Full wwPDB X-ray Structure Validation Report ⓘ

Apr 28, 2024 – 09:54 pm BST

PDB ID	:	2XVJ
Title	:	Crystal structure of the mutant bacterial flavin containing monooxygenase in complex with indole
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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

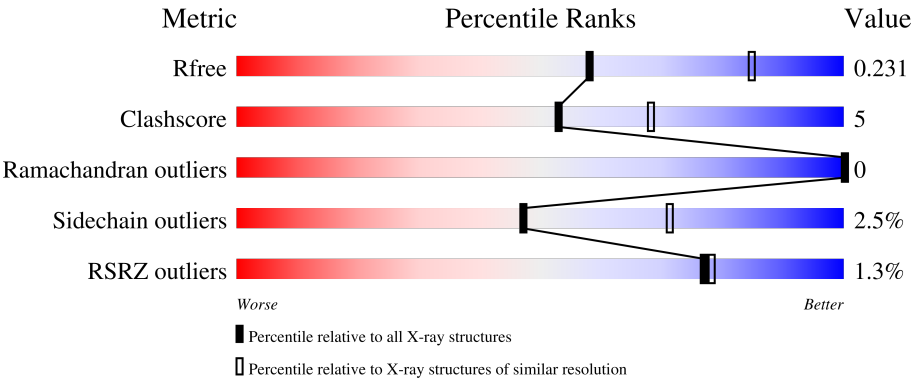
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.48 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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	464	3%, 81%, 15%, .
1	B	464	81%, 14%, .
1	C	464	84%, 12%, .

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
3	NO3	A	1449	-	-	X	-
3	NO3	B	1448	-	-	X	-

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 11214 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
1	A	446	Total	C	N	O	S	0	1	0
			3634	2330	598	686	20			
1	B	445	Total	C	N	O	S	0	1	0
			3611	2317	596	678	20			
1	C	446	Total	C	N	O	S	0	0	0
			3612	2317	595	680	20			

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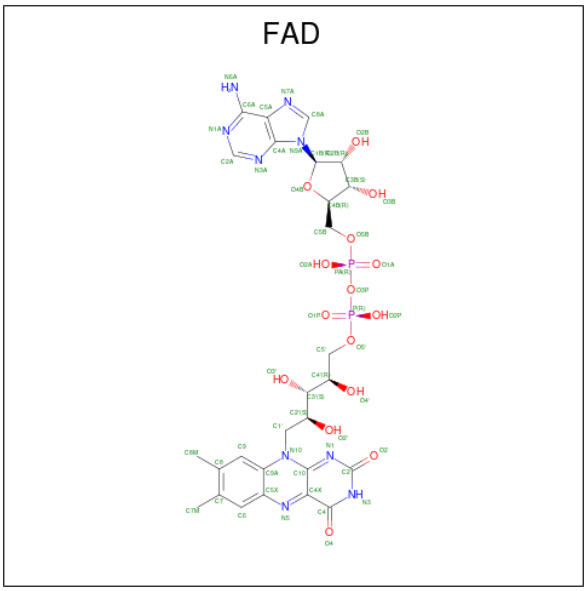
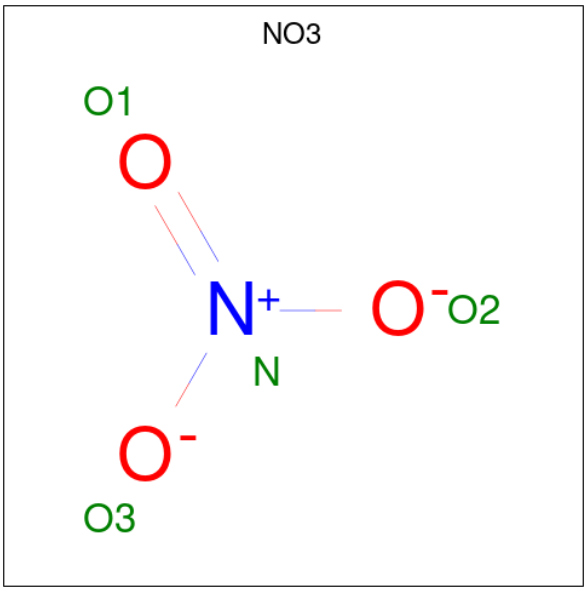
Chain	Residue	Modelled	Actual	Comment	Reference
C	462	HIS	-	expression tag	UNP Q83XK4
C	463	HIS	-	expression tag	UNP Q83XK4
C	464	HIS	-	expression tag	UNP Q83XK4
C	207	SER	TYR	engineered mutation	UNP Q83XK4

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



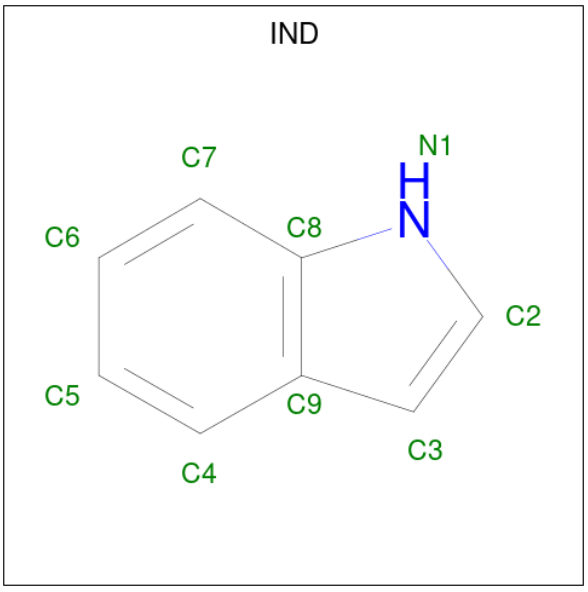
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

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



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

- Molecule 4 is INDOLE (three-letter code: IND) (formula: C_8H_7N).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N 9 8 1	0	0

