



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 12:00 pm BST

PDB ID : 2XVI
Title : Crystal structure of the mutant bacterial flavin containing monooxygenase (Y207S)
Authors : Cho, H.J.; Kang, B.S.
Deposited on : 2010-10-26
Resolution : 2.48 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The 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

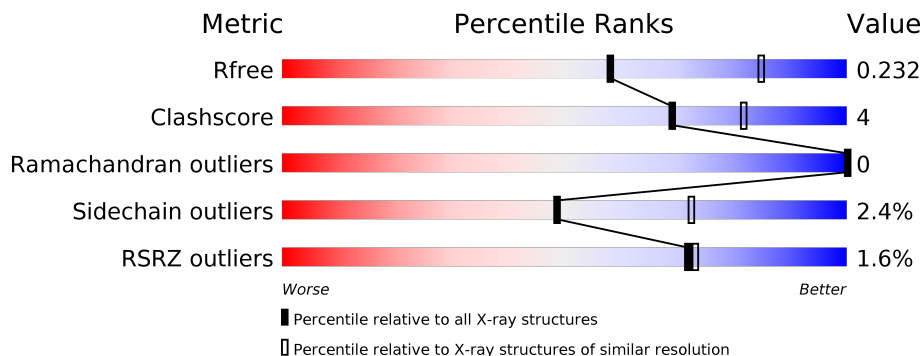
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.48 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 ≥ 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	464	
1	B	464	
1	C	464	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	OXY	C	1450	-	-	-	X

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 11137 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-CONTAINING MONOOXYGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	445	Total 3594	C 2305	N 591	O 678	S 20	0	0	0
1	B	446	Total 3636	C 2333	N 603	O 680	S 20	0	1	0
1	C	446	Total 3593	C 2305	N 594	O 674	S 20	0	0	0

There are 27 discrepancies between the modelled and reference sequences:

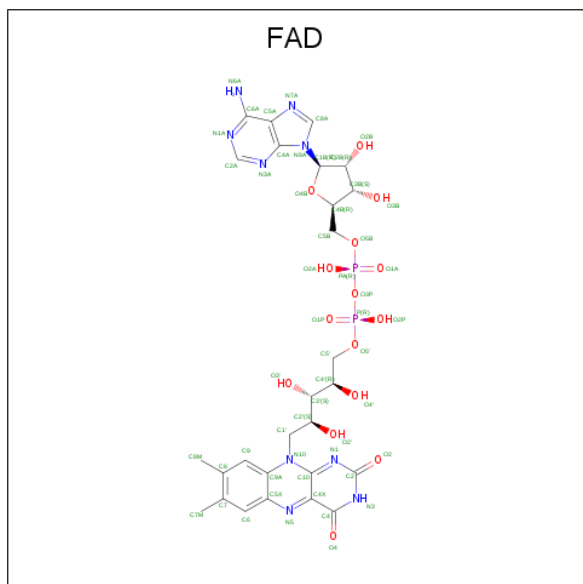
Chain	Residue	Modelled	Actual	Comment	Reference
A	457	LEU	-	expression tag	UNP Q83XK4
A	458	GLU	-	expression tag	UNP Q83XK4
A	459	HIS	-	expression tag	UNP Q83XK4
A	460	HIS	-	expression tag	UNP Q83XK4
A	461	HIS	-	expression tag	UNP Q83XK4
A	462	HIS	-	expression tag	UNP Q83XK4
A	463	HIS	-	expression tag	UNP Q83XK4
A	464	HIS	-	expression tag	UNP Q83XK4
A	207	SER	TYR	engineered mutation	UNP Q83XK4
B	457	LEU	-	expression tag	UNP Q83XK4
B	458	GLU	-	expression tag	UNP Q83XK4
B	459	HIS	-	expression tag	UNP Q83XK4
B	460	HIS	-	expression tag	UNP Q83XK4
B	461	HIS	-	expression tag	UNP Q83XK4
B	462	HIS	-	expression tag	UNP Q83XK4
B	463	HIS	-	expression tag	UNP Q83XK4
B	464	HIS	-	expression tag	UNP Q83XK4
B	207	SER	TYR	engineered mutation	UNP Q83XK4
C	457	LEU	-	expression tag	UNP Q83XK4
C	458	GLU	-	expression tag	UNP Q83XK4
C	459	HIS	-	expression tag	UNP Q83XK4
C	460	HIS	-	expression tag	UNP Q83XK4
C	461	HIS	-	expression tag	UNP Q83XK4

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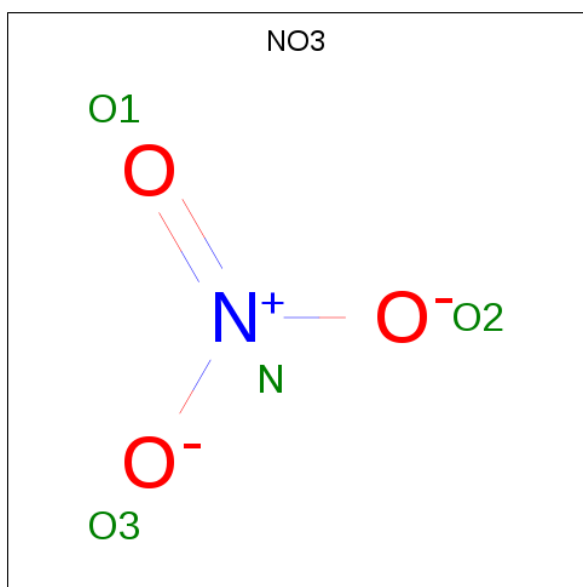
Chain	Residue	Modelled	Actual	Comment	Reference
C	462	HIS	-	expression tag	UNP Q83XK4
C	463	HIS	-	expression tag	UNP Q83XK4
C	464	HIS	-	expression tag	UNP Q83XK4
C	207	SER	TYR	engineered mutation	UNP Q83XK4

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



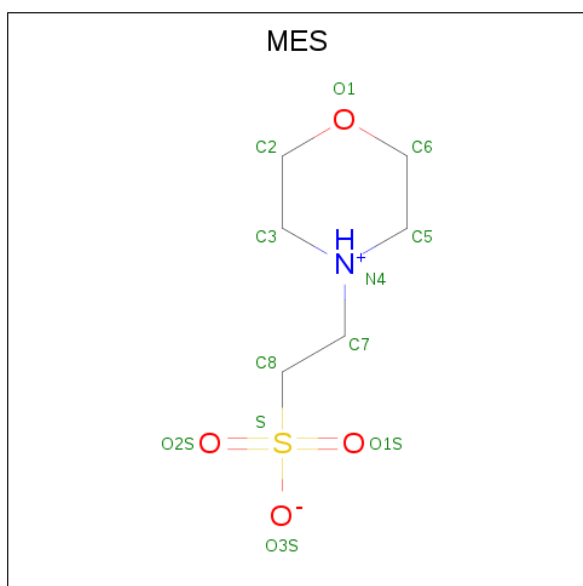
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	N	O			P
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is NITRATE ION (three-letter code: NO3) (formula: NO_3).



