



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 22, 2020 – 04:19 pm BST

PDB ID : 3AX7  
Title : Bovine Xanthine Oxidase, protease cleaved form  
Authors : Ishikita, H.; Eger, B.T.; Pai, E.F.; Okamoto, K.; Nishino, T.  
Deposited on : 2011-03-30  
Resolution : 2.34 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](mailto: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.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

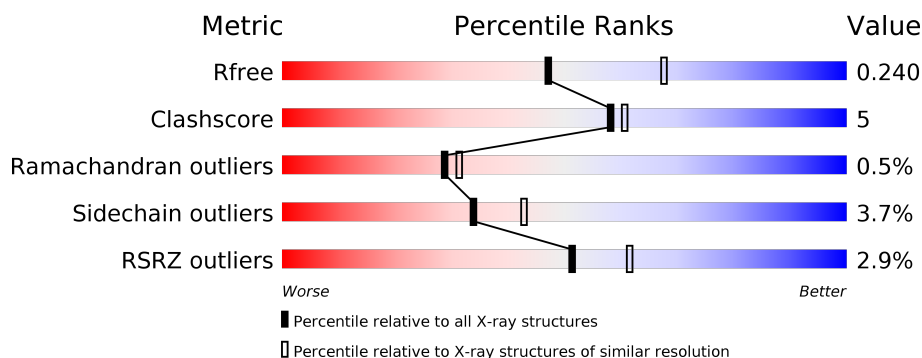
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.34 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	2096 (2.36-2.32)
Clashscore	141614	2193 (2.36-2.32)
Ramachandran outliers	138981	2159 (2.36-2.32)
Sidechain outliers	138945	2160 (2.36-2.32)
RSRZ outliers	127900	2067 (2.36-2.32)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	1332	<div> <div>2%</div> <div> <div></div> <div>81%</div> <div>10%</div> <div>8%</div> </div> </div>
1	B	1332	<div> <div>3%</div> <div> <div></div> <div>80%</div> <div>10%</div> <div>9%</div> </div> </div>

## 2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 19746 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.

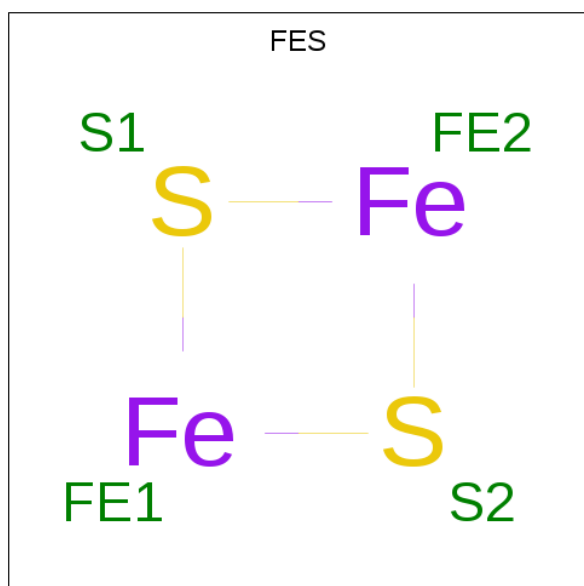
- Molecule 1 is a protein called Xanthine dehydrogenase/oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1225	Total	C	N	O	S	0	0	0
			9476	6013	1632	1771	60			
1	B	1218	Total	C	N	O	S	0	0	0
			9433	5988	1623	1763	59			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	552	HIS	ASP	CONFLICT	UNP P80457
B	552	HIS	ASP	CONFLICT	UNP P80457

- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe<sub>2</sub>S<sub>2</sub>).



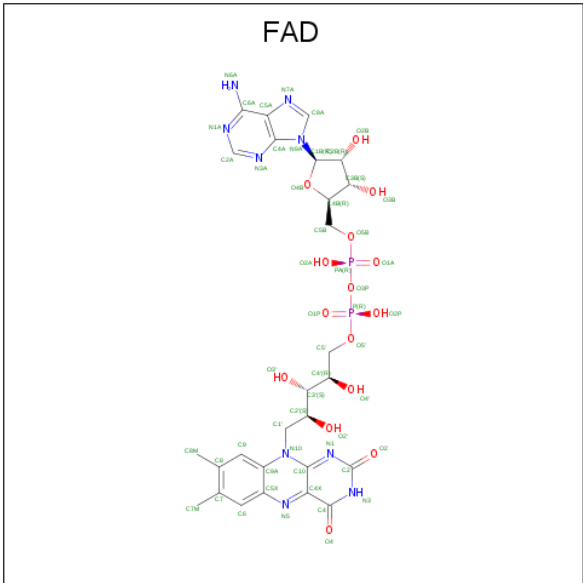
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			4	2	2		

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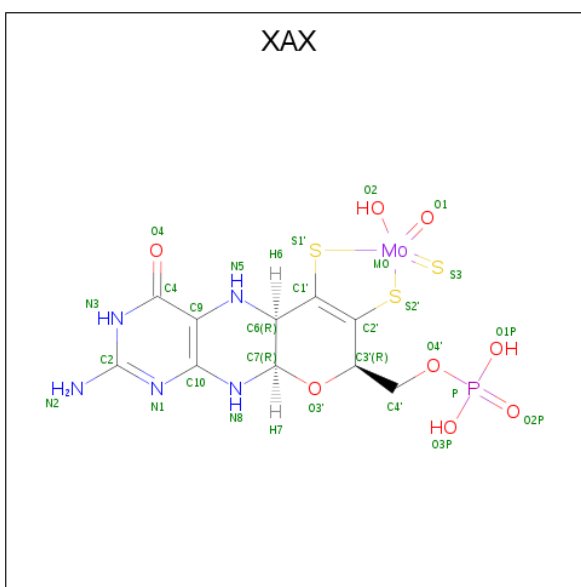
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	Fe	S	0	0
			4	2	2		
2	B	1	Total	Fe	S	0	0
			4	2	2		
2	B	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C<sub>27</sub>H<sub>33</sub>N<sub>9</sub>O<sub>15</sub>P<sub>2</sub>).



