



# Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 09:47 pm BST

PDB ID : 3MKH  
Title : Podospora anserina Nitroalkane Oxidase  
Authors : Tormos, J.R.; Taylor, A.B.; Daubner, S.C.; Hart, P.J.; Fitzpatrick, P.F.  
Deposited on : 2010-04-14  
Resolution : 2.00 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

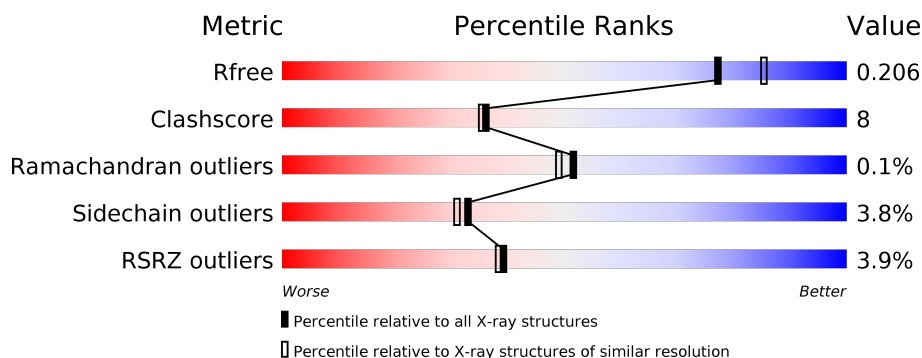
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	438	<div> <div>2%</div> <div> <div></div> <div>87%</div> <div>9%</div> <div>..</div> </div> </div>
1	B	438	<div> <div>%</div> <div> <div></div> <div>87%</div> <div>10%</div> <div>..</div> </div> </div>
1	C	438	<div> <div>5%</div> <div> <div></div> <div>79%</div> <div>17%</div> <div>..</div> </div> </div>
1	D	438	<div> <div>7%</div> <div> <div></div> <div>75%</div> <div>20%</div> <div>..</div> </div> </div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 14315 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 NITROALKANE OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	426	Total	C	N	O	S	0	0	0
			3165	1995	558	597	15			
1	B	426	Total	C	N	O	S	0	0	0
			3165	1995	558	597	15			
1	C	426	Total	C	N	O	S	0	0	0
			3165	1995	558	597	15			
1	D	426	Total	C	N	O	S	0	0	0
			3165	1995	558	597	15			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	ALA	PRO	SEE REMARK 999	UNP B2AM55
A	431	GLY	-	EXPRESSION TAG	UNP B2AM55
A	432	SER	-	EXPRESSION TAG	UNP B2AM55
A	433	HIS	-	EXPRESSION TAG	UNP B2AM55
A	434	HIS	-	EXPRESSION TAG	UNP B2AM55
A	435	HIS	-	EXPRESSION TAG	UNP B2AM55
A	436	HIS	-	EXPRESSION TAG	UNP B2AM55
A	437	HIS	-	EXPRESSION TAG	UNP B2AM55
A	438	HIS	-	EXPRESSION TAG	UNP B2AM55
B	2	ALA	PRO	SEE REMARK 999	UNP B2AM55
B	431	GLY	-	EXPRESSION TAG	UNP B2AM55
B	432	SER	-	EXPRESSION TAG	UNP B2AM55
B	433	HIS	-	EXPRESSION TAG	UNP B2AM55
B	434	HIS	-	EXPRESSION TAG	UNP B2AM55
B	435	HIS	-	EXPRESSION TAG	UNP B2AM55
B	436	HIS	-	EXPRESSION TAG	UNP B2AM55
B	437	HIS	-	EXPRESSION TAG	UNP B2AM55
B	438	HIS	-	EXPRESSION TAG	UNP B2AM55
C	2	ALA	PRO	SEE REMARK 999	UNP B2AM55
C	431	GLY	-	EXPRESSION TAG	UNP B2AM55
C	432	SER	-	EXPRESSION TAG	UNP B2AM55

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Chain	Residue	Modelled	Actual	Comment	Reference
C	433	HIS	-	EXPRESSION TAG	UNP B2AM55
C	434	HIS	-	EXPRESSION TAG	UNP B2AM55
C	435	HIS	-	EXPRESSION TAG	UNP B2AM55
C	436	HIS	-	EXPRESSION TAG	UNP B2AM55
C	437	HIS	-	EXPRESSION TAG	UNP B2AM55
C	438	HIS	-	EXPRESSION TAG	UNP B2AM55
D	2	ALA	PRO	SEE REMARK 999	UNP B2AM55
D	431	GLY	-	EXPRESSION TAG	UNP B2AM55
D	432	SER	-	EXPRESSION TAG	UNP B2AM55
D	433	HIS	-	EXPRESSION TAG	UNP B2AM55
D	434	HIS	-	EXPRESSION TAG	UNP B2AM55
D	435	HIS	-	EXPRESSION TAG	UNP B2AM55
D	436	HIS	-	EXPRESSION TAG	UNP B2AM55
D	437	HIS	-	EXPRESSION TAG	UNP B2AM55
D	438	HIS	-	EXPRESSION TAG	UNP B2AM55

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:  $C_{27}H_{33}N_9O_{15}P_2$ ).



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

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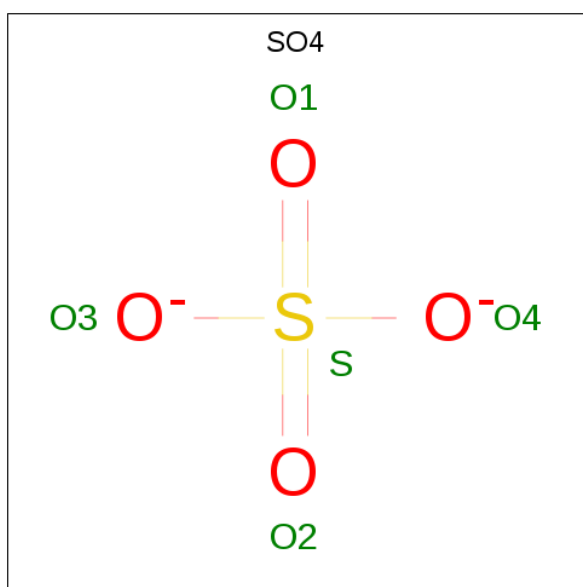
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	479	Total	O	0	0
			479	479		
5	B	448	Total	O	0	0
			448	448		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	265	Total 265	O 265	0	0
5	D	239	Total 239	O 239	0	0



- Molecule 1: NITROALKANE OXIDASE







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	137.50Å 137.50Å 131.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.88 – 2.00 44.88 – 2.00	Depositor EDS
% Data completeness (in resolution range)	92.0 (44.88-2.00) 92.0 (44.88-2.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.73 (at 2.00Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6 _289)	Depositor
R, $R_{free}$	0.176 , 0.209 0.173 , 0.206	Depositor DCC
$R_{free}$ test set	7600 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.0	Xtriage
Anisotropy	0.220	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 48.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.010 for -h,-l,-k 0.001 for -h,l,k 0.001 for l,-k,h 0.012 for -l,-k,-h 0.025 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	14315	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.24% 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: MG, SO4, 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	0/3233	0.55	0/4398
1	B	0.40	0/3233	0.55	0/4398
1	C	0.35	0/3233	0.51	0/4398
1	D	0.33	0/3233	0.51	1/4398 (0.0%)
All	All	0.38	0/12932	0.53	1/17592 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	328	ARG	NE-CZ-NH2	-5.45	117.58	120.30

There are no chirality outliers.

There are no planarity outliers.

