



# Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 05:30 am BST

PDB ID : 5KSW  
Title : DHODB-I74D mutant  
Authors : Pompeu, Y.A.; Stewart, J.D.  
Deposited on : 2016-07-10  
Resolution : 2.47 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

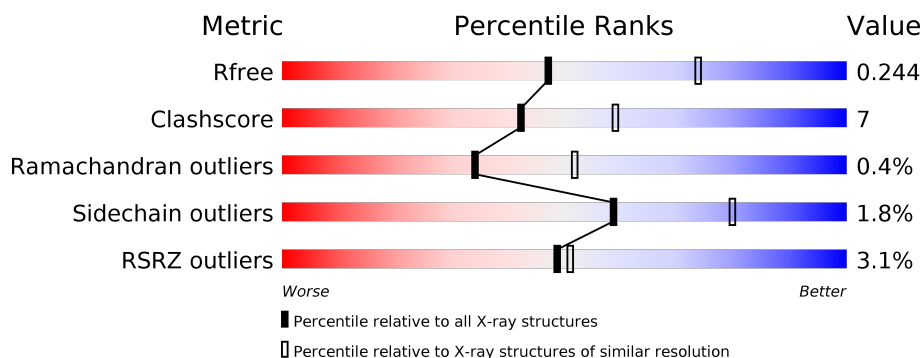
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.47 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	5857 (2.50-2.46)
Clashscore	141614	6594 (2.50-2.46)
Ramachandran outliers	138981	6469 (2.50-2.46)
Sidechain outliers	138945	6471 (2.50-2.46)
RSRZ outliers	127900	5738 (2.50-2.46)

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	306	<div> <div>2%</div> <div> <div></div> <div>87%</div> <div>13%</div> </div> </div>
1	C	306	<div> <div>6%</div> <div> <div></div> <div>80%</div> <div>19%</div> </div> </div>
2	B	262	<div> <div>3%</div> <div> <div></div> <div>83%</div> <div>16%</div> </div> </div>
2	D	262	<div> <div></div> <div> <div></div> <div>82%</div> <div>17%</div> </div> </div>

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 9054 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 Dihydroorotate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	306	Total	C	N	O	S	0	0	0
			2260	1445	374	427	14			
1	C	305	Total	C	N	O	S	0	1	0
			2246	1439	372	421	14			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	67	ALA	SER	conflict	UNP A0A089ZD72
A	74	ASP	ILE	ENGINEERED MUTATION	UNP A0A089ZD72
A	128	VAL	ALA	conflict	UNP A0A089ZD72
A	187	ALA	HIS	conflict	UNP A0A089ZD72
A	189	GLY	CYS	conflict	UNP A0A089ZD72
A	308	ASN	LYS	conflict	UNP A0A089ZD72
C	67	ALA	SER	conflict	UNP A0A089ZD72
C	74	ASP	ILE	ENGINEERED MUTATION	UNP A0A089ZD72
C	128	VAL	ALA	conflict	UNP A0A089ZD72
C	187	ALA	HIS	conflict	UNP A0A089ZD72
C	189	GLY	CYS	conflict	UNP A0A089ZD72
C	308	ASN	LYS	conflict	UNP A0A089ZD72

- Molecule 2 is a protein called Dihydroorotate dehydrogenase B (NAD(+)), electron transfer subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	262	Total	C	N	O	S	0	0	0
			1988	1266	321	381	20			
2	D	262	Total	C	N	O	S	0	0	0
			1985	1265	321	379	20			

- Molecule 3 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: C<sub>17</sub>H<sub>21</sub>N<sub>4</sub>O<sub>9</sub>P).

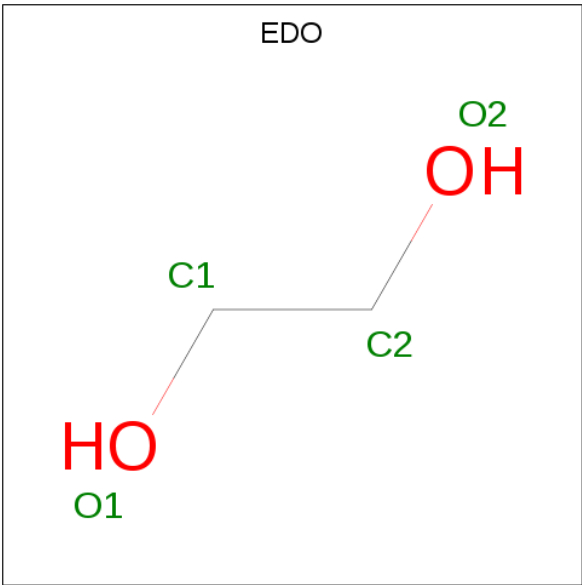


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
3	C	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

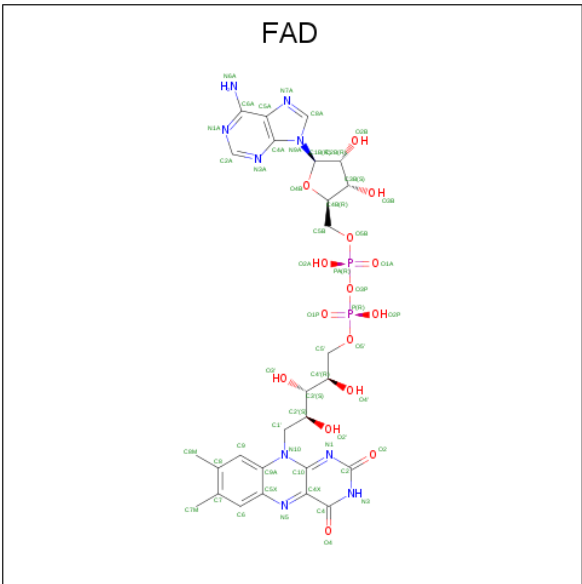
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Cl	0	0
			1	1		
4	A	1	Total	Cl	0	0
			1	1		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C<sub>27</sub>H<sub>33</sub>N<sub>9</sub>O<sub>15</sub>P<sub>2</sub>).



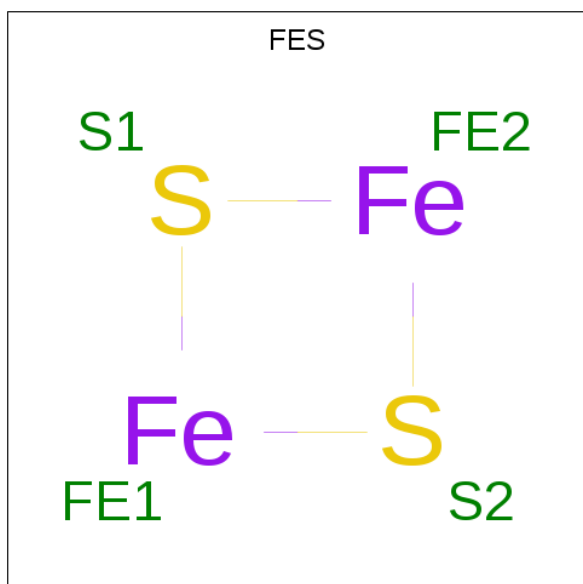
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

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

- Molecule 7 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe<sub>2</sub>S<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	Fe	S	0	0
			4	2	2		
7	D	1	Total	Fe	S	0	0
			4	2	2		