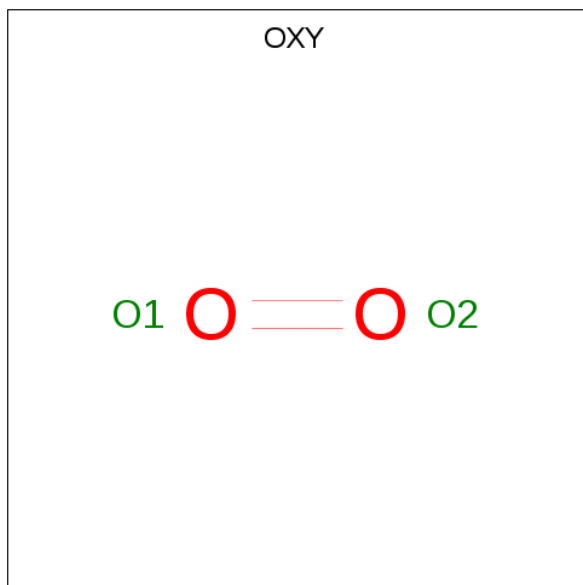
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	N O	0	0
			4	1 3		
3	B	1	Total	N O	0	0
			4	1 3		
3	C	1	Total	N O	0	0
			4	1 3		

- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
4	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 5 is OXYGEN MOLECULE (three-letter code: OXY) (formula: O₂).



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

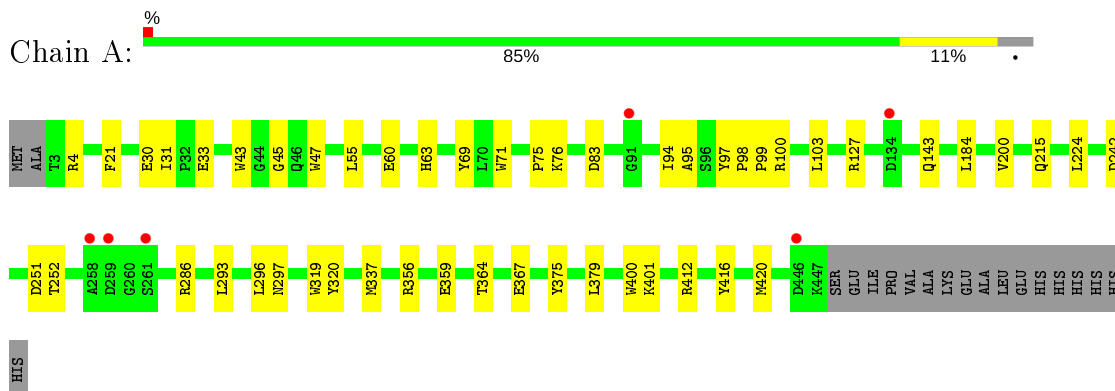
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	49	Total	O	0	0
			49	49		
6	B	44	Total	O	0	0
			44	44		
6	C	20	Total	O	0	0
			20	20		

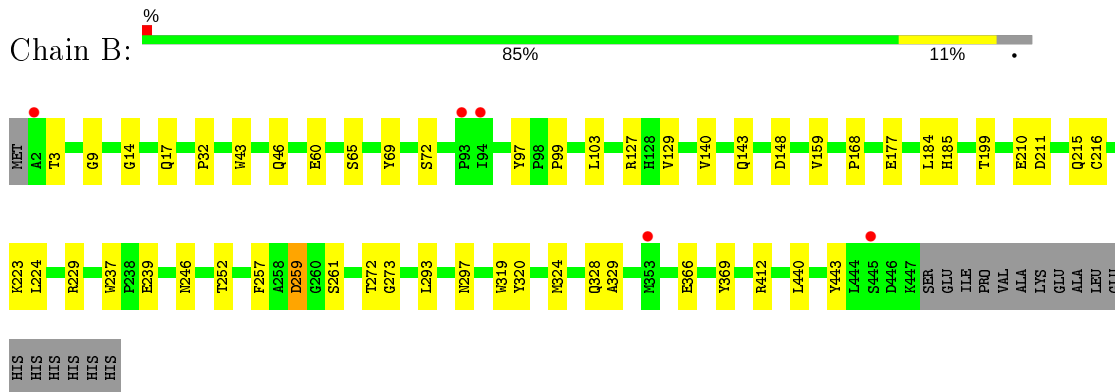
3 Residue-property plots [i](#)

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.

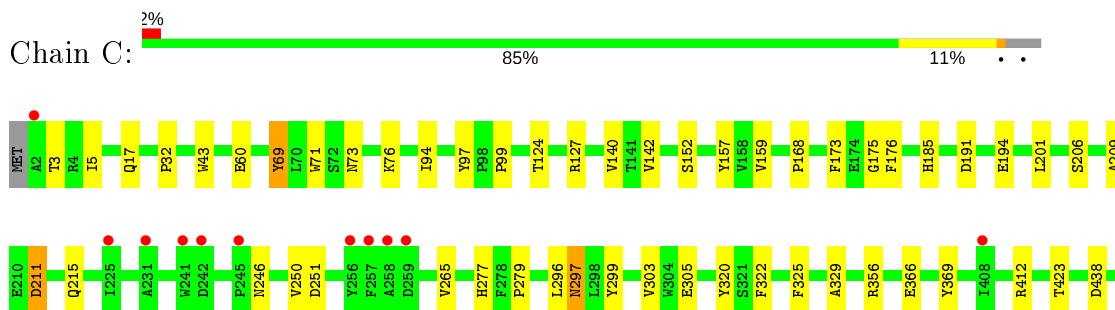
- Molecule 1: FLAVIN-CONTAINING MONOOXYGENASE



