

# wwPDB X-ray Structure Validation Summary Report (i)

May 10, 2022 – 01:11 pm BST

PDB ID : 7P4F

Title: Crystal Structure of Monoamine Oxidase B in complex with inhibitor 1

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Deposited on : 2021-07-11

Resolution : 2.30 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 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 : 2.28.1buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0267$ 

CCP4 : 7.1.010 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

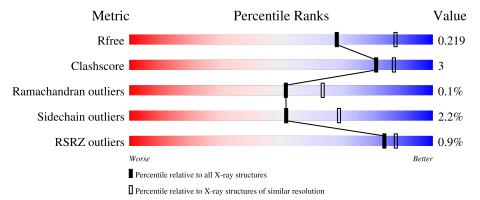
Validation Pipeline (wwPDB-VP) : 2.28.1

## 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 2.30 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\#  ext{Entries},  ext{ resolution range}( ext{Å}))$
$R_{free}$	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	520	90%	5% • •				
1	В	520	87%	7% • 5%				



# 2 Entry composition (i)

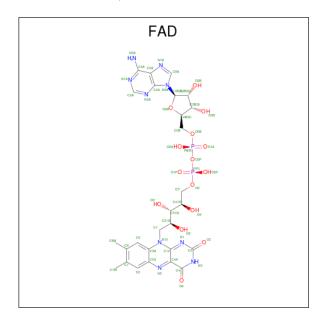
There are 5 unique types of molecules in this entry. The entry contains 8490 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Amine oxidase [flavin-containing] B.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	Λ	499	Total	С	N	О	S	0	0	0
1	Λ	499	3974	2541	681	728	24			
1	P	494	Total	С	N	О	S	0	0	0
1	Б	494	3940	2519	676	721	24	0		

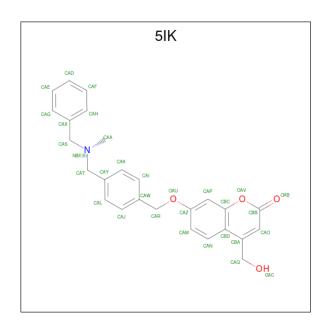
• Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:  $C_{27}H_{33}N_9O_{15}P_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	Λ	1	Total	С	N	О	Р	0	0
		1	53	27	9	15	2		
9	D	1	Total	С	N	О	Р	0	0
	Б	1	53	27	9	15	2		

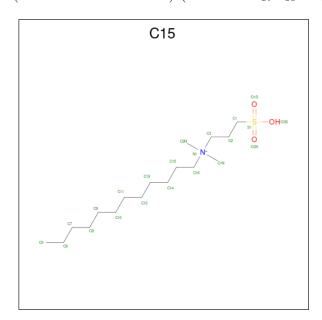
• Molecule 3 is 4-(hydroxymethyl)-7-[[4-[[methyl-(phenylmethyl)amino]methyl]phenyl]methox y]chromen-2-one (three-letter code: 5IK) (formula: C<sub>26</sub>H<sub>25</sub>NO<sub>4</sub>).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	A	1	Total				0	0	
	11	1	31	26	1	4	O O		
2	D	1	Total	С	N	O	0	0	
)	Ъ	1	31	26	1	4	U	U	

• Molecule 4 is N-DODECYL-N,N-DIMETHYL-3-AMMONIO-1-PROPANESULFONATE (three-letter code: C15) (formula: C<sub>17</sub>H<sub>38</sub>NO<sub>3</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf		
4	A	1	Total 14	C 9	N 1	O 3	S 1	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
4	D	1	Total	С	N	О	S	0	0
4	Б	1	11	6	1	3	1	0	0

#### • Molecule 5 is water.

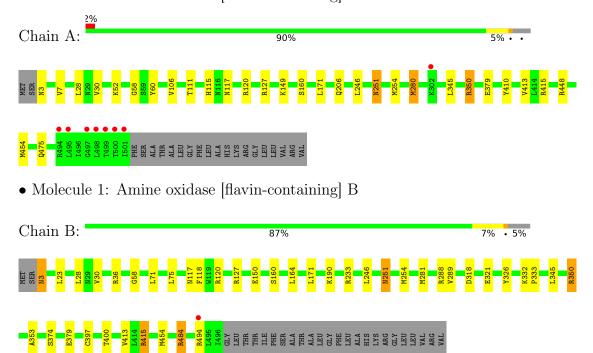
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	181	Total O 181 181	0	0
5	В	202	Total O 202 202	0	0