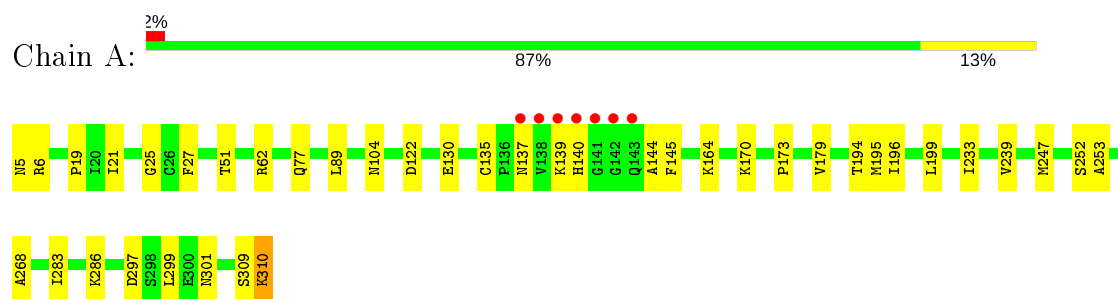
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	122	Total	O	0	0
			122	122		
8	B	86	Total	O	0	0
			86	86		
8	D	107	Total	O	0	0
			107	107		
8	C	74	Total	O	0	0
			74	74		

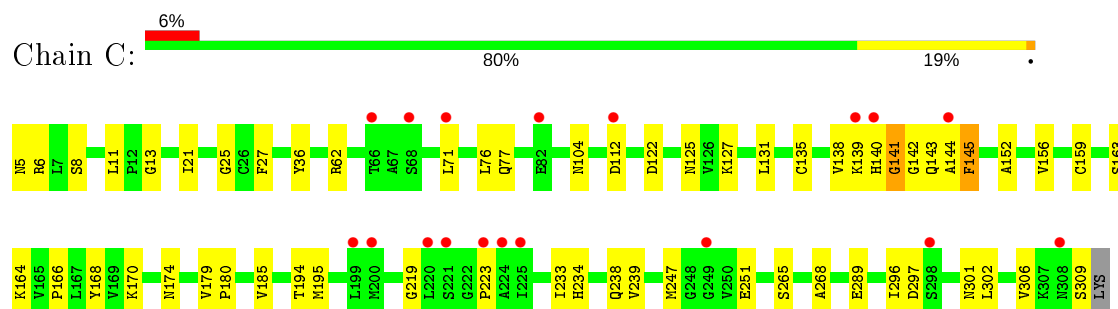
### 3 Residue-property plots [i](#)

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

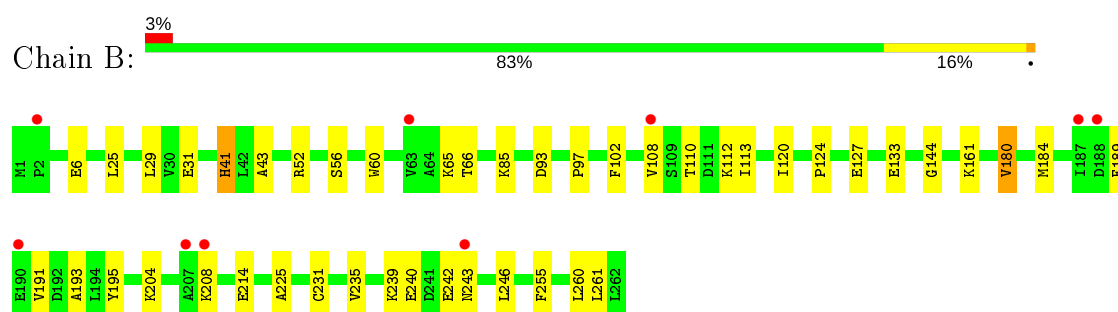
- Molecule 1: Dihydroorotate dehydrogenase



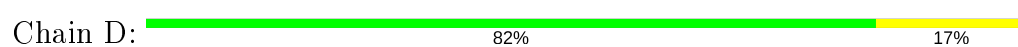
- Molecule 1: Dihydroorotate dehydrogenase

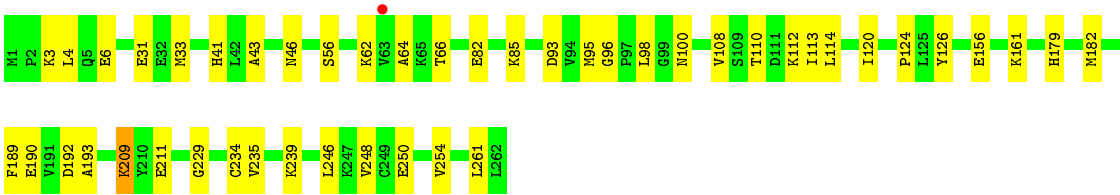


- Molecule 2: Dihydroorotate dehydrogenase B (NAD(+)), electron transfer subunit



- Molecule 2: Dihydroorotate dehydrogenase B (NAD(+)), electron transfer subunit







## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.99Å 151.72Å 214.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.92 – 2.47 41.92 – 2.47	Depositor EDS
% Data completeness (in resolution range)	100.0 (41.92-2.47) 95.4 (41.92-2.47)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.91 (at 2.48Å)	Xtriage
Refinement program	PHENIX 1.10.1 _2155	Depositor
R, $R_{free}$	0.176 , 0.244 0.177 , 0.244	Depositor DCC
$R_{free}$ test set	2369 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	41.1	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 53.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	9054	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FMN, FAD, EDO, FES, CL

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	0.43	0/2302	0.58	0/3134
1	C	0.42	0/2291	0.59	0/3119
2	B	0.45	0/2018	0.64	0/2718
2	D	0.43	0/2015	0.62	0/2714
All	All	0.43	0/8626	0.61	0/11685

There are no bond length outliers.

There are no bond angle outliers.

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	2260	0	2314	28	0
1	C	2246	0	2299	39	0
2	B	1988	0	2004	28	0
2	D	1985	0	2002	30	0
3	A	31	0	19	1	0
3	C	31	0	19	1	0
4	A	1	0	0	1	0
4	B	1	0	0	0	0
5	A	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	4	0	6	0	0
6	B	53	0	31	1	0
6	D	53	0	31	1	0
7	B	4	0	0	0	0
7	D	4	0	0	0	0
8	A	122	0	0	4	1
8	B	86	0	0	3	1
8	C	74	0	0	4	0
8	D	107	0	0	6	0
All	All	9054	0	8731	120	1

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 7.