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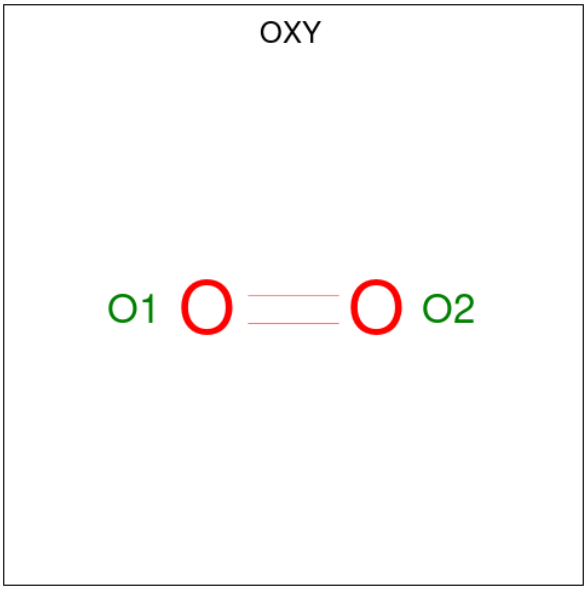
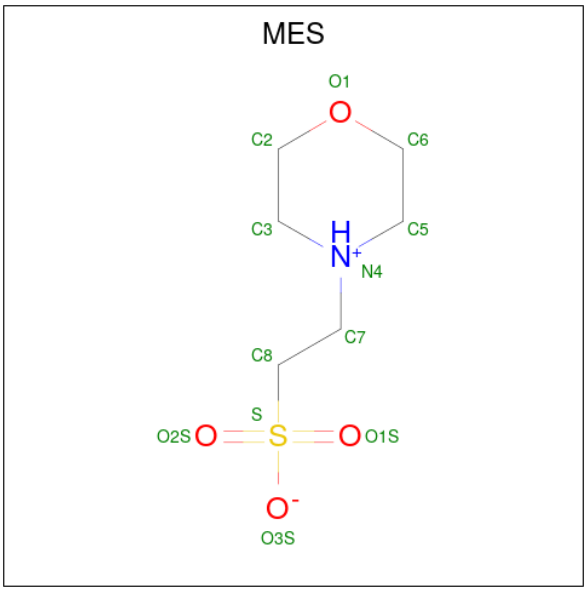
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C N 9 8 1	0	0
4	C	1	Total C N 9 8 1	0	0

- Molecule 5 is OXYGEN MOLECULE (three-letter code: OXY) (formula: O_2).



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

- Molecule 6 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



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

- Molecule 7 is water.

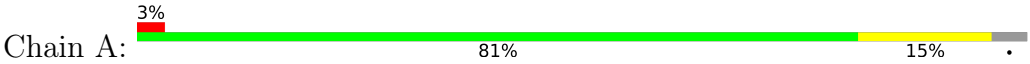
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	32	Total O 32 32	0	0
7	B	54	Total O 54 54	0	0
7	C	55	Total O 55 55	0	0

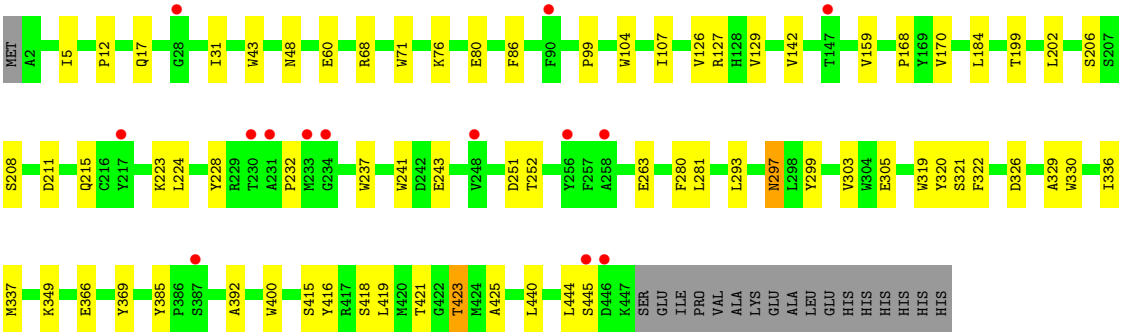
3 Residue-property plots (i)

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

• Molecule 1: FLAVIN-CONTAINING MONOOXYGENASE

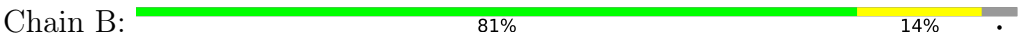
Chain A:

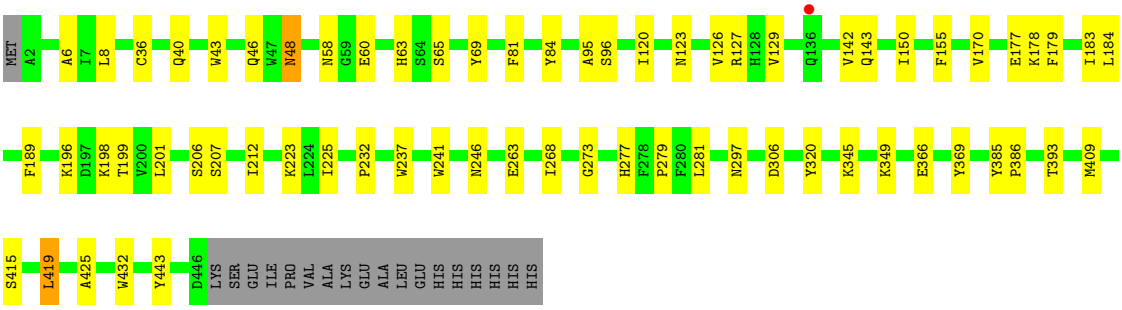




• Molecule 1: FLAVIN-CONTAINING MONOOXYGENASE

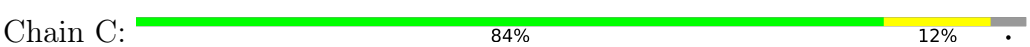
Chain B:

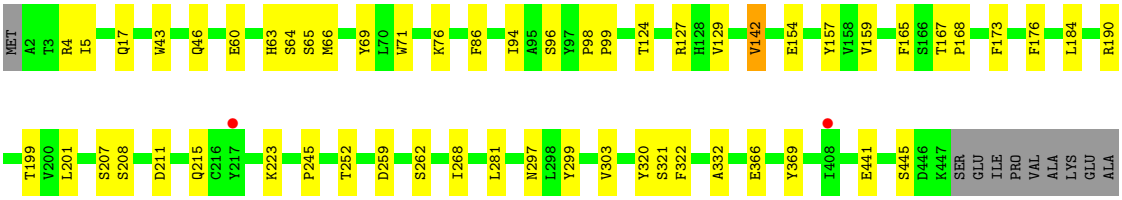




• Molecule 1: FLAVIN-CONTAINING MONOOXYGENASE

Chain C:





LEU
GLU
HIS
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	159.12Å 67.85Å 138.90Å 90.00° 90.62° 90.00°	Depositor
Resolution (Å)	50.01 – 2.48 46.32 – 2.48	Depositor EDS
% Data completeness (in resolution range)	97.3 (50.01-2.48) 84.7 (46.32-2.48)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	223.90 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.199 , 0.237 0.194 , 0.231	Depositor DCC
R_{free} test set	2316 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å^2)	41.9	Xtriage
Anisotropy	0.078	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 36.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	11214	wwPDB-VP
Average B, all atoms (Å^2)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.55	4/3751 (0.1%)	0.55	0/5100
1	B	0.50	2/3728 (0.1%)	0.53	0/5072
1	C	0.51	2/3725 (0.1%)	0.54	0/5068
All	All	0.52	8/11204 (0.1%)	0.54	0/15240

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	320	TYR	CD2-CE2	-7.70	1.27	1.39
1	A	320	TYR	CD1-CE1	-7.63	1.27	1.39
1	C	320	TYR	CD1-CE1	-6.78	1.29	1.39
1	B	320	TYR	CD1-CE1	-6.36	1.29	1.39
1	B	320	TYR	CD2-CE2	-6.09	1.30	1.39
1	A	320	TYR	CD2-CE2	-6.00	1.30	1.39
1	A	416	TYR	CE1-CZ	-5.79	1.31	1.38
1	A	416	TYR	CD1-CE1	-5.56	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