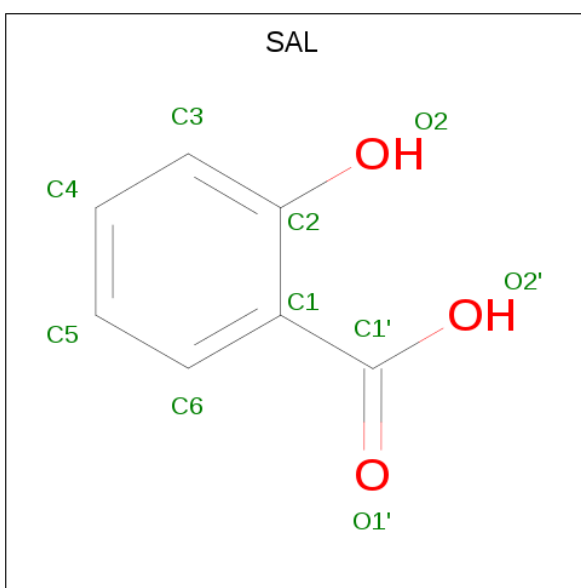
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
3	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 4 is {[ (5aR,8R,9aR)-2-amino-4-oxo-6,7-di(sulfanyl-kappaS)-3,5,5a,8,9a,10-hexahydro-4H-pyran[3,2-g]pteridin-8-yl)methyl dihydrogenato(2-) phosphate}(hydroxy)oxo(thioxo)molybdenum (three-letter code: XAX) (formula: C<sub>10</sub>H<sub>13</sub>MoN<sub>5</sub>O<sub>8</sub>PS<sub>3</sub>).



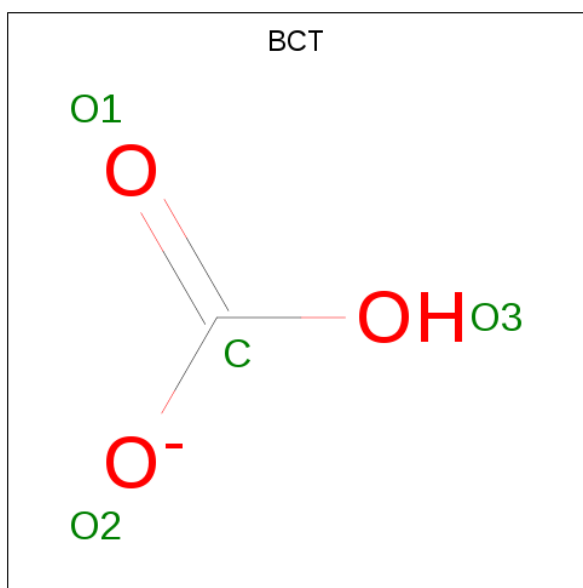
Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
4	A	1	Total	C	Mo	N	O	P	S	0	0
			28	10	1	5	8	1	3		
4	B	1	Total	C	Mo	N	O	P	S	0	0
			28	10	1	5	8	1	3		

- Molecule 5 is 2-HYDROXYBENZOIC ACID (three-letter code: SAL) (formula: C<sub>7</sub>H<sub>6</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	7	3		
5	B	1	Total	C	O	0	0
			10	7	3		

- Molecule 6 is BICARBONATE ION (three-letter code: BCT) (formula:  $\text{CHO}_3$ ).



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

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula:  $\text{C}_3\text{H}_8\text{O}_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			6	3	3		

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

- Molecule 8 is CALCIUM ION (three-letter code: CA) (formula: Ca).

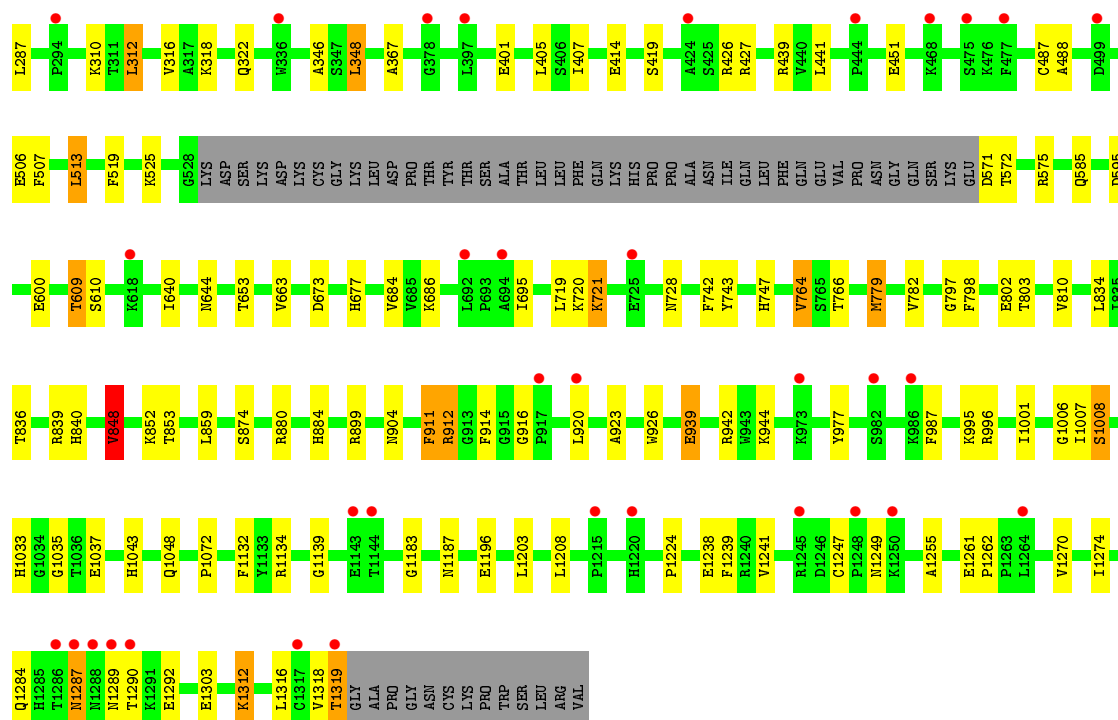
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	1	Total	Ca	0	0
			1	1		
8	A	1	Total	Ca	0	0
			1	1		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	316	Total	O	0	0
			316	316		
9	B	283	Total	O	0	0
			283	283		