### 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	3165	0	3134	33	0
1	B	3165	0	3134	40	0
1	C	3165	0	3134	62	0
1	D	3165	0	3134	78	0
2	A	53	0	31	7	0
2	B	53	0	31	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	53	0	31	5	0
2	D	53	0	31	8	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
4	A	5	0	0	0	0
4	B	5	0	0	1	0
5	A	479	0	0	2	0
5	B	448	0	0	7	0
5	C	265	0	0	6	0
5	D	239	0	0	3	0
All	All	14315	0	12660	204	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 8.

All (204) 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:D:152:GLN:HE22	1:D:167:GLU:H	0.97	0.95
1:C:111:GLY:H	1:C:114:HIS:HD2	1.05	0.94
1:A:111:GLY:H	1:A:114:HIS:HD2	1.04	0.94
1:A:142:ASN:HD21	2:A:500:FAD:H61A	1.16	0.92
1:B:106:ILE:HG21	1:B:118:LEU:HD11	1.49	0.92
1:A:376:ILE:HD11	2:B:500:FAD:HM81	1.53	0.90
1:B:142:ASN:HD21	2:B:500:FAD:H61A	1.19	0.88
2:A:500:FAD:HM81	1:B:376:ILE:HD11	1.60	0.84
1:C:142:ASN:HD21	2:C:500:FAD:H61A	1.25	0.82
1:C:111:GLY:H	1:C:114:HIS:CD2	1.95	0.82
1:D:111:GLY:H	1:D:114:HIS:HD2	1.24	0.81
1:D:353:LEU:HD23	1:D:406:ARG:HD2	1.62	0.81
1:A:111:GLY:H	1:A:114:HIS:CD2	1.96	0.78
2:C:500:FAD:HM81	1:D:376:ILE:HD11	1.64	0.78
1:C:152:GLN:HE22	1:C:167:GLU:H	1.30	0.77
1:D:152:GLN:HE22	1:D:167:GLU:N	1.80	0.77
2:D:500:FAD:O2P	2:D:500:FAD:H52A	1.84	0.77
1:A:55:VAL:HG22	1:A:60:LEU:HD22	1.66	0.75
1:D:401:GLY:HA3	5:D:683:HOH:O	1.87	0.74
1:D:134:PHE:HB3	2:D:500:FAD:O2	1.86	0.74
1:B:111:GLY:H	1:B:114:HIS:HD2	1.34	0.74
1:C:376:ILE:HD11	2:D:500:FAD:HM81	1.69	0.73
1:A:376:ILE:HD11	2:B:500:FAD:C8M	2.19	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:55:VAL:HG22	1:C:60:LEU:HD22	1.72	0.72
1:C:379:TYR:CE2	1:D:232:THR:HB	2.25	0.71
1:D:27:MET:HB2	1:D:28:PRO:HD3	1.75	0.69
1:A:163:VAL:HG22	1:A:249:ARG:HG2	1.75	0.68
1:B:409:HIS:HD2	5:B:474:HOH:O	1.76	0.67
1:D:40:SER:OG	1:D:42:LEU:HD22	1.94	0.67
1:B:55:VAL:HG22	1:B:60:LEU:HD22	1.75	0.67
1:D:353:LEU:CD2	1:D:406:ARG:HD2	2.26	0.65
1:A:39:ASN:HB3	5:A:1295:HOH:O	1.96	0.64
1:A:349:ARG:NH2	1:A:409:HIS:ND1	2.43	0.64
1:D:111:GLY:H	1:D:114:HIS:CD2	2.12	0.63
2:A:500:FAD:C8M	1:B:376:ILE:HD11	2.29	0.63
1:D:172:THR:O	1:D:173:ASN:HB2	1.99	0.63
1:D:169:MET:HE2	1:D:242:HIS:HB2	1.82	0.62
1:D:402:ASN:HA	1:D:406:ARG:HB2	1.81	0.62
1:B:300:ASN:C	1:B:300:ASN:HD22	2.03	0.61
1:C:95:THR:HG23	1:C:173:ASN:HD21	1.65	0.61
1:D:300:ASN:HD22	1:D:300:ASN:C	2.02	0.61
1:A:142:ASN:ND2	2:A:500:FAD:H61A	1.94	0.61
1:D:226:VAL:HG11	1:D:229:HIS:CE1	2.35	0.61
1:B:311:GLN:HE21	1:D:312:ALA:H	1.50	0.59
1:D:141:ALA:HA	1:D:405:ILE:HD11	1.84	0.59
1:A:300:ASN:C	1:A:300:ASN:HD22	2.04	0.59
1:A:312:ALA:H	1:C:311:GLN:NE2	2.01	0.59
1:C:152:GLN:NE2	1:C:167:GLU:H	2.00	0.58
1:C:383:ARG:HH11	1:C:383:ARG:HG2	1.68	0.58
1:A:296:ALA:HB2	1:A:313:PHE:CZ	2.38	0.58
1:D:21:LEU:HD12	1:D:25:LEU:HB2	1.85	0.58
1:C:349:ARG:NH2	1:C:409:HIS:ND1	2.52	0.57
1:A:95:THR:HG23	1:A:173:ASN:HD21	1.69	0.57
1:C:378:ALA:HA	5:C:1387:HOH:O	2.04	0.57
2:B:500:FAD:O2P	2:B:500:FAD:H52A	2.05	0.57
1:D:409:HIS:HD2	5:D:1375:HOH:O	1.88	0.57
1:C:111:GLY:N	1:C:114:HIS:HD2	1.88	0.57
1:D:42:LEU:N	1:D:42:LEU:HD13	2.18	0.57
1:D:144:LEU:HD11	1:D:203:GLU:HB2	1.87	0.56
1:B:42:LEU:HD23	1:B:45:GLN:HE21	1.72	0.55
1:C:383:ARG:HG2	1:C:383:ARG:NH1	2.21	0.55
1:D:55:VAL:HG22	1:D:60:LEU:HD22	1.87	0.55
1:D:119:ALA:HB3	1:D:120:PRO:HD3	1.88	0.55
1:A:311:GLN:NE2	1:C:312:ALA:H	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:500:FAD:H51A	1:B:301:ARG:HD2	1.89	0.55
1:B:356:LYS:NZ	1:B:402:ASN:ND2	2.54	0.54
1:C:136:GLU:HG2	1:C:168:LYS:HD3	1.88	0.54
1:B:106:ILE:CG2	1:B:118:LEU:HD11	2.31	0.54
1:A:408:ARG:HH11	1:A:411:GLN:NE2	2.05	0.54
1:D:165:ASN:HD21	1:D:247:ASN:ND2	2.05	0.54
1:D:169:MET:O	1:D:170:TRP:HB2	2.07	0.54
1:B:156:ARG:NH2	4:B:439:SO4:O2	2.39	0.54
1:B:134:PHE:HB3	2:B:500:FAD:O2	2.08	0.54
1:D:142:ASN:HD21	2:D:500:FAD:H61A	1.56	0.54
1:A:312:ALA:H	1:C:311:GLN:HE21	1.56	0.53
1:D:343:GLY:HA3	1:D:347:ALA:HB3	1.91	0.53
2:C:500:FAD:C8M	1:D:376:ILE:HD11	2.35	0.53
1:D:292:ALA:HB3	1:D:317:LEU:HD21	1.91	0.53
1:A:161:GLU:OE2	1:A:249:ARG:HD3	2.09	0.53
1:B:401:GLY:HA3	5:B:1374:HOH:O	2.09	0.52
1:C:301:ARG:HA	1:C:301:ARG:NE	2.25	0.52
1:C:368:THR:HG23	5:C:587:HOH:O	2.11	0.51
1:B:42:LEU:HD23	1:B:45:GLN:NE2	2.25	0.51
1:C:379:TYR:CD2	1:D:232:THR:HB	2.45	0.51
1:D:207:MET:HG2	1:D:259:ALA:N	2.26	0.51
1:B:136:GLU:HG2	1:B:168:LYS:HD3	1.93	0.50
1:B:200:GLN:HE21	1:B:204:ASN:HD22	1.59	0.50
1:D:206:VAL:HB	5:D:767:HOH:O	2.11	0.50
1:C:292:ALA:HB3	1:C:317:LEU:HD21	1.94	0.50
1:C:24:ASN:C	1:C:25:LEU:HD12	2.32	0.50
1:A:172:THR:O	1:A:173:ASN:HB2	2.11	0.50
1:C:39:ASN:HB3	5:C:1225:HOH:O	2.13	0.49
1:C:303:GLY:HA2	1:D:142:ASN:HB3	1.95	0.49
1:A:65:SER:H	1:A:107:ASN:ND2	2.10	0.49
1:D:29:ALA:HB2	1:D:50:THR:OG1	2.12	0.49
1:D:121:PHE:CE2	1:D:184:LEU:HD12	2.48	0.48
1:B:274:ALA:HB1	1:B:356:LYS:HB2	1.95	0.48
1:B:38:PRO:HD2	5:B:530:HOH:O	2.13	0.48
1:C:106:ILE:HG23	1:C:110:ALA:HB2	1.94	0.48