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	3634	0	3374	46	0
1	B	3611	0	3351	37	0
1	C	3612	0	3343	32	0
2	A	53	0	31	1	0
2	B	53	0	31	2	0
2	C	53	0	31	1	0
3	A	4	0	0	2	0
3	B	4	0	0	4	0
3	C	4	0	0	0	0
4	A	9	0	7	0	0
4	B	9	0	7	0	0
4	C	9	0	7	1	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
5	C	2	0	0	0	0
6	B	12	0	12	1	0
7	A	32	0	0	0	0
7	B	54	0	0	0	0
7	C	55	0	0	0	0
All	All	11214	0	10194	115	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (115) 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:C:46:GLN:NE2	1:C:65:SER:H	1.61	0.97
1:C:46:GLN:HE22	1:C:65:SER:N	1.65	0.94
1:C:46:GLN:HE22	1:C:65:SER:H	0.82	0.78
1:B:273:GLY:N	3:B:1448:NO3:O3	2.24	0.70
1:A:228:TYR:HE1	1:A:232:PRO:HA	1.56	0.69
1:A:293:LEU:HD21	1:A:319:TRP:CD1	2.28	0.68
1:A:60:GLU:OE2	1:A:127:ARG:NH2	2.30	0.63
1:B:60:GLU:OE2	1:B:127:ARG:NH2	2.27	0.63
1:B:46:GLN:OE1	1:B:65:SER:N	2.29	0.62
1:A:251:ASP:OD1	1:A:252:THR:N	2.33	0.62
1:B:179:PHE:HD2	1:B:183:ILE:HD11	1.65	0.61
1:B:199:THR:HG23	1:B:223:LYS:HB3	1.83	0.61
1:A:5:ILE:HD11	1:A:336:ILE:HG21	1.82	0.60
1:C:321:SER:OG	1:C:322:PHE:N	2.34	0.60
1:A:319:TRP:HB3	1:A:369[B]:TYR:CE1	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:60:GLU:OE2	1:C:127:ARG:NH2	2.29	0.59
1:A:369[B]:TYR:OH	1:A:400:TRP:HZ3	1.85	0.59
1:A:321:SER:OG	1:A:322:PHE:N	2.37	0.57
1:A:251:ASP:OD1	1:A:251:ASP:C	2.43	0.57
1:A:71:TRP:HA	1:A:99:PRO:HA	1.87	0.55
1:A:369[B]:TYR:HH	1:A:400:TRP:HZ3	1.55	0.55
1:A:211:ASP:O	1:A:215:GLN:HG2	2.06	0.55
1:A:104:TRP:HA	1:A:107:ILE:HD12	1.89	0.55
1:A:319:TRP:HE3	1:A:369[B]:TYR:CZ	2.25	0.54
1:B:409:MET:HE3	1:B:432:TRP:CZ3	2.43	0.54
1:B:409:MET:HE3	1:B:432:TRP:HZ3	1.73	0.54
1:B:206:SER:HB2	3:B:1448:NO3:O2	2.08	0.54
1:A:326:ASP:HB3	1:A:385:TYR:CE1	2.42	0.53
1:B:127:ARG:HH21	1:B:143:GLN:NE2	2.05	0.53
1:A:330:TRP:CD1	1:A:385:TYR:HB2	2.44	0.53
1:B:225:ILE:HD11	1:B:263:GLU:HG2	1.91	0.52
1:A:17:GLN:HA	1:A:329:ALA:HB1	1.91	0.52
1:A:202:LEU:HD11	1:A:224:LEU:HD22	1.92	0.51
1:C:124:THR:CG2	1:C:142:VAL:HG13	2.40	0.51
1:A:293:LEU:HD21	1:A:319:TRP:HD1	1.74	0.51
1:B:366:GLU:HA	1:B:369:TYR:CZ	2.46	0.51
1:C:17:GLN:HE22	1:C:332:ALA:HB3	1.75	0.51
1:C:71:TRP:HA	1:C:99:PRO:HA	1.93	0.50
1:C:199:THR:HG23	1:C:223:LYS:HB3	1.94	0.49
1:C:173:PHE:HB2	1:C:176:PHE:CE1	2.48	0.49
1:B:81:PHE:HB2	1:B:84:TYR:O	2.12	0.49
1:A:319:TRP:CE3	1:A:369[B]:TYR:CZ	3.00	0.49
1:B:415:SER:HB3	1:B:425:ALA:O	2.13	0.48
1:C:299:TYR:HB3	1:C:303:VAL:HB	1.96	0.48
1:B:273:GLY:CA	3:B:1448:NO3:O3	2.62	0.48
1:A:444:LEU:O	1:A:445:SER:OG	2.29	0.47
1:C:4:ARG:NH1	1:C:154:GLU:O	2.47	0.47
6:B:1450:MES:H82	6:B:1450:MES:H51	1.69	0.47
1:B:237:TRP:HB3	1:B:241:TRP:HB3	1.95	0.47
1:C:5:ILE:HD12	1:C:157:TYR:HB2	1.96	0.47
1:C:441:GLU:O	1:C:445:SER:HB2	2.15	0.47
1:A:297:ASN:HA	1:A:305:GLU:OE1	2.14	0.47
1:B:177:GLU:H	1:B:177:GLU:CD	2.18	0.47
1:A:366:GLU:HA	1:A:369[A]:TYR:CZ	2.49	0.47
1:C:76:LYS:HE2	1:C:94:ILE:HG13	1.96	0.46
1:B:277:HIS:CE1	1:B:279:PRO:HG3	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:211:ASP:O	1:C:215:GLN:HG2	2.16	0.46
1:B:8:LEU:HD22	1:B:126:VAL:HG21	1.97	0.46
1:B:65:SER:HB2	2:B:1447:FAD:HM82	1.97	0.46
1:B:385:TYR:CD1	1:B:386:PRO:HD2	2.50	0.46
1:A:68:ARG:HH12	1:B:178:LYS:HA	1.81	0.45
1:A:80:GLU:OE1	1:A:418:SER:OG	2.30	0.45
1:B:201:LEU:HD23	1:B:268:ILE:HG12	1.99	0.45
1:C:64:SER:O	1:C:190:ARG:NH2	2.49	0.45
1:C:201:LEU:HD23	1:C:268:ILE:HG12	1.99	0.45
1:B:129:VAL:HB	1:B:281:LEU:HD23	1.99	0.45
1:A:421:THR:OG1	1:A:423:THR:OG1	2.28	0.45
1:B:40:GLN:HE22	1:B:48:ASN:HB3	1.81	0.45
1:C:46:GLN:NE2	1:C:65:SER:N	2.43	0.45
1:C:46:GLN:HE21	1:C:66:MET:HG3	1.82	0.45
1:B:127:ARG:HH21	1:B:143:GLN:HE21	1.64	0.44
1:A:206:SER:HB2	3:A:1449:NO3:O1	2.18	0.44
1:A:415:SER:OG	1:A:425:ALA:O	2.29	0.44
1:B:225:ILE:CD1	1:B:263:GLU:HG2	2.48	0.44
1:A:126:VAL:HA	1:A:142:VAL:HG12	1.98	0.44
1:A:129:VAL:HB	1:A:281:LEU:HD23	1.98	0.44
1:A:168:PRO:HB2	1:A:170:VAL:HG22	2.00	0.44
1:B:189:PHE:CD2	1:B:212:ILE:HG12	2.53	0.44
1:A:392:ALA:CB	1:A:419:LEU:HD22	2.48	0.43
1:A:392:ALA:HB3	1:A:419:LEU:HD22	2.00	0.43
1:B:273:GLY:HA2	3:B:1448:NO3:O3	2.19	0.43
1:A:31:ILE:HG22	1:A:337:MET:CE	2.49	0.43
1:B:345:LYS:O	1:B:349:LYS:HG2	2.18	0.43
1:B:393:THR:OG1	1:B:419:LEU:HD13	2.19	0.43
1:B:95:ALA:HB3	1:B:443:TYR:OH	2.19	0.42
1:A:237:TRP:HB3	1:A:241:TRP:HB3	2.01	0.42
1:C:96:SER:C	1:C:98:PRO:HD3	2.39	0.42
1:C:98:PRO:HA	1:C:99:PRO:HD3	1.93	0.42
1:A:76:LYS:HB2	1:A:86:PHE:CZ	2.54	0.42
1:B:58:ASN:ND2	1:B:123:ASN:OD1	2.52	0.42
1:A:17:GLN:HE21	1:A:159:VAL:HG11	1.85	0.42
1:C:63:HIS:NE2	2:C:1448:FAD:H51A	2.34	0.42
1:A:223:LYS:HD3	1:A:263:GLU:OE2	2.20	0.41
1:C:165:PHE:CE2	4:C:1450:IND:H4	2.54	0.41
1:A:280:PHE:CD1	1:A:281:LEU:HG	2.55	0.41
1:A:319:TRP:HB3	1:A:369[B]:TYR:CD1	2.55	0.41
1:B:63:HIS:NE2	2:B:1447:FAD:H51A	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:THR:HG23	1:A:223:LYS:HB3	2.02	0.41
1:C:167:THR:HA	1:C:168:PRO:HD3	1.87	0.41
1:C:17:GLN:HG3	1:C:159:VAL:HG11	2.03	0.41
1:C:76:LYS:HD2	1:C:86:PHE:CZ	2.55	0.41
1:C:129:VAL:HB	1:C:281:LEU:HD23	2.03	0.41
1:C:76:LYS:HB2	1:C:86:PHE:CZ	2.56	0.41
1:A:12:PRO:HG2	2:A:1448:FAD:H4'	2.03	0.41
1:A:208:SER:HB2	3:A:1449:NO3:O1	2.20	0.41
1:A:305:GLU:OE1	1:A:349:LYS:HE3	2.21	0.41
1:B:142:VAL:O	1:B:150:ILE:HG23	2.21	0.41
1:B:6:ALA:HB2	1:B:155:PHE:CG	2.56	0.40
1:B:196:LYS:O	1:B:198:LYS:HD3	2.20	0.40
1:C:245:PRO:HG3	1:C:259:ASP:HB3	2.03	0.40
1:C:366:GLU:HA	1:C:369:TYR:CZ	2.56	0.40
1:A:60:GLU:CD	1:A:127:ARG:HH22	2.23	0.40
1:A:299:TYR:HB3	1:A:303:VAL:HB	2.03	0.40
1:B:36:CYS:HB3	1:B:120:ILE:HA	2.04	0.40
1:C:124:THR:HG23	1:C:142:VAL:HG13	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	445/464 (96%)	430 (97%)	15 (3%)	0	100	100
1	B	444/464 (96%)	428 (96%)	16 (4%)	0	100	100
1	C	444/464 (96%)	427 (96%)	17 (4%)	0	100	100
All	All	1333/1392 (96%)	1285 (96%)	48 (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	381/405 (94%)	374 (98%)	7 (2%)	59	80
1	B	377/405 (93%)	365 (97%)	12 (3%)	39	63
1	C	375/405 (93%)	366 (98%)	9 (2%)	49	72
All	All	1133/1215 (93%)	1105 (98%)	28 (2%)	47	71

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	TRP
1	A	48	ASN
1	A	184	LEU
1	A	243	GLU
1	A	297	ASN
1	A	423	THR
1	A	440	LEU
1	B	43	TRP
1	B	48	ASN
1	B	69	TYR
1	B	96	SER
1	B	170	VAL
1	B	184	LEU
1	B	207	SER
1	B	232	PRO
1	B	246	ASN
1	B	297	ASN
1	B	306	ASP
1	B	419	LEU
1	C	43	TRP
1	C	69	TYR
1	C	142	VAL
1	C	184	LEU
1	C	207	SER
1	C	208	SER
1	C	252	THR