All (120) 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)
2:B:85:LYS:NZ	8:B:401:HOH:O	2.03	0.90
1:A:286:LYS:NZ	8:A:501:HOH:O	2.06	0.88
2:D:235:VAL:HG11	1:C:77:GLN:HE22	1.45	0.82
2:D:192:ASP:OD1	8:D:401:HOH:O	2.01	0.79
2:D:190:GLU:OE1	8:D:401:HOH:O	2.03	0.76
1:A:310:LYS:NZ	8:A:502:HOH:O	2.20	0.73
1:A:253:ALA:H	1:A:286:LYS:HE3	1.54	0.73
1:A:283:ILE:HA	1:A:286:LYS:HE2	1.70	0.73
1:C:5:ASN:N	1:C:8:SER:HG	1.88	0.72
1:A:179:VAL:HG13	1:A:239:VAL:HG11	1.74	0.70
1:A:309:SER:N	1:A:310:LYS:HA	2.07	0.69
1:C:141:GLY:O	1:C:143:GLN:N	2.28	0.66
1:C:159:CYS:O	1:C:163:SER:OG	2.17	0.63
2:B:239:LYS:HG3	2:B:261:LEU:HG	1.81	0.61
1:A:77:GLN:HE22	2:B:235:VAL:HG11	1.67	0.60
2:B:65:LYS:HA	8:B:403:HOH:O	2.01	0.59
2:D:113:ILE:HD12	2:D:193:ALA:HB3	1.85	0.59
2:D:4:LEU:HD22	1:C:36:TYR:CG	2.38	0.58
1:A:122:ASP:OD1	1:A:164:LYS:HE2	2.03	0.58
1:C:306:VAL:O	1:C:309:SER:OG	2.19	0.57
2:B:97:PRO:O	8:B:402:HOH:O	2.18	0.57
1:C:247:MET:HB3	1:C:268:ALA:HB3	1.88	0.56
2:D:43:ALA:HB3	2:D:93:ASP:HB3	1.87	0.56
1:A:62:ARG:HH12	1:A:77:GLN:NE2	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:156:VAL:HG21	1:C:185:VAL:HG22	1.88	0.55
2:B:43:ALA:HB3	2:B:93:ASP:HB3	1.87	0.55
1:C:5:ASN:CG	1:C:6:ARG:H	2.10	0.54
1:C:138:VAL:HG23	1:C:141:GLY:HA2	1.90	0.54
1:C:143:GLN:HE22	1:C:174:ASN:HD22	1.56	0.54
2:D:108:VAL:HG12	2:D:110:THR:HG23	1.89	0.54
2:B:133:GLU:OE2	2:B:161:LYS:HG2	2.08	0.53
1:C:25:GLY:HA3	3:C:401:FMN:N5	2.24	0.53
2:D:96:GLY:HA2	2:D:98:LEU:HG	1.92	0.52
1:A:21:ILE:O	1:A:268:ALA:HA	2.10	0.52
1:A:5:ASN:CG	1:A:6:ARG:H	2.13	0.52
2:D:4:LEU:O	2:D:95:MET:HA	2.10	0.52
1:A:25:GLY:HA3	3:A:401:FMN:N5	2.24	0.51
1:A:247:MET:HB3	1:A:268:ALA:HB3	1.92	0.51
2:B:242:GLU:O	2:B:243:ASN:HB2	2.10	0.50
1:A:137:ASN:ND2	1:A:140:HIS:HB2	2.27	0.50
1:A:252:SER:HB2	8:A:566:HOH:O	2.12	0.50
2:B:133:GLU:OE2	2:B:161:LYS:N	2.43	0.49
1:C:11:LEU:HD21	1:C:265:SER:HB2	1.93	0.49
1:C:195:MET:HG3	1:C:233:ILE:HG12	1.94	0.49
1:C:223:PRO:HG3	1:C:251[B]:GLU:HG3	1.94	0.49
2:D:56:SER:HB3	2:D:124:PRO:HB3	1.93	0.49
2:B:102:PHE:CD1	2:B:195:TYR:HD2	2.31	0.48
1:C:234:HIS:O	1:C:238:GLN:HG2	2.14	0.48
2:D:239:LYS:HG3	2:D:261:LEU:HG	1.95	0.48
1:C:127:LYS:O	1:C:166:PRO:HD2	2.14	0.48
2:D:235:VAL:HA	2:D:246:LEU:O	2.14	0.47
1:C:143:GLN:NE2	1:C:174:ASN:HB2	2.29	0.47
2:B:204:LYS:O	2:B:208:LYS:HG3	2.13	0.47
1:C:104:ASN:O	8:C:501:HOH:O	2.21	0.47
1:C:289:GLU:HG3	8:C:510:HOH:O	2.15	0.47
2:B:189:PHE:CE1	2:B:191:VAL:HG12	2.50	0.47
1:C:143:GLN:NE2	1:C:174:ASN:HD22	2.13	0.47
1:C:62:ARG:HH12	1:C:77:GLN:NE2	2.13	0.47
2:D:120:ILE:HD12	6:D:301:FAD:N1	2.30	0.47
1:C:179:VAL:HG13	1:C:239:VAL:HG11	1.98	0.46
2:D:179:HIS:HB3	8:D:491:HOH:O	2.15	0.46
2:B:214:GLU:OE1	2:D:182:MET:HB2	2.14	0.46
2:B:235:VAL:HA	2:B:246:LEU:O	2.16	0.46
2:B:240:GLU:HB2	2:D:209:LYS:HE2	1.97	0.46
2:B:108:VAL:CG1	2:B:110:THR:HG22	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:152:ALA:O	1:C:156:VAL:HG23	2.15	0.46
2:B:144:GLY:HA3	2:B:180:VAL:HG13	1.97	0.46
2:B:113:ILE:HD12	2:B:193:ALA:HB3	1.97	0.46
2:D:3:LYS:NZ	8:D:403:HOH:O	2.31	0.46
1:A:135:CYS:O	1:A:144:ALA:HA	2.15	0.46
1:A:195:MET:HA	1:A:196:ILE:HA	1.69	0.46
1:A:195:MET:HG3	1:A:233:ILE:HG12	1.98	0.46
1:A:19:PRO:HG3	1:A:299:LEU:HD13	1.98	0.46
2:B:108:VAL:HG13	2:B:110:THR:HG22	1.98	0.46
1:C:125:ASN:OD1	1:C:125:ASN:N	2.49	0.45
1:A:104:ASN:ND2	1:A:130:GLU:OE1	2.37	0.45
1:C:143:GLN:HE22	1:C:174:ASN:HB2	1.81	0.45
2:D:112:LYS:HE2	2:D:189:PHE:CE1	2.51	0.45
1:C:179:VAL:HB	1:C:180:PRO:HD3	1.98	0.45
1:A:297:ASP:OD2	1:A:301:ASN:ND2	2.49	0.45
1:A:51:THR:HB	4:A:402:CL:CL	2.55	0.44
2:B:112:LYS:HG2	2:B:189:PHE:CE1	2.52	0.44
1:A:89:LEU:HD23	1:A:89:LEU:HA	1.81	0.44
1:C:141:GLY:O	1:C:144:ALA:N	2.51	0.44
1:C:76:LEU:HD12	1:C:138:VAL:HG11	2.00	0.44
2:D:156:GLU:OE2	2:D:156:GLU:N	2.47	0.44
1:C:21:ILE:O	1:C:268:ALA:HA	2.17	0.44
2:D:229:GLY:HA3	2:D:250:GLU:OE1	2.18	0.43
1:A:137:ASN:HD21	1:A:139:LYS:NZ	2.15	0.43
1:A:173:PRO:HG2	1:A:199:LEU:HD23	2.01	0.43
2:D:126:TYR:CE1	2:D:156:GLU:HB3	2.53	0.43
2:D:64:ALA:O	2:D:66:THR:HG23	2.19	0.43
1:C:297:ASP:CG	1:C:301:ASN:HD22	2.22	0.43
1:C:71:LEU:HA	1:C:219:GLY:O	2.19	0.43
1:A:137:ASN:OD1	1:A:139:LYS:HG2	2.19	0.43
2:B:25:LEU:O	2:B:66:THR:HA	2.19	0.43
2:D:33:MET:O	2:D:62:LYS:NZ	2.50	0.43
1:C:13:GLY:HA3	1:C:166:PRO:HG2	2.00	0.43
2:B:60:TRP:CZ3	2:B:127:GLU:HG2	2.54	0.42
2:B:120:ILE:HD12	6:B:301:FAD:N1	2.35	0.42
2:B:41:HIS:CD2	2:B:52:ARG:HB2	2.54	0.42
1:C:139:LYS:HE2	8:C:560:HOH:O	2.18	0.42
2:B:6:GLU:OE1	2:B:29:LEU:HG	2.20	0.42
1:A:139:LYS:HE3	8:A:525:HOH:O	2.19	0.42
2:D:234:CYS:O	2:D:248:VAL:HG23	2.20	0.42
1:C:170:LYS:HA	1:C:194:THR:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:31:GLU:HA	2:D:62:LYS:HE3	2.01	0.42
2:D:190:GLU:OE1	8:D:402:HOH:O	2.22	0.41
1:C:296:ILE:HD13	1:C:302:LEU:HD22	2.02	0.41
1:A:170:LYS:HA	1:A:194:THR:O	2.20	0.41
2:B:255:PHE:HB2	2:B:260:LEU:HD11	2.02	0.41
1:C:112:ASP:HB3	8:C:538:HOH:O	2.20	0.41
2:D:100:ASN:O	2:D:254:VAL:HG13	2.20	0.41
2:D:85:LYS:HE2	8:D:437:HOH:O	2.21	0.41
2:D:3:LYS:HE3	2:D:6:GLU:HB2	2.03	0.41
1:C:122:ASP:OD1	1:C:164:LYS:NZ	2.41	0.40
1:C:144:ALA:O	1:C:145:PHE:CG	2.74	0.40
2:D:46:ASN:HB2	2:D:82:GLU:OE2	2.21	0.40
2:B:225:ALA:HB3	2:B:231:CYS:SG	2.62	0.40
2:B:56:SER:HB3	2:B:124:PRO:HB3	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:A:566:HOH:O	8:B:407:HOH:O[4_555]	1.91	0.29