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	133.48Å 73.15Å 146.27Å 90.00° 98.18° 90.00°	Depositor
Resolution (Å)	36.37 – 2.34 36.37 – 2.34	Depositor EDS
% Data completeness (in resolution range)	93.8 (36.37-2.34) 93.8 (36.37-2.34)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.40 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, $R_{free}$	0.194 , 0.240 0.195 , 0.240	Depositor DCC
$R_{free}$ test set	5554 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.4	Xtriage
Anisotropy	0.221	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 41.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	19746	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.32% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: XAX, GOL, SAL, CA, FES, BCT, 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	0.64	0/9675	0.69	6/13085 (0.0%)
1	B	0.62	0/9631	0.68	6/13026 (0.0%)
All	All	0.63	0/19306	0.68	12/26111 (0.0%)

There are no bond length outliers.

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	839	ARG	NE-CZ-NH2	-7.26	116.67	120.30
1	B	942	ARG	NE-CZ-NH2	-6.35	117.12	120.30
1	A	779	MET	CG-SD-CE	-5.94	90.69	100.20
1	A	1325	CYS	N-CA-C	5.65	126.26	111.00
1	A	439	ARG	NE-CZ-NH2	-5.51	117.54	120.30

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	9476	0	9483	91	0
1	B	9433	0	9439	97	0
2	A	8	0	0	0	0
2	B	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	53	0	31	1	0
3	B	53	0	31	3	0
4	A	28	0	10	0	0
4	B	28	0	10	1	0
5	A	10	0	5	1	0
5	B	10	0	4	1	0
6	A	4	0	0	0	0
6	B	4	0	0	0	0
7	A	18	0	24	0	0
7	B	12	0	16	0	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
9	A	316	0	0	3	0
9	B	283	0	0	9	0
All	All	19746	0	19053	185	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.

The worst 5 of 185 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:779:MET:SD	9:B:1452:HOH:O	2.09	1.10
1:A:884:HIS:HE1	1:A:1006:GLY:H	1.17	0.92
1:A:720:LYS:O	1:A:721:LYS:HB2	1.70	0.90
1:A:131:GLN:HE21	1:A:133:GLU:H	1.14	0.88
1:A:995:LYS:NZ	1:A:1284:GLN:HE21	1.72	0.86

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	1219/1332 (92%)	1172 (96%)	39 (3%)	8 (1%)	22	22
1	B	1212/1332 (91%)	1162 (96%)	45 (4%)	5 (0%)	34	38
All	All	2431/2664 (91%)	2334 (96%)	84 (4%)	13 (0%)	29	31

5 of 13 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	721	LYS
1	A	1008	SER
1	A	1325	CYS
1	B	1008	SER
1	B	721	LYS

### 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	1030/1128 (91%)	990 (96%)	40 (4%)	32	41
1	B	1026/1128 (91%)	990 (96%)	36 (4%)	36	45
All	All	2056/2256 (91%)	1980 (96%)	76 (4%)	34	43

5 of 76 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1001	ILE
1	B	129	ARG
1	B	1238	GLU
1	A	1203	LEU
1	A	1287	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 32 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1284	GLN

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Mol	Chain	Res	Type
1	B	131	GLN
1	B	1048	GLN
1	A	1324	ASN
1	B	146	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

Of 19 ligands modelled in this entry, 2 are monoatomic - leaving 17 for Mogul analysis.

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	$\# Z  > 2$
7	GOL	A	1338	-	5,5,5	0.61	0	5,5,5	0.59	0
4	XAX	A	3003	-	20,31,31	2.48	5 (25%)	22,52,52	2.00	7 (31%)
6	BCT	B	1337	-	0,3,3	0.00	-	0,3,3	0.00	-
2	FES	B	1334	1	0,4,4	0.00	-	-	-	-
2	FES	A	1334	1	0,4,4	0.00	-	-	-	-
4	XAX	B	3003	-	20,31,31	2.13	5 (25%)	22,52,52	2.35	11 (50%)
3	FAD	B	1335	-	51,58,58	1.40	6 (11%)	60,89,89	1.69	9 (15%)
2	FES	A	1333	1	0,4,4	0.00	-	-	-	-
6	BCT	A	1337	-	0,3,3	0.00	-	0,3,3	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
7	GOL	A	1340	-	5,5,5	0.33	0	5,5,5	0.56	0
7	GOL	B	1338	-	5,5,5	0.36	0	5,5,5	0.42	0
5	SAL	B	1336	-	8,10,10	1.70	2 (25%)	9,13,13	1.44	2 (22%)
5	SAL	A	1336	-	8,10,10	1.47	2 (25%)	9,13,13	1.55	3 (33%)
7	GOL	A	1339	-	5,5,5	0.35	0	5,5,5	0.37	0
7	GOL	B	1339	-	5,5,5	0.35	0	5,5,5	0.44	0
2	FES	B	1333	1	0,4,4	0.00	-	-	-	-
3	FAD	A	1335	-	51,58,58	1.34	6 (11%)	60,89,89	1.81	10 (16%)