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Mol	Chain	Res	Type
1	C	262	SER
1	C	297	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	40	GLN
1	A	48	ASN
1	A	128	HIS
1	A	254	ASN
1	A	282	ASN
1	A	297	ASN
1	A	377	GLN
1	A	407	ASN
1	B	17	GLN
1	B	40	GLN
1	B	48	ASN
1	B	132	ASN
1	B	143	GLN
1	B	277	HIS
1	B	414	HIS
1	C	17	GLN
1	C	46	GLN
1	C	89	HIS
1	C	128	HIS
1	C	164	HIS
1	C	414	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

13 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	FAD	A	1448	-	53,58,58	1.12	4 (7%)	68,89,89	1.38	11 (16%)
4	IND	B	1449	-	8,10,10	1.01	0	9,13,13	2.28	2 (22%)
2	FAD	C	1448	-	53,58,58	1.12	4 (7%)	68,89,89	1.40	11 (16%)
2	FAD	B	1447	-	53,58,58	1.10	4 (7%)	68,89,89	1.46	8 (11%)
5	OXY	A	1451	-	1,1,1	0.13	0	-		
5	OXY	C	1451	-	1,1,1	0.15	0	-		
4	IND	A	1450	-	8,10,10	1.05	0	9,13,13	2.32	2 (22%)
3	NO3	A	1449	-	1,3,3	3.49	1 (100%)	0,3,3	-	-
6	MES	B	1450	-	12,12,12	2.20	1 (8%)	14,16,16	2.37	6 (42%)
3	NO3	C	1449	-	1,3,3	3.53	1 (100%)	0,3,3	-	-
4	IND	C	1450	-	8,10,10	1.02	0	9,13,13	2.32	2 (22%)
3	NO3	B	1448	-	1,3,3	2.29	1 (100%)	0,3,3	-	-
5	OXY	B	1451	-	1,1,1	0.13	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
2	FAD	A	1448	-	-	4/30/50/50	0/6/6/6
4	IND	B	1449	-	-	-	0/2/2/2
2	FAD	C	1448	-	-	3/30/50/50	0/6/6/6
2	FAD	B	1447	-	-	3/30/50/50	0/6/6/6
4	IND	A	1450	-	-	-	0/2/2/2
6	MES	B	1450	-	-	1/6/14/14	0/1/1/1
4	IND	C	1450	-	-	-	0/2/2/2

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	1450	MES	C8-S	-7.28	1.67	1.77
2	A	1448	FAD	C4X-N5	4.22	1.39	1.30
2	B	1447	FAD	C4X-N5	3.90	1.38	1.30
2	C	1448	FAD	C2A-N3A	3.89	1.38	1.32
2	C	1448	FAD	C4X-N5	3.83	1.38	1.30
2	B	1447	FAD	C2A-N3A	3.54	1.37	1.32
3	C	1449	NO3	O1-N	3.53	1.40	1.24
3	A	1449	NO3	O1-N	3.49	1.40	1.24
2	A	1448	FAD	C2A-N3A	3.34	1.37	1.32
2	B	1447	FAD	C10-N1	2.84	1.39	1.33
2	A	1448	FAD	C10-N1	2.67	1.38	1.33
2	C	1448	FAD	C2A-N1A	2.54	1.38	1.33
2	C	1448	FAD	C10-N1	2.39	1.38	1.33
2	A	1448	FAD	C2A-N1A	2.33	1.38	1.33
3	B	1448	NO3	O1-N	2.29	1.34	1.24
2	B	1447	FAD	C2A-N1A	2.10	1.37	1.33

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	1447	FAD	N3A-C2A-N1A	-6.70	118.21	128.68
4	C	1450	IND	C3-C9-C8	6.28	111.56	106.20
4	A	1450	IND	C3-C9-C8	6.26	111.55	106.20
4	B	1449	IND	C3-C9-C8	6.12	111.43	106.20
2	C	1448	FAD	N3A-C2A-N1A	-5.63	119.88	128.68
2	A	1448	FAD	N3A-C2A-N1A	-5.38	120.27	128.68
6	B	1450	MES	C5-N4-C3	4.80	119.63	108.83
6	B	1450	MES	C7-N4-C5	4.12	121.78	111.23
2	A	1448	FAD	C4-N3-C2	-3.18	119.77	125.64
2	B	1447	FAD	P-O3P-PA	-3.14	122.05	132.83
6	B	1450	MES	O3S-S-C8	3.02	110.66	105.77
2	B	1447	FAD	C4X-C10-N10	3.00	120.87	116.48
6	B	1450	MES	C7-N4-C3	2.88	118.59	111.23
2	A	1448	FAD	C4X-C4-N3	2.87	120.48	113.19
2	C	1448	FAD	C4X-C10-N10	2.80	120.58	116.48
2	C	1448	FAD	C4-N3-C2	-2.75	120.57	125.64
2	A	1448	FAD	C9A-C5X-N5	-2.69	119.51	122.43
2	C	1448	FAD	P-O3P-PA	-2.64	123.75	132.83
2	A	1448	FAD	C4X-C10-N10	2.61	120.29	116.48
2	C	1448	FAD	C4X-C4-N3	2.59	119.78	113.19
2	B	1447	FAD	C10-C4X-N5	-2.55	119.45	124.86
6	B	1450	MES	C2-C3-N4	-2.51	106.30	110.10
4	A	1450	IND	C3-C9-C4	-2.49	128.03	136.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	C	1450	IND	C3-C9-C4	-2.47	128.10	136.72
4	B	1449	IND	C3-C9-C4	-2.46	128.14	136.72
2	C	1448	FAD	C9A-C5X-N5	-2.43	119.79	122.43
2	A	1448	FAD	C10-C4X-N5	-2.38	119.80	124.86
2	A	1448	FAD	P-O3P-PA	-2.38	124.67	132.83
2	B	1447	FAD	C9A-C5X-N5	-2.37	119.86	122.43
2	C	1448	FAD	O4-C4-C4X	-2.37	120.32	126.60
2	C	1448	FAD	C1B-N9A-C4A	-2.36	122.50	126.64
2	B	1447	FAD	C4X-C4-N3	2.36	119.17	113.19
2	B	1447	FAD	C4-N3-C2	-2.32	121.36	125.64
2	C	1448	FAD	C10-C4X-N5	-2.31	119.95	124.86
2	A	1448	FAD	C4X-C10-N1	-2.23	119.55	124.73
2	A	1448	FAD	O4-C4-C4X	-2.22	120.72	126.60
2	C	1448	FAD	C10-N1-C2	2.19	121.28	116.90
2	A	1448	FAD	C5X-C9A-N10	2.14	120.16	117.95
2	C	1448	FAD	C4X-C10-N1	-2.11	119.83	124.73
6	B	1450	MES	O2S-S-C8	2.11	109.46	106.92
2	B	1447	FAD	C5'-C4'-C3'	2.01	116.09	112.20
2	A	1448	FAD	C4A-C5A-N7A	-2.01	107.31	109.40

