



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 12, 2024 – 11:22 pm BST

PDB ID : 2XVF
Title : Crystal structure of bacterial flavin-containing monooxygenase
Authors : Cho, H.J.; Kang, B.S.
Deposited on : 2010-10-26
Resolution : 2.40 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 : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

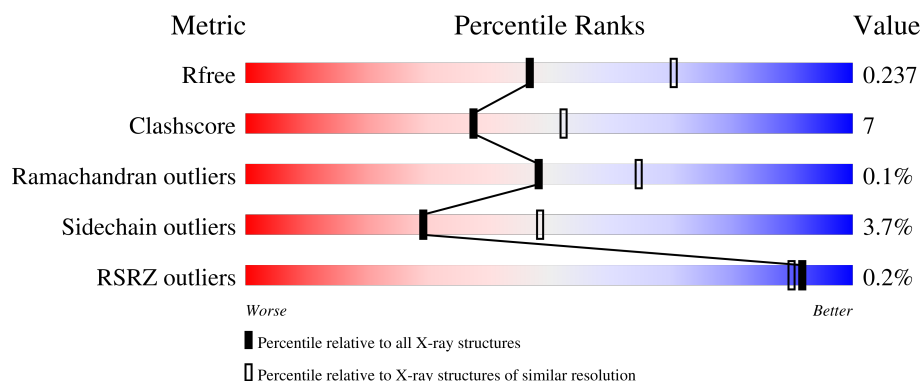
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION




The reported resolution of this entry is 2.40 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}	164625	4642 (2.40-2.40)
Clashscore	180529	5218 (2.40-2.40)
Ramachandran outliers	177936	5158 (2.40-2.40)
Sidechain outliers	177891	5159 (2.40-2.40)
RSRZ outliers	164620	4642 (2.40-2.40)

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 $\leq 5\%$. 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	 82% 13% . .
1	B	464	 81% 13% . .
1	C	464	 77% 18% . .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11219 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	445	Total	C	N	O	S	Se	0	0	0
			3608	2321	597	670	7	13			
1	B	444	Total	C	N	O	S	Se	0	2	0
			3626	2330	604	672	7	13			
1	C	445	Total	C	N	O	S	Se	0	0	0
			3587	2298	591	678	7	13			

There are 24 discrepancies between the modelled and reference sequences:

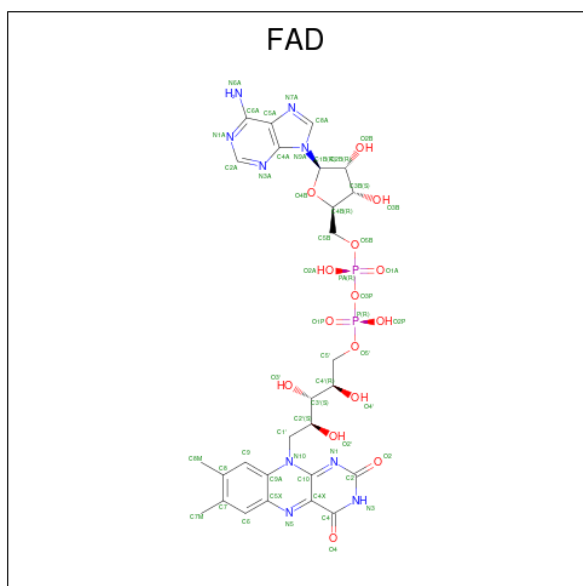
Chain	Residue	Modelled	Actual	Comment	Reference
A	457	LEU	-	expression tag	UNP Q83XK4
A	458	GLU	-	expression tag	UNP Q83XK4
A	459	HIS	-	expression tag	UNP Q83XK4
A	460	HIS	-	expression tag	UNP Q83XK4
A	461	HIS	-	expression tag	UNP Q83XK4
A	462	HIS	-	expression tag	UNP Q83XK4
A	463	HIS	-	expression tag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

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Chain	Residue	Modelled	Actual	Comment	Reference
C	464	HIS	-	expression tag	UNP Q83XK4

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $\text{C}_{27}\text{H}_{33}\text{N}_9\text{O}_{15}\text{P}_2$).



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

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



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		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		

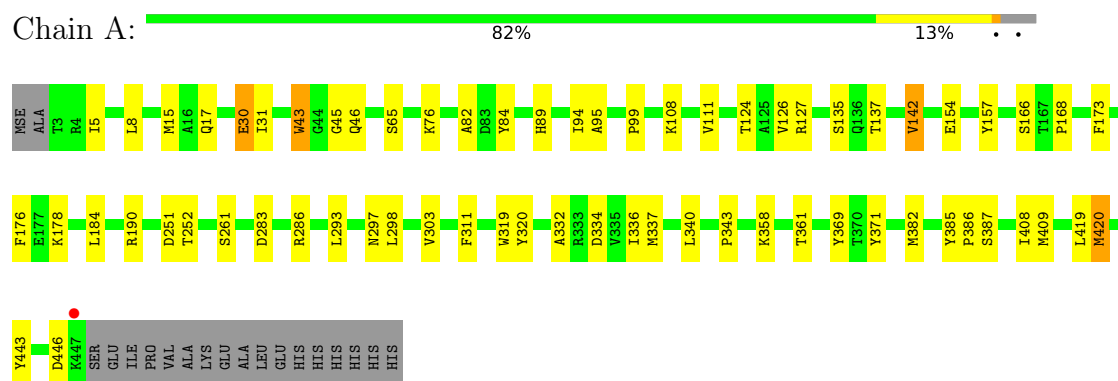
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	75	Total 75	O 75	0	0
5	B	108	Total 108	O 108	0	0
5	C	42	Total 42	O 42	0	0