## 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	304/306 (99%)	296 (97%)	7 (2%)	1 (0%)	41 59
1	C	304/306 (99%)	291 (96%)	10 (3%)	3 (1%)	15 26
2	B	260/262 (99%)	253 (97%)	7 (3%)	0	100 100
2	D	260/262 (99%)	252 (97%)	8 (3%)	0	100 100
All	All	1128/1136 (99%)	1092 (97%)	32 (3%)	4 (0%)	34 52

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	142	GLY
1	C	145	PHE
1	C	141	GLY
1	A	145	PHE

### 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	244/246 (99%)	242 (99%)	2 (1%)	81	92
1	C	240/246 (98%)	235 (98%)	5 (2%)	53	76
2	B	213/222 (96%)	209 (98%)	4 (2%)	57	78
2	D	213/222 (96%)	208 (98%)	5 (2%)	50	74
All	All	910/936 (97%)	894 (98%)	16 (2%)	59	80

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	PHE
1	A	310	LYS
2	B	31	GLU
2	B	41	HIS
2	B	180	VAL
2	B	184	MET
2	D	41	HIS
2	D	114	LEU
2	D	161	LYS
2	D	209	LYS
2	D	211	GLU
1	C	27	PHE
1	C	131	LEU
1	C	135	CYS
1	C	140	HIS
1	C	168	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	95	ASN
1	A	137	ASN
1	A	140	HIS
2	B	41	HIS
2	B	87	GLN
2	B	237	HIS
2	D	237	HIS
1	C	77	GLN
1	C	95	ASN
1	C	143	GLN
1	C	301	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 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
6	FAD	D	301	-	51,58,58	1.78	8 (15%)	60,89,89	2.04	10 (16%)
5	EDO	C	402	-	3,3,3	0.47	0	2,2,2	0.18	0
5	EDO	A	403	-	3,3,3	0.43	0	2,2,2	0.28	0
3	FMN	A	401	-	31,33,33	2.05	5 (16%)	40,50,50	2.01	9 (22%)
3	FMN	C	401	-	31,33,33	2.25	6 (19%)	40,50,50	2.65	11 (27%)
7	FES	B	302	2	0,4,4	0.00	-	-	-	-
7	FES	D	302	2	0,4,4	0.00	-	-	-	-
6	FAD	B	301	-	51,58,58	1.85	6 (11%)	60,89,89	2.00	10 (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
6	FAD	D	301	-	-	0/30/50/50	0/6/6/6
5	EDO	C	402	-	-	0/1/1/1	-
5	EDO	A	403	-	-	0/1/1/1	-
3	FMN	A	401	-	-	5/18/18/18	0/3/3/3
3	FMN	C	401	-	-	6/18/18/18	0/3/3/3
7	FES	B	302	2	-	-	0/1/1/1
7	FES	D	302	2	-	-	0/1/1/1
6	FAD	B	301	-	-	2/30/50/50	0/6/6/6

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	401	FMN	C4A-C10	9.32	1.48	1.38
6	B	301	FAD	C4X-C10	9.07	1.47	1.38
6	D	301	FAD	C4X-C10	8.78	1.47	1.38
3	A	401	FMN	C4A-C10	8.60	1.47	1.38
6	B	301	FAD	C4-C4X	4.12	1.48	1.41
3	C	401	FMN	C9A-C5A	3.86	1.50	1.42
6	B	301	FAD	C8-C7	3.75	1.50	1.40
6	D	301	FAD	C4-C4X	3.71	1.47	1.41
6	D	301	FAD	C9A-C5X	3.70	1.50	1.42
6	B	301	FAD	C9A-C5X	3.66	1.49	1.42
3	C	401	FMN	C4-C4A	3.59	1.47	1.41
3	A	401	FMN	C4-C4A	3.38	1.47	1.41
6	D	301	FAD	C8-C7	3.32	1.49	1.40
3	A	401	FMN	C8-C7	3.26	1.49	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	401	FMN	C8-C7	3.22	1.48	1.40
3	A	401	FMN	C9A-C5A	3.16	1.48	1.42
6	B	301	FAD	C5A-C4A	2.69	1.48	1.40
6	D	301	FAD	C5A-C4A	2.61	1.47	1.40
3	A	401	FMN	C9A-N10	2.34	1.41	1.38
3	C	401	FMN	C9A-N10	2.21	1.41	1.38
3	C	401	FMN	C1'-N10	-2.21	1.46	1.48
6	D	301	FAD	C2A-N3A	2.19	1.35	1.32
6	B	301	FAD	C2A-N3A	2.05	1.35	1.32
6	D	301	FAD	C2B-C1B	-2.05	1.50	1.53
6	D	301	FAD	C9A-N10	2.04	1.41	1.38

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	301	FAD	C4-N3-C2	8.40	122.24	115.14
3	C	401	FMN	C4-N3-C2	8.26	122.12	115.14
6	D	301	FAD	C4-N3-C2	7.97	121.88	115.14
3	A	401	FMN	C4-N3-C2	7.52	121.49	115.14
6	D	301	FAD	C1'-N10-C9A	6.73	123.59	118.29
6	B	301	FAD	C4-C4X-C10	-6.68	115.53	119.95
3	C	401	FMN	C5'-C4'-C3'	-6.51	99.63	112.20
6	B	301	FAD	C1'-N10-C9A	5.86	122.91	118.29
3	C	401	FMN	C1'-N10-C9A	5.74	122.81	118.29
3	C	401	FMN	O2P-P-O5'	-5.28	92.67	106.73
6	D	301	FAD	C4-C4X-C10	-5.20	116.51	119.95
3	A	401	FMN	C4-C4A-C10	-4.67	116.86	119.95
3	C	401	FMN	C4-C4A-C10	-4.46	117.00	119.95
6	B	301	FAD	C4-C4X-N5	4.28	123.49	118.60
3	A	401	FMN	C4A-N5-C5A	3.99	120.76	116.77
3	C	401	FMN	C4A-C4-N3	-3.86	118.15	123.43
6	D	301	FAD	C5X-C9A-N10	3.81	120.48	117.72
6	D	301	FAD	C4X-N5-C5X	3.77	120.54	116.77
6	B	301	FAD	C4X-C4-N3	-3.44	118.73	123.43
6	D	301	FAD	C9A-N10-C10	-3.38	117.49	121.91
3	C	401	FMN	C4A-N5-C5A	3.35	120.12	116.77
6	D	301	FAD	C4X-C4-N3	-3.33	118.88	123.43
3	C	401	FMN	P-O5'-C5'	3.24	127.22	118.30
6	B	301	FAD	C4X-N5-C5X	3.23	120.00	116.77
3	A	401	FMN	C1'-N10-C10	3.20	121.28	118.41
3	C	401	FMN	C5A-C9A-N10	2.95	119.85	117.72
6	B	301	FAD	C9A-N10-C10	-2.75	118.31	121.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	FMN	C9A-N10-C10	-2.71	118.36	121.91
3	A	401	FMN	C4A-C4-N3	-2.65	119.81	123.43
6	B	301	FAD	N3A-C2A-N1A	-2.60	124.61	128.68
6	B	301	FAD	C4A-C5A-N7A	-2.53	106.76	109.40
6	D	301	FAD	N3A-C2A-N1A	-2.49	124.78	128.68
3	C	401	FMN	C9A-N10-C10	-2.49	118.64	121.91
3	A	401	FMN	C1'-N10-C9A	2.49	120.25	118.29
6	D	301	FAD	C4-C4X-N5	2.46	121.41	118.60
6	D	301	FAD	O4B-C1B-C2B	-2.23	103.66	106.93
3	C	401	FMN	O3P-P-O2P	2.06	115.52	107.64
3	A	401	FMN	C5A-C9A-N10	2.06	119.21	117.72
3	A	401	FMN	C4-C4A-N5	2.05	120.94	118.60
6	B	301	FAD	C5X-C9A-N10	2.04	119.19	117.72

There are no chirality outliers.