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
7	GOL	A	1338	-	-	2/4/4/4	-
2	FES	B	1334	1	-	-	0/1/1/1
2	FES	A	1333	1	-	-	0/1/1/1
3	FAD	B	1335	-	-	0/30/50/50	0/6/6/6
2	FES	B	1333	1	-	-	0/1/1/1
4	XAX	B	3003	-	-	3/6/46/46	0/4/4/4
2	FES	A	1334	1	-	-	0/1/1/1
7	GOL	A	1340	-	-	0/4/4/4	-
7	GOL	B	1338	-	-	2/4/4/4	-
5	SAL	B	1336	-	-	0/0/4/4	0/1/1/1
5	SAL	A	1336	-	-	0/0/4/4	0/1/1/1
7	GOL	A	1339	-	-	0/4/4/4	-
7	GOL	B	1339	-	-	2/4/4/4	-
4	XAX	A	3003	-	-	2/6/46/46	0/4/4/4
3	FAD	A	1335	-	-	1/30/50/50	0/6/6/6

The worst 5 of 26 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	3003	XAX	C4-C9	5.48	1.48	1.41
4	A	3003	XAX	C9-N5	5.29	1.48	1.38
4	B	3003	XAX	C4-C9	5.27	1.48	1.41
4	A	3003	XAX	C9-C10	4.87	1.50	1.41
3	B	1335	FAD	C2A-N3A	4.66	1.39	1.32

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1335	FAD	C1'-N10-C9A	6.87	123.70	118.29
3	A	1335	FAD	C4-N3-C2	6.47	120.60	115.14
3	A	1335	FAD	N3A-C2A-N1A	-6.32	118.80	128.68
3	A	1335	FAD	C1'-N10-C9A	5.48	122.60	118.29
4	B	3003	XAX	C4-N3-C2	5.23	124.24	115.93

There are no chirality outliers.

5 of 12 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	3003	XAX	C4'-O4'-P-O2P
7	A	1338	GOL	C1-C2-C3-O3
7	B	1338	GOL	O1-C1-C2-C3
4	A	3003	XAX	C3'-C4'-O4'-P
4	B	3003	XAX	C3'-C4'-O4'-P

There are no ring outliers.

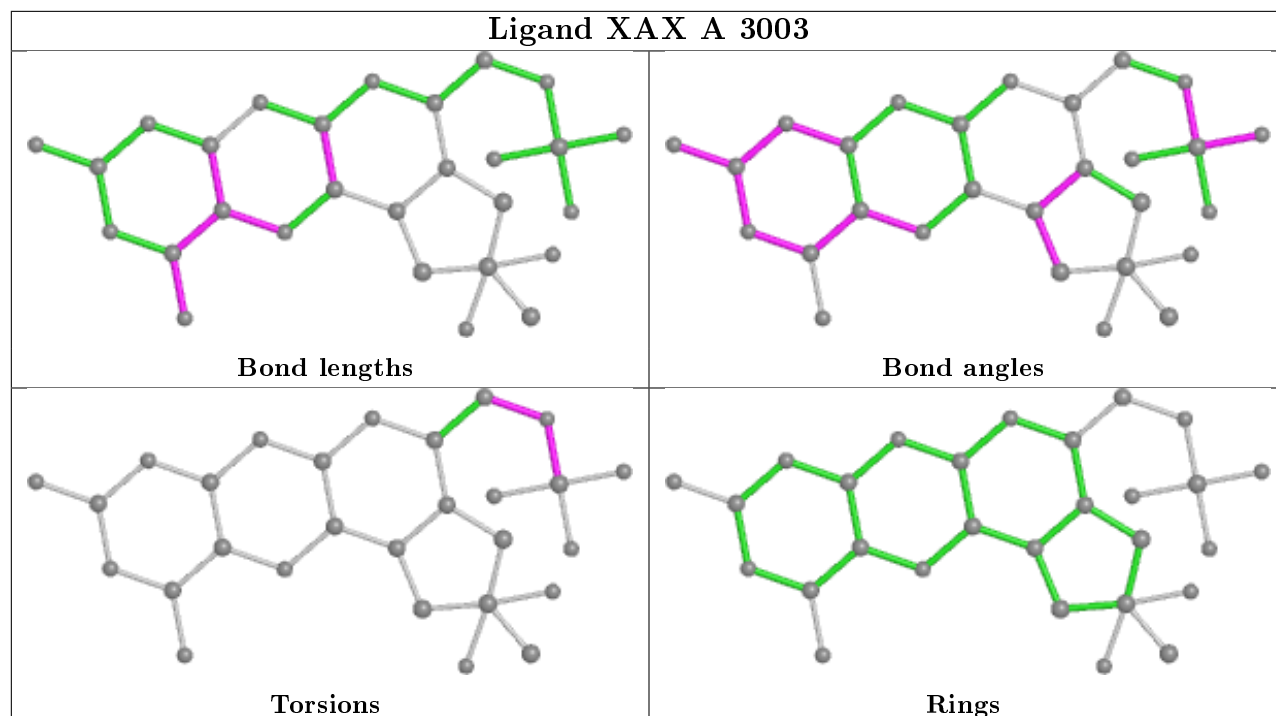
5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	3003	XAX	1	0
3	B	1335	FAD	3	0
5	B	1336	SAL	1	0
5	A	1336	SAL	1	0
3	A	1335	FAD	1	0