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1448	FAD	N10-C1'-C2'-O2'
2	A	1448	FAD	N10-C1'-C2'-C3'
2	A	1448	FAD	PA-O3P-P-O5'
2	B	1447	FAD	N10-C1'-C2'-O2'
2	B	1447	FAD	N10-C1'-C2'-C3'
2	C	1448	FAD	N10-C1'-C2'-O2'
2	C	1448	FAD	N10-C1'-C2'-C3'
6	B	1450	MES	C8-C7-N4-C5
2	C	1448	FAD	O4B-C4B-C5B-O5B
2	B	1447	FAD	O4B-C4B-C5B-O5B
2	A	1448	FAD	O4B-C4B-C5B-O5B

There are no ring outliers.

7 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1448	FAD	1	0

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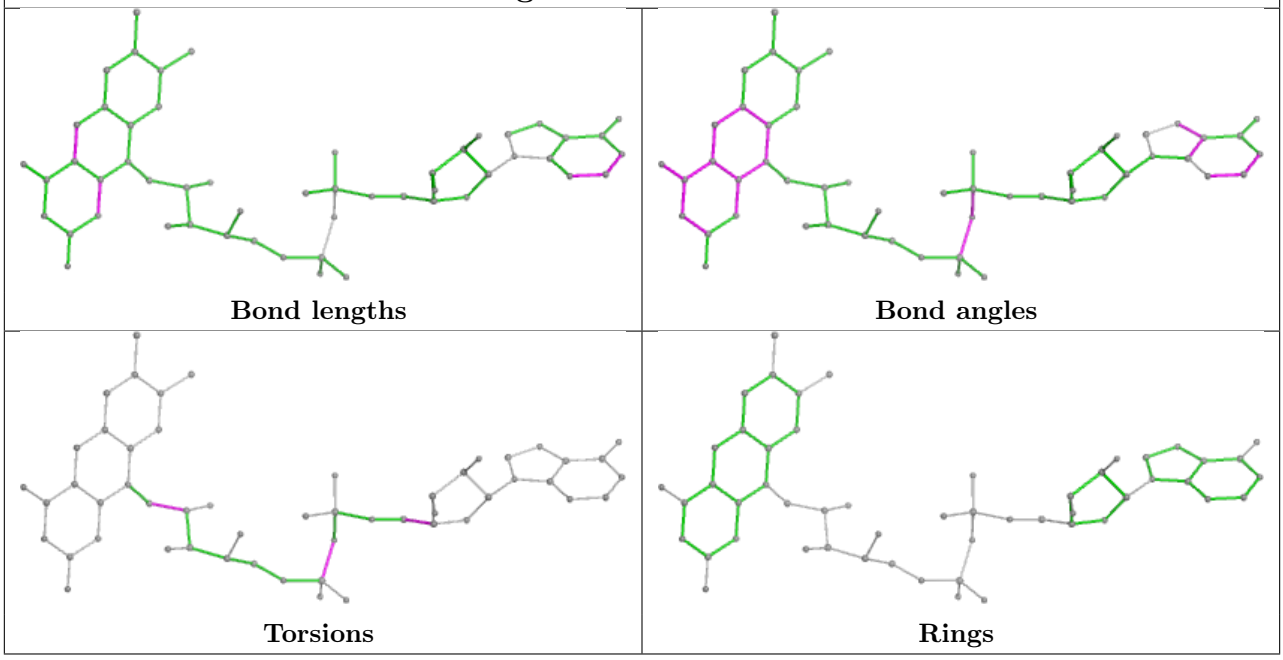
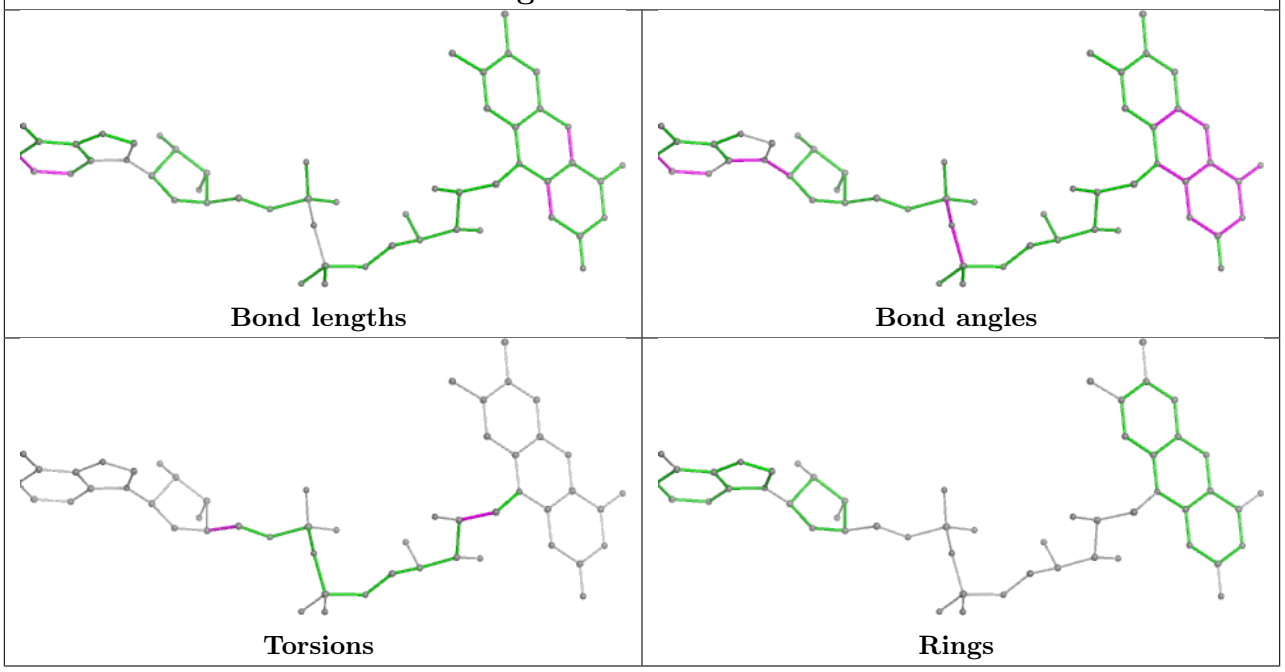
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1448	FAD	1	0
2	B	1447	FAD	2	0
3	A	1449	NO3	2	0
6	B	1450	MES	1	0
4	C	1450	IND	1	0
3	B	1448	NO3	4	0

