



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 05:12 am BST

PDB ID : 3ALL
Title : Crystal structure of 2-methyl-3-hydroxypyridine-5-carboxylic acid oxygenase, mutant Y270A
Authors : Kobayashi, J.; Yoshida, H.; Yoshikane, Y.; Kamitori, S.; Yagi, T.
Deposited on : 2010-08-04
Resolution : 1.78 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 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

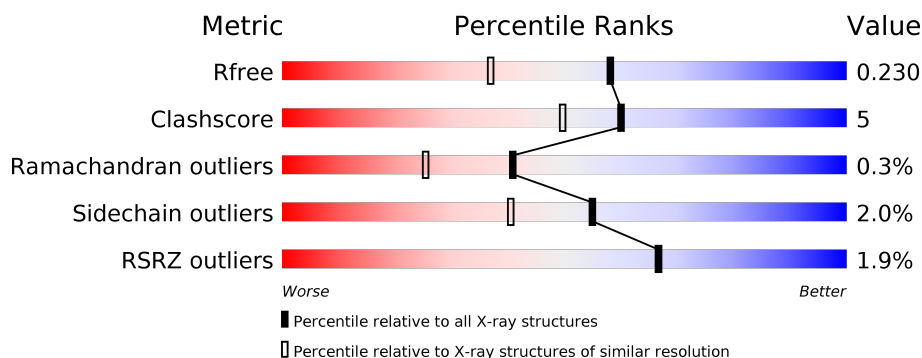
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.78 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	9185 (1.80-1.76)
Clashscore	141614	10184 (1.80-1.76)
Ramachandran outliers	138981	10051 (1.80-1.76)
Sidechain outliers	138945	10050 (1.80-1.76)
RSRZ outliers	127900	9032 (1.80-1.76)

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 ≥ 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	379	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 85%, yellow 12%, grey 3%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 85% 12% .. </div> </div>
1	B	379	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 2%, green 90%, yellow 7%, grey 3%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 2% 90% 7% .. </div> </div>

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	BME	B	381	-	X	X	-
4	GOL	B	382	-	X	-	-

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.

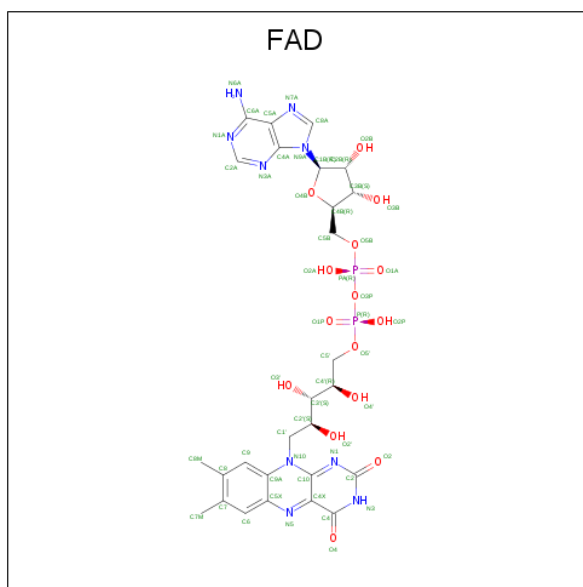
- Molecule 1 is a protein called 2-methyl-3-hydroxypyridine-5-carboxylic acid oxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	370	Total 2901	C 1828	N 520	O 539	S 14	0	7	0
1	B	371	Total 2892	C 1824	N 520	O 534	S 14	0	3	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	270	ALA	TYR	ENGINEERED MUTATION	UNP Q988D3
B	270	ALA	TYR	ENGINEERED MUTATION	UNP Q988D3

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



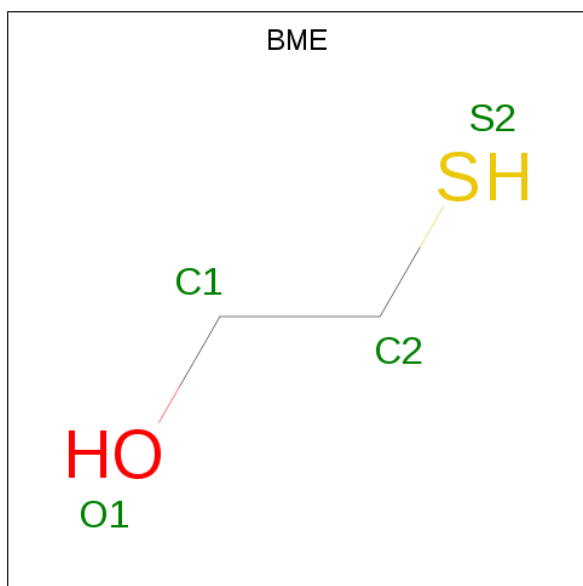
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