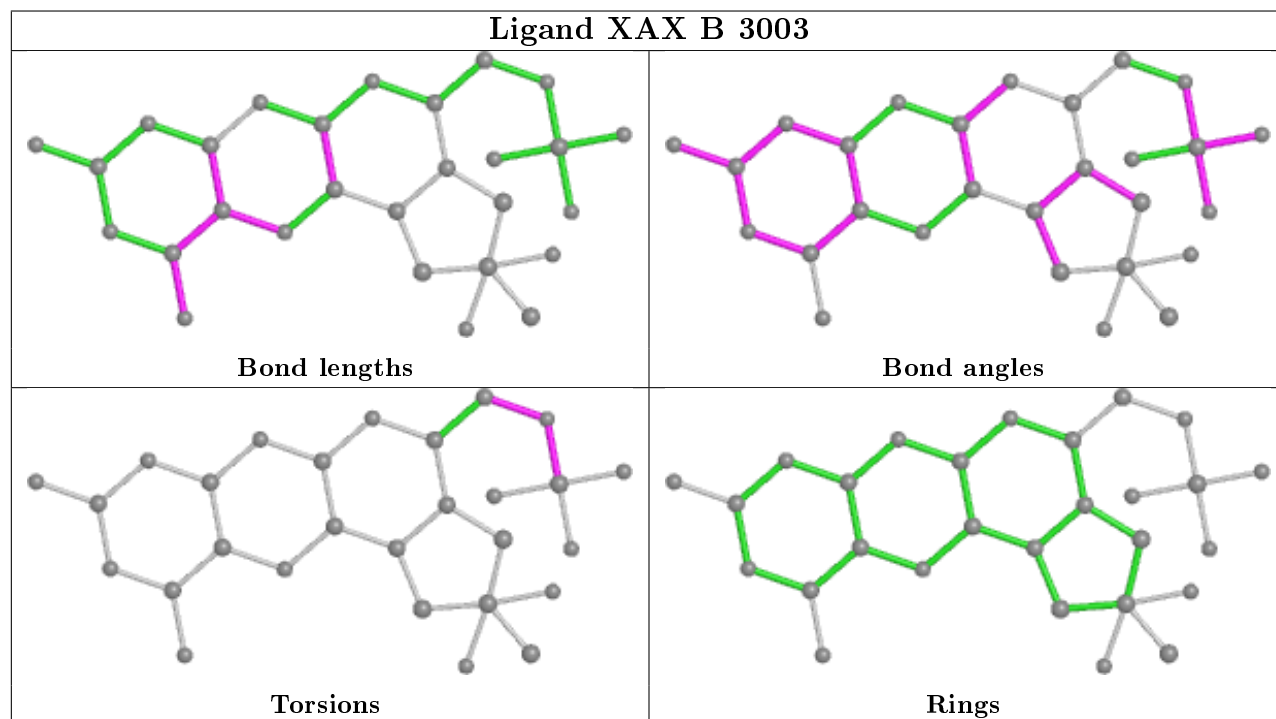
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

