

wwPDB X-ray Structure Validation Summary Report (i)

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 wwPDB X-ray Structure Validation Summary 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 (i) 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 (1)) were used in the production of this report:

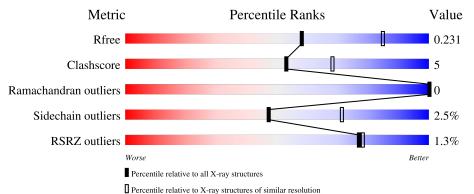
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
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: $X\text{-}RAY \, DIFFRACTION$

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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
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	А	464	81%	15%	·
1	В	464	81%	14%	·
1	С	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	А	1449	-	-	Х	-
3	NO3	В	1448	-	-	Х	-



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.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	А	446	Total	С	Ν	0	\mathbf{S}	0	1	0
	A	440	3634	2330	598	686	20	0		
1	Р	445	Total	С	Ν	0	S	0	1	0
			3611	2317	596	678	20			
1	1 C	446	Total	С	Ν	0	S	0	0	0
			3612	2317	595	680	20			0

• Molecule 1 is a protein called FLAVIN-CONTAINING MONOOXYGENASE.

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

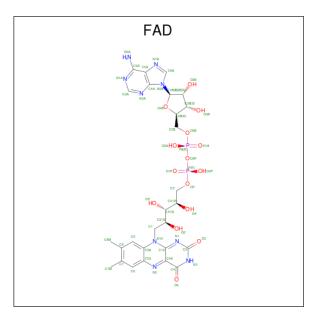
There are 27 discrepancies between the modelled and reference sequences:



Continu	Continued from previous page									
Chain	Residue	Modelled	Actual	Comment	Reference					
С	462	HIS	-	expression tag	UNP Q83XK4					
С	463	HIS	-	expression tag	UNP Q83XK4					
С	464	HIS	-	expression tag	UNP Q83XK4					
С	207	SER	TYR	engineered mutation	UNP Q83XK4					

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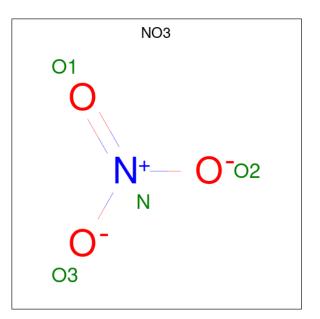
• Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	2 A	1	Total	С	Ν	Ο	Р	0	0
2		1	53	27	9	15	2	0	
2	2 B	1	Total	С	Ν	Ο	Р	0	0
2			53	27	9	15	2		
2	2 C	1 1	Total	С	Ν	Ο	Р	0	0
	U	1	53	27	9	15	2	0	0

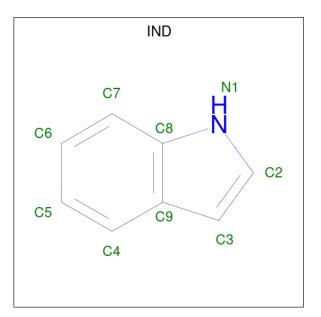
• Molecule 3 is NITRATE ION (three-letter code: NO3) (formula: NO_3).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	Total N O 4 1 3	0	0
3	В	1	$\begin{array}{ccc} \text{Total} & \text{N} & \text{O} \\ 4 & 1 & 3 \end{array}$	0	0
3	С	1	Total N O 4 1 3	0	0

 $\bullet\,$ Molecule 4 is INDOLE (three-letter code: IND) (formula: $\rm C_8H_7N).$



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

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

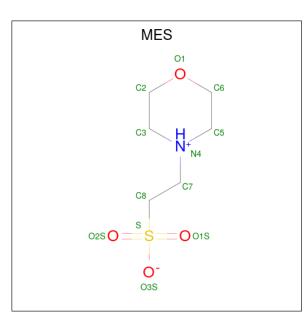
• Molecule 5 is OXYGEN MOLECULE (three-letter code: OXY) (formula: O_2).