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

- Molecule 3 is BETA-MERCAPTOETHANOL (three-letter code: BME) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	S	0	0
			4	2	1	1		
3	B	1	Total	C	O	S	0	0
			4	2	1	1		

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



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

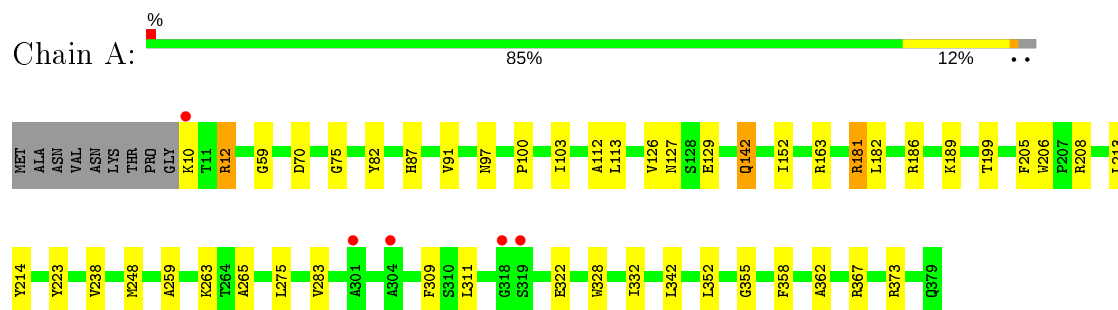
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	216	Total	O	0	0
			216	216		
5	B	126	Total	O	0	0
			126	126		

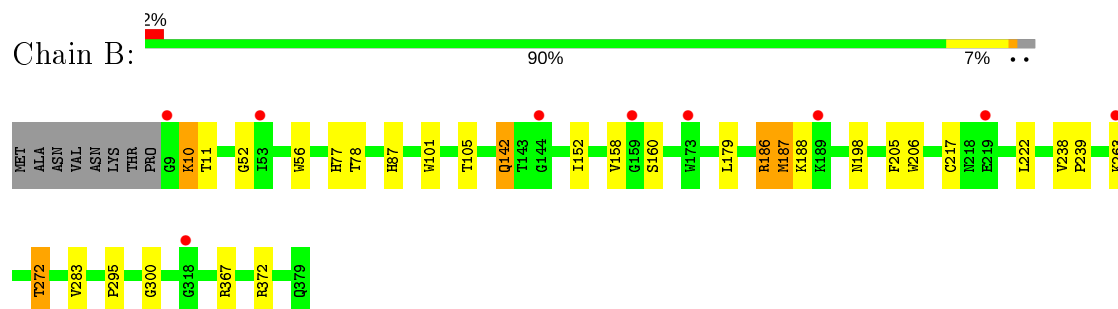
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: 2-methyl-3-hydroxypyridine-5-carboxylic acid oxygenase



- Molecule 1: 2-methyl-3-hydroxypyridine-5-carboxylic acid oxygenase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	57.09Å 284.98Å 123.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.78 – 1.78 33.78 – 1.78	Depositor EDS
% Data completeness (in resolution range)	98.4 (33.78-1.78) 98.4 (33.78-1.78)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.70 (at 1.78Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.196 , 0.230 0.196 , 0.230	Depositor DCC
R_{free} test set	4767 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	20.3	Xtriage
Anisotropy	0.237	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6279	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% 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: GOL, FAD, BME

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	1.43	17/2987 (0.6%)	1.17	13/4052 (0.3%)
1	B	1.16	1/2966 (0.0%)	1.02	4/4022 (0.1%)
All	All	1.30	18/5953 (0.3%)	1.10	17/8074 (0.2%)