1:C:119:ALA:HB3	1:C:120:PRO:HD3	1.94	0.48
1:B:356:LYS:NZ	1:B:402:ASN:HD21	2.11	0.48
1:A:235:HIS:CE1	1:A:391:THR:HG23	2.49	0.47
1:D:408:ARG:N	1:D:408:ARG:HD2	2.30	0.47
1:A:398:PHE:C	1:A:398:PHE:CD1	2.87	0.47
1:C:106:ILE:HG21	1:C:118:LEU:HD13	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:398:PHE:CD1	1:B:398:PHE:C	2.88	0.47
1:C:172:THR:O	1:C:173:ASN:HB2	2.15	0.47
1:B:406:ARG:NH1	5:B:1174:HOH:O	2.48	0.47
1:D:118:LEU:N	1:D:118:LEU:HD12	2.30	0.47
1:C:251:PRO:HB2	1:C:253:LYS:HG2	1.97	0.47
1:B:156:ARG:NH1	5:B:457:HOH:O	2.45	0.46
1:B:402:ASN:HA	1:B:406:ARG:HB2	1.97	0.46
1:D:144:LEU:CD1	1:D:203:GLU:HB2	2.45	0.46
1:B:409:HIS:CD2	5:B:474:HOH:O	2.60	0.46
1:A:136:GLU:HG2	1:A:168:LYS:HD3	1.97	0.46
1:A:27:MET:HB2	1:A:28:PRO:HD3	1.97	0.46
1:A:165:ASN:HD21	1:A:247:ASN:ND2	2.13	0.46
1:D:106:ILE:HG23	1:D:110:ALA:HB2	1.98	0.46
1:D:292:ALA:CB	1:D:317:LEU:HD21	2.45	0.46
1:D:83:GLU:OE1	1:D:328:ARG:HD3	2.15	0.46
1:D:172:THR:HG23	1:D:238:VAL:HG13	1.97	0.46
1:A:48:GLN:HB3	1:A:49:PRO:HD3	1.97	0.45
1:D:278:GLY:HA2	1:D:359:CYS:HB2	1.97	0.45
1:C:303:GLY:HA2	1:D:142:ASN:CB	2.47	0.45
1:B:311:GLN:NE2	1:D:312:ALA:H	2.15	0.45
1:C:398:PHE:CD1	1:C:398:PHE:C	2.89	0.45
1:D:394:VAL:CG2	2:D:500:FAD:HM83	2.46	0.45
1:C:157:LEU:C	1:C:157:LEU:HD13	2.37	0.45
1:D:114:HIS:HE1	1:D:255:VAL:O	2.00	0.45
1:D:408:ARG:HA	1:D:408:ARG:NE	2.32	0.45
1:A:308:LEU:HB2	5:A:592:HOH:O	2.17	0.45
1:B:356:LYS:HZ3	1:B:402:ASN:HD21	1.65	0.45
1:C:408:ARG:HH11	1:C:411:GLN:NE2	2.15	0.45
1:C:408:ARG:HA	1:C:408:ARG:HD3	1.75	0.44
1:B:48:GLN:HG2	5:B:1169:HOH:O	2.17	0.44
1:C:376:ILE:HA	1:C:376:ILE:HD12	1.85	0.44
1:B:408:ARG:HA	1:B:408:ARG:HD3	1.79	0.44
1:B:172:THR:O	1:B:173:ASN:HB2	2.18	0.44
1:C:25:LEU:N	1:C:25:LEU:HD12	2.32	0.44
1:D:155:ALA:HB2	1:D:164:ILE:HG12	1.99	0.44
1:C:383:ARG:HB2	5:C:1387:HOH:O	2.17	0.44
1:C:278:GLY:HA2	1:C:359:CYS:HB2	2.00	0.44
1:C:89:GLU:O	1:C:89:GLU:HG3	2.18	0.43
1:C:292:ALA:CB	1:C:317:LEU:HD21	2.48	0.43
1:D:36:HIS:CD2	1:D:36:HIS:N	2.87	0.43
1:B:142:ASN:ND2	2:B:500:FAD:H61A	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:GLN:HB2	1:C:311:GLN:HB2	2.00	0.43
1:C:25:LEU:C	1:C:28:PRO:HD2	2.39	0.43
1:B:356:LYS:HZ2	1:B:402:ASN:ND2	2.15	0.43
1:C:165:ASN:HD21	1:C:247:ASN:ND2	2.16	0.43
1:C:169:MET:O	1:C:170:TRP:HB2	2.19	0.43
1:C:29:ALA:HB2	1:C:50:THR:HG21	2.00	0.43
1:B:135:SER:CB	2:B:500:FAD:HO4'	2.31	0.43
1:D:95:THR:HG23	1:D:173:ASN:HD21	1.83	0.43
1:C:379:TYR:CE1	1:D:233:PRO:HG2	2.54	0.43
1:C:38:PRO:HA	1:C:43:ARG:CZ	2.48	0.43
2:D:500:FAD:O2P	2:D:500:FAD:C5B	2.63	0.43
1:D:62:GLY:O	1:D:72:GLY:HA2	2.18	0.42
1:D:317:LEU:HD23	1:D:317:LEU:HA	1.80	0.42
1:D:35:GLN:HB2	1:D:36:HIS:CD2	2.53	0.42
1:D:101:LEU:HD11	1:D:132:LEU:HD23	2.02	0.42
1:D:37:PRO:HA	1:D:38:PRO:HD3	1.77	0.42
1:D:207:MET:HE3	1:D:209:ILE:HD11	2.01	0.42
1:D:399:ASP:H	2:D:500:FAD:H2'	1.83	0.42
1:C:106:ILE:CG2	1:C:110:ALA:HB2	2.49	0.42
1:C:158:GLU:OE1	1:C:163:VAL:HG21	2.20	0.42
1:C:379:TYR:CE1	1:D:394:VAL:HB	2.55	0.42
1:D:169:MET:O	1:D:170:TRP:CB	2.68	0.42
1:C:134:PHE:HB3	2:C:500:FAD:O2	2.20	0.42
1:C:377:SER:O	1:C:380:ASP:HB3	2.20	0.42
1:C:235:HIS:CE1	1:C:391:THR:HG23	2.55	0.42
1:D:106:ILE:HG21	1:D:118:LEU:HD11	2.01	0.42
1:B:300:ASN:C	1:B:300:ASN:ND2	2.72	0.41
1:D:111:GLY:N	1:D:114:HIS:HD2	2.03	0.41
1:D:251:PRO:HB2	1:D:253:LYS:HG2	2.02	0.41
1:A:312:ALA:HB1	1:B:403:VAL:HG13	2.00	0.41
1:A:278:GLY:HA2	1:A:359:CYS:HB2	2.01	0.41
1:A:408:ARG:HA	1:A:408:ARG:HD3	1.82	0.41
1:D:356:LYS:NZ	1:D:402:ASN:ND2	2.68	0.41
2:A:500:FAD:H51A	1:B:301:ARG:CD	2.49	0.41
2:C:500:FAD:H52A	2:C:500:FAD:O2P	2.20	0.41
1:C:65:SER:H	1:C:107:ASN:ND2	2.18	0.41
1:D:37:PRO:HB2	1:D:40:SER:CB	2.51	0.41
1:A:133:VAL:HG12	1:A:168:LYS:HG3	2.03	0.41
1:B:169:MET:O	1:B:170:TRP:HB2	2.20	0.41
1:C:310:ARG:HD3	5:C:453:HOH:O	2.20	0.41
1:A:311:GLN:HE21	1:C:312:ALA:H	1.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:170:TRP:CH2	1:D:230:VAL:HG11	2.55	0.41
1:D:400:GLY:O	1:D:406:ARG:HG2	2.21	0.41
2:A:500:FAD:H9	2:A:500:FAD:H1'2	1.91	0.41
1:C:175:ALA:HB3	5:C:915:HOH:O	2.20	0.41
1:B:89:GLU:O	1:B:89:GLU:HG3	2.22	0.40
1:C:376:ILE:HD11	2:D:500:FAD:C8M	2.45	0.40
1:D:270:PHE:HA	1:D:273:SER:OG	2.21	0.40
1:D:318:SER:O	1:D:322:ILE:HG13	2.21	0.40
1:C:207:MET:HG2	1:C:259:ALA:HA	2.04	0.40
1:D:121:PHE:HE2	1:D:184:LEU:HD12	1.85	0.40
1:D:391:THR:O	1:D:394:VAL:HG12	2.21	0.40
1:D:33:TYR:HB2	1:D:43:ARG:HG3	2.03	0.40
1:C:395:LEU:HB2	1:C:396:PRO:HD3	2.03	0.40
1:D:152:GLN:HG2	1:D:152:GLN:O	2.22	0.40
1:D:41:PRO:O	1:D:44:PHE:HB3	2.20	0.40
1:C:164:ILE:HD11	1:C:209:ILE:HG21	2.02	0.40
1:D:175:ALA:O	1:D:179:PHE:HA	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	424/438 (97%)	414 (98%)	10 (2%)	0	100	100
1	B	424/438 (97%)	415 (98%)	8 (2%)	1 (0%)	47	44
1	C	424/438 (97%)	411 (97%)	12 (3%)	1 (0%)	47	44
1	D	424/438 (97%)	413 (97%)	11 (3%)	0	100	100
All	All	1696/1752 (97%)	1653 (98%)	41 (2%)	2 (0%)	51	49