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

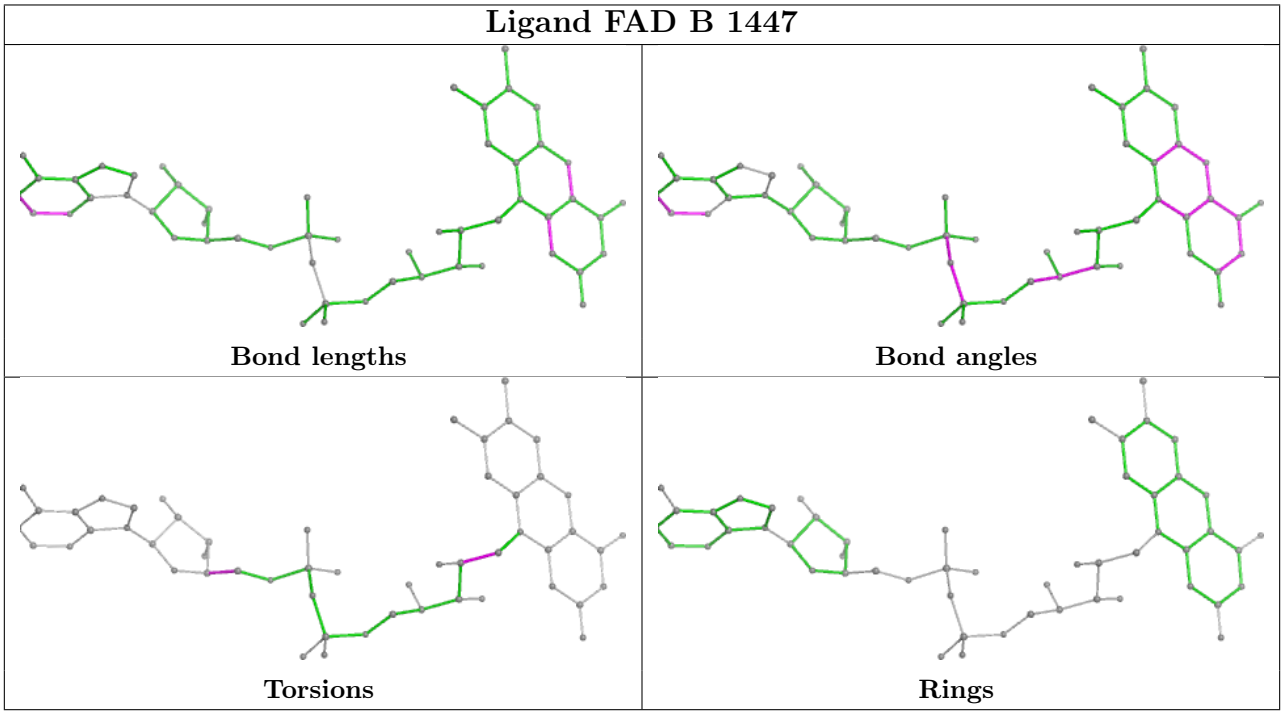
Ligand FAD A 1448



Ligand FAD C 1448



Ligand FAD B 1447



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^{th} 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	446/464 (96%)	0.07	14 (3%) 49 51	23, 45, 68, 86	0
1	B	445/464 (95%)	-0.19	1 (0%) 95 95	23, 38, 58, 67	0
1	C	446/464 (96%)	-0.15	2 (0%) 92 93	20, 37, 65, 83	0
All	All	1337/1392 (96%)	-0.09	17 (1%) 77 78	20, 40, 65, 86	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	231	ALA	5.5
1	A	230	THR	3.8
1	A	446	ASP	3.7
1	A	233	MET	3.4
1	A	217	TYR	2.9
1	A	234	GLY	2.7
1	C	217	TYR	2.6
1	A	90	PHE	2.5
1	A	258	ALA	2.4
1	A	256	TYR	2.4
1	A	147	THR	2.4
1	A	248	VAL	2.4
1	C	408	ILE	2.3
1	A	387	SER	2.2
1	A	28	GLY	2.2
1	A	445	SER	2.1
1	B	136	GLN	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 (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

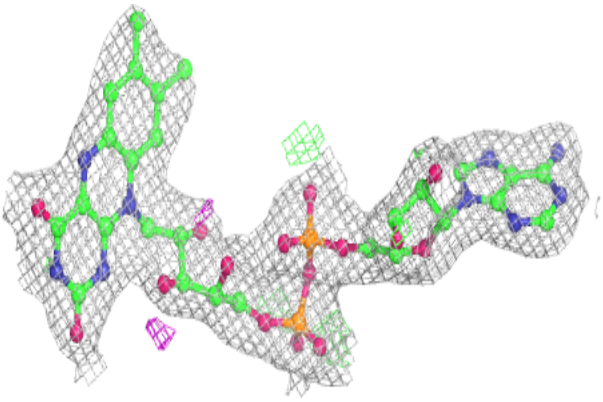
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

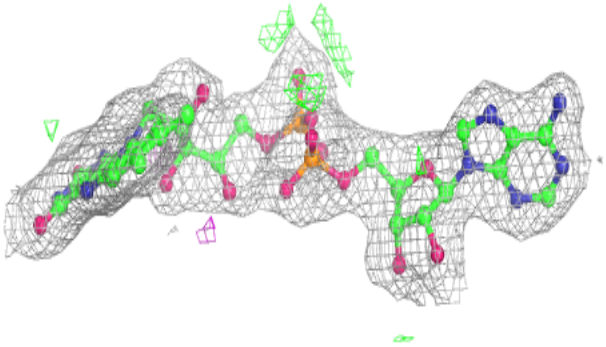
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	IND	C	1450	9/9	0.68	0.34	94,94,95,95	0
6	MES	B	1450	12/12	0.71	0.31	102,102,103,103	0
4	IND	B	1449	9/9	0.85	0.22	59,59,59,60	0
3	NO3	B	1448	4/4	0.86	0.30	71,72,72,72	0
5	OXY	C	1451	2/2	0.87	0.40	69,69,69,71	0
3	NO3	A	1449	4/4	0.89	0.31	73,73,73,74	0
5	OXY	B	1451	2/2	0.93	0.32	69,69,69,69	0
3	NO3	C	1449	4/4	0.94	0.22	66,66,67,67	0
4	IND	A	1450	9/9	0.95	0.20	80,80,81,81	0
5	OXY	A	1451	2/2	0.97	0.16	55,55,55,56	0
2	FAD	A	1448	53/53	0.97	0.12	23,26,37,38	0
2	FAD	B	1447	53/53	0.98	0.12	15,24,27,28	0
2	FAD	C	1448	53/53	0.98	0.13	18,25,28,29	0