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 mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	91	VAL	CB-CG2	6.79	1.67	1.52
1	A	309	PHE	CE1-CZ	6.61	1.49	1.37
1	A	181	ARG	CB-CG	6.45	1.70	1.52
1	A	358	PHE	CD1-CE1	6.40	1.52	1.39
1	A	91	VAL	N-CA	6.37	1.59	1.46
1	B	295	PRO	C-O	6.35	1.35	1.23
1	A	126	VAL	CB-CG2	5.92	1.65	1.52
1	A	214	TYR	CE1-CZ	5.82	1.46	1.38
1	A	238	VAL	CB-CG1	5.81	1.65	1.52
1	A	82	TYR	CE1-CZ	5.70	1.46	1.38
1	A	223	TYR	CG-CD2	5.52	1.46	1.39
1	A	112	ALA	CA-CB	5.46	1.64	1.52
1	A	355	GLY	C-O	5.34	1.32	1.23
1	A	322	GLU	CD-OE1	5.34	1.31	1.25
1	A	248	MET	N-CA	5.30	1.56	1.46
1	A	332	ILE	CA-CB	5.17	1.66	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	59	GLY	N-CA	5.15	1.53	1.46
1	A	223	TYR	CD1-CE1	5.10	1.47	1.39

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	208	ARG	NE-CZ-NH1	9.40	125.00	120.30
1	A	275	LEU	CB-CG-CD2	7.84	124.34	111.00
1	B	367	ARG	NE-CZ-NH2	-7.59	116.51	120.30
1	A	163	ARG	NE-CZ-NH2	-6.80	116.90	120.30
1	A	208	ARG	NE-CZ-NH2	-6.36	117.12	120.30
1	A	12[A]	ARG	NE-CZ-NH2	-6.35	117.12	120.30
1	A	12[B]	ARG	NE-CZ-NH2	-6.35	117.12	120.30
1	B	372[A]	ARG	NE-CZ-NH2	-5.85	117.38	120.30
1	B	372[B]	ARG	NE-CZ-NH2	-5.85	117.38	120.30
1	B	186	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	A	367[A]	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	A	367[B]	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	A	12[A]	ARG	NE-CZ-NH1	5.13	122.87	120.30
1	A	12[B]	ARG	NE-CZ-NH1	5.13	122.87	120.30
1	A	373	ARG	NE-CZ-NH1	-5.04	117.78	120.30
1	A	352	LEU	CB-CG-CD1	-5.04	102.44	111.00
1	A	186	ARG	NE-CZ-NH1	5.03	122.81	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	198	ASN	Peptide

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	2901	0	2882	29	0
1	B	2892	0	2874	28	0
2	A	53	0	31	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	53	0	31	1	0
3	A	4	0	5	0	0
3	B	4	0	5	4	0
4	A	18	0	24	2	0
4	B	12	0	16	2	0
5	A	216	0	0	2	0
5	B	126	0	0	2	0
All	All	6279	0	5868	54	0