# 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: Amine oxidase [flavin-containing] B





# 4 Data and refinement statistics (i)

Property	Value	Source	
Space group	C 2 2 2	Depositor	
Cell constants	131.24Å 222.43Å 86.21Å	D : t	
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor	
Resolution (Å)	47.31 - 2.30	Depositor	
Resolution (A)	47.26 - 2.30	EDS	
% Data completeness	99.6 (47.31-2.30)	Depositor	
(in resolution range)	99.6 (47.26-2.30)	EDS	
$R_{merge}$	(Not available)	Depositor	
$R_{sym}$	(Not available)	Depositor	
$< I/\sigma(I) > 1$	3.49 (at 2.29Å)	Xtriage	
Refinement program	REFMAC 5.8.0230	Depositor	
D D	0.161 , 0.213	Depositor	
$R, R_{free}$	0.169 , $0.219$	DCC	
$R_{free}$ test set	1455 reflections $(2.59%)$	wwPDB-VP	
Wilson B-factor $(Å^2)$	26.2	Xtriage	
Anisotropy	0.118	Xtriage	
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	(Not available), (Not available)	EDS	
L-test for twinning <sup>2</sup>	$< L > = 0.49, < L^2> = 0.32$	Xtriage	
Estimated twinning fraction	0.013  for  1/2*h-1/2*k,-3/2*h-1/2*k,-l	Vtriego	
Estimated twinning fraction	0.020  for  1/2 *h + 1/2 *k, 3/2 *h - 1/2 *k, -1	Xtriage	
$F_o, F_c$ correlation	0.95	EDS	
Total number of atoms	8490	wwPDB-VP	
Average B, all atoms $(\mathring{A}^2)$	29.0	wwPDB-VP	

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.01% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>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.



<sup>&</sup>lt;sup>1</sup>Intensities 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: C15, FAD, 5IK

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		
Moi Chain		RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.57	0/4071	0.69	2/5526~(0.0%)	
1	В	0.60	1/4037 (0.0%)	0.71	$2/5479 \ (0.0\%)$	
All	All	0.58	1/8108 (0.0%)	0.70	4/11005~(0.0%)	

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a maintain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	В	0	4
All	All	0	7

#### All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(A)	Ideal(Å)
1	В	150	GLU	CD-OE2	5.33	1.31	1.25

#### All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$Ideal(^{o})$
1	A	448	ARG	NE-CZ-NH2	-5.37	117.62	120.30
1	В	288	ARG	NE-CZ-NH2	-5.33	117.63	120.30
1	В	36	ARG	NE-CZ-NH2	-5.20	117.70	120.30
1	A	448	ARG	NE-CZ-NH1	5.07	122.84	120.30

There are no chirality outliers.

5 of 7 planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	A	127	ARG	Sidechain
1	A	350	ARG	Sidechain
1	A	415	ARG	Sidechain
1	В	127	ARG	Sidechain
1	В	350	ARG	Sidechain

#### 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	3974	0	3976	20	1
1	В	3940	0	3937	24	0
2	A	53	0	29	1	0
2	В	53	0	29	1	0
3	A	31	0	0	0	0
3	В	31	0	0	0	0
4	A	14	0	19	0	0
4	В	11	0	12	0	0
5	A	181	0	0	2	1
5	В	202	0	0	3	1
All	All	8490	0	8002	42	2

The all-atom clash score is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clash score for this structure is 3.

The worst 5 of 42 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:A:280:MET:HE1	1:B:353:ALA:HB1	1.40	1.02
1:B:117:ASN:HD22	1:B:120:ARG:HH21	1.17	0.93
1:A:280:MET:HE1	1:B:353:ALA:CB	2.05	0.85
1:B:494:ARG:NH2	5:B:701:HOH:O	2.11	0.84
1:A:117:ASN:HD22	1:A:120:ARG:HH21	1.35	0.74

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



Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
5:A:816:HOH:O	5:B:897:HOH:O[6_565]	1.85	0.35
1:A:475:GLN:NE2	1:A:475:GLN:NE2[3_656]	1.99	0.21

### 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	Percentil	$\overline{\mathbf{es}}$
1	A	497/520~(96%)	482 (97%)	14 (3%)	1 (0%)	47 58	
1	В	492/520 (95%)	475 (96%)	17 (4%)	0	100 100	0
All	All	989/1040 (95%)	957 (97%)	31 (3%)	1 (0%)	51 64	