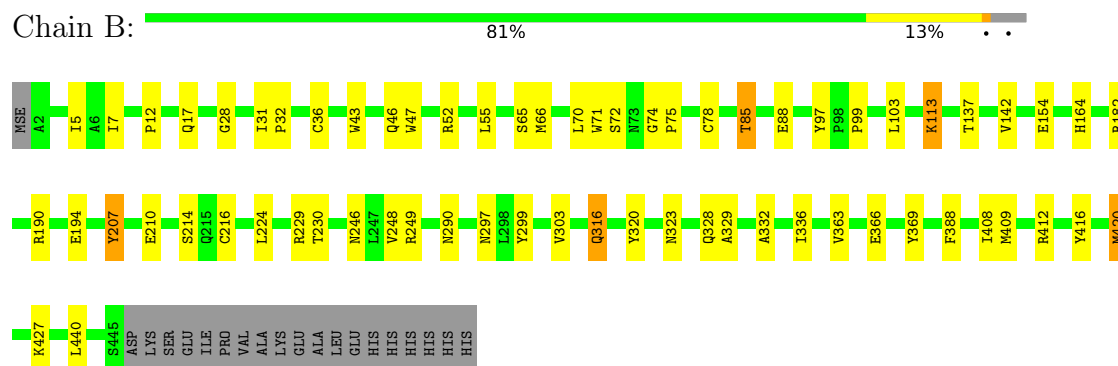
3 Residue-property plots

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



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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	158.62Å 70.08Å 140.96Å 90.00° 90.71° 90.00°	Depositor
Resolution (Å)	34.38 – 2.40 34.38 – 2.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (34.38-2.40) 84.7 (34.38-2.40)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	44.13 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.184 , 0.236 0.187 , 0.237	Depositor DCC
R_{free} test set	2627 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	41.2	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 33.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.021 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11219	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% 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, GOL, FAD

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.54	1/3709 (0.0%)	0.59	1/5023 (0.0%)
1	B	0.51	0/3734	0.59	0/5055
1	C	0.50	0/3688	0.55	0/5004
All	All	0.52	1/11131 (0.0%)	0.58	1/15082 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	320	TYR	CD1-CE1	-6.02	1.30	1.39

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	190	ARG	NE-CZ-NH2	-5.41	117.60	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3608	0	3356	43	0
1	B	3626	0	3393	49	0
1	C	3587	0	3274	56	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	53	0	31	1	0
2	B	53	0	31	1	0
2	C	53	0	31	3	0
3	A	4	0	0	0	0
3	B	4	0	0	0	0
4	B	6	0	8	0	0
5	A	75	0	0	0	0
5	B	108	0	0	3	0
5	C	42	0	0	1	0
All	All	11219	0	10124	146	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 7.