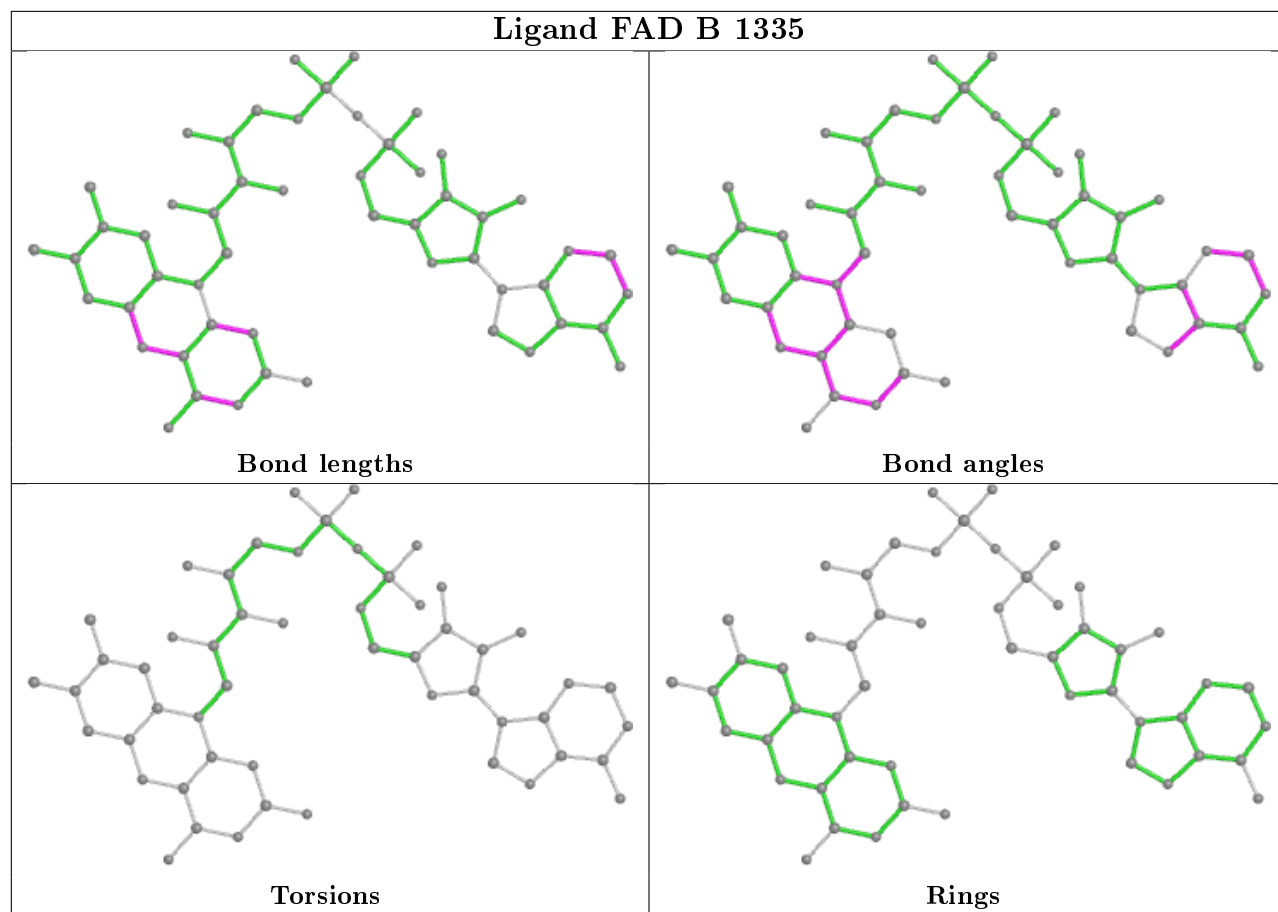


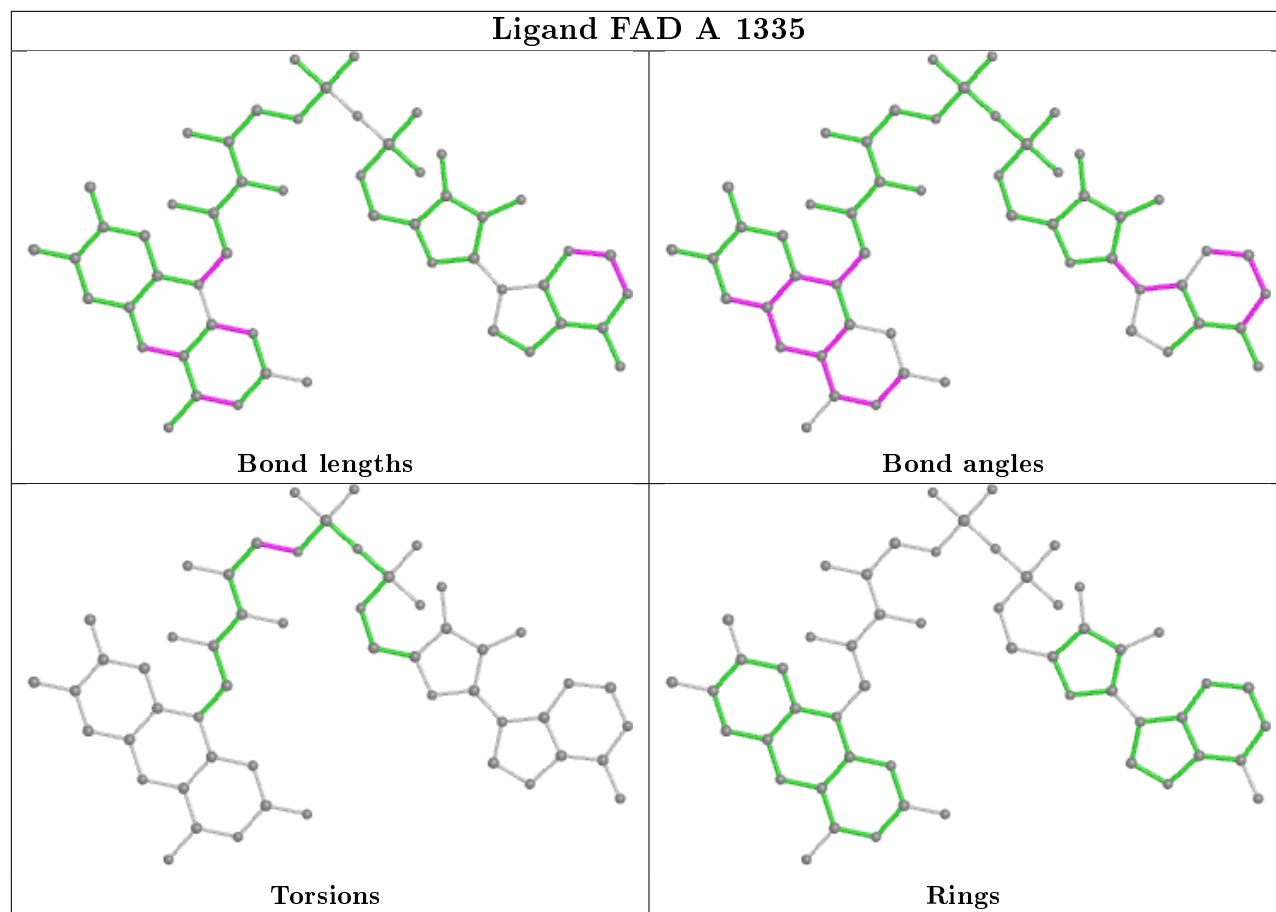
## Ligand XAX A 3003



## Ligand XAX B 3003







## 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	1225/1332 (91%)	-0.01	24 (1%) 65 74	17, 27, 43, 62	0
1	B	1218/1332 (91%)	0.07	46 (3%) 40 51	17, 28, 44, 64	0
All	All	2443/2664 (91%)	0.03	70 (2%) 51 61	17, 28, 43, 64	0