- Molecule 1: FLAVIN-CONTAINING MONOOXYGENASE



- Molecule 1: FLAVIN-CONTAINING MONOOXYGENASE



4437
SER
GLU
ILE
PRO
VAL
ALA
LYS
GLU
ALA
LEU
GLU
HIS
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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	159.57Å 70.01Å 140.69Å 90.00° 90.19° 90.00°	Depositor
Resolution (Å)	50.01 – 2.48 40.49 – 2.48	Depositor EDS
% Data completeness (in resolution range)	96.9 (50.01-2.48) 83.9 (40.49-2.48)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	26.75 (at 2.48Å)	Xtrriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.201 , 0.240 0.196 , 0.232	Depositor DCC
R_{free} test set	2344 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	59.4	Xtrriage
Anisotropy	0.049	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.018 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11137	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.08% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: NO3, FAD, MES, OXY

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	0.50	1/3707 (0.0%)	0.51	0/5046
1	B	0.48	1/3752 (0.0%)	0.52	0/5099
1	C	0.47	2/3706 (0.1%)	0.51	0/5047
All	All	0.48	4/11165 (0.0%)	0.52	0/15192

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	320	TYR	CD1-CE1	-6.48	1.29	1.39
1	C	320	TYR	CD1-CE1	-5.50	1.31	1.39
1	C	320	TYR	CD2-CE2	-5.42	1.31	1.39
1	B	320	TYR	CD1-CE1	-5.41	1.31	1.39

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	3594	0	3306	23	0
1	B	3636	0	3397	30	0
1	C	3593	0	3308	28	0
2	A	53	0	31	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	53	0	31	1	0
2	C	53	0	31	0	0
3	A	4	0	0	0	0
3	B	4	0	0	1	0
3	C	4	0	0	0	0
4	A	12	0	12	0	0
4	B	12	0	12	0	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
5	C	2	0	0	0	0
6	A	49	0	0	0	0
6	B	44	0	0	0	0
6	C	20	0	0	0	0
All	All	11137	0	10128	79	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 4.

All (79) 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:229:ARG:HG2	1:B:246:ASN:HB3	1.65	0.77
1:B:46:GLN:HE22	1:B:65:SER:H	1.34	0.73
1:B:293:LEU:HD21	1:B:319:TRP:HD1	1.58	0.69
1:A:76:LYS:HE2	1:A:94:ILE:HG13	1.76	0.68
1:C:60:GLU:OE2	1:C:127:ARG:NH2	2.25	0.66
1:C:76:LYS:HE2	1:C:94:ILE:HG13	1.82	0.62
1:A:286:ARG:NH2	1:B:148:ASP:OD1	2.31	0.62
1:B:293:LEU:HD21	1:B:319:TRP:CD1	2.34	0.61
1:B:3:THR:HB	1:B:32:PRO:HB3	1.83	0.60
1:C:71:TRP:HA	1:C:99:PRO:HA	1.83	0.60
1:A:45:GLY:HA2	2:A:1448:FAD:O3B	2.02	0.60
1:A:98:PRO:HG2	1:A:103:LEU:HG	1.84	0.59
1:A:76:LYS:HE3	1:A:95:ALA:O	2.01	0.59
1:B:60:GLU:OE2	1:B:127:ARG:NH2	2.36	0.59
1:B:46:GLN:NE2	1:B:65:SER:H	2.00	0.58
1:A:71:TRP:HA	1:A:99:PRO:HA	1.87	0.55
1:C:299:TYR:HB3	1:C:303:VAL:HB	1.88	0.55
1:B:210:GLU:HB2	1:B:237:TRP:HH2	1.72	0.55
1:A:375:TYR:CZ	1:A:379:LEU:HD11	2.41	0.55
1:A:127:ARG:HH21	1:A:143:GLN:NE2	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:ARG:HH21	1:A:143:GLN:HE21	1.56	0.53
1:B:199:THR:HG23	1:B:223:LYS:HB3	1.91	0.53
1:C:251:ASP:OD1	1:C:251:ASP:N	2.42	0.52
1:C:124:THR:HG23	1:C:142:VAL:HG13	1.93	0.51
1:C:69:TYR:HA	1:C:438:ASP:HB3	1.93	0.51
1:C:201:LEU:HB2	1:C:265:VAL:HG21	1.92	0.50
1:B:17:GLN:HA	1:B:329:ALA:HB1	1.94	0.50
1:B:273:GLY:N	3:B:1449:NO3:O3	2.39	0.50
1:B:366:GLU:HA	1:B:369:TYR:CZ	2.47	0.50
1:A:293:LEU:HD21	1:A:319:TRP:CD1	2.47	0.49
1:B:324:MET:O	1:B:328:GLN:HG3	2.13	0.49
1:C:3:THR:HB	1:C:32:PRO:HB3	1.95	0.49
1:A:21:PHE:O	1:A:31:ILE:HG21	2.13	0.48
1:A:200:VAL:HB	1:A:224:LEU:HD23	1.95	0.48
1:A:364:THR:HG23	1:A:367:GLU:H	1.79	0.48
1:C:191:ASP:O	1:C:194:GLU:HG2	2.14	0.48
1:B:97:TYR:CD1	1:B:412:ARG:HA	2.49	0.47
1:B:257:PHE:C	1:B:259:ASP:H	2.17	0.47
1:A:4:ARG:HE	1:A:33:GLU:CD	2.18	0.47
1:A:47:TRP:O	1:A:100:ARG:NH1	2.48	0.47
1:B:72:SER:HB2	1:B:103:LEU:HD11	1.97	0.47
1:C:5:ILE:HD13	1:C:32:PRO:HB2	1.97	0.46
1:C:5:ILE:HG13	1:C:157:TYR:HB2	1.97	0.46
1:A:63:HIS:CD2	2:A:1448:FAD:H3B	2.51	0.46
1:B:127:ARG:NH2	1:B:143:GLN:OE1	2.46	0.46
1:C:97:TYR:CD1	1:C:412:ARG:HG2	2.50	0.46
1:B:257:PHE:C	1:B:259:ASP:N	2.69	0.46
1:B:259:ASP:HB3	1:B:261:SER:H	1.81	0.46
1:C:277:HIS:CD2	1:C:279:PRO:HD3	2.52	0.45
1:B:211:ASP:O	1:B:215:GLN:HG2	2.17	0.45
1:A:60:GLU:OE2	1:A:127:ARG:NH2	2.40	0.44
1:C:124:THR:CG2	1:C:142:VAL:HG13	2.47	0.44
1:C:168:PRO:HB3	1:C:185:HIS:CE1	2.52	0.44
1:C:296:LEU:HG	1:C:356:ARG:NH2	2.33	0.44
1:B:9:GLY:O	1:B:14:GLY:HA3	2.18	0.44
1:C:17:GLN:HE21	1:C:159:VAL:HG11	1.83	0.44
1:C:140:VAL:O	1:C:152:SER:HA	2.18	0.43
1:A:400:TRP:CZ3	1:A:401:LYS:HE2	2.54	0.43
1:C:211:ASP:O	1:C:215:GLN:HG2	2.19	0.43
1:A:296:LEU:HG	1:A:356:ARG:NH2	2.34	0.43
1:B:177:GLU:H	1:B:177:GLU:CD	2.22	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:366:GLU:HA	1:C:369:TYR:CZ	2.54	0.43
1:C:322:PHE:HA	1:C:325:PHE:CD2	2.54	0.43
1:C:173:PHE:HB2	1:C:176:PHE:CE1	2.54	0.42
1:B:129:VAL:HG22	1:B:140:VAL:HG13	2.01	0.42
1:B:216:CYS:HB2	1:B:224:LEU:HD21	2.02	0.42
1:A:55:LEU:HD12	1:B:168:PRO:CD	2.49	0.42
1:C:206:SER:H	1:C:209:ALA:HB3	1.85	0.42
1:A:97:TYR:CD2	1:A:412:ARG:HG2	2.55	0.41
1:B:99:PRO:HD3	1:B:443:TYR:CD1	2.55	0.41
1:C:175:GLY:HA3	1:C:250:VAL:O	2.21	0.41
1:C:17:GLN:HA	1:C:329:ALA:HB1	2.02	0.41
1:B:17:GLN:NE2	1:B:159:VAL:HG11	2.35	0.41
1:A:251:ASP:OD1	1:A:252:THR:N	2.54	0.41
1:B:168:PRO:HB3	1:B:185:HIS:CE1	2.55	0.41
1:C:124:THR:CG2	1:C:142:VAL:CG1	2.99	0.41
1:A:75:PRO:HB3	1:A:416:TYR:CD1	2.56	0.40
1:B:65:SER:HB2	2:B:1448:FAD:HM82	2.03	0.40
1:C:297:ASN:HA	1:C:305:GLU:OE1	2.21	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	443/464 (96%)	425 (96%)	18 (4%)	0	100	100
1	B	445/464 (96%)	425 (96%)	20 (4%)	0	100	100
1	C	444/464 (96%)	428 (96%)	16 (4%)	0	100	100
All	All	1332/1392 (96%)	1278 (96%)	54 (4%)	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	372/405 (92%)	361 (97%)	11 (3%)	41	65
1	B	380/405 (94%)	371 (98%)	9 (2%)	49	72
1	C	370/405 (91%)	363 (98%)	7 (2%)	57	78
All	All	1122/1215 (92%)	1095 (98%)	27 (2%)	49	72