All (1) Ramachandran outliers are listed below:

]	Mol	Chain	$\operatorname{Res}$	Type
	1	A	52	LYS

### 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	428/444 (96%)	422 (99%)	6 (1%)	67 81		
1	В	424/444 (96%)	411 (97%)	13 (3%)	40 55		
All	All	852/888 (96%)	833 (98%)	19 (2%)	52 69		

5 of 19 residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	В	350	ARG
1	В	397	CYS
1	В	415	ARG
1	В	379	GLU
1	В	160	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
1	В	251	ASN
1	В	464	GLN
1	A	452	HIS
1	A	475	GLN
1	В	3	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)

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	e Chain Res		Link	Bond lengths			Bond angles		
MIOI	туре	Chain	nes	Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	5IK	В	602	-	32,34,34	2.34	4 (12%)	42,46,46	1.22	3 (7%)
4	C15	A	603	-	13,13,21	2.39	2 (15%)	17,18,26	1.13	1 (5%)
2	FAD	В	601	1	51,58,58	1.08	4 (7%)	60,89,89	1.84	13 (21%)
3	5IK	A	602	-	32,34,34	2.54	5 (15%)	42,46,46	1.15	4 (9%)
2	FAD	A	601	1	51,58,58	1.02	2 (3%)	60,89,89	1.66	15 (25%)
4	C15	В	603	-	10,10,21	3.22	2 (20%)	14,15,26	1.36	3 (21%)

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
3	5IK	В	602	-	-	1/15/15/15	0/4/4/4
4	C15	A	603	-	-	1/13/13/21	-
2	FAD	В	601	1	-	2/30/50/50	0/6/6/6
3	5IK	A	602	-	-	0/15/15/15	0/4/4/4
2	FAD	A	601	1	-	1/30/50/50	0/6/6/6
4	C15	В	603	-	-	1/8/8/21	-

The worst 5 of 19 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(A)	$Ideal(\AA)$
3	A	602	5IK	CAQ-CBA	-10.92	1.41	1.51
3	В	602	5IK	CAQ-CBA	-10.00	1.42	1.51
4	В	603	C15	C1-S1	-7.32	1.67	1.77
4	В	603	C15	O2S-S1	6.68	1.64	1.45
3	A	602	5IK	CAT-CAY	-6.51	1.39	1.51

The worst 5 of 39 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$Ideal(^{o})$
2	В	601	FAD	C2-N3-C4	6.83	120.91	115.14
2	В	601	FAD	C1'-N10-C9A	5.14	122.34	118.29
3	В	602	5IK	CAT-NBE-CAS	4.76	119.63	110.97
2	A	601	FAD	C9A-N10-C10	-4.67	115.79	121.91
2	A	601	FAD	N3A-C2A-N1A	-4.38	121.83	128.68

There are no chirality outliers.

5 of 6 torsion outliers are listed below:



Mol	Chain	Res	Type	Atoms
4	В	603	C15	S1-C1-C2-C3
4	A	603	C15	C13-C14-C15-C16
2	В	601	FAD	PA-O3P-P-O5'
3	В	602	5IK	OAC-CAQ-CBA-CBD
2	A	601	FAD	O4B-C4B-C5B-O5B

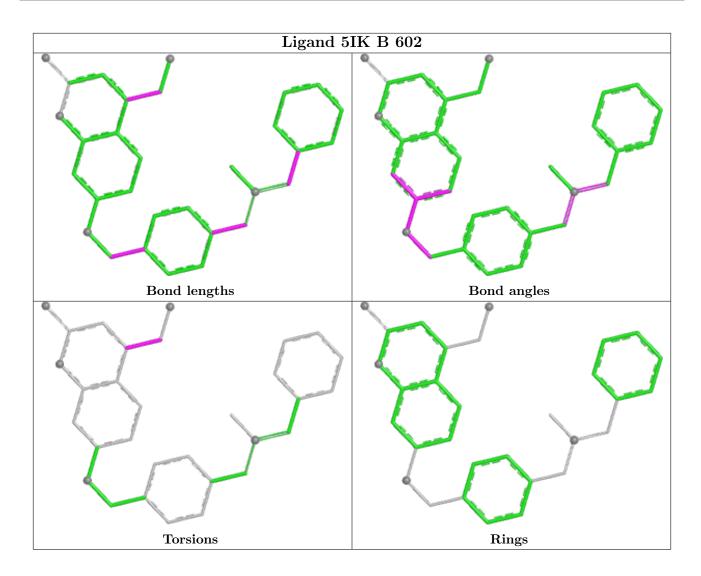
There are no ring outliers.

