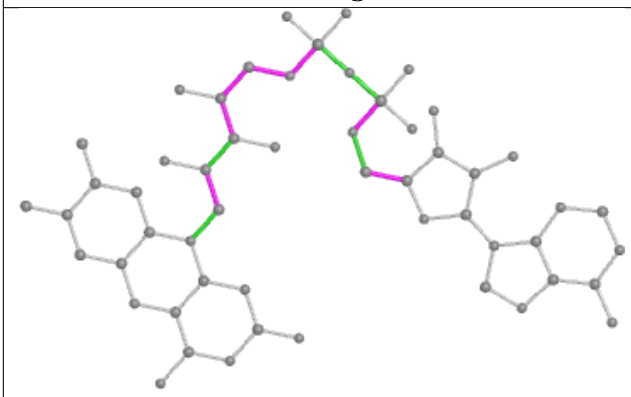
Ligand FAD A 601



Bond lengths



Bond angles



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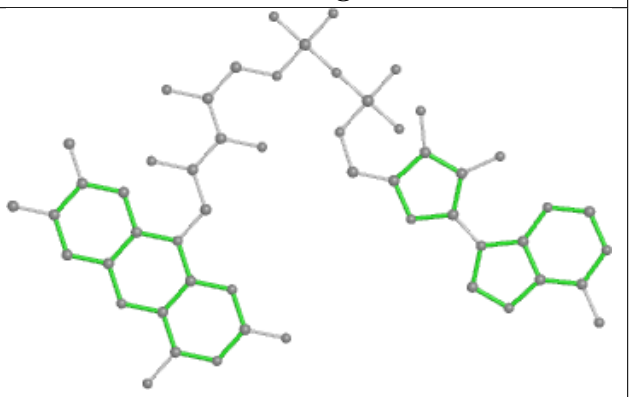
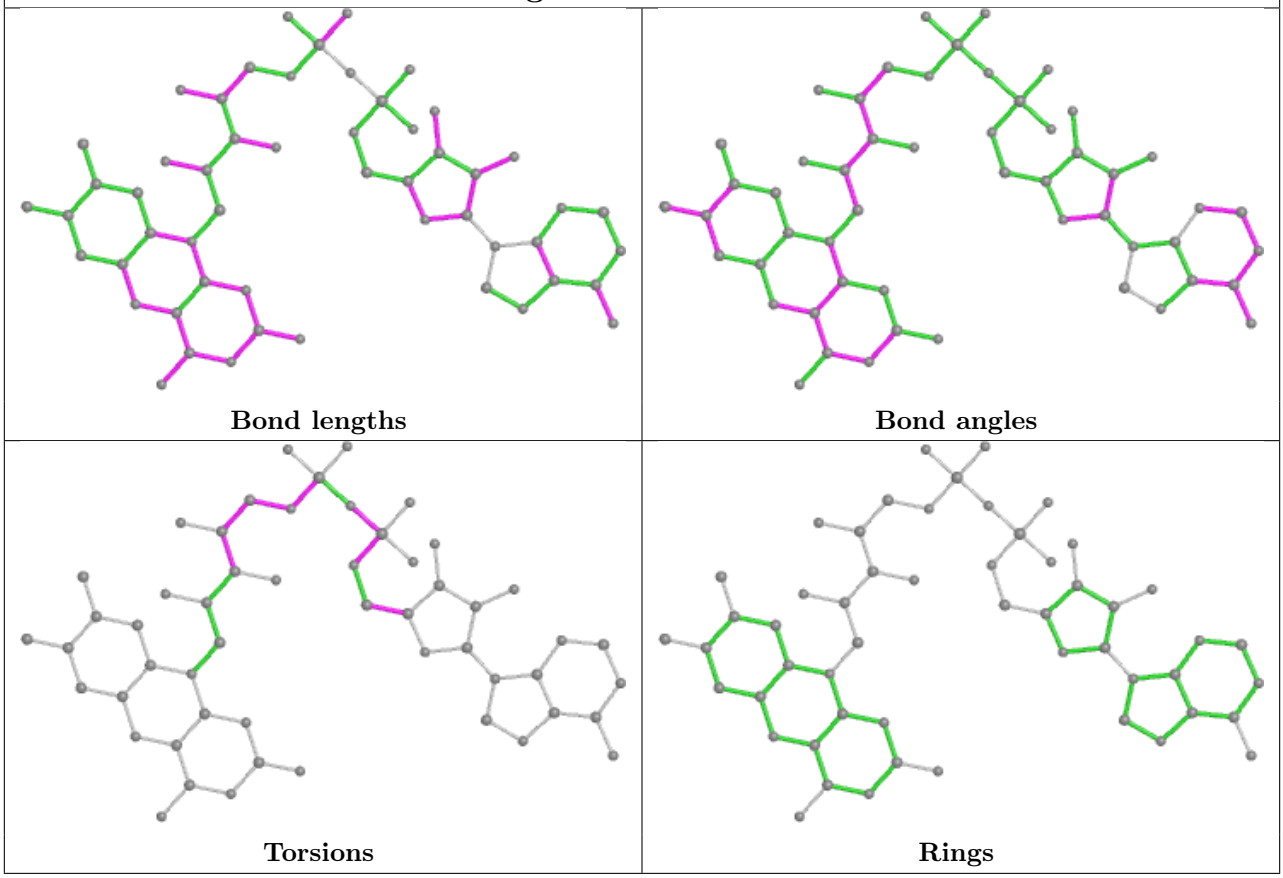