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	GLU
1	A	43	TRP
1	A	69	TYR
1	A	83	ASP
1	A	184	LEU
1	A	215	GLN
1	A	242	ASP
1	A	297	ASN
1	A	337	MET
1	A	359	GLU
1	A	420	MET
1	B	43	TRP
1	B	69	TYR
1	B	184	LEU
1	B	239	GLU
1	B	252	THR
1	B	259	ASP
1	B	272	THR
1	B	297	ASN
1	B	440	LEU
1	C	43	TRP
1	C	69	TYR
1	C	73	ASN
1	C	211	ASP
1	C	246	ASN
1	C	297	ASN

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Mol	Chain	Res	Type
1	C	423	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	GLN
1	A	48	ASN
1	A	128	HIS
1	A	143	GLN
1	A	215	GLN
1	A	282	ASN
1	A	297	ASN
1	B	17	GLN
1	B	40	GLN
1	B	46	GLN
1	B	48	ASN
1	B	132	ASN
1	B	428	HIS
1	C	17	GLN
1	C	22	GLN
1	C	25	GLN
1	C	40	GLN
1	C	48	ASN
1	C	128	HIS
1	C	164	HIS
1	C	277	HIS
1	C	428	HIS
1	C	429	HIS

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

11 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
4	MES	B	1450	-	12,12,12	2.16	1 (8%)	14,16,16	2.39	6 (42%)
5	OXY	B	1451	-	1,1,1	0.14	0	-		
5	OXY	C	1450	-	1,1,1	0.14	0	-		
2	FAD	B	1448	-	51,58,58	1.30	6 (11%)	60,89,89	1.70	9 (15%)
4	MES	A	1450	-	12,12,12	2.15	1 (8%)	14,16,16	2.48	6 (42%)
2	FAD	A	1448	-	51,58,58	1.46	6 (11%)	60,89,89	1.59	8 (13%)
2	FAD	C	1448	-	51,58,58	1.36	7 (13%)	60,89,89	1.64	8 (13%)
3	NO3	B	1449	-	1,3,3	3.38	1 (100%)	0,3,3	0.00	-
3	NO3	C	1449	-	1,3,3	3.49	1 (100%)	0,3,3	0.00	-
3	NO3	A	1449	-	1,3,3	3.45	1 (100%)	0,3,3	0.00	-
5	OXY	A	1451	-	1,1,1	0.12	0	-		

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
4	MES	B	1450	-	-	4/6/14/14	0/1/1/1
2	FAD	A	1448	-	-	4/30/50/50	0/6/6/6
2	FAD	C	1448	-	-	8/30/50/50	0/6/6/6
2	FAD	B	1448	-	-	3/30/50/50	0/6/6/6
4	MES	A	1450	-	-	4/6/14/14	0/1/1/1

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1450	MES	C8-S	-7.19	1.67	1.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1450	MES	C8-S	-7.17	1.67	1.77
2	A	1448	FAD	C10-N1	4.50	1.39	1.33
2	C	1448	FAD	C10-N1	4.21	1.38	1.33
2	B	1448	FAD	C10-N1	4.18	1.38	1.33
2	A	1448	FAD	C2A-N3A	4.14	1.38	1.32
2	A	1448	FAD	C4X-N5	3.92	1.38	1.33
2	C	1448	FAD	C2A-N3A	3.91	1.38	1.32
2	C	1448	FAD	C4X-N5	3.64	1.38	1.33
2	B	1448	FAD	C4X-N5	3.62	1.38	1.33
2	B	1448	FAD	C2A-N3A	3.52	1.37	1.32
3	C	1449	NO3	O1-N	3.49	1.40	1.24
3	A	1449	NO3	O1-N	3.45	1.40	1.24
3	B	1449	NO3	O1-N	3.38	1.39	1.24
2	A	1448	FAD	C4-N3	3.16	1.38	1.33
2	A	1448	FAD	C1'-N10	3.06	1.51	1.48
2	B	1448	FAD	C4-N3	2.97	1.38	1.33
2	C	1448	FAD	C4-N3	2.82	1.37	1.33
2	C	1448	FAD	C1'-N10	2.55	1.50	1.48
2	C	1448	FAD	C2A-N1A	2.47	1.38	1.33
2	A	1448	FAD	C2A-N1A	2.42	1.38	1.33
2	B	1448	FAD	C1'-N10	2.10	1.50	1.48
2	C	1448	FAD	C5X-N5	2.02	1.38	1.35
2	B	1448	FAD	C2A-N1A	2.01	1.37	1.33