All (146) 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:388:PHE:HB3	1:B:420:MSE:HE1	1.46	0.96
1:C:305:GLU:HG2	1:C:348:MSE:HE2	1.48	0.96
1:B:388:PHE:CB	1:B:420:MSE:HE1	1.96	0.95
1:A:46:GLN:NE2	1:A:65:SER:H	1.65	0.94
1:A:31:ILE:HG22	1:A:337:MSE:CE	1.99	0.93
1:C:299:TYR:CB	1:C:348:MSE:HE3	1.98	0.92
1:B:46:GLN:HE22	1:B:65:SER:H	1.13	0.92
1:C:46:GLN:NE2	1:C:65:SER:H	1.67	0.90
1:A:31:ILE:HG22	1:A:337:MSE:HE1	1.54	0.89
1:A:46:GLN:HE22	1:A:65:SER:N	1.69	0.89
1:A:82:ALA:HA	1:A:420:MSE:HE3	1.54	0.89
1:B:85:THR:HG22	1:B:88:GLU:H	1.40	0.87
1:C:299:TYR:HB3	1:C:348:MSE:HE3	1.54	0.87
1:C:293:LEU:HD11	1:C:368:MSE:HG3	1.62	0.81
1:A:15:MSE:HE2	1:A:111:VAL:HG11	1.64	0.79
1:B:46:GLN:NE2	1:B:65:SER:H	1.84	0.74
1:A:30:GLU:O	1:A:30:GLU:HG3	1.86	0.73
1:C:251:ASP:OD1	1:C:253:GLU:N	2.23	0.70
1:A:46:GLN:HE22	1:A:65:SER:H	0.84	0.70
1:B:316:GLN:NE2	1:B:328:GLN:HE22	1.90	0.70
1:C:46:GLN:HE21	1:C:65:SER:H	1.37	0.69
1:C:364:THR:HG22	1:C:366:GLU:H	1.59	0.68
1:B:5:ILE:HD11	1:B:336:ILE:HG21	1.75	0.68
1:C:299:TYR:HB2	1:C:348:MSE:HE3	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:LYS:HA	1:A:111:VAL:HG12	1.77	0.66
1:C:251:ASP:OD1	1:C:251:ASP:C	2.32	0.65
1:A:8:LEU:HD22	1:A:126:VAL:HG21	1.78	0.65
1:C:251:ASP:OD1	1:C:252:THR:N	2.29	0.65
1:C:43:TRP:CH2	1:C:111:VAL:HG12	2.32	0.65
1:A:76:LYS:HE2	1:A:94:ILE:HG13	1.79	0.63
1:B:72:SER:HB2	1:B:103:LEU:HD11	1.80	0.63
1:B:440:LEU:C	1:B:440:LEU:HD23	2.19	0.62
1:C:65:SER:OG	1:C:187:HIS:HD2	1.82	0.61
1:B:388:PHE:HB2	1:B:420:MSE:HE1	1.82	0.60
1:B:52:ARG:O	1:B:190:ARG:NH1	2.34	0.59
1:C:104:TRP:O	1:C:108:LYS:HG2	2.03	0.59
1:C:43:TRP:CZ3	1:C:111:VAL:HG12	2.38	0.59
1:C:80:GLU:HG2	1:C:86:PHE:CD2	2.38	0.59
1:A:31:ILE:HG22	1:A:337:MSE:HE3	1.82	0.59
1:A:251:ASP:OD1	1:A:252:THR:N	2.36	0.59
1:A:45:GLY:HA2	2:A:1448:FAD:O3B	2.03	0.58
1:A:283:ASP:OD1	1:A:286:ARG:NH1	2.37	0.58
1:B:113:LYS:N	1:B:113:LYS:HD3	2.19	0.57
1:C:135:SER:OG	1:C:137:THR:HG22	2.04	0.57
1:C:358:LYS:O	1:C:361:THR:HB	2.04	0.57
1:B:229:ARG:HG2	1:B:246:ASN:HB3	1.86	0.56
1:B:5:ILE:CD1	1:B:336:ILE:HG21	2.35	0.56
1:A:5:ILE:HD11	1:A:336:ILE:HG21	1.87	0.56
1:C:22:GLN:HG2	1:C:114:ALA:HB1	1.88	0.56
1:C:15:MSE:HB3	1:C:111:VAL:HG21	1.88	0.56
1:C:80:GLU:HG2	1:C:86:PHE:CE2	2.40	0.56
1:C:75:PRO:HB3	1:C:416:TYR:CD2	2.41	0.55
1:A:15:MSE:HE3	1:A:43:TRP:CG	2.42	0.55
1:A:124:THR:CG2	1:A:142:VAL:HG13	2.37	0.55
1:A:251:ASP:OD1	1:A:251:ASP:C	2.43	0.55
1:C:201:LEU:HD11	1:C:227:CYS:SG	2.47	0.54
1:C:73:ASN:HB2	2:C:1448:FAD:O4	2.08	0.54
1:A:17:GLN:HE22	1:A:332:ALA:HB3	1.72	0.54
1:B:408:ILE:HG23	1:B:409:MSE:HE3	1.90	0.54
1:A:168:PRO:HD2	1:B:55:LEU:HD12	1.90	0.53
1:C:299:TYR:HB2	1:C:348:MSE:CE	2.37	0.53
1:C:43:TRP:CZ3	1:C:111:VAL:CG1	2.92	0.52
1:C:85:THR:HG23	1:C:88:GLU:H	1.74	0.52
1:B:71:TRP:HA	1:B:99:PRO:HA	1.92	0.52
1:A:124:THR:CG2	1:A:142:VAL:CG1	2.89	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:ILE:HG23	1:A:409:MSE:HE2	1.92	0.51
1:C:208:SER:HB3	1:C:212:ILE:HG13	1.92	0.51
1:A:334:ASP:HB3	1:A:340:LEU:HD13	1.93	0.51
1:C:46:GLN:HE22	1:C:65:SER:H	1.56	0.51
1:C:376:ILE:HG22	1:C:380:ILE:HD12	1.93	0.51
1:A:293:LEU:HD21	1:A:319:TRP:CD1	2.45	0.50
1:A:135:SER:OG	1:A:137:THR:HG22	2.11	0.50
1:C:73:ASN:HB2	2:C:1448:FAD:C4	2.43	0.49
1:B:207:TYR:CD1	1:B:207:TYR:C	2.86	0.49
1:C:324:MSE:O	1:C:328:GLN:HG3	2.13	0.49