All (13) torsion outliers are listed below:

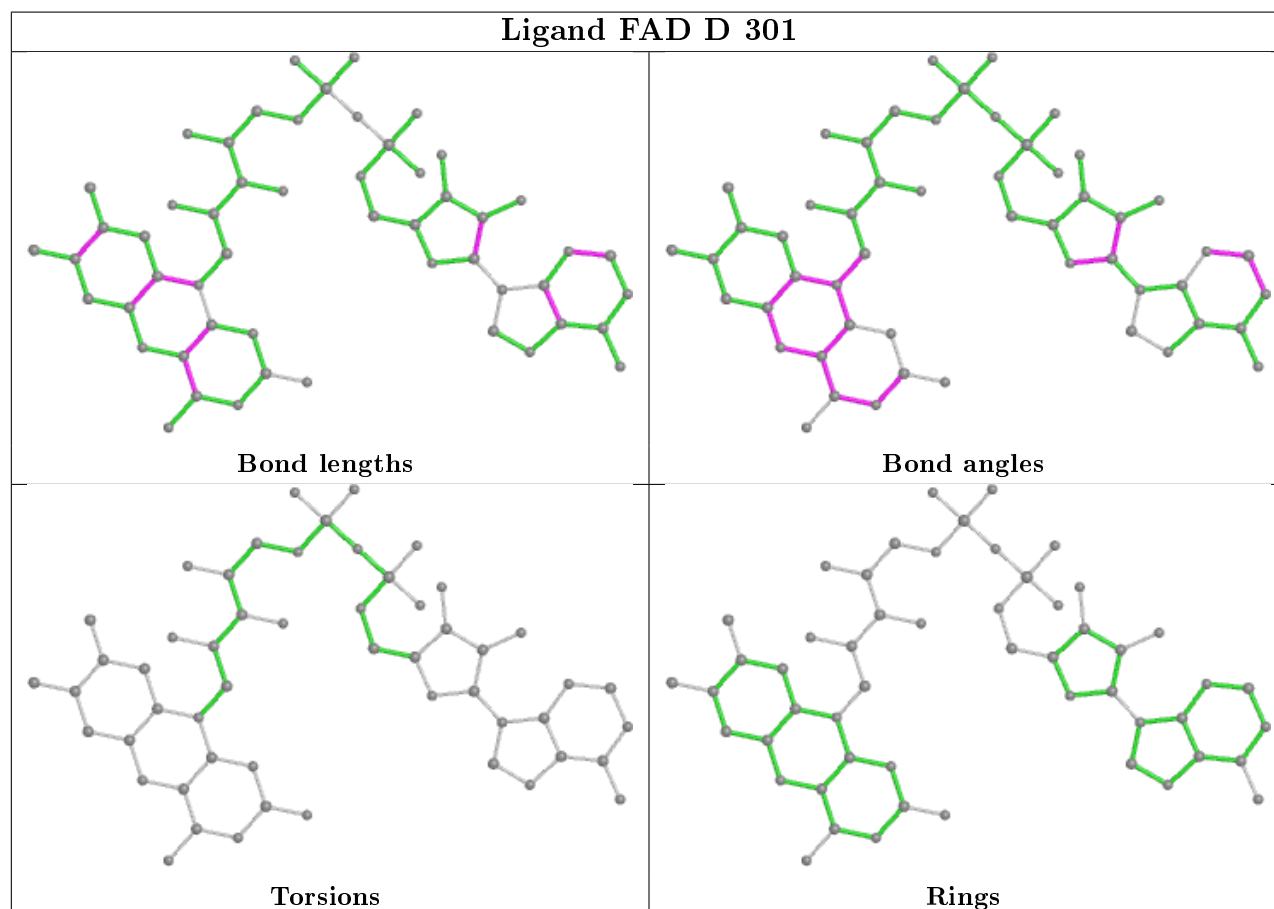
Mol	Chain	Res	Type	Atoms
3	C	401	FMN	C2'-C3'-C4'-O4'
3	C	401	FMN	C2'-C3'-C4'-C5'
3	C	401	FMN	O3'-C3'-C4'-O4'
3	C	401	FMN	O3'-C3'-C4'-C5'
6	B	301	FAD	C5B-O5B-PA-O3P
3	A	401	FMN	O3'-C3'-C4'-C5'
3	A	401	FMN	C2'-C3'-C4'-C5'
3	A	401	FMN	C2'-C3'-C4'-O4'
3	A	401	FMN	O3'-C3'-C4'-O4'
3	A	401	FMN	C4'-C5'-O5'-P
3	C	401	FMN	C4'-C5'-O5'-P
6	B	301	FAD	C5B-O5B-PA-O2A
3	C	401	FMN	O4'-C4'-C5'-O5'

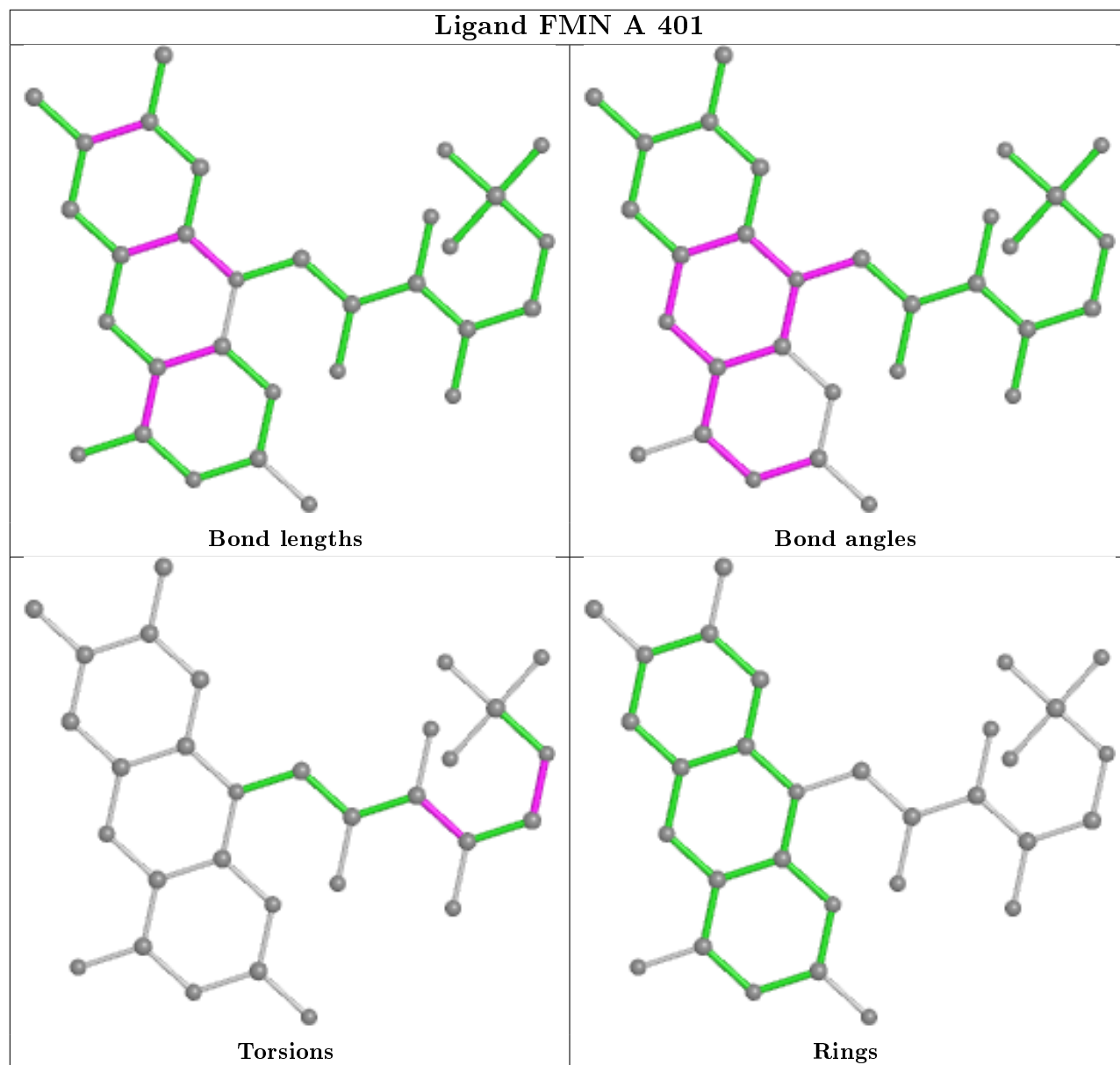
There are no ring outliers.