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

All (54) 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:A:12[B]:ARG:HH11	1:A:12[B]:ARG:HG3	1.05	1.15
1:B:272:THR:HG23	3:B:381:BME:H11	1.23	1.14
1:B:272:THR:CG2	3:B:381:BME:H11	1.96	0.94
1:A:12[B]:ARG:NH1	1:A:12[B]:ARG:HG3	1.85	0.90
1:B:272:THR:HG23	3:B:381:BME:C1	2.01	0.90
1:B:187:MET:HA	1:B:187:MET:HE2	1.56	0.88
1:B:10[B]:LYS:HE3	1:B:11:THR:N	1.90	0.86
1:B:77:HIS:CE1	4:B:382:GOL:H11	2.13	0.84
1:A:12[B]:ARG:CG	1:A:12[B]:ARG:HH11	1.93	0.78
1:A:142:GLN:HE21	1:A:142:GLN:H	1.33	0.77
1:B:187:MET:HA	1:B:187:MET:CE	2.17	0.74
1:A:87:HIS:CE1	1:A:205:PHE:H	2.05	0.74
1:A:87:HIS:HE1	1:A:205:PHE:H	1.39	0.71
1:B:238:VAL:HA	1:B:239:PRO:C	2.09	0.70
1:A:181:ARG:O	1:A:182:LEU:HD23	1.92	0.69
1:B:77:HIS:NE2	4:B:382:GOL:H11	2.08	0.67
1:A:129:GLU:H	1:A:142:GLN:NE2	1.93	0.67
1:A:70:ASP:OD1	5:A:554:HOH:O	2.13	0.65
1:A:129:GLU:H	1:A:142:GLN:HE22	1.46	0.63
1:B:217:CYS:O	5:B:473:HOH:O	2.15	0.63
1:A:87:HIS:HE1	1:A:206:TRP:H	1.49	0.59
1:A:97:ASN:HD21	1:B:77:HIS:HA	1.68	0.58
1:B:87:HIS:CE1	1:B:205:PHE:H	2.23	0.56
1:A:182:LEU:HD21	1:A:265:ALA:CB	2.38	0.54
1:B:10[B]:LYS:CE	1:B:11:THR:N	2.69	0.53
1:B:87:HIS:HE1	1:B:206:TRP:H	1.55	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:187:MET:CA	1:B:187:MET:CE	2.86	0.53
1:A:182:LEU:HD12	1:A:259:ALA:HB1	1.90	0.53
1:A:75:GLY:HA2	4:A:382:GOL:H12	1.90	0.53
1:A:152:ILE:HB	1:A:283:VAL:HG22	1.91	0.52
1:A:12[A]:ARG:NH2	5:A:445:HOH:O	2.28	0.52
1:A:142:GLN:HE21	1:A:142:GLN:N	2.05	0.52
1:A:87:HIS:CE1	1:A:206:TRP:H	2.28	0.51
1:B:152:ILE:HB	1:B:283:VAL:HG22	1.95	0.48
1:B:300:GLY:HA3	2:B:380:FAD:H1'2	1.95	0.48
1:B:52:GLY:HA2	1:B:105:THR:HA	1.99	0.44
1:A:97:ASN:ND2	1:B:78:THR:H	2.15	0.44
1:A:182:LEU:HD21	1:A:265:ALA:HB1	2.00	0.44
1:B:187:MET:H	1:B:187:MET:HE3	1.82	0.44
1:A:97:ASN:HD21	1:B:77:HIS:CA	2.29	0.44
1:A:181:ARG:C	1:A:182:LEU:HD23	2.40	0.43
1:B:186:ARG:NH2	1:B:188:LYS:HD2	2.34	0.42
1:B:186:ARG:HD2	1:B:222:LEU:HB2	2.01	0.42
1:A:213:LEU:HD22	4:A:384:GOL:H31	2.00	0.42
1:B:56:TRP:HA	1:B:101:TRP:HA	2.02	0.41
1:A:311:LEU:HA	1:A:328:TRP:CD1	2.56	0.41
1:B:142:GLN:HE21	1:B:142:GLN:HB3	1.66	0.41
1:A:113:LEU:HD23	1:A:113:LEU:HA	1.88	0.41
1:A:342:LEU:HD23	1:A:362:ALA:HA	2.03	0.41
1:B:272:THR:CG2	5:B:451:HOH:O	2.68	0.41
1:B:272:THR:HG23	3:B:381:BME:C2	2.51	0.40
1:A:182:LEU:HD21	1:A:265:ALA:HB2	2.02	0.40
1:A:103:ILE:HD11	1:A:199:THR:OG1	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	375/379 (99%)	365 (97%)	10 (3%)	0	100	100
1	B	372/379 (98%)	362 (97%)	7 (2%)	3 (1%)	19	7
All	All	747/758 (98%)	727 (97%)	17 (2%)	3 (0%)	41	19

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	160	SER
1	B	10[A]	LYS
1	B	10[B]	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	304/304 (100%)	298 (98%)	6 (2%)	55	40
1	B	300/304 (99%)	294 (98%)	6 (2%)	55	40
All	All	604/608 (99%)	592 (98%)	12 (2%)	55	40

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	LYS
1	A	100	PRO
1	A	127	ASN
1	A	142	GLN
1	A	189	LYS
1	A	263	LYS
1	B	142	GLN
1	B	158	VAL
1	B	179	LEU
1	B	187	MET
1	B	263	LYS
1	B	272	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	ASN
1	A	87	HIS
1	A	97	ASN
1	A	115	ASN
1	A	127	ASN
1	A	142	GLN
1	A	198	ASN
1	A	307	ASN
1	B	33	ASN
1	B	87	HIS
1	B	115	ASN
1	B	127	ASN
1	B	142	GLN
1	B	170	GLN
1	B	198	ASN
1	B	307	ASN
1	B	379	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