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1448	FAD	C4-N3-C2	5.93	120.15	115.14
2	B	1448	FAD	N3A-C2A-N1A	-5.92	119.42	128.68
2	A	1448	FAD	N3A-C2A-N1A	-5.73	119.73	128.68
2	C	1448	FAD	N3A-C2A-N1A	-5.64	119.87	128.68
4	B	1450	MES	C5-N4-C3	5.63	121.51	108.83
2	B	1448	FAD	C4-N3-C2	5.54	119.82	115.14
2	B	1448	FAD	C1'-N10-C9A	5.48	122.61	118.29
2	A	1448	FAD	C4-N3-C2	5.48	119.77	115.14
4	A	1450	MES	C5-N4-C3	4.97	120.03	108.83
2	C	1448	FAD	C1'-N10-C9A	4.58	121.90	118.29
2	A	1448	FAD	C4X-N5-C5X	4.03	120.80	116.77
4	A	1450	MES	O3S-S-C8	3.89	112.06	105.77
2	B	1448	FAD	C4X-N5-C5X	3.72	120.49	116.77
4	A	1450	MES	C7-N4-C3	3.47	120.11	111.23
4	B	1450	MES	C7-N4-C5	3.41	119.95	111.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1448	FAD	C4X-N5-C5X	3.39	120.16	116.77
4	A	1450	MES	C7-N4-C5	3.37	119.86	111.23
2	A	1448	FAD	C1'-N10-C9A	3.30	120.89	118.29
2	A	1448	FAD	C5X-C9A-N10	3.29	120.10	117.72
4	A	1450	MES	C2-C3-N4	-2.96	105.62	110.10
4	B	1450	MES	C7-N4-C3	2.85	118.52	111.23
4	B	1450	MES	O3S-S-C8	2.77	110.25	105.77
2	B	1448	FAD	C5X-C9A-N10	2.68	119.66	117.72
2	A	1448	FAD	C4X-C4-N3	-2.53	119.97	123.43
2	C	1448	FAD	C4X-C4-N3	-2.50	120.01	123.43
2	C	1448	FAD	P-O3P-PA	-2.50	124.26	132.83
2	B	1448	FAD	C10-C4X-N5	-2.44	119.57	121.26
2	A	1448	FAD	C10-C4X-N5	-2.40	119.60	121.26
2	A	1448	FAD	P-O3P-PA	-2.39	124.61	132.83
4	B	1450	MES	O2S-S-C8	2.38	109.78	106.92
2	B	1448	FAD	C4X-C4-N3	-2.33	120.24	123.43
2	C	1448	FAD	C5X-C9A-N10	2.32	119.40	117.72
2	B	1448	FAD	P-O3P-PA	-2.29	124.96	132.83
4	B	1450	MES	C2-C3-N4	-2.25	106.69	110.10
4	A	1450	MES	O2S-S-C8	2.18	109.53	106.92
2	B	1448	FAD	C4A-C5A-N7A	-2.07	107.24	109.40
2	C	1448	FAD	C4A-C5A-N7A	-2.02	107.30	109.40

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1450	MES	C7-C8-S-O1S
2	B	1448	FAD	N10-C1'-C2'-O2'
2	B	1448	FAD	N10-C1'-C2'-C3'
2	A	1448	FAD	N10-C1'-C2'-O2'
2	C	1448	FAD	N10-C1'-C2'-O2'
2	C	1448	FAD	C5'-O5'-P-O2P
2	C	1448	FAD	C5'-O5'-P-O3P
4	B	1450	MES	C7-C8-S-O3S
2	C	1448	FAD	C3'-C4'-C5'-O5'
4	A	1450	MES	C7-C8-S-O3S
4	B	1450	MES	C8-C7-N4-C3
4	A	1450	MES	C8-C7-N4-C3
2	A	1448	FAD	PA-O3P-P-O5'
2	C	1448	FAD	C5'-O5'-P-O1P
4	B	1450	MES	C7-C8-S-O2S