2 monomers are involved in 2 short contacts:

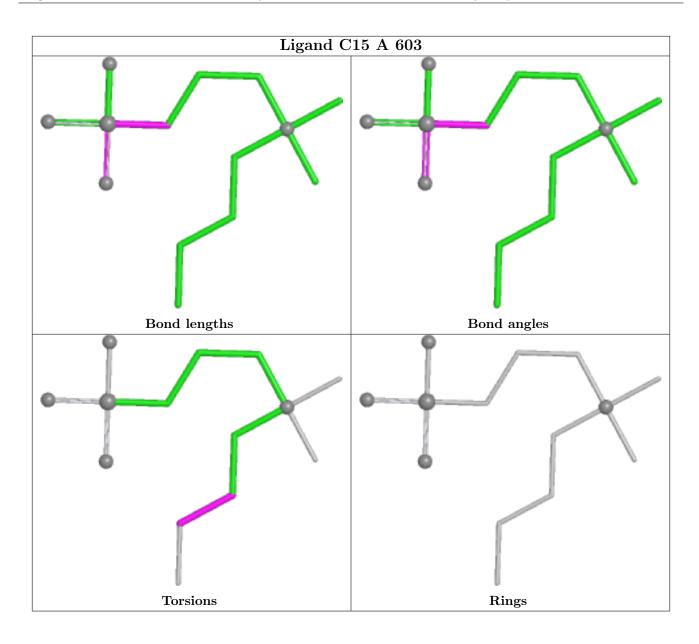
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	В	601	FAD	1	0
2	A	601	FAD	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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.