9 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	A	384	-	5,5,5	0.48	0	5,5,5	1.30	0
3	BME	B	381	-	3,3,3	1.66	1 (33%)	1,2,2	2.15	1 (100%)
3	BME	A	381	-	3,3,3	0.72	0	1,2,2	2.08	1 (100%)
4	GOL	A	383	-	5,5,5	0.42	0	5,5,5	1.23	1 (20%)
2	FAD	A	380	-	51,58,58	2.13	12 (23%)	60,89,89	2.49	15 (25%)
2	FAD	B	380	-	51,58,58	2.39	17 (33%)	60,89,89	2.05	16 (26%)
4	GOL	B	382	-	5,5,5	0.65	0	5,5,5	1.77	2 (40%)
4	GOL	A	382	-	5,5,5	0.67	0	5,5,5	0.93	0
4	GOL	B	383	-	5,5,5	0.58	0	5,5,5	1.03	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
4	GOL	A	384	-	-	4/4/4/4	-
3	BME	B	381	-	-	1/1/1/1	-
3	BME	A	381	-	-	1/1/1/1	-
4	GOL	A	383	-	-	1/4/4/4	-
2	FAD	A	380	-	-	5/30/50/50	0/6/6/6
2	FAD	B	380	-	-	1/30/50/50	0/6/6/6
4	GOL	B	382	-	-	4/4/4/4	-
4	GOL	A	382	-	-	4/4/4/4	-
4	GOL	B	383	-	-	2/4/4/4	-

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	380	FAD	O4B-C1B	9.03	1.53	1.41
2	B	380	FAD	O4B-C1B	6.61	1.50	1.41
2	B	380	FAD	C10-N1	5.86	1.40	1.33
2	B	380	FAD	C4X-N5	5.69	1.41	1.33
2	B	380	FAD	C1'-N10	-5.42	1.42	1.48
2	B	380	FAD	C9A-N10	5.20	1.45	1.38
2	A	380	FAD	C4X-N5	4.13	1.39	1.33
2	A	380	FAD	O4'-C4'	4.04	1.51	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	380	FAD	O3'-C3'	3.86	1.52	1.43
2	B	380	FAD	O4'-C4'	3.75	1.51	1.43
2	A	380	FAD	C8M-C8	3.70	1.58	1.51
2	B	380	FAD	C2A-N3A	3.58	1.37	1.32
2	B	380	FAD	C5X-N5	3.47	1.41	1.35
2	B	380	FAD	O3'-C3'	3.18	1.50	1.43
2	A	380	FAD	C1'-N10	2.91	1.51	1.48
2	A	380	FAD	C4X-C10	2.75	1.41	1.38
2	B	380	FAD	C9-C8	2.65	1.44	1.37
2	B	380	FAD	C7M-C7	2.58	1.56	1.51
2	B	380	FAD	C2-N3	2.49	1.43	1.38
2	B	380	FAD	C8M-C8	2.45	1.55	1.51
3	B	381	BME	C2-S2	2.42	1.89	1.80
2	A	380	FAD	C9A-C5X	-2.42	1.37	1.42
2	B	380	FAD	C9A-C5X	-2.26	1.38	1.42
2	A	380	FAD	O3B-C3B	2.21	1.48	1.43
2	A	380	FAD	C2B-C1B	-2.20	1.50	1.53
2	B	380	FAD	C4X-C10	2.13	1.41	1.38
2	B	380	FAD	O2B-C2B	2.13	1.48	1.43
2	A	380	FAD	C4A-N3A	2.11	1.38	1.35
2	A	380	FAD	C4'-C3'	-2.06	1.49	1.53
2	B	380	FAD	O2'-C2'	2.04	1.47	1.43