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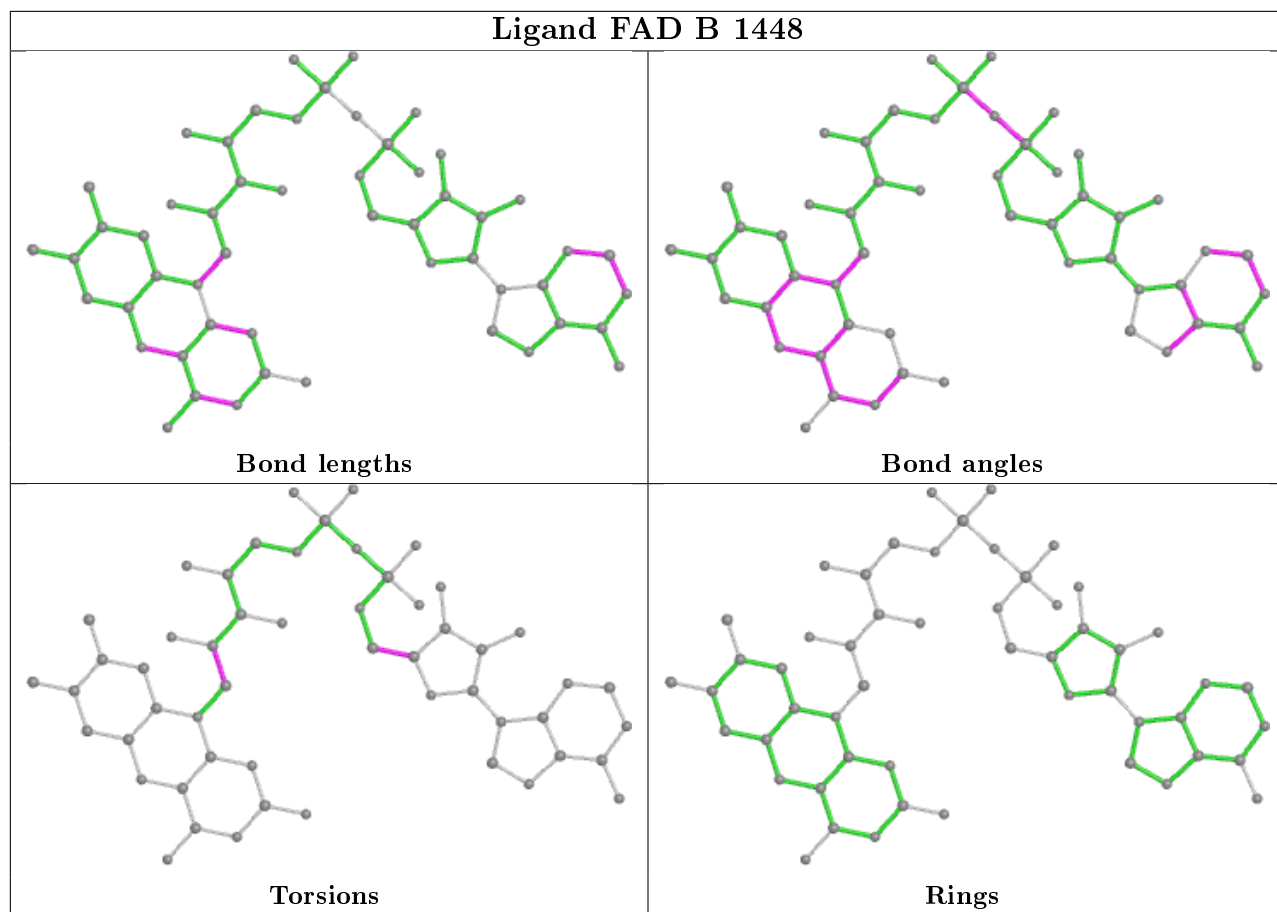
Mol	Chain	Res	Type	Atoms
4	A	1450	MES	C7-C8-S-O2S
2	A	1448	FAD	N10-C1'-C2'-C3'
2	C	1448	FAD	N10-C1'-C2'-C3'
2	C	1448	FAD	O4B-C4B-C5B-O5B
2	C	1448	FAD	O4'-C4'-C5'-O5'
2	B	1448	FAD	O4B-C4B-C5B-O5B
2	A	1448	FAD	O4B-C4B-C5B-O5B
4	A	1450	MES	C7-C8-S-O1S

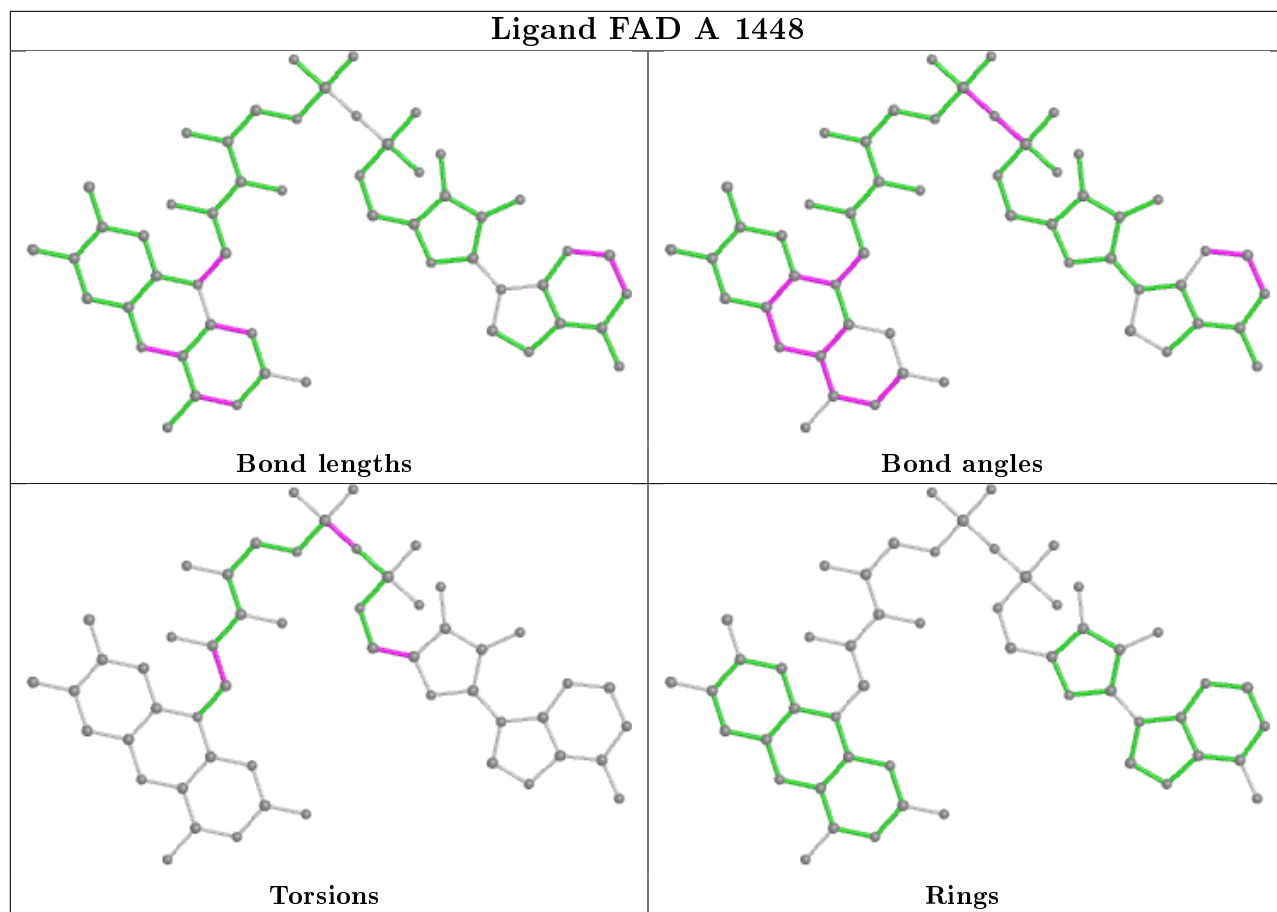
There are no ring outliers.