The worst 5 of 70 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1287	ASN	6.3
1	A	58	TYR	5.8
1	B	1290	THR	5.5
1	B	165	LYS	5.5
1	B	2	THR	5.0

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

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

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

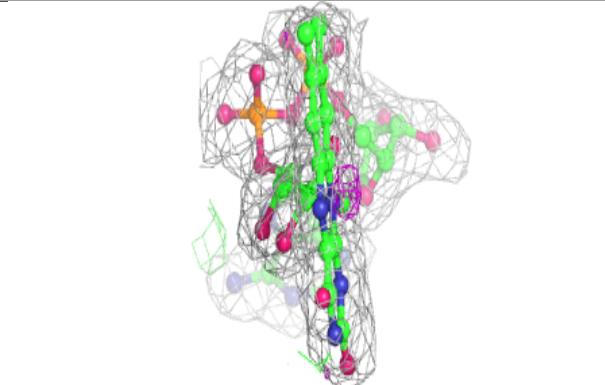
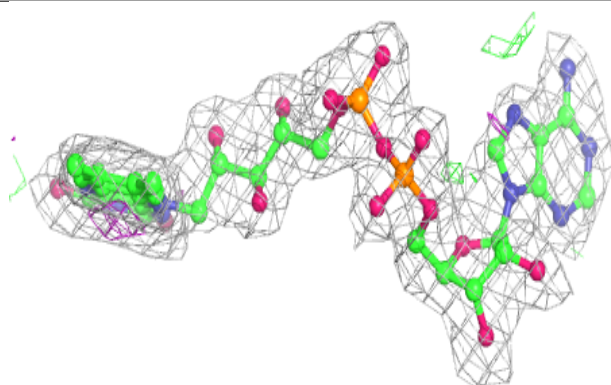
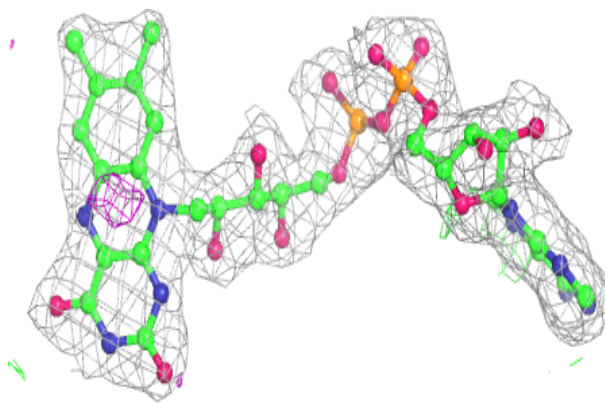
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	SAL	B	1336	10/10	0.87	0.18	35,37,38,40	0
5	SAL	A	1336	10/10	0.92	0.21	27,30,31,31	0
7	GOL	A	1340	6/6	0.95	0.13	40,40,40,41	0
7	GOL	B	1338	6/6	0.96	0.15	29,30,30,31	0
3	FAD	B	1335	53/53	0.96	0.14	23,28,30,31	0
7	GOL	A	1338	6/6	0.96	0.11	27,28,30,30	0
7	GOL	B	1339	6/6	0.96	0.13	19,25,27,27	0
7	GOL	A	1339	6/6	0.97	0.11	26,27,28,29	0
6	BCT	B	1337	4/4	0.97	0.14	17,17,18,20	0
3	FAD	A	1335	53/53	0.97	0.13	20,24,27,30	0
8	CA	B	1340	1/1	0.98	0.13	24,24,24,24	0
6	BCT	A	1337	4/4	0.99	0.21	20,21,22,22	0
4	XAX	B	3003	28/28	0.99	0.15	14,23,29,34	0
4	XAX	A	3003	28/28	0.99	0.15	15,20,25,33	0
2	FES	A	1333	4/4	0.99	0.10	20,20,21,24	0
2	FES	A	1334	4/4	0.99	0.09	19,21,21,22	0
2	FES	B	1333	4/4	0.99	0.11	20,20,21,23	0
8	CA	A	1341	1/1	0.99	0.10	22,22,22,22	0
2	FES	B	1334	4/4	1.00	0.10	20,21,22,23	0