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	380	FAD	C1'-N10-C9A	7.35	124.07	118.29
2	A	380	FAD	C4-N3-C2	7.23	121.25	115.14
2	B	380	FAD	C4-N3-C2	7.09	121.13	115.14
2	A	380	FAD	C10-C4X-N5	-6.14	117.01	121.26
2	A	380	FAD	C5X-C9A-N10	6.08	122.12	117.72
2	A	380	FAD	C4X-N5-C5X	5.82	122.59	116.77
2	B	380	FAD	N3A-C2A-N1A	-5.26	120.46	128.68
2	A	380	FAD	N3A-C2A-N1A	-4.74	121.27	128.68
2	B	380	FAD	C10-C4X-N5	-4.17	118.38	121.26
2	B	380	FAD	C1'-N10-C9A	4.00	121.44	118.29
2	B	380	FAD	O4B-C4B-C5B	-3.98	96.29	109.37
2	B	380	FAD	C4X-N5-C5X	3.91	120.68	116.77
2	A	380	FAD	O4B-C1B-C2B	-3.85	101.30	106.93
2	A	380	FAD	C4X-C4-N3	-3.70	118.38	123.43
2	B	380	FAD	O4B-C1B-C2B	-3.35	102.03	106.93
2	A	380	FAD	C4-C4X-N5	3.23	122.28	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	380	FAD	C4X-C4-N3	-3.17	119.09	123.43
2	A	380	FAD	C6-C5X-C9A	3.17	123.21	119.05
2	A	380	FAD	C7-C6-C5X	-3.09	116.84	121.22
2	A	380	FAD	C9A-C5X-N5	-2.93	117.78	122.36
2	B	380	FAD	P-O3P-PA	-2.92	122.80	132.83
2	B	380	FAD	O4'-C4'-C5'	-2.90	103.40	109.92
2	B	380	FAD	C5'-C4'-C3'	-2.78	106.83	112.20
2	A	380	FAD	C9A-N10-C10	-2.67	118.41	121.91
2	B	380	FAD	N6A-C6A-N1A	2.56	123.89	118.57
4	B	382	GOL	O1-C1-C2	-2.45	98.45	110.20
4	B	382	GOL	O2-C2-C1	-2.38	98.65	109.12
2	A	380	FAD	O2A-PA-O1A	2.34	123.81	112.24
2	B	380	FAD	C4-C4X-N5	2.27	121.19	118.60
4	A	383	GOL	C3-C2-C1	-2.23	103.05	111.70
2	B	380	FAD	C2A-N1A-C6A	2.22	122.54	118.75
2	A	380	FAD	O2B-C2B-C1B	-2.21	102.69	110.85
2	B	380	FAD	O2A-PA-O1A	2.16	122.94	112.24
3	B	381	BME	O1-C1-C2	2.15	119.31	110.83
2	B	380	FAD	C1B-N9A-C4A	-2.12	122.92	126.64
3	A	381	BME	O1-C1-C2	-2.08	102.65	110.83

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	384	GOL	O1-C1-C2-C3
4	A	384	GOL	C1-C2-C3-O3
2	A	380	FAD	C1'-C2'-C3'-C4'
4	B	382	GOL	O1-C1-C2-C3
4	A	382	GOL	O2-C2-C3-O3
4	B	383	GOL	C1-C2-C3-O3
2	A	380	FAD	O2'-C2'-C3'-C4'
4	B	382	GOL	C1-C2-C3-O3
4	A	382	GOL	C1-C2-C3-O3
4	A	384	GOL	O1-C1-C2-O2
4	A	384	GOL	O2-C2-C3-O3
4	B	382	GOL	O1-C1-C2-O2
4	B	383	GOL	O2-C2-C3-O3
4	B	382	GOL	O2-C2-C3-O3
4	A	383	GOL	O1-C1-C2-C3
2	A	380	FAD	O2'-C2'-C3'-O3'
3	B	381	BME	O1-C1-C2-S2