OXY	
01 () () 02	

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	1	Total O 2 2	0	0
5	В	1	Total O 2 2	0	0
5	С	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	В	1	Total	С	Ν	0	S	0	0
0	D	1	12	6	1	4	1	0	0

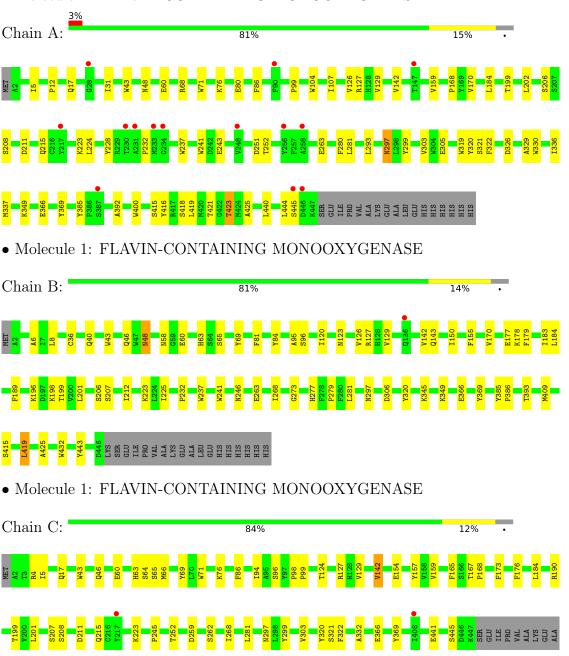
• Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	А	32	$\begin{array}{cc} \text{Total} & \text{O} \\ 32 & 32 \end{array}$	0	0
7	В	54	$\begin{array}{cc} \text{Total} & \text{O} \\ 54 & 54 \end{array}$	0	0
7	С	55	$\begin{array}{cc} \text{Total} & \text{O} \\ 55 & 55 \end{array}$	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

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4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	159.12Å 67.85 Å 138.90 Å	Depositor
a, b, c, α , β , γ	90.00° 90.62° 90.00°	Depositor
Resolution (Å)	50.01 - 2.48	Depositor
Resolution (A)	46.32 - 2.48	EDS
% Data completeness	97.3 (50.01-2.48)	Depositor
(in resolution range)	84.7(46.32-2.48)	EDS
R _{merge}	0.08	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$223.90 (at 2.48 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
P. P.	0.199 , 0.237	Depositor
R, R_{free}	0.194 , 0.231	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.

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.



¹Intensities estimated from amplitudes.

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: 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		Bo	nd lengths	Bond	angles
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	0.55	4/3751~(0.1%)	0.55	0/5100
1	В	0.50	2/3728~(0.1%)	0.53	0/5072
1	С	0.51	2/3725~(0.1%)	0.54	0/5068
All	All	0.52	8/11204 (0.1%)	0.54	0/15240

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\operatorname{Observed}(\operatorname{\AA})$	$\mathrm{Ideal}(\mathrm{\AA})$
1	С	320	TYR	CD2-CE2	-7.70	1.27	1.39
1	А	320	TYR	CD1-CE1	-7.63	1.27	1.39
1	С	320	TYR	CD1-CE1	-6.78	1.29	1.39
1	В	320	TYR	CD1-CE1	-6.36	1.29	1.39
1	В	320	TYR	CD2-CE2	-6.09	1.30	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	А	3634	0	3374	46	0
1	В	3611	0	3351	37	0
1	С	3612	0	3343	32	0



Mol	Chain	Non-H		H(added)	Clashes	Symm-Clashes
2	А	53	0	31	1	0
2	В	53	0	31	2	0
2	С	53	0	31	1	0
3	А	4	0	0	2	0
3	В	4	0	0	4	0
3	С	4	0	0	0	0
4	А	9	0	7	0	0
4	В	9	0	7	0	0
4	С	9	0	7	1	0
5	А	2	0	0	0	0
5	В	2	0	0	0	0
5	С	2	0	0	0	0
6	В	12	0	12	1	0
7	А	32	0	0	0	0
7	В	54	0	0	0	0
7	С	55	0	0	0	0
All	All	11214	0	10194	115	0

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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.

The worst 5 of 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

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.



Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	А	445/464~(96%)	430 (97%)	15 (3%)	0	100 100
1	В	444/464~(96%)	428 (96%)	16 (4%)	0	100 100
1	С	444/464 (96%)	427 (96%)	17 (4%)	0	100 100
All	All	1333/1392~(96%)	1285 (96%)	48 (4%)	0	100 100

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

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	А	381/405~(94%)	374~(98%)	7 (2%)	59 80		
1	В	377/405~(93%)	365~(97%)	12 (3%)	39 63		
1	С	375/405~(93%)	366~(98%)	9(2%)	49 72		
All	All	1133/1215 (93%)	1105 (98%)	28 (2%)	47 71		

5 of 28 residues with a non-rotameric side chain are listed below:

Mol	Chain	Res	Type
1	В	232	PRO
1	С	297	ASN
1	В	306	ASP
1	С	208	SER
1	В	297	ASN

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 22 such side chains are listed below:

Mol	Chain	Res	Type
1	В	277	HIS
1	С	46	GLN
1	С	17	GLN



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Mol	Chain	Res	Type
1	С	89	HIS
1	А	297	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 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	Turne	Chain	Res	Link	В	ond leng	gths	E	ond ang	gles
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2
2	FAD	А	1448	-	$53,\!58,\!58$	1.12	4 (7%)	68,89,89	1.38	11 (16%)
4	IND	В	1449	-	8,10,10	1.01	0	9,13,13	2.28	2 (22%)
2	FAD	С	1448	-	$53,\!58,\!58$	1.12	4 (7%)	68,89,89	1.40	11 (16%)
2	FAD	В	1447	-	$53,\!58,\!58$	1.10	4 (7%)	68,89,89	1.46	8 (11%)
5	OXY	А	1451	-	$1,\!1,\!1$	0.13	0	-		
5	OXY	С	1451	-	$1,\!1,\!1$	0.15	0	-		
4	IND	А	1450	-	8,10,10	1.05	0	9,13,13	2.32	2 (22%)
3	NO3	А	1449	-	$1,\!3,\!3$	3.49	1 (100%)	0,3,3	-	-
6	MES	В	1450	-	12,12,12	2.20	1 (8%)	14,16,16	2.37	6 (42%)
3	NO3	С	1449	-	$1,\!3,\!3$	<mark>3.53</mark>	1 (100%)	0,3,3	-	-



Mal	Mol Type Chain		Res Linl		Bond lengths			Bond angles		
IVIOI	Type	Chain	nes	Link	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
4	IND	С	1450	-	8,10,10	1.02	0	9,13,13	2.32	2 (22%)
3	NO3	В	1448	-	$1,\!3,\!3$	2.29	1 (100%)	0,3,3	-	-
5	OXY	В	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	А	1448	-	-	4/30/50/50	0/6/6/6
4	IND	В	1449	-	-	-	0/2/2/2
2	FAD	С	1448	-	-	3/30/50/50	0/6/6/6
2	FAD	В	1447	-	-	3/30/50/50	0/6/6/6
4	IND	А	1450	-	-	-	0/2/2/2
6	MES	В	1450	-	-	1/6/14/14	0/1/1/1
4	IND	С	1450	-	-	_	0/2/2/2

The worst 5 of 16 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\operatorname{Observed}(\operatorname{\AA})$	$\mathrm{Ideal}(\mathrm{\AA})$
6	В	1450	MES	C8-S	-7.28	1.67	1.77
2	А	1448	FAD	C4X-N5	4.22	1.39	1.30
2	В	1447	FAD	C4X-N5	3.90	1.38	1.30
2	С	1448	FAD	C2A-N3A	3.89	1.38	1.32
2	С	1448	FAD	C4X-N5	3.83	1.38	1.30

The worst 5 of 42 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	В	1447	FAD	N3A-C2A-N1A	-6.70	118.21	128.68
4	С	1450	IND	C3-C9-C8	6.28	111.56	106.20
4	А	1450	IND	C3-C9-C8	6.26	111.55	106.20
4	В	1449	IND	C3-C9-C8	6.12	111.43	106.20
2	С	1448	FAD	N3A-C2A-N1A	-5.63	119.88	128.68

There are no chirality outliers.