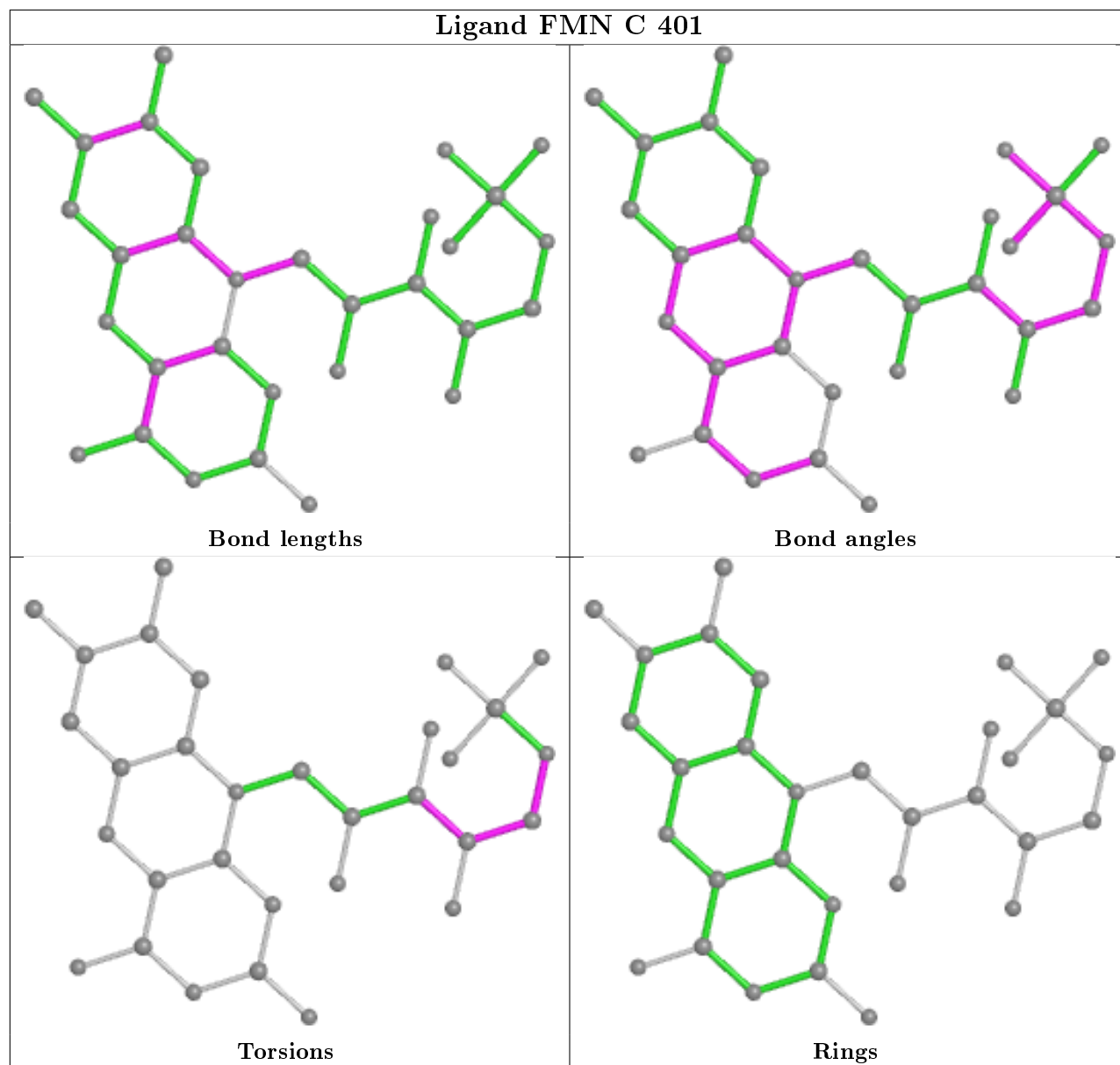
4 monomers are involved in 4 short contacts:

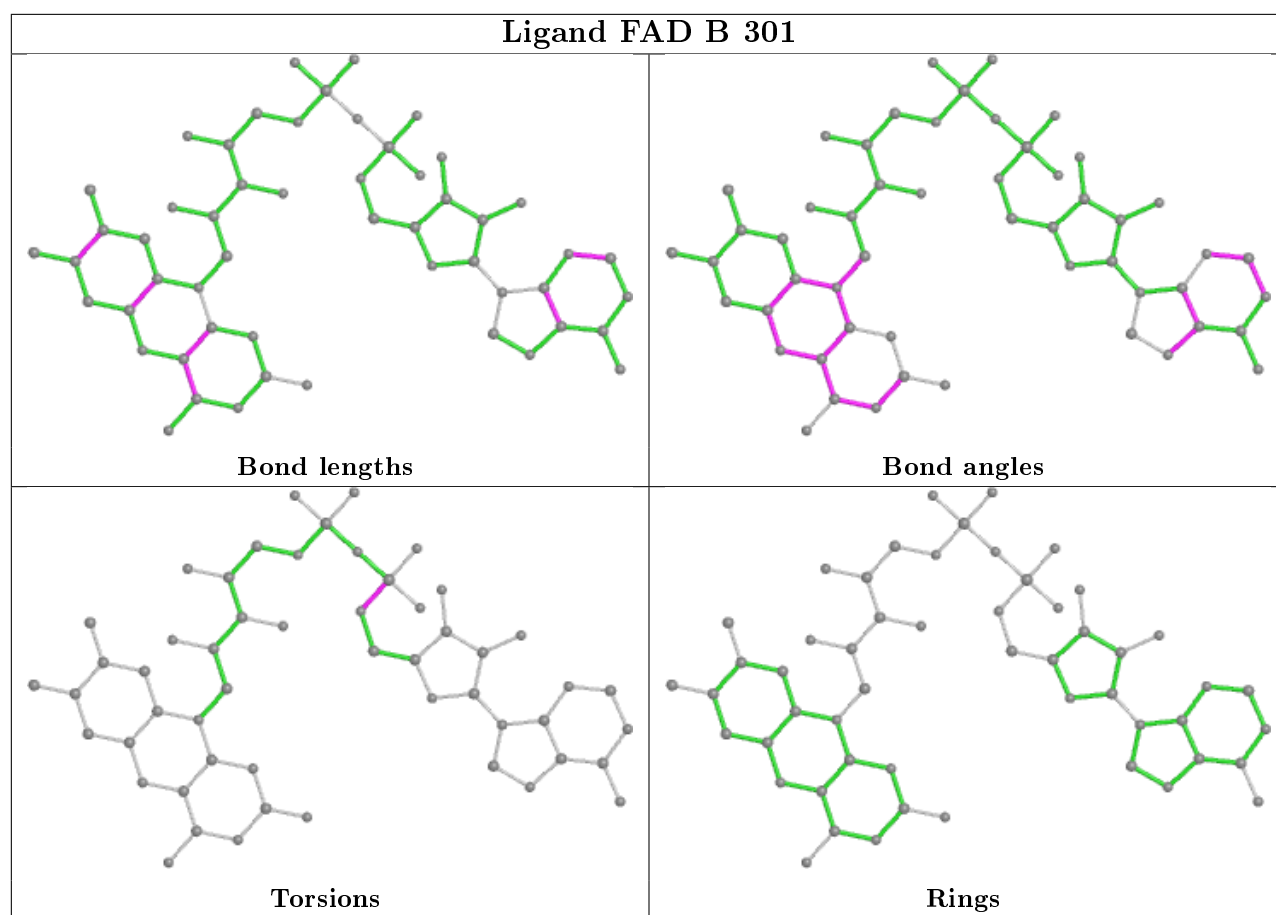
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	301	FAD	1	0
3	A	401	FMN	1	0
3	C	401	FMN	1	0
6	B	301	FAD	1	0