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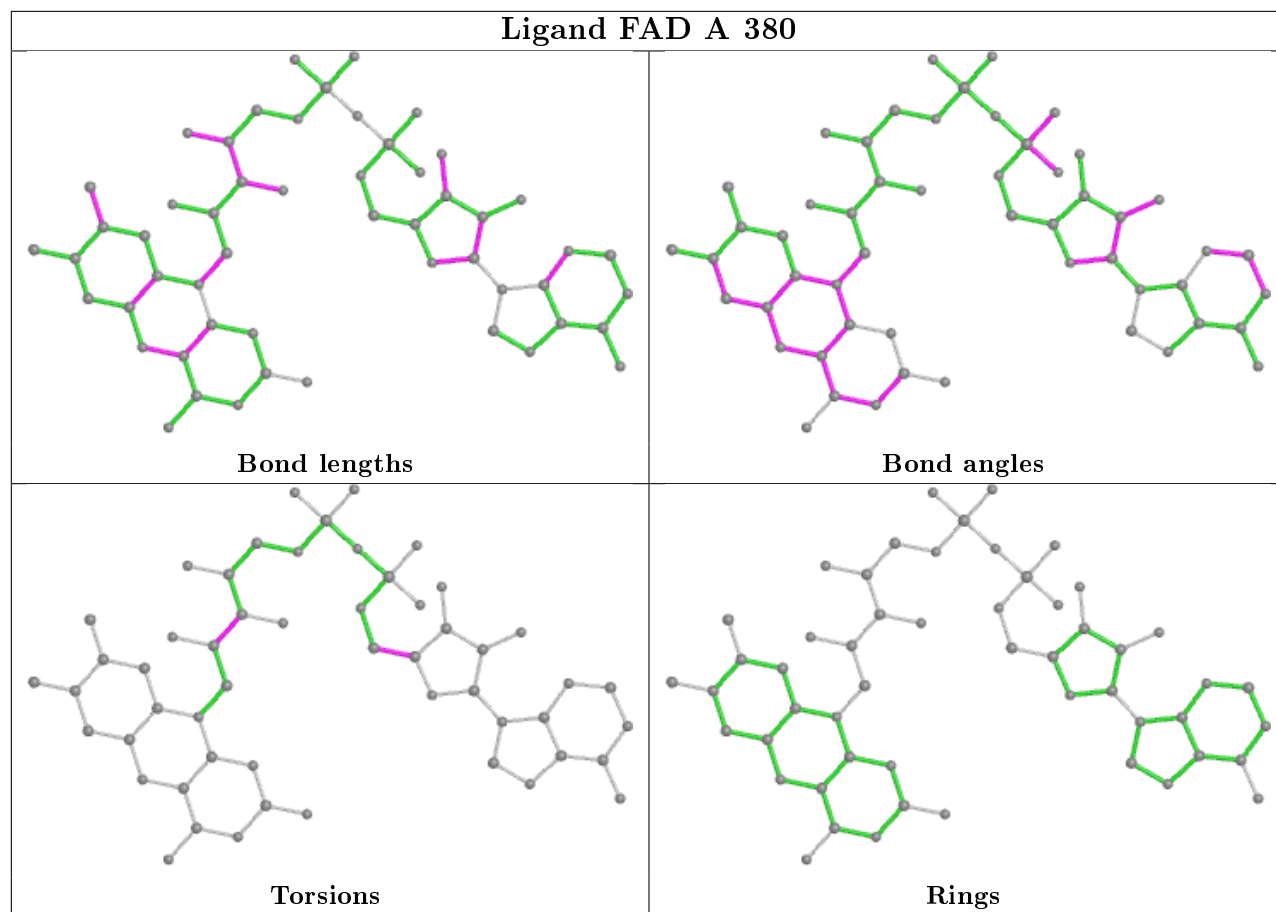
Mol	Chain	Res	Type	Atoms
4	A	382	GOL	O1-C1-C2-O2
4	A	382	GOL	O1-C1-C2-C3
2	A	380	FAD	C1'-C2'-C3'-O3'
2	A	380	FAD	O4B-C4B-C5B-O5B
2	B	380	FAD	O4B-C4B-C5B-O5B
3	A	381	BME	O1-C1-C2-S2

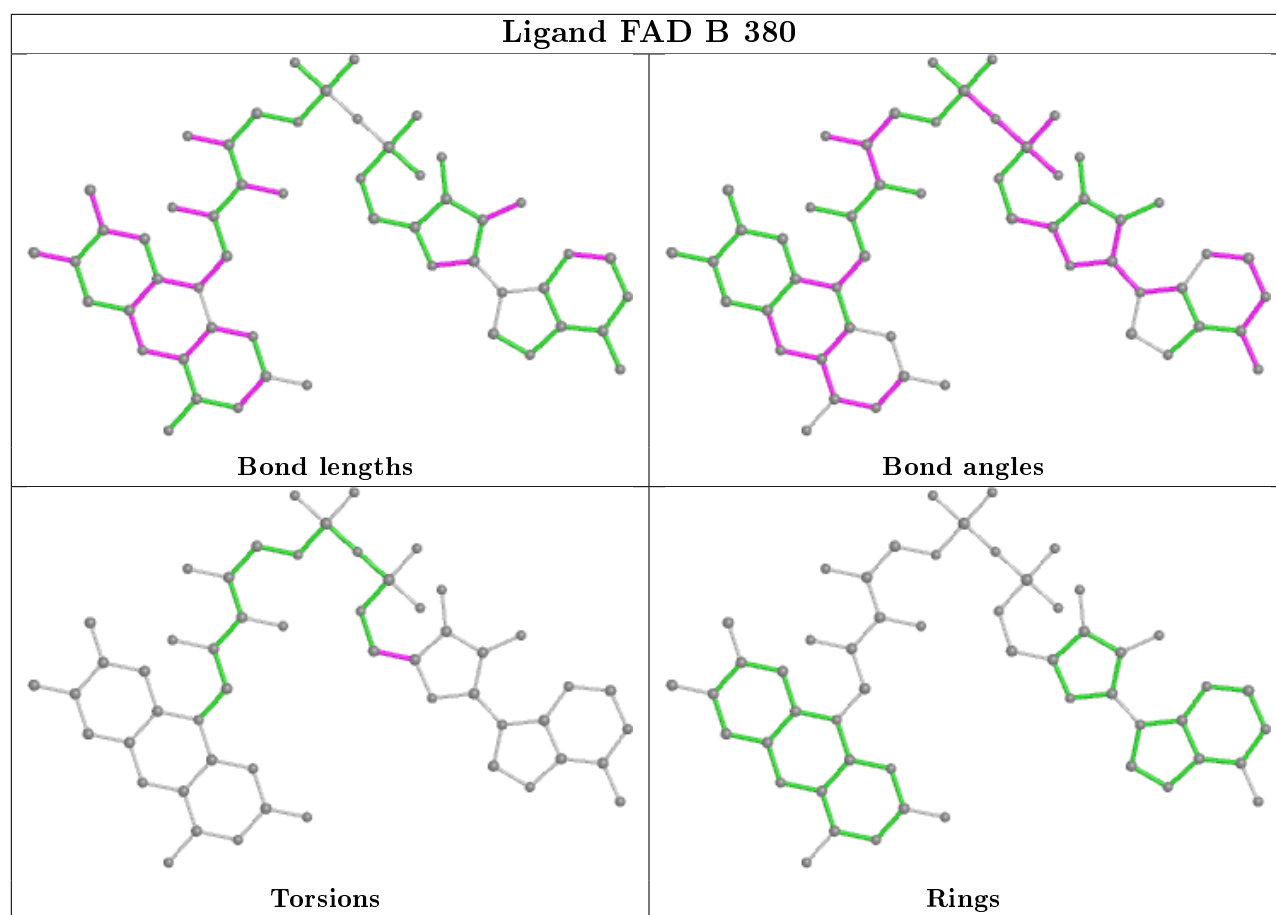
There are no ring outliers.