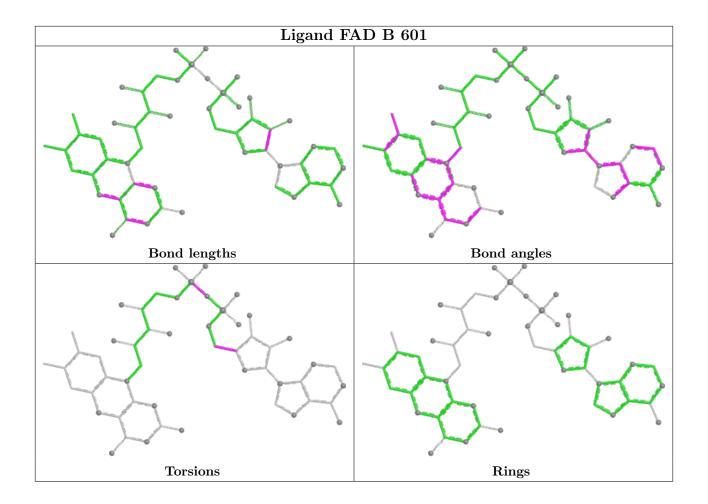




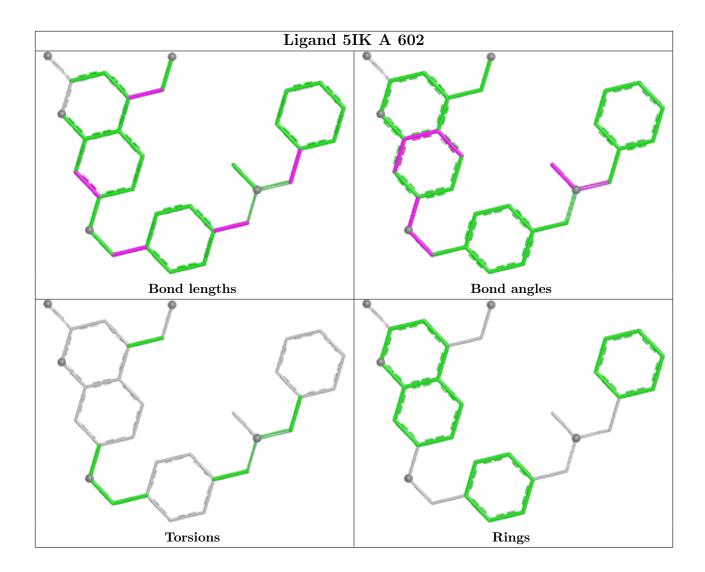




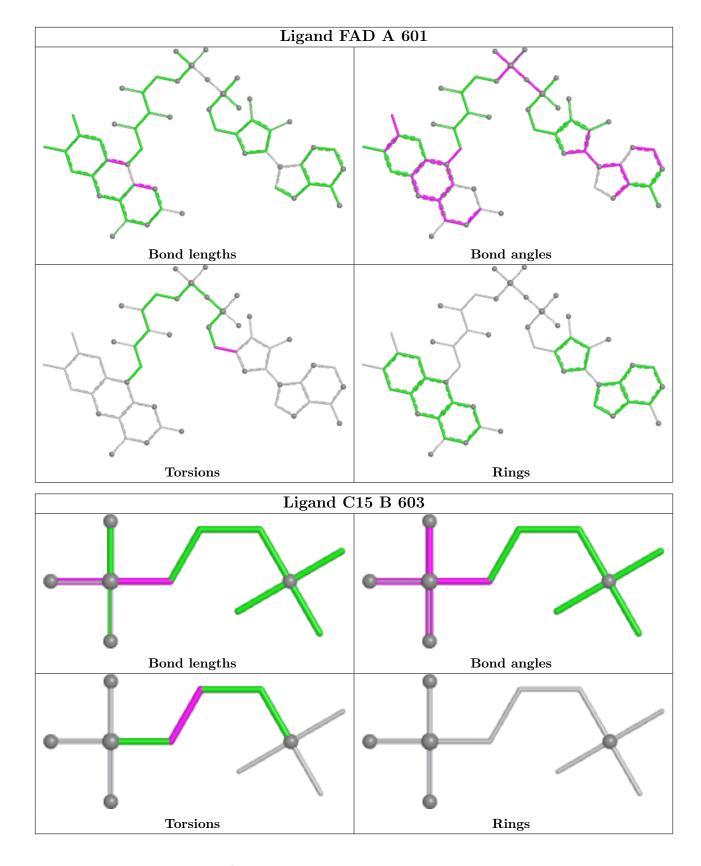












## 5.7 Other polymers (i)

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	$oxed{ Analysed} egin{array}{c c} <& RSRZ> & \#RSRZ \end{array}$		$\# \mathbf{RSRZ} >$	2	$\mathbf{OWAB}(\mathbf{\mathring{A}}^2)$	Q < 0.9
1	A	499/520 (95%)	-0.32	8 (1%) 72	77	18, 27, 50, 92	0
1	В	494/520~(95%)	-0.34	1 (0%) 95	96	16, 25, 47, 84	0
All	All	993/1040 (95%)	-0.33	9 (0%) 84	88	16, 26, 49, 92	0

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	500	THR	5.0
1	A	498	LEU	4.5
1	A	501	ILE	4.0
1	A	499	THR	3.4
1	A	495	LEU	3.1

## 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.

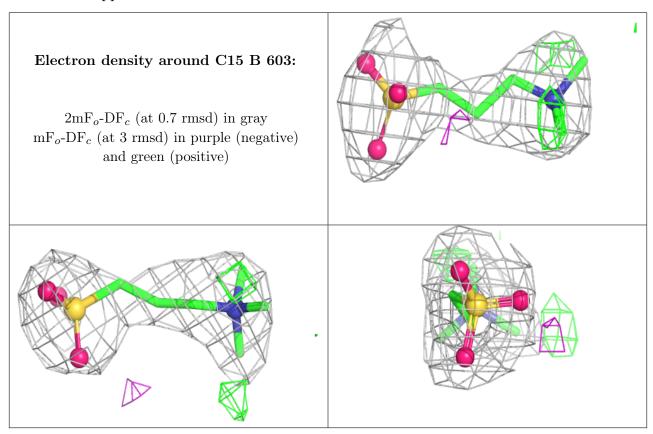
### 6.4 Ligands (i)

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.