5 of 11 torsion outliers are listed below:



Mol	Chain	Res	Type	Atoms
2	А	1448	FAD	N10-C1'-C2'-O2'
2	А	1448	FAD	N10-C1'-C2'-C3'
2	А	1448	FAD	PA-O3P-P-O5'
2	В	1447	FAD	N10-C1'-C2'-O2'
2	В	1447	FAD	N10-C1'-C2'-C3'

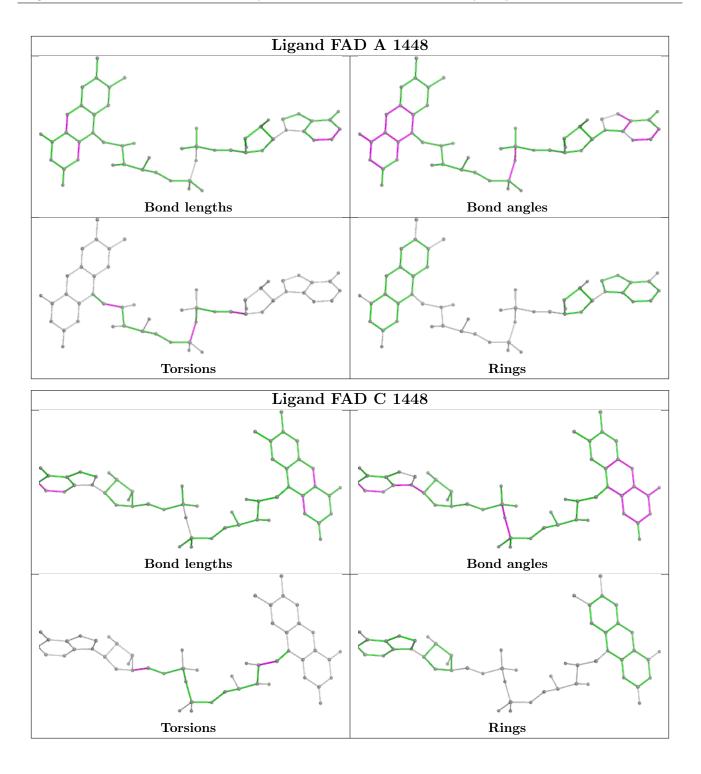
There are no ring outliers.

7 monomers are involved in 12 short contacts:

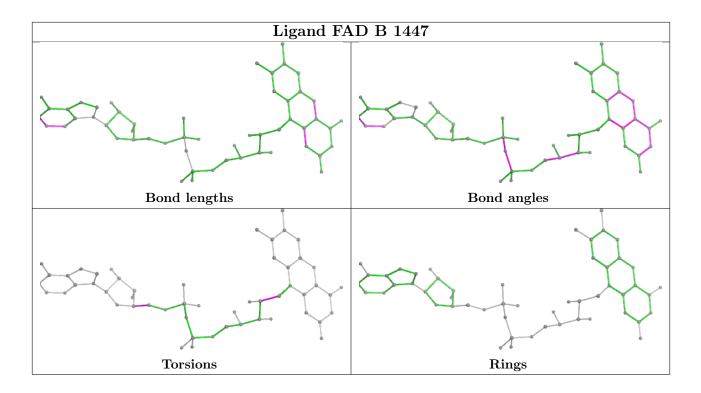
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	А	1448	FAD	1	0
2	С	1448	FAD	1	0
2	В	1447	FAD	2	0
3	А	1449	NO3	2	0
6	В	1450	MES	1	0
4	С	1450	IND	1	0
3	В	1448	NO3	4	0

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.









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, 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	А	446/464~(96%)	0.07	14 (3%) 49 51	23, 45, 68, 86	0
1	В	445/464~(95%)	-0.19	1 (0%) 95 95	23, 38, 58, 67	0
1	С	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

The worst 5 of 17 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	231	ALA	5.5
1	А	230	THR	3.8
1	А	446	ASP	3.7
1	А	233	MET	3.4
1	А	217	TYR	2.9

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, 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.