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	306/306 (100%)	-0.13	7 (2%) 60 62	26, 41, 64, 105	0
1	C	305/306 (99%)	0.28	18 (5%) 22 22	28, 50, 70, 90	0
2	B	262/262 (100%)	0.07	9 (3%) 45 47	32, 50, 76, 91	1 (0%)
2	D	262/262 (100%)	-0.22	1 (0%) 92 93	30, 47, 66, 82	2 (0%)
All	All	1135/1136 (99%)	0.00	35 (3%) 49 51	26, 47, 71, 105	3 (0%)

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	140	HIS	5.5
1	A	141	GLY	4.3
1	C	224	ALA	3.9
2	B	188	ASP	3.7
1	A	140	HIS	3.7
1	A	143	GLN	3.7
1	A	137	ASN	3.5
1	A	139	LYS	3.5
1	A	142	GLY	3.0
2	B	187	ILE	2.8
2	B	108	VAL	2.7
1	C	199	LEU	2.7
1	C	68	SER	2.7
1	C	139	LYS	2.7
1	C	223	PRO	2.7
2	B	207	ALA	2.6
2	B	63	VAL	2.6
1	C	220	LEU	2.6
1	A	138	VAL	2.5
1	C	71	LEU	2.5
1	C	225	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	190	GLU	2.4
2	B	243	ASN	2.4
2	B	2	PRO	2.4
1	C	144	ALA	2.3
1	C	298	SER	2.3
1	C	82	GLU	2.3
1	C	308	ASN	2.2
1	C	66	THR	2.2
1	C	249	GLY	2.2
2	B	208	LYS	2.1
1	C	221	SER	2.1
1	C	112	ASP	2.1
2	D	63	VAL	2.1
1	C	200	MET	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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.

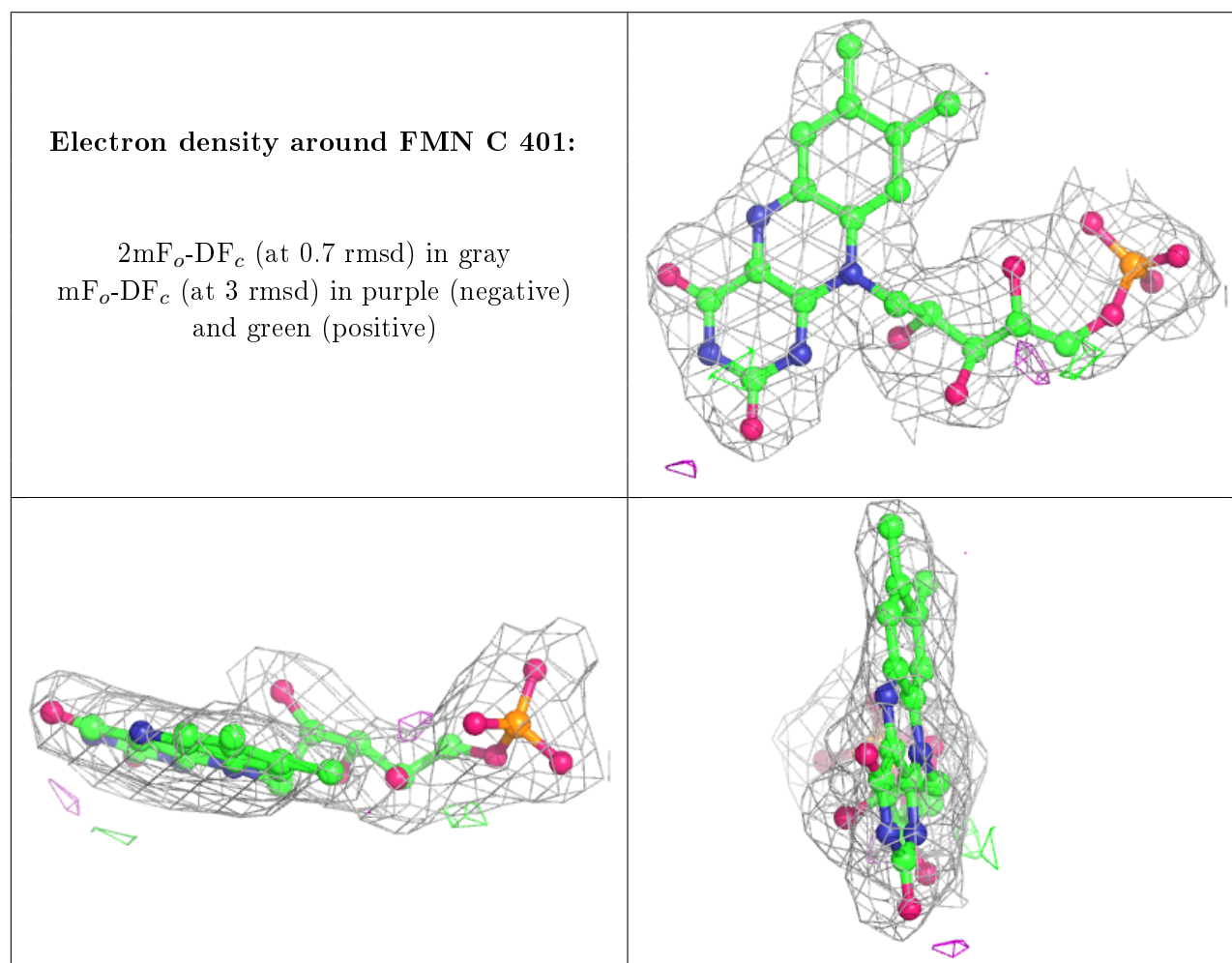
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	CL	B	303	1/1	0.81	0.21	85,85,85,85	0
3	FMN	C	401	31/31	0.96	0.24	12,37,47,51	0
6	FAD	D	301	53/53	0.97	0.14	20,38,50,53	0
4	CL	A	402	1/1	0.97	0.06	65,65,65,65	0
5	EDO	C	402	4/4	0.97	0.22	25,33,48,57	0
3	FMN	A	401	31/31	0.98	0.20	17,31,44,48	0
5	EDO	A	403	4/4	0.98	0.25	30,33,37,48	0
6	FAD	B	301	53/53	0.98	0.15	28,43,53,65	0
7	FES	B	302	4/4	0.99	0.14	38,39,40,41	0