Rings

Ligand FAD B 601



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	518/550 (94%)	-0.23	11 (2%)	63	65	27, 40, 68, 109	0
1	B	519/550 (94%)	-0.16	17 (3%)	46	46	26, 42, 71, 113	0
All	All	1037/1100 (94%)	-0.20	28 (2%)	54	55	26, 41, 71, 113	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	529	GLY	4.1
1	A	460	GLU	3.5
1	B	1	MET	3.4
1	A	356	LEU	3.4
1	A	458	PHE	3.3
1	B	356	LEU	3.1
1	B	446	PRO	3.1
1	A	457	ASN	3.0
1	A	446	PRO	2.9
1	B	2	SER	2.7
1	B	100	LEU	2.7
1	B	501	GLU	2.7
1	A	82	ILE	2.5
1	B	102	GLY	2.5
1	A	447	ALA	2.4
1	B	460	GLU	2.4
1	B	82	ILE	2.4
1	B	457	ASN	2.4
1	B	275	ALA	2.3
1	A	225	VAL	2.2
1	B	425	ARG	2.2
1	B	344	VAL	2.2
1	A	2	SER	2.2
1	B	101	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	28	LEU	2.1
1	A	42	ILE	2.1
1	A	355	PRO	2.1
1	B	273	THR	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 monosaccharides in this entry.

6.4 Ligands (i)

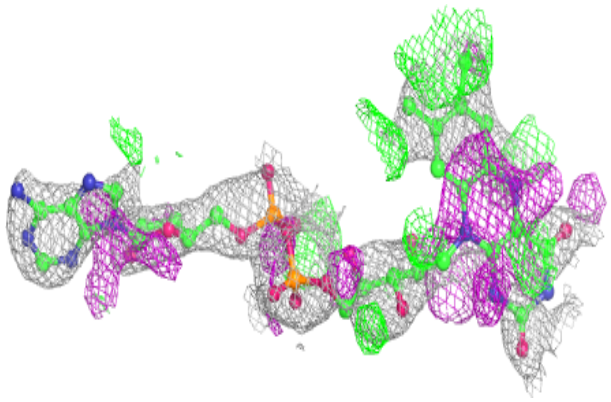
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

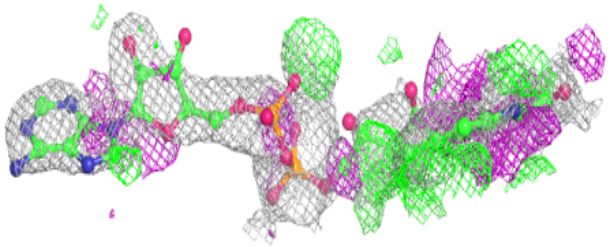
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	FAD	A	601	53/53	0.74	0.29	66,94,112,129	0
2	FAD	B	601	53/53	0.78	0.25	55,91,101,105	0

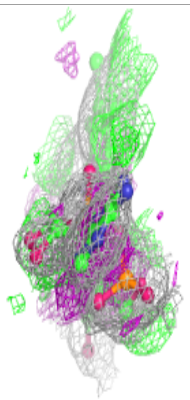
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD A 601:

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

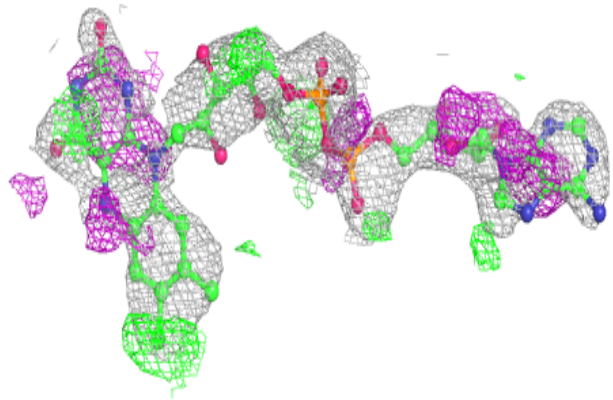


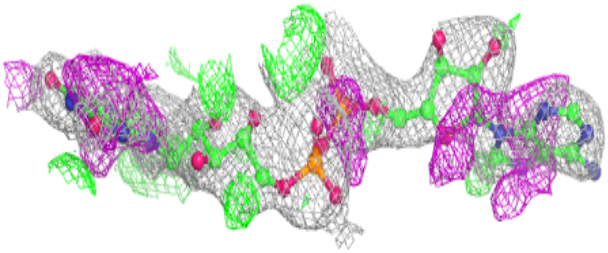


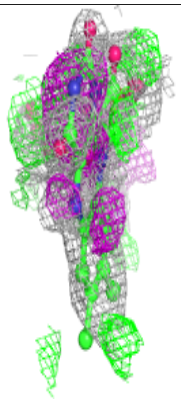


Electron density around FAD B 601:

$2mF_o$-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.