All (2) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	B	72	GLY
1	C	344	ASP

### 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	323/334 (97%)	312 (97%)	11 (3%)	37	36
1	B	323/334 (97%)	312 (97%)	11 (3%)	37	36
1	C	323/334 (97%)	309 (96%)	14 (4%)	29	26
1	D	323/334 (97%)	310 (96%)	13 (4%)	31	29
All	All	1292/1336 (97%)	1243 (96%)	49 (4%)	33	31

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	LEU
1	A	60	LEU
1	A	65	SER
1	A	118	LEU
1	A	276	LEU
1	A	300	ASN
1	A	316	LEU
1	A	344	ASP
1	A	394	VAL
1	A	398	PHE
1	A	408	ARG
1	B	34	LEU
1	B	184	LEU
1	B	247	ASN
1	B	300	ASN
1	B	316	LEU
1	B	317	LEU
1	B	349	ARG
1	B	388	LEU
1	B	394	VAL

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Mol	Chain	Res	Type
1	B	398	PHE
1	B	408	ARG
1	C	30	ARG
1	C	34	LEU
1	C	60	LEU
1	C	152	GLN
1	C	184	LEU
1	C	276	LEU
1	C	300	ASN
1	C	316	LEU
1	C	388	LEU
1	C	393	VAL
1	C	394	VAL
1	C	398	PHE
1	C	408	ARG
1	C	412	GLN
1	D	23	ARG
1	D	34	LEU
1	D	42	LEU
1	D	132	LEU
1	D	144	LEU
1	D	184	LEU
1	D	276	LEU
1	D	300	ASN
1	D	316	LEU
1	D	388	LEU
1	D	394	VAL
1	D	398	PHE
1	D	408	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (57) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	GLN
1	A	107	ASN
1	A	114	HIS
1	A	142	ASN
1	A	173	ASN
1	A	204	ASN
1	A	247	ASN
1	A	254	ASN
1	A	300	ASN

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Mol	Chain	Res	Type
1	A	311	GLN
1	A	382	GLN
1	A	402	ASN
1	A	411	GLN
1	B	35	GLN
1	B	45	GLN
1	B	107	ASN
1	B	114	HIS
1	B	142	ASN
1	B	173	ASN
1	B	204	ASN
1	B	247	ASN
1	B	254	ASN
1	B	300	ASN
1	B	311	GLN
1	B	402	ASN
1	B	409	HIS
1	B	411	GLN
1	C	45	GLN
1	C	107	ASN
1	C	114	HIS
1	C	142	ASN
1	C	152	GLN
1	C	173	ASN
1	C	204	ASN
1	C	247	ASN
1	C	254	ASN
1	C	300	ASN
1	C	311	GLN
1	C	382	GLN
1	C	402	ASN
1	C	411	GLN
1	C	412	GLN
1	D	36	HIS
1	D	45	GLN
1	D	107	ASN
1	D	114	HIS
1	D	142	ASN
1	D	152	GLN
1	D	173	ASN
1	D	229	HIS
1	D	247	ASN

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Mol	Chain	Res	Type
1	D	254	ASN
1	D	300	ASN
1	D	402	ASN
1	D	409	HIS
1	D	411	GLN
1	D	412	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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 for Mogul analysis.

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	500	-	51,58,58	1.56	5 (9%)	60,89,89	1.54	10 (16%)
4	SO4	A	440	-	4,4,4	0.14	0	6,6,6	0.19	0
4	SO4	B	439	-	4,4,4	0.16	0	6,6,6	0.07	0
2	FAD	A	500	-	51,58,58	1.50	5 (9%)	60,89,89	1.57	9 (15%)
2	FAD	C	500	-	51,58,58	1.49	5 (9%)	60,89,89	1.50	8 (13%)
2	FAD	D	500	-	51,58,58	1.55	5 (9%)	60,89,89	1.51	9 (15%)

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	500	-	-	3/30/50/50	0/6/6/6
2	FAD	A	500	-	-	5/30/50/50	0/6/6/6
2	FAD	D	500	-	-	10/30/50/50	0/6/6/6
2	FAD	C	500	-	-	8/30/50/50	0/6/6/6