1:B:290[B]:ASN:HD22	1:B:290[B]:ASN:H	1.61	0.48
1:A:99:PRO:HD3	1:A:443:TYR:CE1	2.47	0.48
1:A:358:LYS:HB3	1:A:371:TYR:HE1	1.78	0.48
1:B:366:GLU:HA	1:B:369:TYR:CE2	2.48	0.48
1:B:17:GLN:HA	1:B:329:ALA:HB1	1.96	0.48
1:B:72:SER:OG	1:B:74:GLY:O	2.30	0.47
1:B:214:SER:OG	1:B:409:MSE:HE1	2.14	0.47
1:C:71:TRP:HA	1:C:99:PRO:HA	1.95	0.47
1:A:5:ILE:CD1	1:A:336:ILE:HG21	2.45	0.47
1:B:12:PRO:HG2	2:B:1446:FAD:H4'	1.97	0.47
1:B:388:PHE:HB3	1:B:420:MSE:CE	2.32	0.47
1:A:124:THR:HG23	1:A:142:VAL:HG13	1.98	0.46
1:C:19:ARG:NH1	1:C:22:GLN:NE2	2.63	0.46
1:B:207:TYR:C	1:B:207:TYR:HD1	2.17	0.46
1:B:299:TYR:HB3	1:B:303:VAL:HB	1.97	0.46
1:C:305:GLU:OE2	1:C:352:SER:OG	2.34	0.46
1:B:31:ILE:HG22	1:B:32:PRO:O	2.15	0.46
1:B:164:HIS:HD2	5:B:2045:HOH:O	1.99	0.46
1:A:76:LYS:HE3	1:A:95:ALA:O	2.16	0.46
1:A:166:SER:O	1:A:168:PRO:HD3	2.16	0.46
1:A:385:TYR:CD1	1:A:386:PRO:HD2	2.51	0.46
1:B:47:TRP:CE2	1:B:66:MSE:HE1	2.51	0.45
1:B:137:THR:HG23	1:B:154:GLU:OE1	2.17	0.45
1:C:201:LEU:HB2	1:C:265:VAL:HG21	1.99	0.45
1:A:127:ARG:HD3	1:A:127:ARG:HA	1.78	0.45
1:C:251:ASP:CG	1:C:252:THR:N	2.70	0.45
1:C:46:GLN:HG3	1:C:66:MSE:HG3	1.99	0.45
1:A:358:LYS:O	1:A:361:THR:HB	2.17	0.45
1:B:412:ARG:NH1	5:B:2097:HOH:O	2.48	0.45
1:C:262:SER:O	1:C:263:GLU:HG2	2.17	0.44
1:C:12:PRO:HG2	2:C:1448:FAD:H4'	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:277:HIS:NE2	1:C:279:PRO:HG3	2.32	0.44
1:B:408:ILE:HG23	1:B:409:MSE:CE	2.47	0.44
1:C:159:VAL:CG1	1:C:313:ILE:HD13	2.47	0.44
1:A:319:TRP:HE3	1:A:369:TYR:CZ	2.36	0.44
1:A:303:VAL:HG22	1:A:311:PHE:CE2	2.53	0.44
1:B:28:GLY:HA2	1:C:296:LEU:HD11	1.98	0.44
1:B:210:GLU:HG2	5:B:2061:HOH:O	2.18	0.44
1:B:17:GLN:HE22	1:B:332:ALA:HB3	1.83	0.43
1:B:248:VAL:HG12	1:B:249:ARG:HG2	2.01	0.43
1:A:298:LEU:HA	1:A:303:VAL:O	2.18	0.43
1:C:251:ASP:OD1	1:C:254:ASN:N	2.38	0.43
1:C:13:SER:OG	1:C:313:ILE:HG22	2.18	0.43
1:C:245:PRO:HB2	1:C:258:ALA:HB3	2.01	0.43
1:B:320:TYR:HD2	1:B:323:ASN:HD22	1.65	0.43
1:A:82:ALA:CA	1:A:420:MSE:HE3	2.39	0.43
1:C:75:PRO:HG2	1:C:78:CYS:SG	2.59	0.43
1:B:31:ILE:HD13	1:B:31:ILE:HA	1.86	0.42
1:B:97:TYR:CD2	1:B:412:ARG:HA	2.53	0.42
1:A:340:LEU:HD21	1:A:382:MSE:HB3	2.01	0.42
1:C:60:GLU:HG3	5:C:2004:HOH:O	2.18	0.42
1:C:417:ARG:HA	1:C:424:MSE:HA	2.01	0.42
1:A:173:PHE:HB2	1:A:176:PHE:CE1	2.55	0.42
1:A:5:ILE:HG12	1:A:157:TYR:HB2	2.01	0.42
1:B:46:GLN:HE22	1:B:65:SER:N	1.97	0.42
1:B:75:PRO:HG2	1:B:78:CYS:SG	2.60	0.41
1:C:3:THR:HA	1:C:156:ASP:OD2	2.20	0.41
1:B:182:ARG:NH2	1:B:194:GLU:OE1	2.41	0.41
1:C:95:ALA:HB3	1:C:443:TYR:OH	2.21	0.41
1:C:111:VAL:HG13	1:C:116:VAL:CG2	2.50	0.41
1:B:7:ILE:O	1:B:36:CYS:HA	2.20	0.41
1:B:97:TYR:CE2	1:B:412:ARG:HA	2.56	0.41
1:C:5:ILE:HG13	1:C:157:TYR:HB2	2.02	0.41
1:C:64:SER:O	1:C:190:ARG:NH2	2.54	0.41
1:B:216:CYS:HB2	1:B:224:LEU:HD21	2.02	0.41
1:B:70:LEU:C	1:B:70:LEU:HD23	2.41	0.41
1:B:75:PRO:HB3	1:B:416:TYR:CD1	2.56	0.41
1:C:124:THR:HG23	1:C:142:VAL:HG13	2.03	0.41
1:B:47:TRP:CZ2	1:B:66:MSE:HE1	2.57	0.40
1:C:94:ILE:HD12	1:C:443:TYR:HE2	1.86	0.40
1:A:84:TYR:CE2	1:A:89:HIS:HB2	2.57	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%)	423 (96%)	19 (4%)	1 (0%)	44	59
1	B	444/464 (96%)	429 (97%)	15 (3%)	0	100	100
1	C	443/464 (96%)	419 (95%)	24 (5%)	0	100	100
All	All	1330/1392 (96%)	1271 (96%)	58 (4%)	1 (0%)	48	65