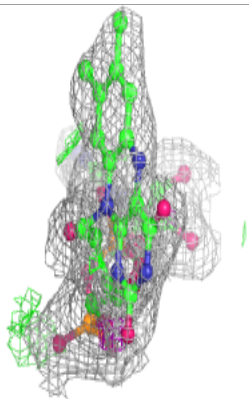
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD A 1448:

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

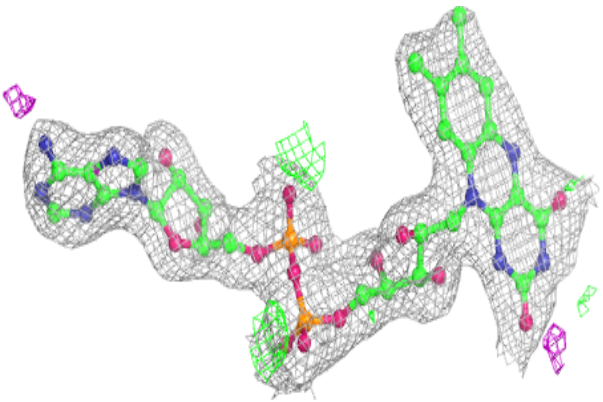


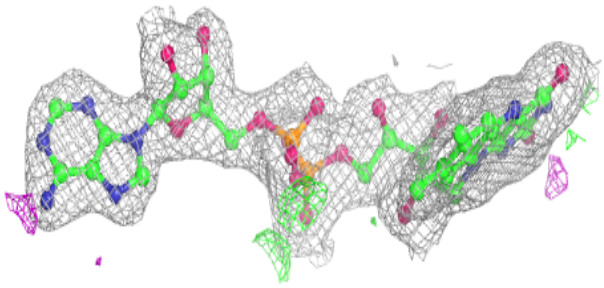


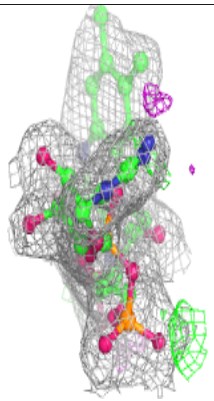


Electron density around FAD B 1447:

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

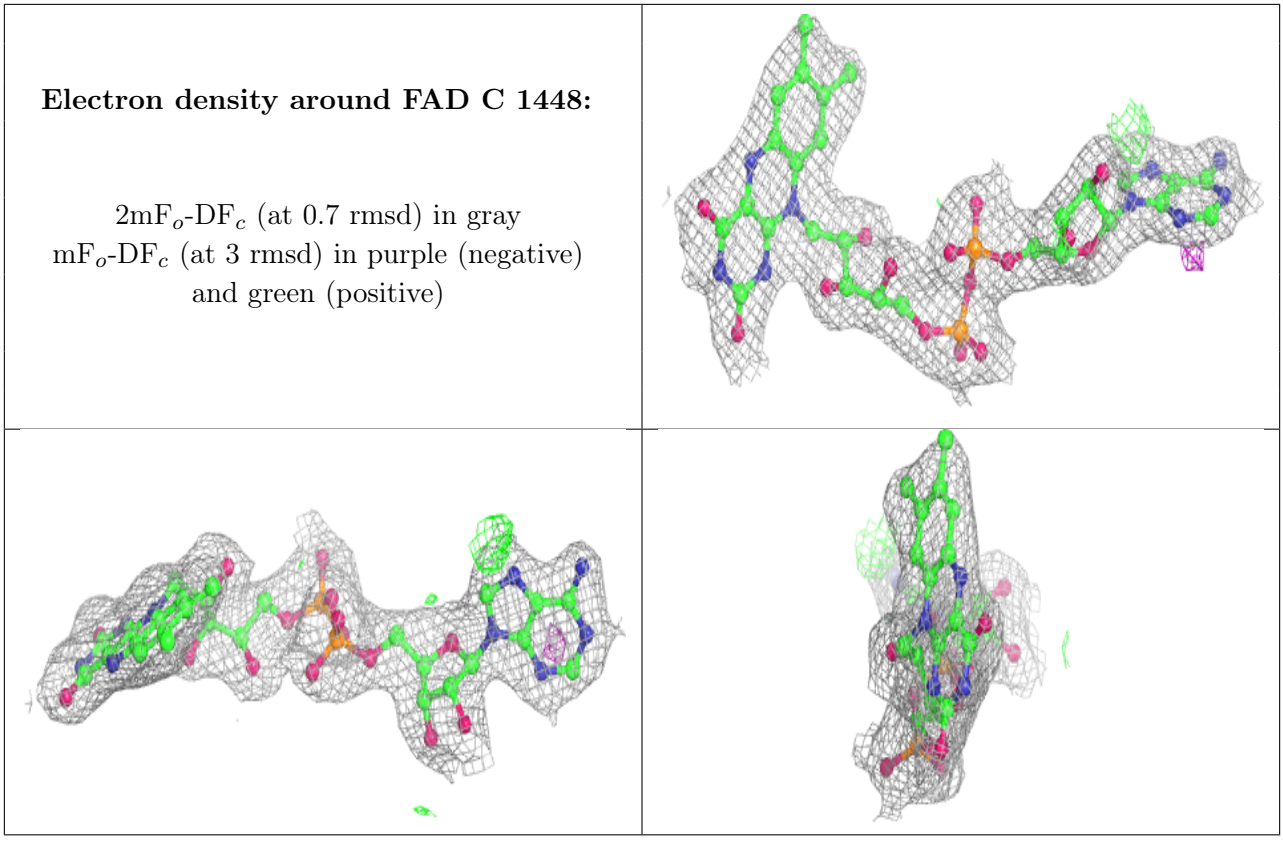






Electron density around FAD C 1448:

$2mF_o$-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.