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	500	FAD	C10-N1	4.95	1.39	1.33
2	B	500	FAD	C10-N1	4.89	1.39	1.33
2	D	500	FAD	C4X-N5	4.84	1.40	1.33
2	D	500	FAD	C10-N1	4.81	1.39	1.33
2	A	500	FAD	C4X-N5	4.81	1.40	1.33
2	C	500	FAD	C4X-N5	4.74	1.40	1.33
2	A	500	FAD	C10-N1	4.73	1.39	1.33
2	B	500	FAD	C4X-N5	4.72	1.40	1.33
2	B	500	FAD	O4B-C1B	4.23	1.47	1.41
2	D	500	FAD	O4B-C1B	4.04	1.46	1.41
2	D	500	FAD	C4-N3	4.02	1.40	1.33
2	B	500	FAD	C4-N3	3.86	1.39	1.33
2	A	500	FAD	C4-N3	3.84	1.39	1.33
2	C	500	FAD	C4-N3	3.71	1.39	1.33
2	C	500	FAD	O4B-C1B	3.30	1.45	1.41
2	A	500	FAD	O4B-C1B	3.13	1.45	1.41
2	A	500	FAD	C5X-N5	3.01	1.40	1.35
2	B	500	FAD	C5X-N5	2.83	1.40	1.35
2	C	500	FAD	C5X-N5	2.76	1.39	1.35
2	D	500	FAD	C5X-N5	2.69	1.39	1.35

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	FAD	C4-N3-C2	6.35	120.50	115.14
2	C	500	FAD	C4-N3-C2	6.10	120.29	115.14
2	B	500	FAD	C4-N3-C2	5.81	120.05	115.14
2	D	500	FAD	C4-N3-C2	5.76	120.01	115.14
2	A	500	FAD	N3A-C2A-N1A	-4.41	121.79	128.68
2	D	500	FAD	N3A-C2A-N1A	-4.41	121.79	128.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	500	FAD	N3A-C2A-N1A	-4.33	121.91	128.68
2	B	500	FAD	N3A-C2A-N1A	-4.17	122.16	128.68
2	B	500	FAD	C1'-N10-C9A	3.77	121.26	118.29
2	A	500	FAD	C1'-N10-C9A	3.53	121.07	118.29
2	C	500	FAD	C1'-N10-C9A	3.51	121.06	118.29
2	A	500	FAD	C10-C4X-N5	-3.13	119.09	121.26
2	D	500	FAD	C1'-N10-C9A	3.04	120.69	118.29
2	A	500	FAD	C4X-C4-N3	-3.04	119.28	123.43
2	D	500	FAD	C4X-N5-C5X	2.94	119.71	116.77
2	D	500	FAD	C5X-C9A-N10	2.90	119.82	117.72
2	B	500	FAD	C10-C4X-N5	-2.86	119.28	121.26
2	B	500	FAD	C4X-C4-N3	-2.79	119.61	123.43
2	C	500	FAD	C4X-C4-N3	-2.76	119.66	123.43
2	A	500	FAD	C5X-C9A-N10	2.75	119.71	117.72
2	C	500	FAD	C5X-C9A-N10	2.67	119.65	117.72
2	B	500	FAD	P-O3P-PA	-2.62	123.82	132.83
2	D	500	FAD	C4X-C4-N3	-2.62	119.85	123.43
2	B	500	FAD	C5X-C9A-N10	2.47	119.50	117.72
2	B	500	FAD	C9A-N10-C10	-2.46	118.68	121.91
2	D	500	FAD	P-O3P-PA	-2.42	124.51	132.83
2	D	500	FAD	C10-C4X-N5	-2.42	119.59	121.26
2	C	500	FAD	C10-C4X-N5	-2.41	119.59	121.26
2	A	500	FAD	C4X-N5-C5X	2.38	119.14	116.77
2	C	500	FAD	C9A-N10-C10	-2.34	118.84	121.91
2	B	500	FAD	C4A-C5A-N7A	-2.32	106.98	109.40
2	D	500	FAD	C3B-C2B-C1B	2.16	104.24	100.98
2	A	500	FAD	C9A-N10-C10	-2.16	119.08	121.91
2	B	500	FAD	C4X-N5-C5X	2.10	118.87	116.77
2	A	500	FAD	C9A-C5X-N5	-2.02	119.19	122.36
2	C	500	FAD	C4X-N5-C5X	2.00	118.77	116.77

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	500	FAD	C5B-O5B-PA-O2A
2	D	500	FAD	O4B-C4B-C5B-O5B
2	D	500	FAD	C1'-C2'-C3'-O3'
2	D	500	FAD	C1'-C2'-C3'-C4'
2	A	500	FAD	C5B-O5B-PA-O2A
2	C	500	FAD	C5B-O5B-PA-O2A
2	D	500	FAD	C3B-C4B-C5B-O5B

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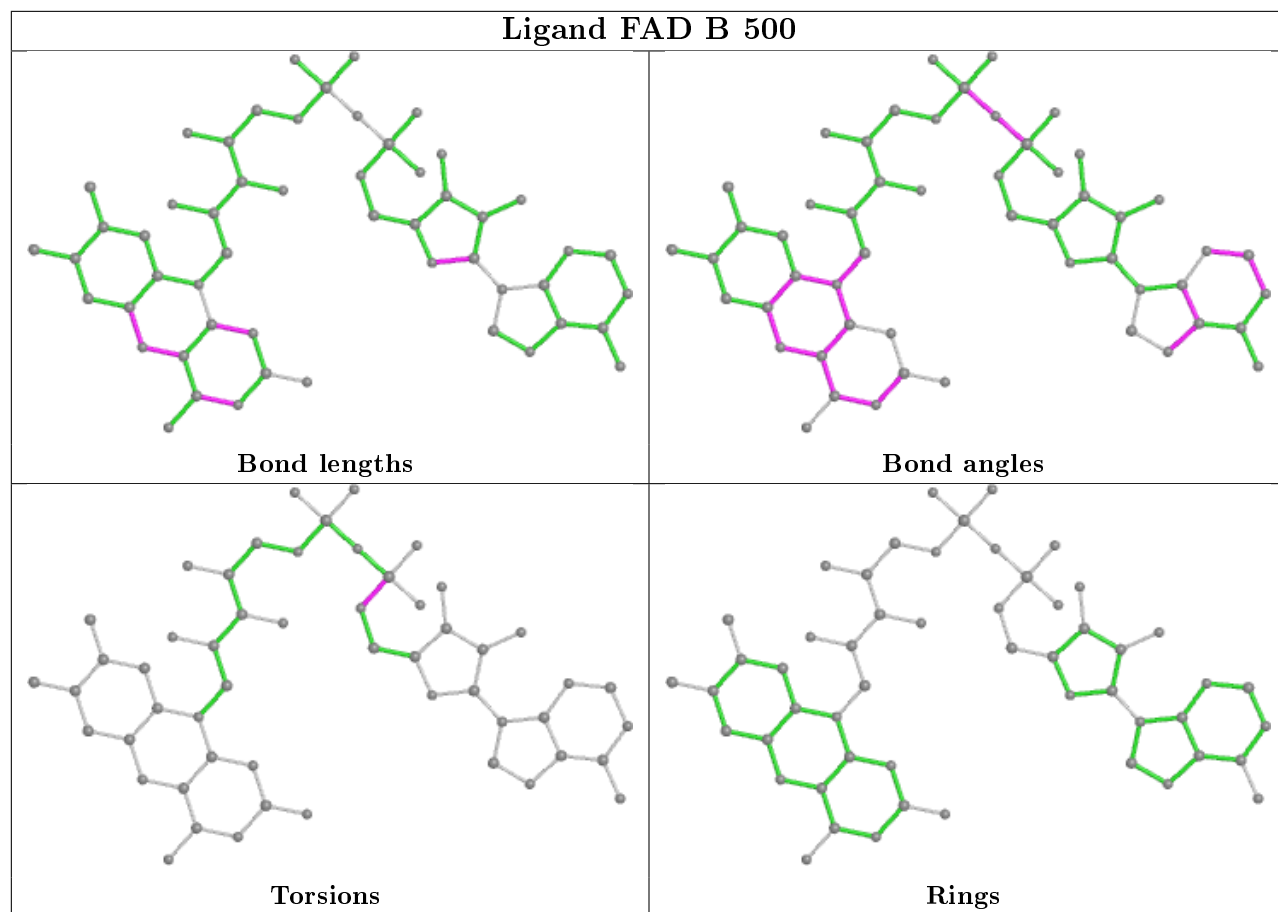
Mol	Chain	Res	Type	Atoms
2	D	500	FAD	O2'-C2'-C3'-O3'
2	C	500	FAD	PA-O3P-P-O1P
2	D	500	FAD	O3'-C3'-C4'-O4'
2	B	500	FAD	C5B-O5B-PA-O3P
2	A	500	FAD	C5B-O5B-PA-O3P
2	C	500	FAD	C5B-O5B-PA-O3P
2	D	500	FAD	O2'-C2'-C3'-C4'
2	D	500	FAD	C2'-C3'-C4'-O4'
2	A	500	FAD	C5B-O5B-PA-O1A
2	C	500	FAD	C5B-O5B-PA-O1A
2	D	500	FAD	O3'-C3'-C4'-C5'
2	C	500	FAD	O3'-C3'-C4'-O4'
2	A	500	FAD	PA-O3P-P-O1P
2	C	500	FAD	O3'-C3'-C4'-C5'
2	D	500	FAD	C2'-C3'-C4'-C5'
2	C	500	FAD	C2'-C3'-C4'-O4'
2	A	500	FAD	PA-O3P-P-O2P
2	C	500	FAD	PA-O3P-P-O2P
2	B	500	FAD	C5B-O5B-PA-O1A