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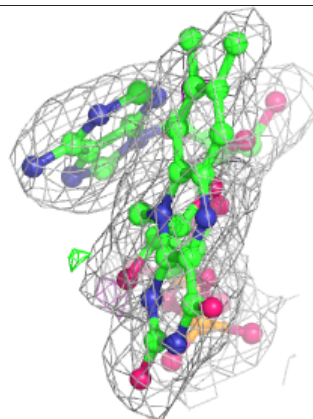
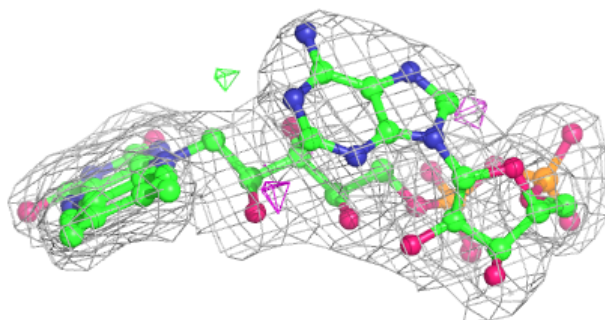
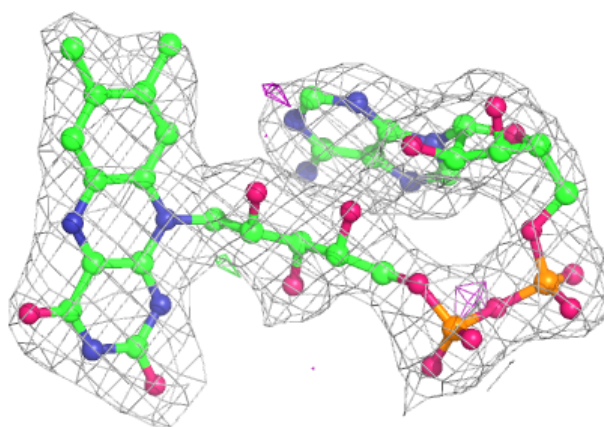
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
7	FES	D	302	4/4	1.00	0.14	32,34,36,37	0

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.



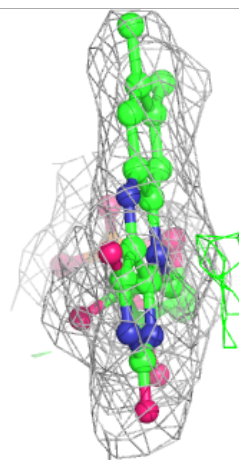
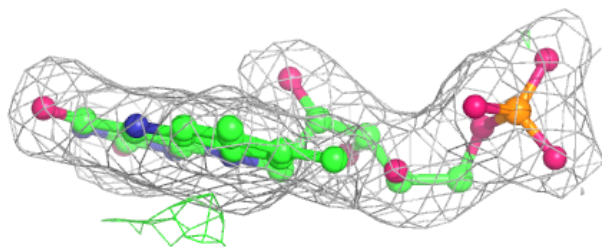
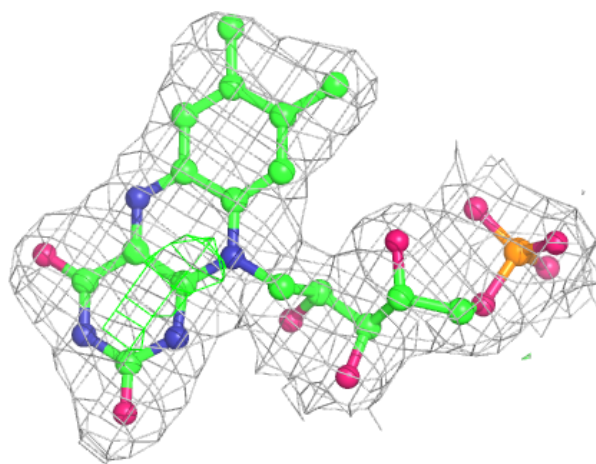
**Electron density around FAD D 301:**

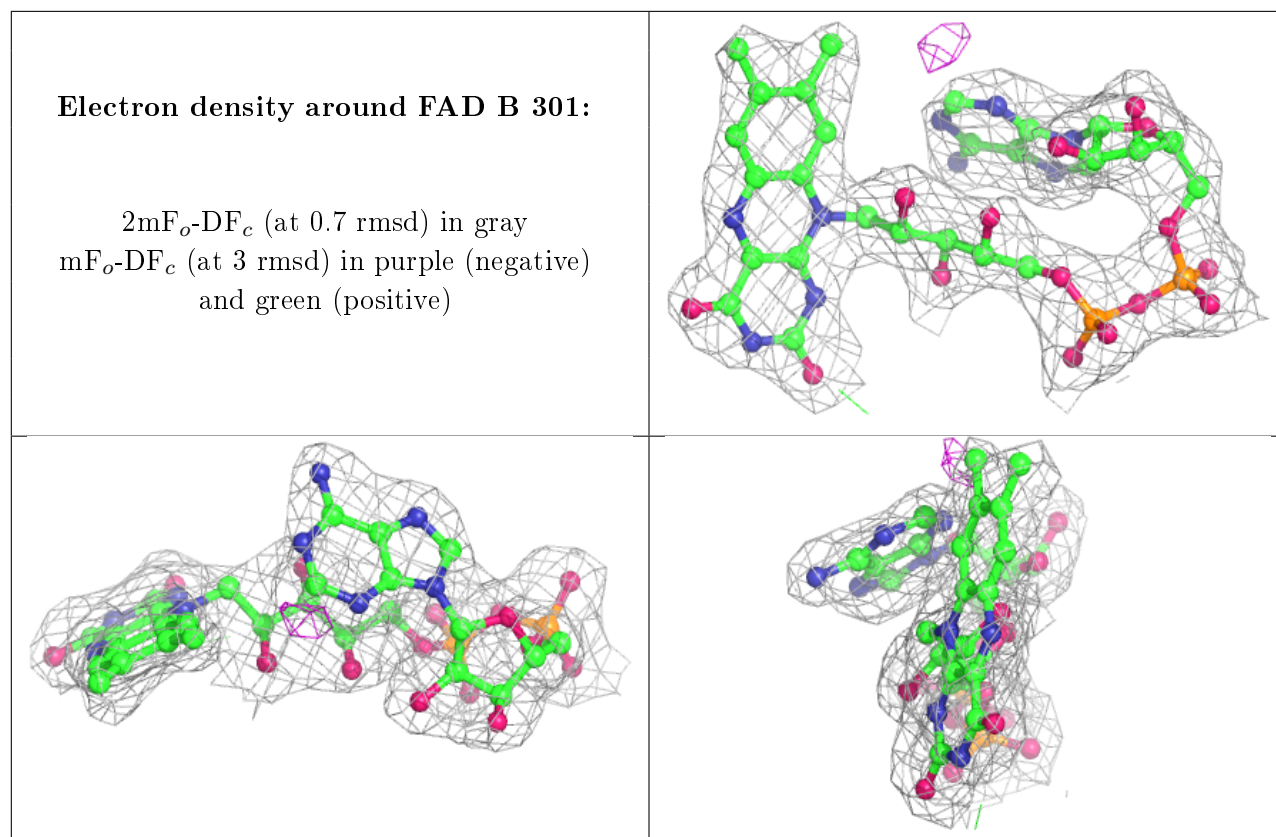
$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around FMN A 401:**

$2mF_o-DF_c$  (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.