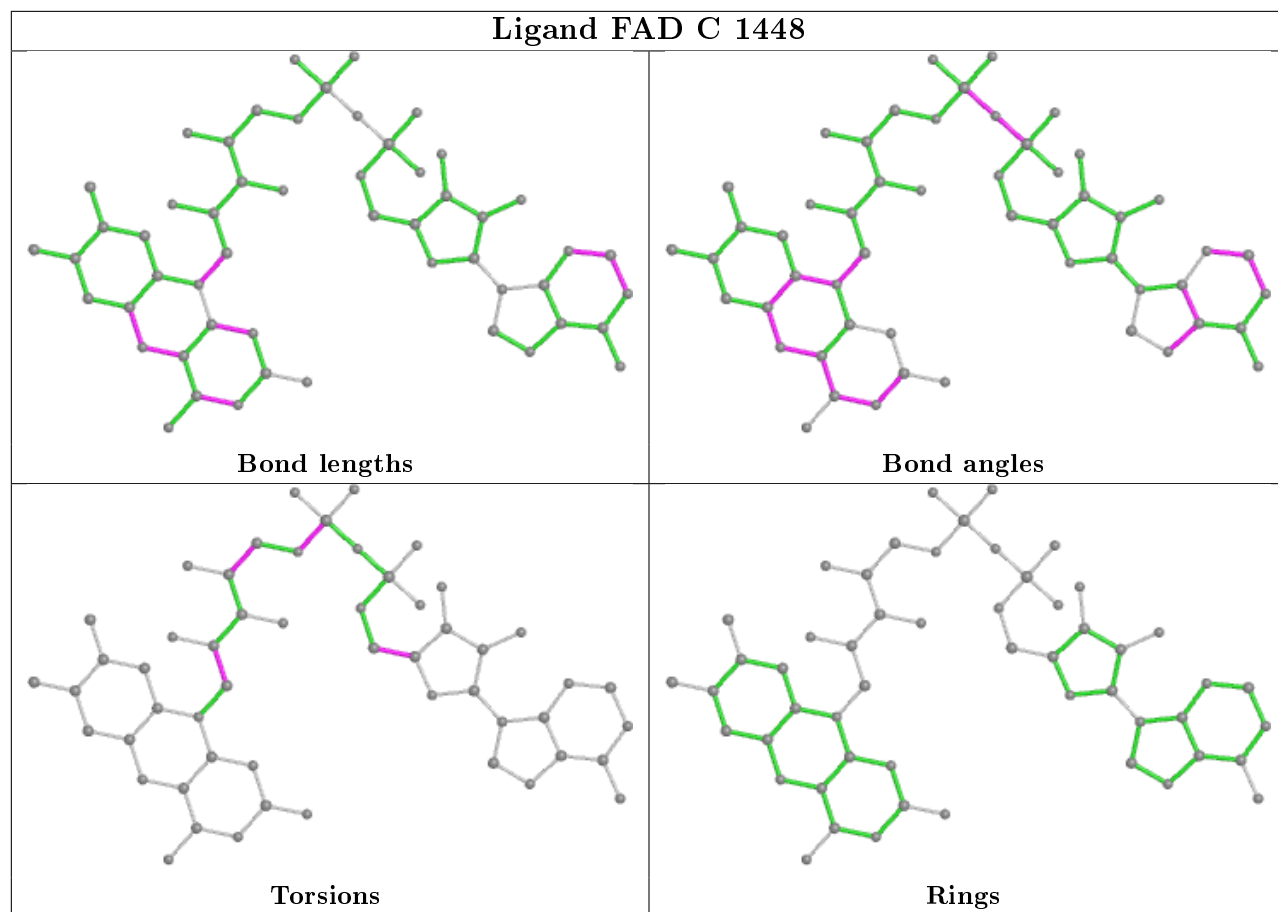
3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1448	FAD	1	0
2	A	1448	FAD	2	0
3	B	1449	NO3	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, 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	445/464 (95%)	0.03	6 (1%) 77 78	33, 62, 86, 95	0
1	B	446/464 (96%)	-0.10	5 (1%) 80 82	33, 56, 74, 84	0
1	C	446/464 (96%)	-0.03	11 (2%) 57 59	42, 64, 94, 112	0
All	All	1337/1392 (96%)	-0.03	22 (1%) 72 73	33, 60, 89, 112	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	91	GLY	3.2
1	A	446	ASP	3.1
1	C	241	TRP	3.1
1	C	2	ALA	2.8
1	A	258	ALA	2.8
1	C	245	PRO	2.7
1	C	258	ALA	2.6
1	C	225	ILE	2.6
1	C	257	PHE	2.4
1	B	445	SER	2.4
1	B	94	ILE	2.4
1	B	353	MET	2.4
1	A	134	ASP	2.3
1	A	259	ASP	2.3
1	C	259	ASP	2.3
1	C	256	TYR	2.3
1	A	261	SER	2.2
1	C	231	ALA	2.2
1	C	242	ASP	2.1
1	B	2	ALA	2.1
1	C	408	ILE	2.0
1	B	93	PRO	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](#)

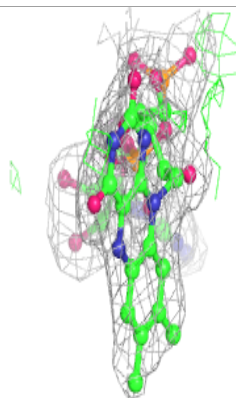
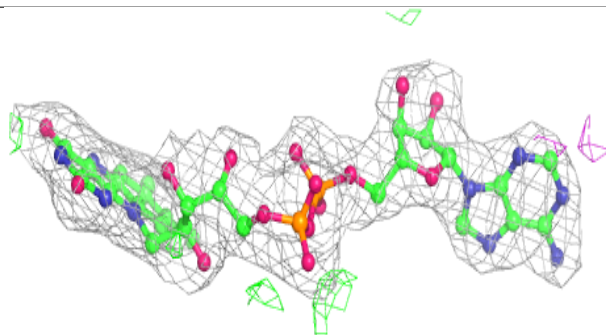
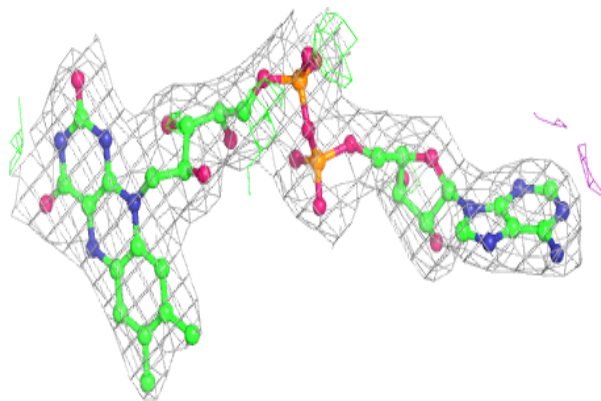
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	OXY	C	1450	2/2	0.75	0.87	93,93,93,93	0
5	OXY	A	1451	2/2	0.82	0.79	73,73,73,73	0
3	NO3	C	1449	4/4	0.87	0.27	87,87,87,87	0
5	OXY	B	1451	2/2	0.88	0.59	66,66,66,67	0
4	MES	B	1450	12/12	0.90	0.19	102,105,106,106	0
3	NO3	B	1449	4/4	0.92	0.30	70,70,70,71	0
4	MES	A	1450	12/12	0.95	0.15	87,88,89,89	0
2	FAD	C	1448	53/53	0.96	0.20	44,49,52,53	0
3	NO3	A	1449	4/4	0.96	0.23	75,75,76,76	0
2	FAD	A	1448	53/53	0.97	0.22	39,42,54,55	0
2	FAD	B	1448	53/53	0.98	0.18	38,41,51,51	0