There are no ring outliers.

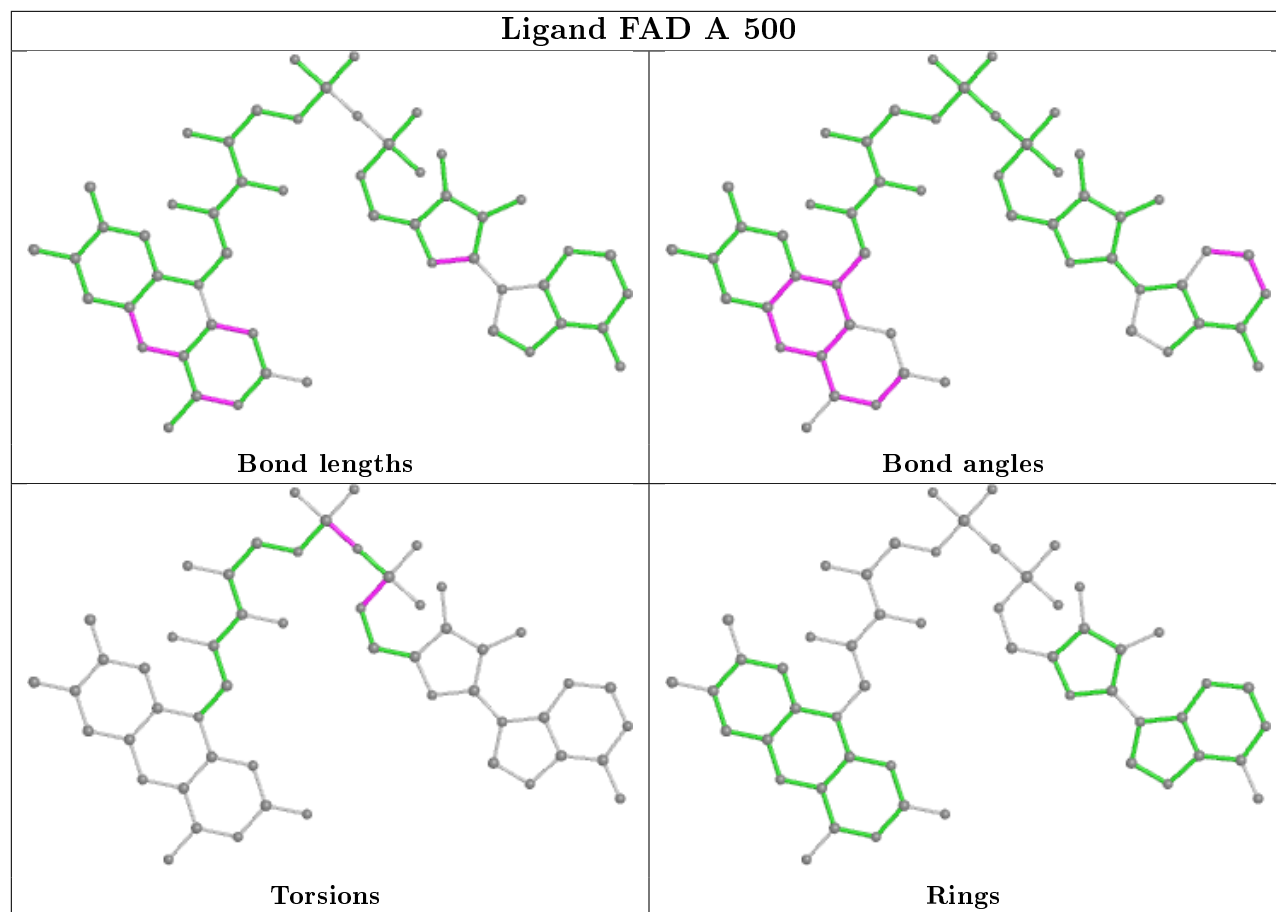
5 monomers are involved in 28 short contacts:

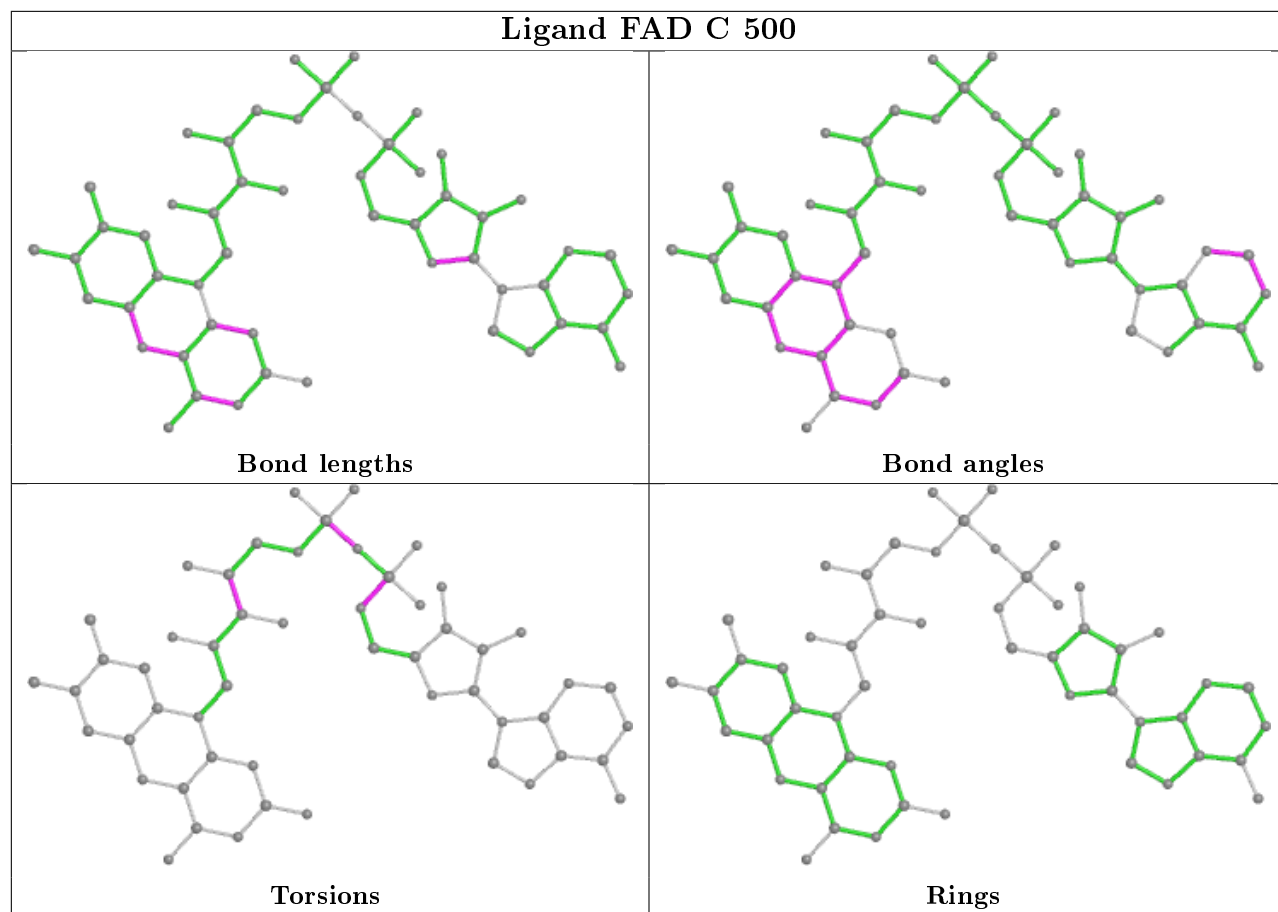
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	500	FAD	7	0
4	B	439	SO4	1	0
2	A	500	FAD	7	0
2	C	500	FAD	5	0
2	D	500	FAD	8	0