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	343	PRO

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	373/391 (95%)	361 (97%)	12 (3%)	34	54
1	B	379/391 (97%)	368 (97%)	11 (3%)	37	58
1	C	369/391 (94%)	351 (95%)	18 (5%)	21	36
All	All	1121/1173 (96%)	1080 (96%)	41 (4%)	29	48

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	GLU
1	A	43	TRP

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Mol	Chain	Res	Type
1	A	142	VAL
1	A	154	GLU
1	A	178	LYS
1	A	184	LEU
1	A	261	SER
1	A	297	ASN
1	A	387	SER
1	A	419	LEU
1	A	420	MSE
1	A	446	ASP
1	B	43	TRP
1	B	85	THR
1	B	113	LYS
1	B	142	VAL
1	B	207	TYR
1	B	230	THR
1	B	297	ASN
1	B	316	GLN
1	B	363	VAL
1	B	420	MSE
1	B	427	LYS
1	C	3	THR
1	C	4	ARG
1	C	43	TRP
1	C	68	ARG
1	C	142	VAL
1	C	184	LEU
1	C	193	LEU
1	C	205	SER
1	C	227	CYS
1	C	246	ASN
1	C	272	THR
1	C	297	ASN
1	C	330	TRP
1	C	346	GLU
1	C	387	SER
1	C	402	HIS
1	C	407	ASN
1	C	419	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	40	GLN
1	A	46	GLN
1	A	48	ASN
1	A	73	ASN
1	A	282	ASN
1	B	17	GLN
1	B	40	GLN
1	B	46	GLN
1	B	48	ASN
1	B	164	HIS
1	B	297	ASN
1	B	316	GLN
1	B	323	ASN
1	B	377	GLN
1	C	17	GLN
1	C	22	GLN
1	C	40	GLN
1	C	46	GLN
1	C	48	ASN
1	C	73	ASN
1	C	132	ASN
1	C	187	HIS
1	C	403	HIS
1	C	407	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	GOL	B	1448	-	5,5,5	0.35	0	5,5,5	0.21	0
3	NO3	A	1449	-	1,3,3	3.59	1 (100%)	0,3,3	-	-
2	FAD	A	1448	-	53,58,58	1.13	3 (5%)	68,89,89	1.43	9 (13%)
2	FAD	B	1446	-	53,58,58	1.11	4 (7%)	68,89,89	1.43	10 (14%)
3	NO3	B	1447	-	1,3,3	3.59	1 (100%)	0,3,3	-	-
2	FAD	C	1448	-	53,58,58	1.14	4 (7%)	68,89,89	1.37	8 (11%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FAD	B	1446	-	-	3/30/50/50	0/6/6/6
2	FAD	C	1448	-	-	3/30/50/50	0/6/6/6
4	GOL	B	1448	-	-	2/4/4/4	-
2	FAD	A	1448	-	-	3/30/50/50	0/6/6/6

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1448	FAD	C4X-N5	4.21	1.38	1.30
2	C	1448	FAD	C4X-N5	3.98	1.38	1.30
2	C	1448	FAD	C2A-N3A	3.77	1.38	1.32
2	B	1446	FAD	C2A-N3A	3.65	1.38	1.32
3	B	1447	NO3	O1-N	3.59	1.40	1.24
3	A	1449	NO3	O1-N	3.59	1.40	1.24
2	B	1446	FAD	C4X-N5	3.35	1.37	1.30
2	A	1448	FAD	C2A-N3A	3.34	1.37	1.32
2	B	1446	FAD	C10-N1	2.79	1.38	1.33
2	C	1448	FAD	C10-N1	2.54	1.38	1.33
2	A	1448	FAD	C10-N1	2.26	1.37	1.33
2	C	1448	FAD	C2A-N1A	2.25	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1446	FAD	C2A-N1A	2.20	1.38	1.33

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1448	FAD	N3A-C2A-N1A	-6.25	118.92	128.68
2	B	1446	FAD	N3A-C2A-N1A	-6.02	119.28	128.68
2	C	1448	FAD	N3A-C2A-N1A	-5.68	119.80	128.68
2	B	1446	FAD	C4-N3-C2	-3.30	119.54	125.64
2	A	1448	FAD	C4X-C10-N10	3.17	121.11	116.48
2	C	1448	FAD	C4-N3-C2	-3.16	119.80	125.64
2	C	1448	FAD	C4X-C10-N10	3.00	120.86	116.48
2	A	1448	FAD	C4-N3-C2	-2.88	120.33	125.64
2	B	1446	FAD	C4X-C4-N3	2.78	120.26	113.19
2	C	1448	FAD	C4X-C4-N3	2.70	120.06	113.19
2	A	1448	FAD	P-O3P-PA	-2.67	123.65	132.83
2	B	1446	FAD	C4X-C10-N1	-2.63	118.62	124.73
2	A	1448	FAD	C10-C4X-N5	-2.58	119.39	124.86
2	B	1446	FAD	C4X-C10-N10	2.49	120.12	116.48
2	C	1448	FAD	C4X-C10-N1	-2.47	119.00	124.73
2	C	1448	FAD	P-O3P-PA	-2.43	124.47	132.83
2	A	1448	FAD	C4X-C4-N3	2.40	119.29	113.19
2	C	1448	FAD	C10-C4X-N5	-2.38	119.81	124.86
2	B	1446	FAD	C10-C4X-N5	-2.38	119.82	124.86
2	B	1446	FAD	P-O3P-PA	-2.36	124.72	132.83
2	A	1448	FAD	C4X-C10-N1	-2.24	119.53	124.73
2	C	1448	FAD	O4-C4-C4X	-2.18	120.83	126.60
2	B	1446	FAD	C4-C4X-C10	2.17	120.44	116.79
2	B	1446	FAD	O4-C4-C4X	-2.14	120.93	126.60
2	B	1446	FAD	O2'-C2'-C1'	-2.13	104.65	109.80
2	A	1448	FAD	C5X-C9A-N10	2.12	120.15	117.95
2	A	1448	FAD	C4A-C5A-N7A	-2.05	107.26	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	B	1446	FAD	N10-C1'-C2'-O2'
2	B	1446	FAD	N10-C1'-C2'-C3'
2	C	1448	FAD	N10-C1'-C2'-O2'