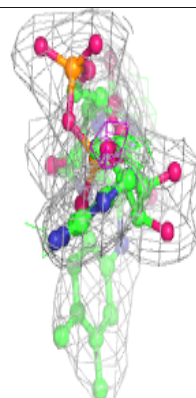
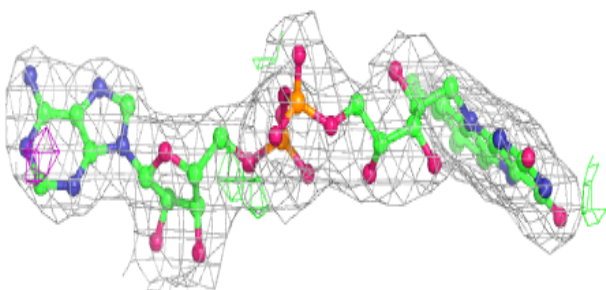
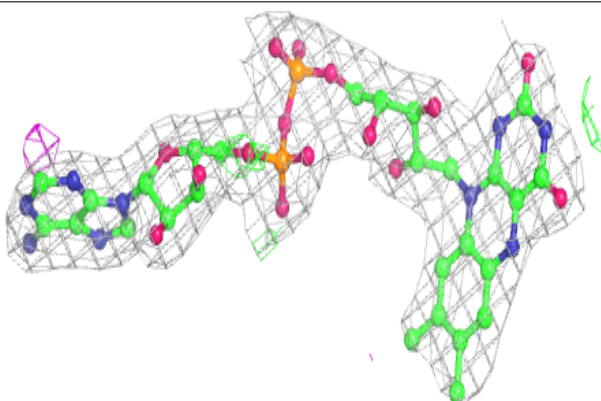
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

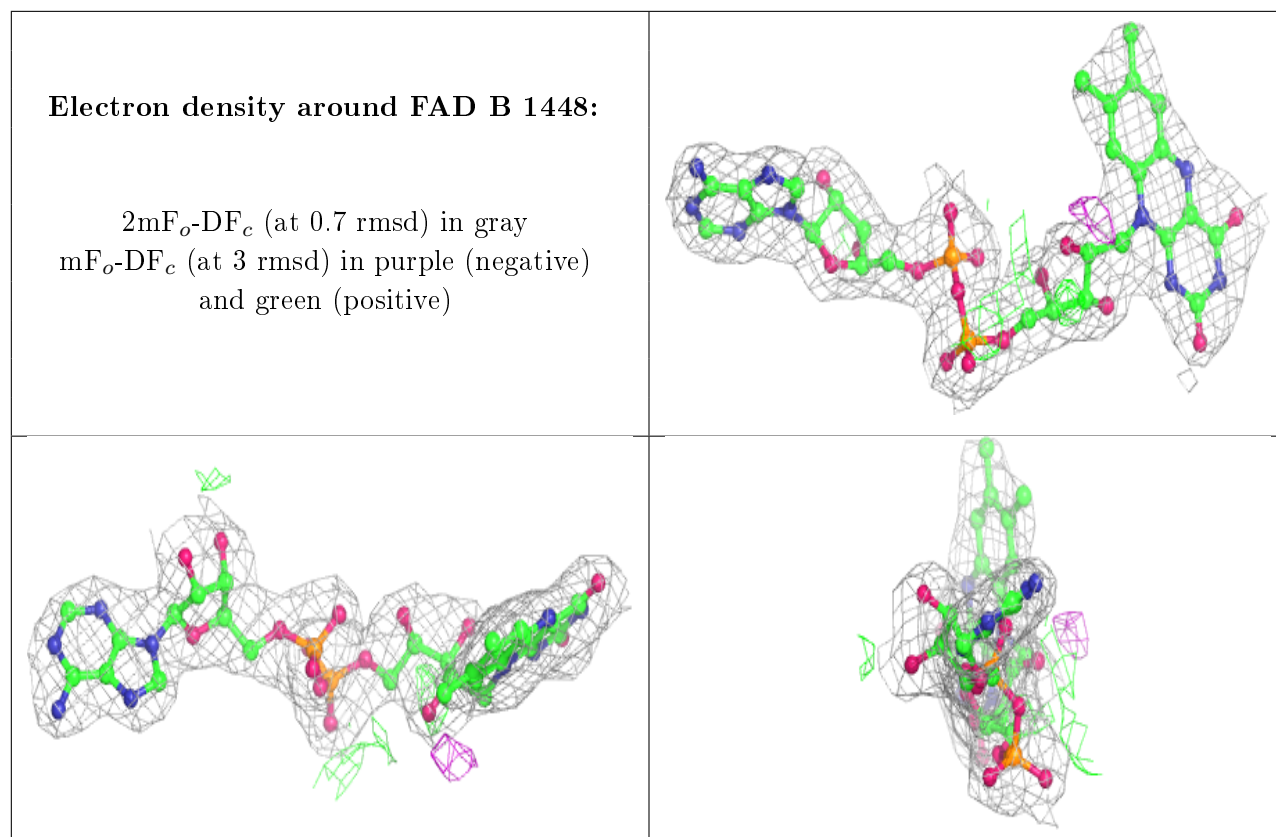
Electron density around FAD C 1448:

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

**Electron density around FAD A 1448:**

$2mF_o-DF_c$ (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.