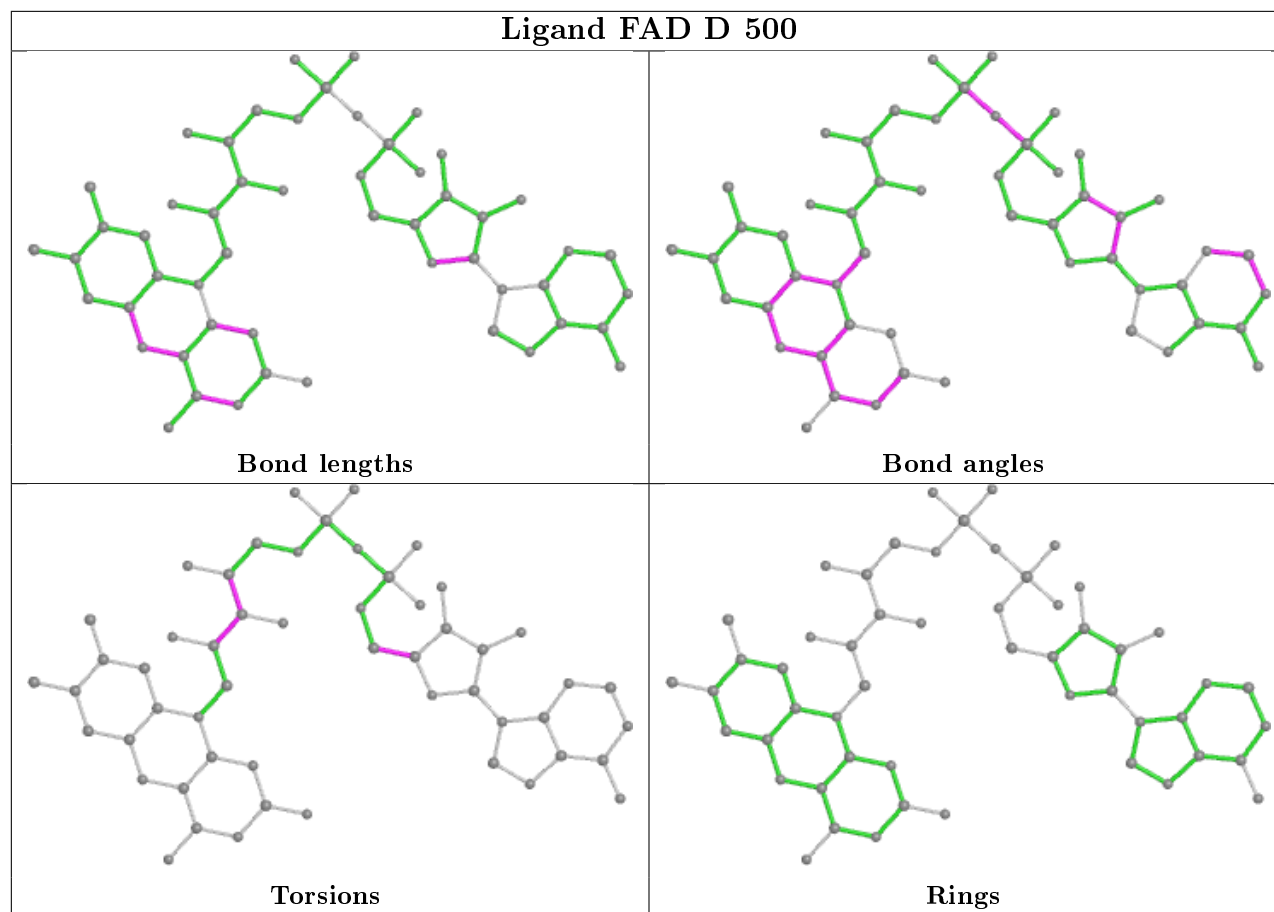
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

### 6.1 Protein, DNA and RNA chains ⓘ

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	426/438 (97%)	-0.16	10 (2%) 60 59	18, 26, 42, 61	0
1	B	426/438 (97%)	-0.04	5 (1%) 79 78	19, 29, 44, 67	0
1	C	426/438 (97%)	0.31	22 (5%) 27 26	23, 38, 60, 75	0
1	D	426/438 (97%)	0.32	29 (6%) 17 16	21, 41, 68, 96	0
All	All	1704/1752 (97%)	0.11	66 (3%) 39 38	18, 32, 59, 96	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	221	GLU	5.9
1	A	302	GLY	5.5
1	C	306	PRO	5.4
1	C	305	VAL	4.8
1	C	157	LEU	4.6
1	D	200	GLN	4.5
1	A	304	ALA	4.3
1	B	221	GLU	3.9
1	D	192	ALA	3.8
1	D	198	GLU	3.8
1	C	304	ALA	3.8
1	C	300	ASN	3.7
1	D	196	LEU	3.6
1	D	221	GLU	3.6
1	C	160	ASP	3.3
1	C	222	GLY	3.3
1	C	199	GLY	3.2
1	C	198	GLU	3.2
1	D	222	GLY	3.1
1	D	161	GLU	3.1
1	A	160	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	199	GLY	3.0
1	D	404	GLY	3.0
1	A	305	VAL	3.0
1	B	405	ILE	3.0
1	A	198	GLU	2.9
1	D	111	GLY	2.9
1	D	141	ALA	2.9
1	D	405	ILE	2.9
1	D	195	PRO	2.9
1	C	23	ARG	2.8
1	D	304	ALA	2.8
1	D	160	ASP	2.8
1	A	300	ASN	2.7
1	C	112	PRO	2.7
1	D	42	LEU	2.7
1	A	306	PRO	2.6
1	A	159	GLY	2.6
1	D	343	GLY	2.6
1	C	194	THR	2.6
1	C	378	ALA	2.5
1	C	303	GLY	2.5
1	D	232	THR	2.5
1	C	196	LEU	2.5
1	B	404	GLY	2.5
1	D	158	GLU	2.5
1	D	247	ASN	2.5
1	C	28	PRO	2.5
1	D	28	PRO	2.5
1	D	398	PHE	2.5
1	B	23	ARG	2.4
1	B	198	GLU	2.3
1	C	308	LEU	2.3
1	C	156	ARG	2.3
1	C	379	TYR	2.3
1	C	197	GLU	2.3
1	A	197	GLU	2.2
1	D	163	VAL	2.2
1	D	157	LEU	2.2
1	A	199	GLY	2.2
1	D	197	GLU	2.1
1	D	143	ALA	2.1
1	C	220	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	31	GLN	2.1
1	D	399	ASP	2.0
1	D	23	ARG	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](#)