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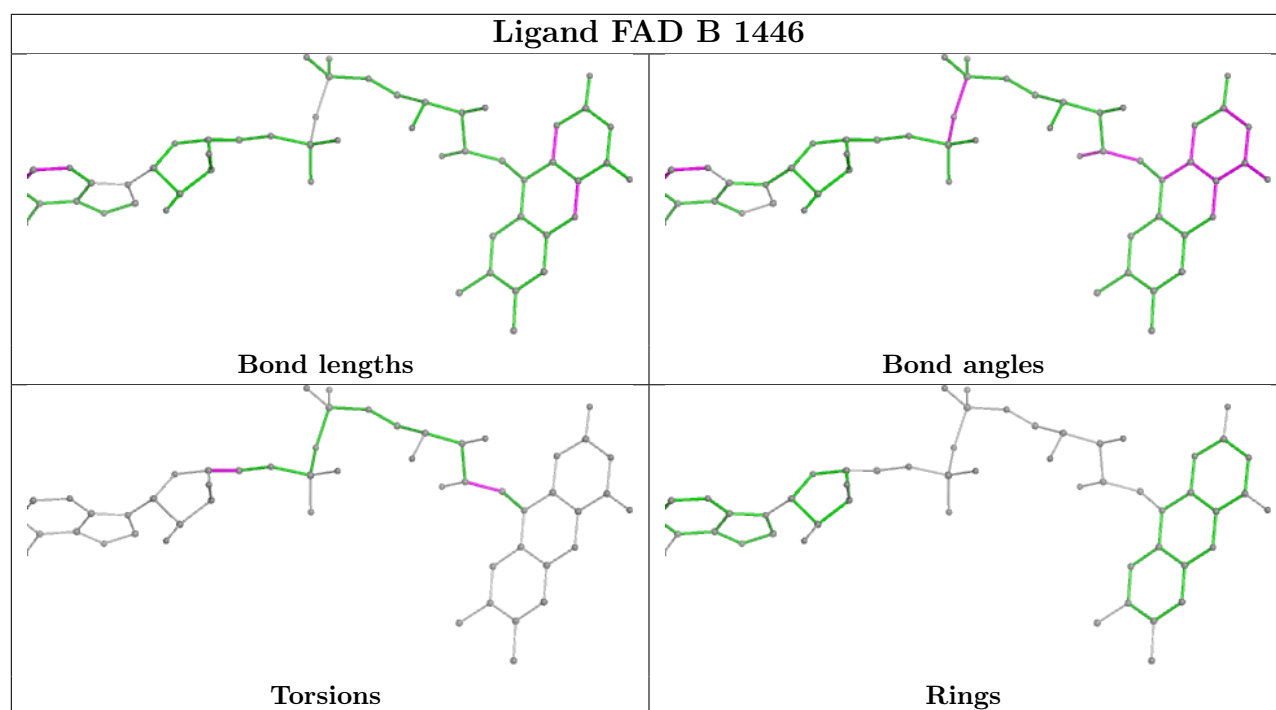
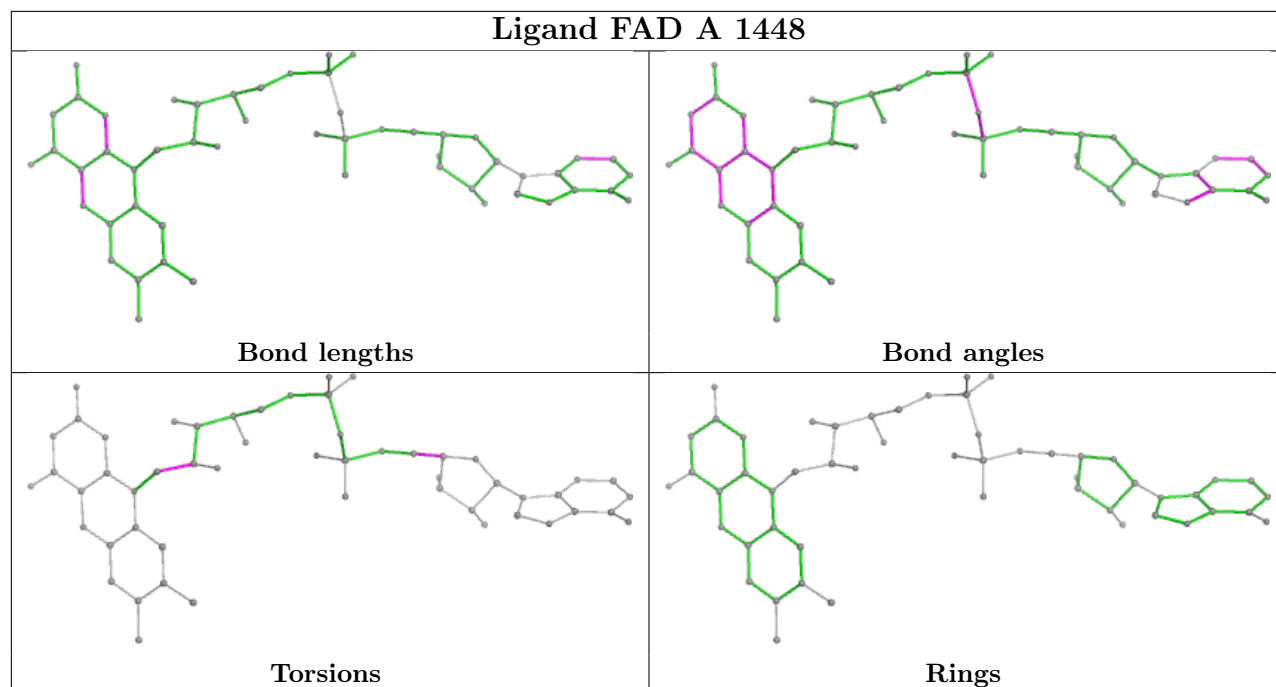
Mol	Chain	Res	Type	Atoms
2	C	1448	FAD	N10-C1'-C2'-C3'
4	B	1448	GOL	O1-C1-C2-C3
4	B	1448	GOL	O1-C1-C2-O2
2	A	1448	FAD	O4B-C4B-C5B-O5B
2	B	1446	FAD	O4B-C4B-C5B-O5B
2	C	1448	FAD	O4B-C4B-C5B-O5B