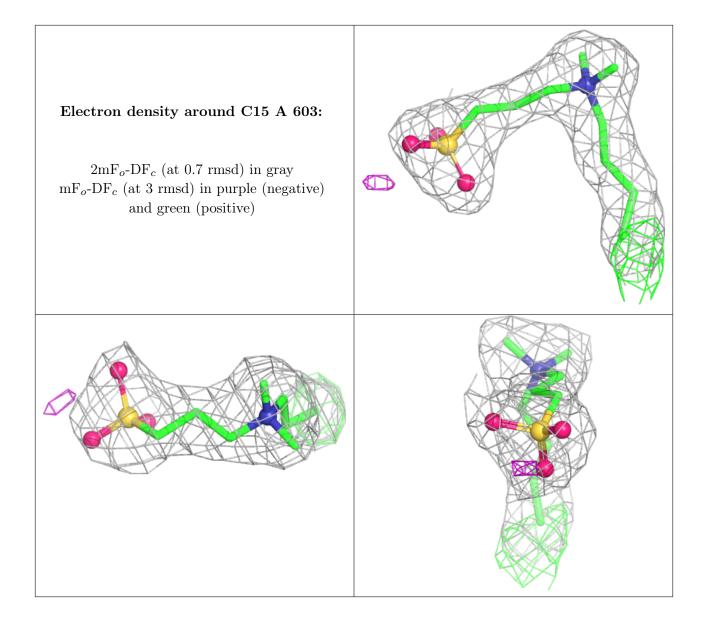


Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
4	C15	В	603	11/22	0.85	0.19	66,70,79,81	0
4	C15	A	603	14/22	0.87	0.15	39,48,68,71	0
3	5IK	A	602	31/31	0.96	0.12	24,29,35,35	0
3	5IK	В	602	31/31	0.96	0.13	25,28,38,40	0
2	FAD	A	601	53/53	0.98	0.10	16,18,21,22	0
2	FAD	В	601	53/53	0.98	0.13	14,17,19,20	0

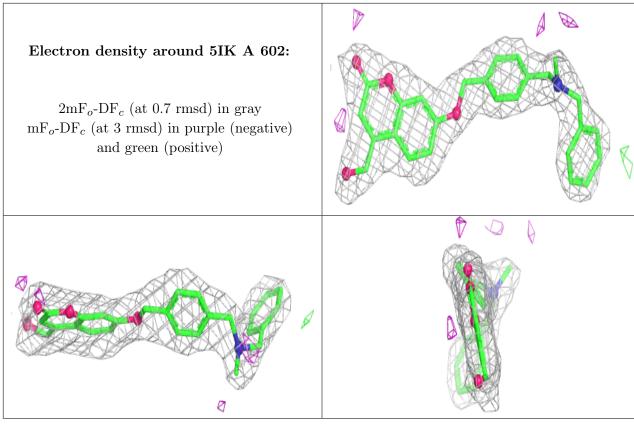
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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









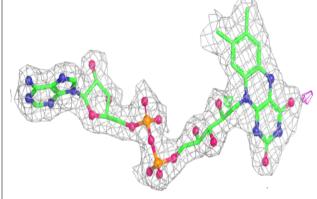


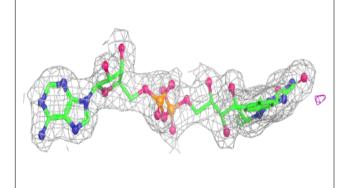
# Electron density around 5IK B 602: 2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative) and green (positive)

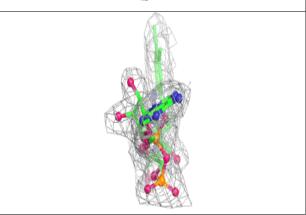


## Electron density around FAD A 601:

 $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$  (at 0.7 rmsd) in gray  $\mathrm{mF}_o\text{-}\mathrm{DF}_c$  (at 3 rmsd) in purple (negative) and green (positive)

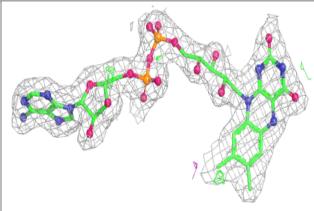


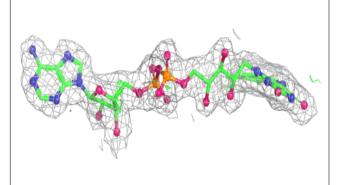


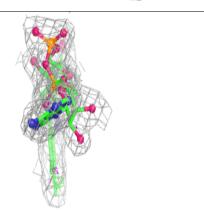


#### Electron density around FAD B 601:

 $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$  (at 0.7 rmsd) in gray  $\mathrm{mF}_o\text{-}\mathrm{DF}_c$  (at 3 rmsd) in purple (negative) and green (positive)









# 6.5 Other polymers (i)

There are no such residues in this entry.