FAD

С

1448

53/53

2

Page 2	1	WV	wwPDB X-ray Structure Validation Summary Report							
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$B-factors(Å^2)$	Q<0.9		
4	IND	С	1450	9/9	0.68	0.34	94,94,95,95	0		
6	MES	В	1450	12/12	0.71	0.31	102,102,103,103	0		
4	IND	В	1449	9/9	0.85	0.22	59,59,59,60	0		
3	NO3	В	1448	4/4	0.86	0.30	71,72,72,72	0		
5	OXY	С	1451	2/2	0.87	0.40	69,69,69,71	0		
3	NO3	А	1449	4/4	0.89	0.31	73,73,73,74	0		
5	OXY	В	1451	2/2	0.93	0.32	69,69,69,69	0		
3	NO3	С	1449	4/4	0.94	0.22	66,66,67,67	0		
4	IND	А	1450	9/9	0.95	0.20	80,80,81,81	0		
5	OXY	А	1451	2/2	0.97	0.16	$55,\!55,\!55,\!56$	0		
2	FAD	А	1448	53/53	0.97	0.12	23,26,37,38	0		
2	FAD	В	1447	53/53	0.98	0.12	15,24,27,28	0		

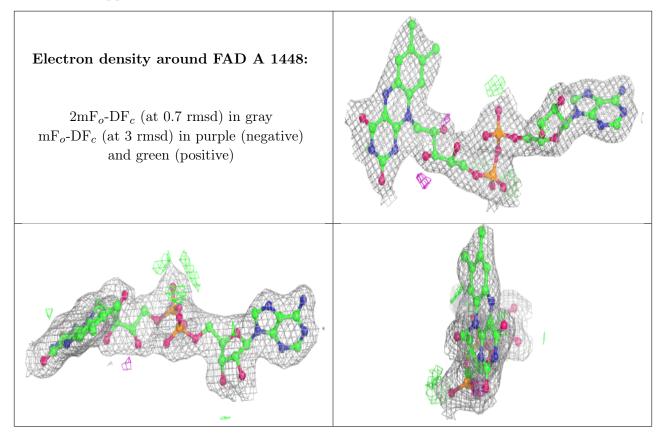
0.98

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.

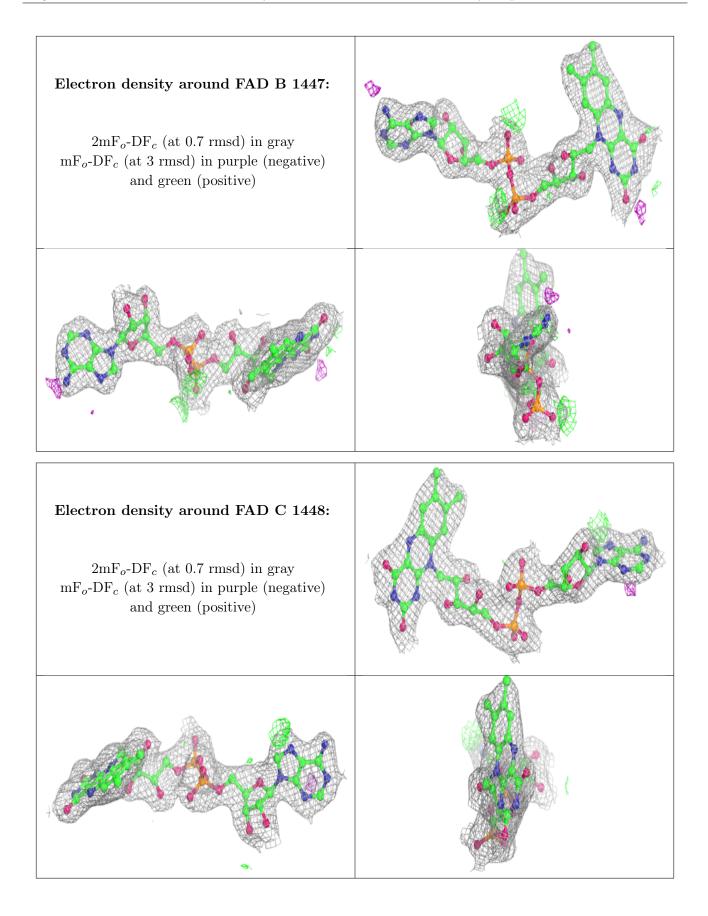
0.13

18,25,28,29

0









6.5 Other polymers (i)

There are no such residues in this entry.