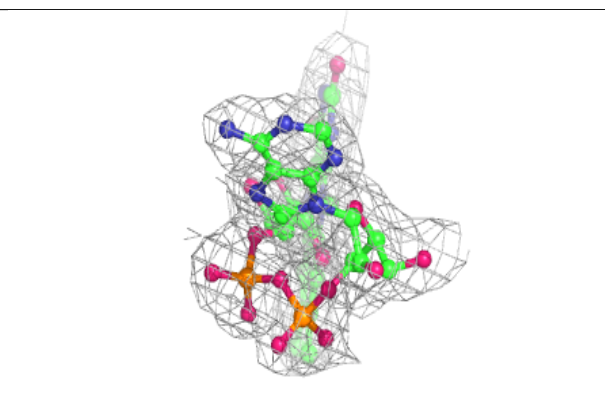
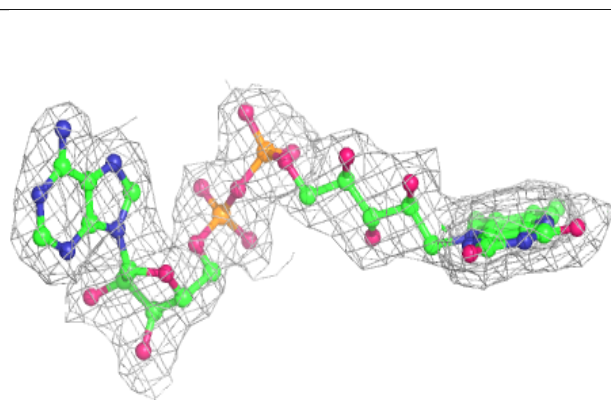
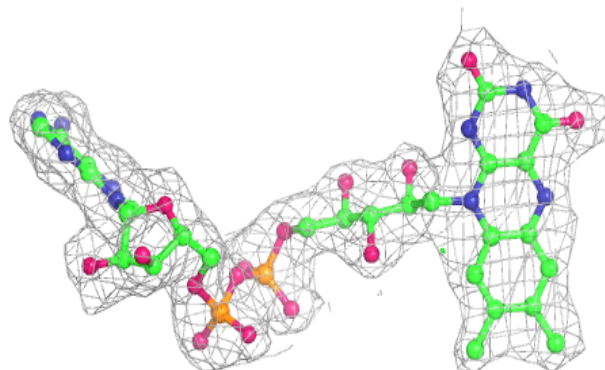
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around FAD B 1335:**

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

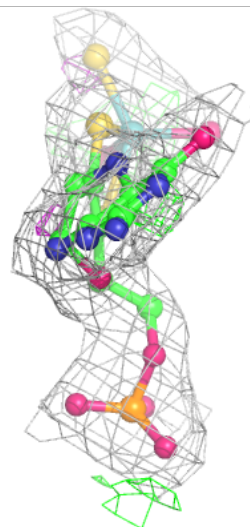
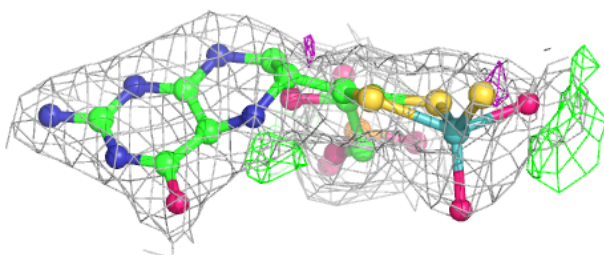
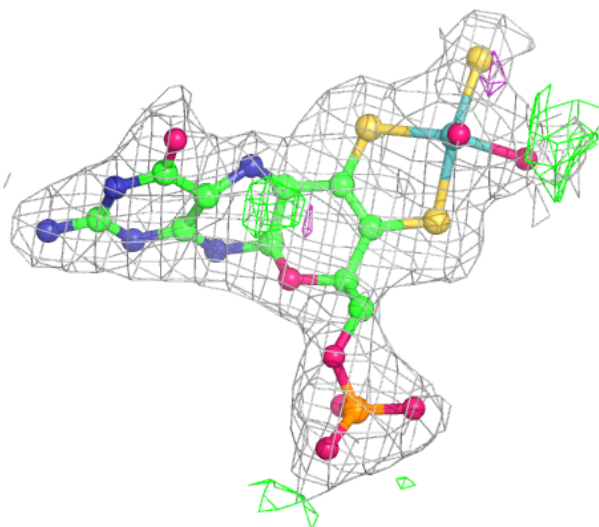
**Electron density around FAD A 1335:**

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



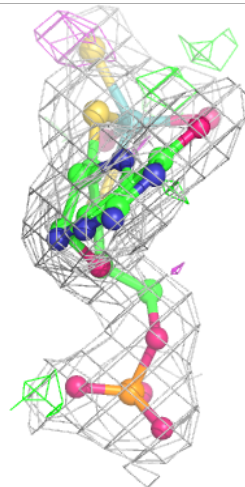
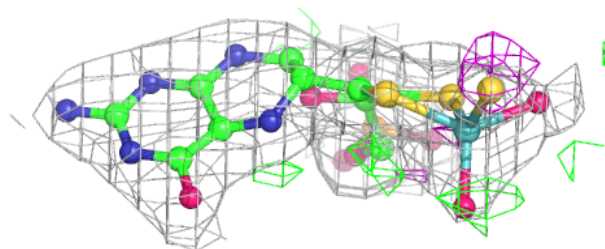
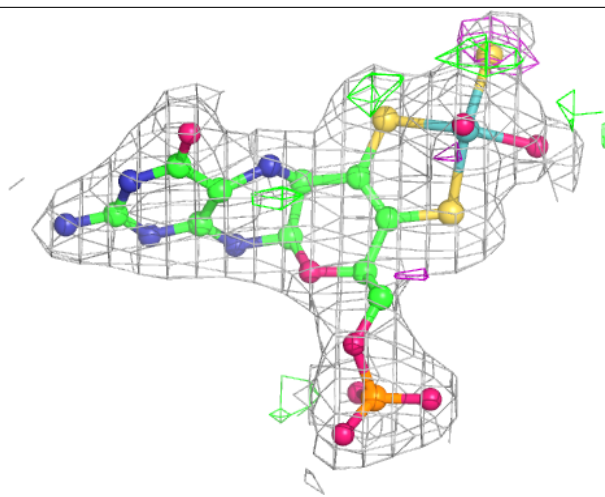
**Electron density around XAX B 3003:**

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



**Electron density around XAX A 3003:**

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



## 6.5 Other polymers ⓘ

There are no such residues in this entry.