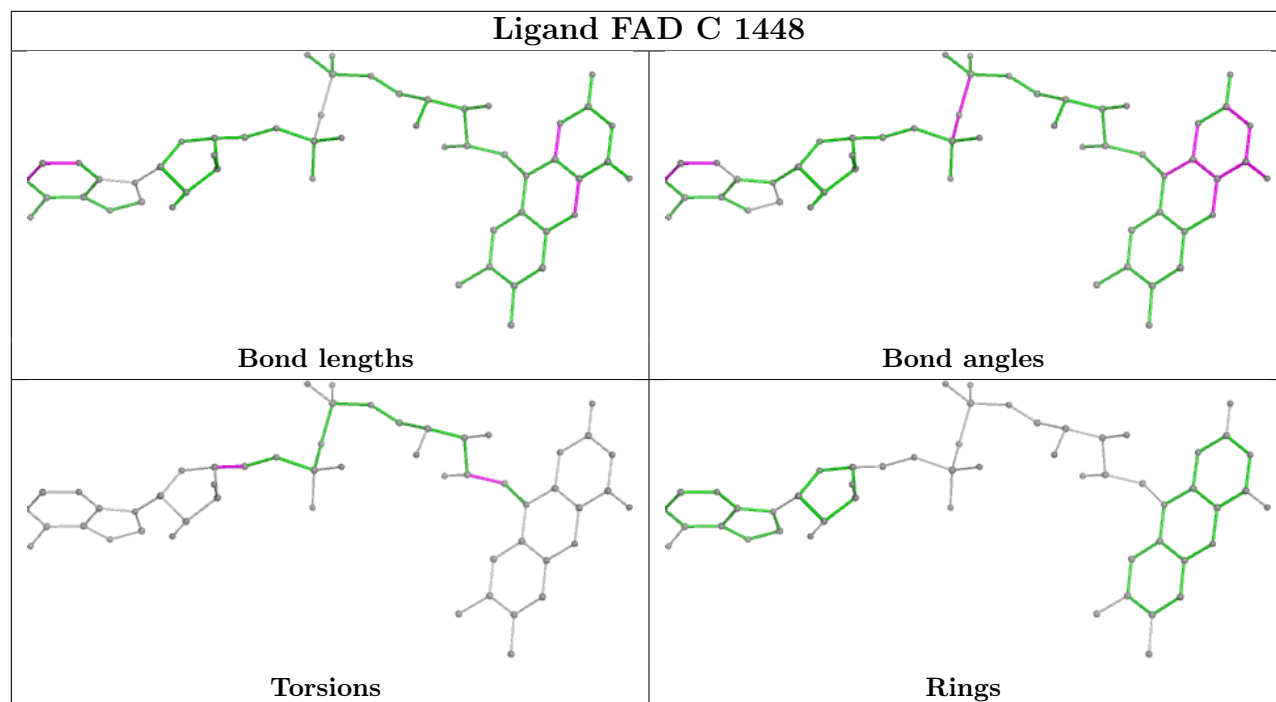
There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1448	FAD	1	0
2	B	1446	FAD	1	0
2	C	1448	FAD	3	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	432/464 (93%)	-0.46	1 (0%) 92 90	24, 45, 68, 85	0
1	B	431/464 (92%)	-0.58	0 100 100	21, 40, 60, 76	2 (0%)
1	C	432/464 (93%)	-0.18	1 (0%) 92 90	34, 55, 95, 113	0
All	All	1295/1392 (93%)	-0.41	2 (0%) 92 90	21, 46, 83, 113	2 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	447	LYS	4.1
1	A	447	LYS	2.6