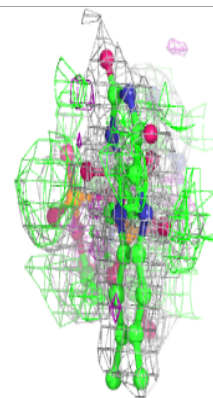
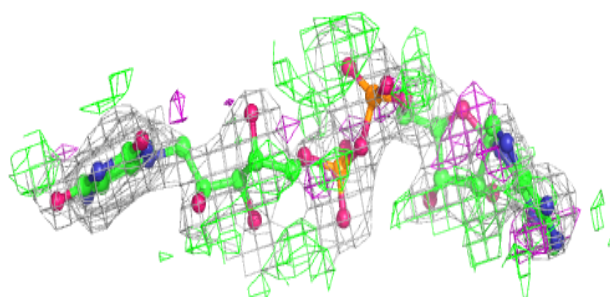
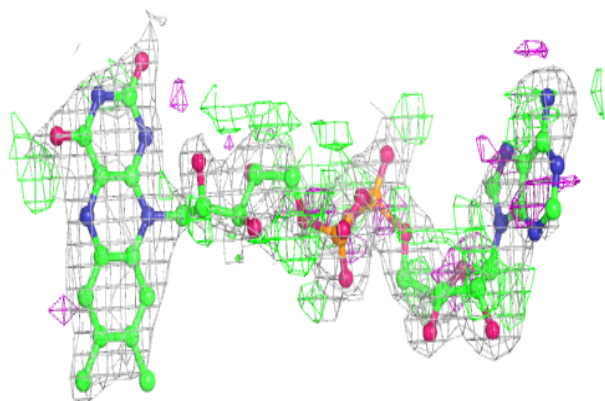
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	FAD	D	500	53/53	0.76	0.24	38,50,58,58	53
2	FAD	B	500	53/53	0.89	0.19	26,32,41,44	53
2	FAD	C	500	53/53	0.93	0.14	24,32,37,37	53
3	MG	A	439	1/1	0.93	0.07	44,44,44,44	0
3	MG	C	439	1/1	0.95	0.10	53,53,53,53	0
2	FAD	A	500	53/53	0.95	0.12	21,26,30,33	53
4	SO4	A	440	5/5	0.96	0.21	48,56,61,73	0
4	SO4	B	439	5/5	0.96	0.15	52,53,61,62	0

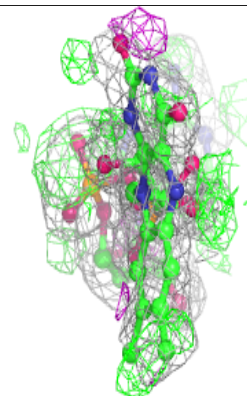
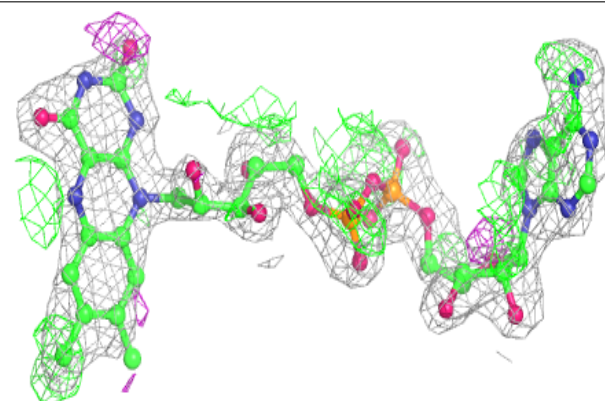
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around FAD D 500:**

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

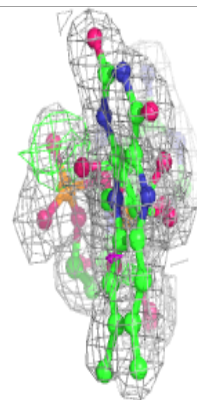
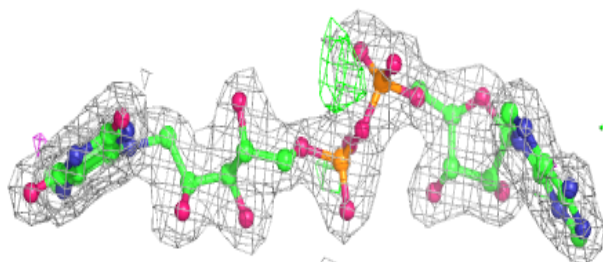
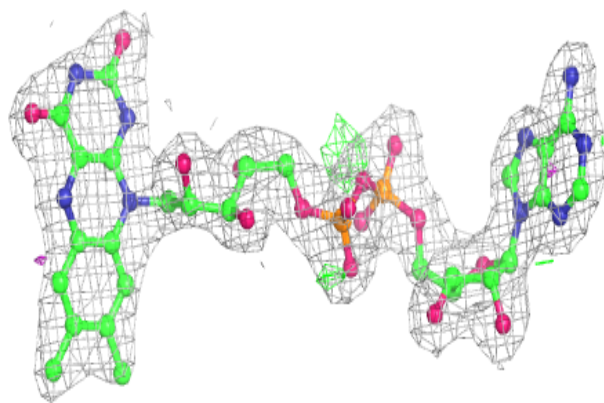
**Electron density around FAD B 500:**

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

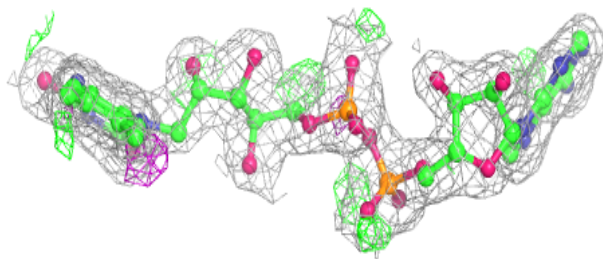
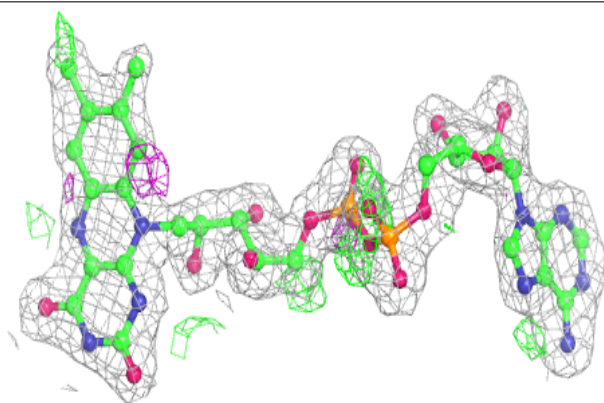


**Electron density around FAD C 500:**

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

**Electron density around FAD A 500:**

$2mF_o-DF_c$  (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.