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

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

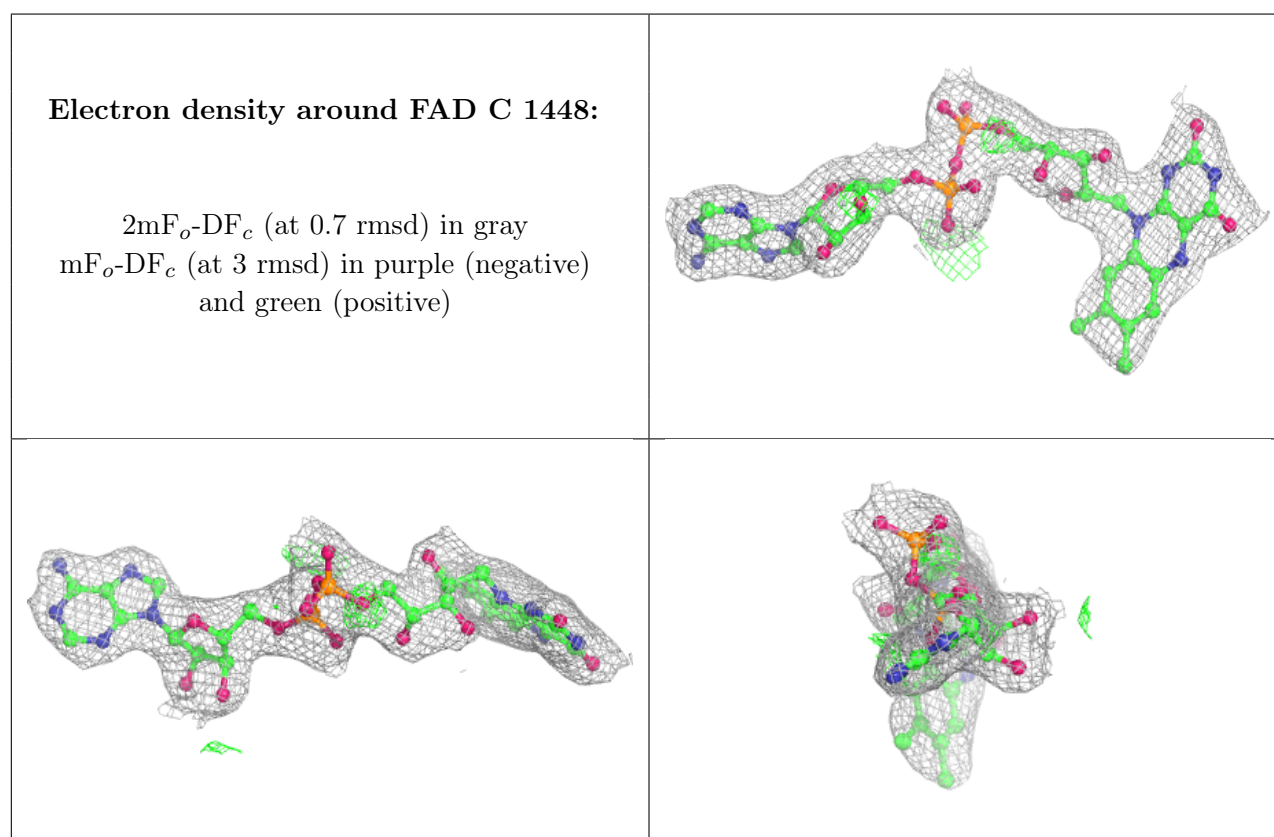
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	B	1448	6/6	0.85	0.25	93,93,94,94	0

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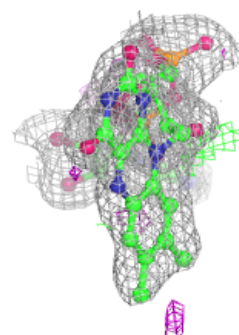
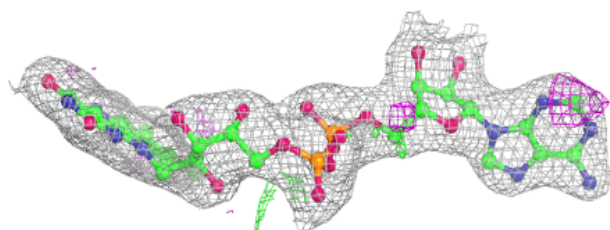
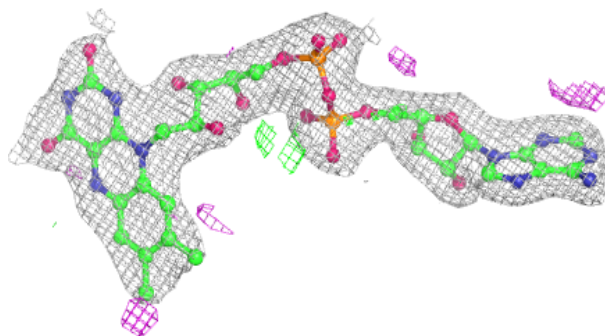
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NO3	A	1449	4/4	0.92	0.18	74,75,75,75	0
3	NO3	B	1447	4/4	0.93	0.12	59,60,60,60	0
2	FAD	C	1448	53/53	0.96	0.07	40,44,51,52	0
2	FAD	A	1448	53/53	0.98	0.07	31,35,39,42	0
2	FAD	B	1446	53/53	0.98	0.06	31,33,34,36	0

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.

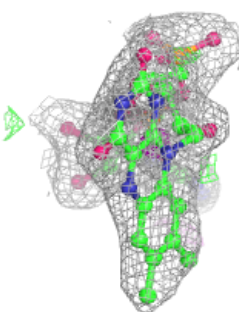
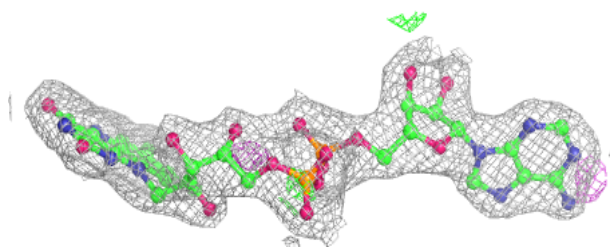
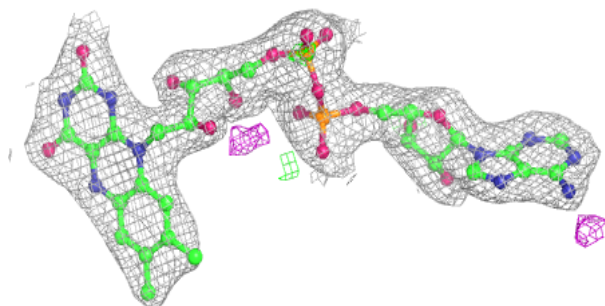


Electron density around FAD A 1448:

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

**Electron density around FAD B 1446:**

$2mF_o-DF_c$ (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.