5 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	384	GOL	1	0
3	B	381	BME	4	0
2	B	380	FAD	1	0
4	B	382	GOL	2	0
4	A	382	GOL	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.





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	370/379 (97%)	-0.21	5 (1%) 75 75	11, 18, 29, 52	0
1	B	371/379 (97%)	0.02	9 (2%) 59 58	15, 24, 39, 51	0
All	All	741/758 (97%)	-0.09	14 (1%) 66 67	11, 20, 36, 52	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	10	LYS	4.1
1	B	9	GLY	3.3
1	B	144	GLY	3.2
1	B	189[A]	LYS	3.1
1	A	318	GLY	3.0
1	B	263	LYS	2.7
1	A	319	SER	2.6
1	A	301	ALA	2.6
1	B	53	ILE	2.4
1	B	318	GLY	2.4
1	B	219	GLU	2.3
1	B	173	TRP	2.0
1	A	304	ALA	2.0
1	B	159	GLY	2.0

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](#)

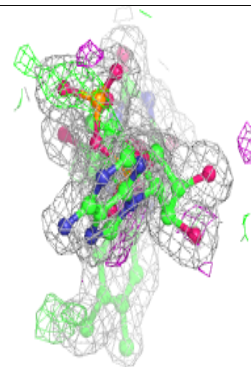
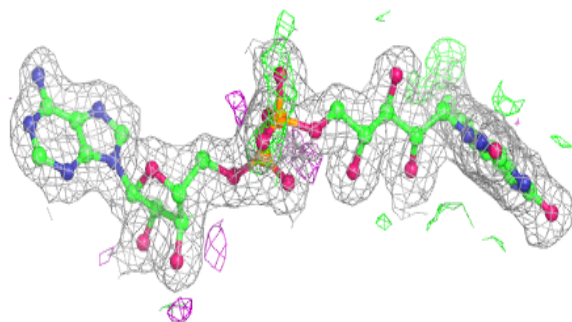
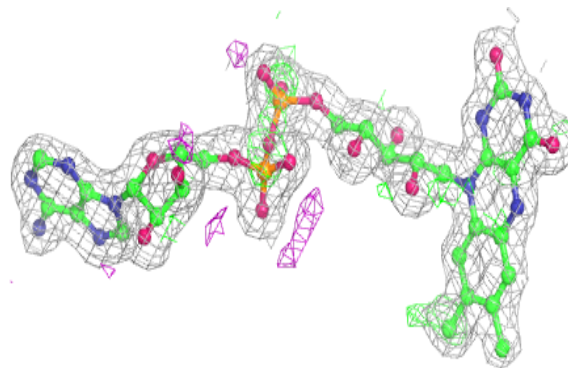
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

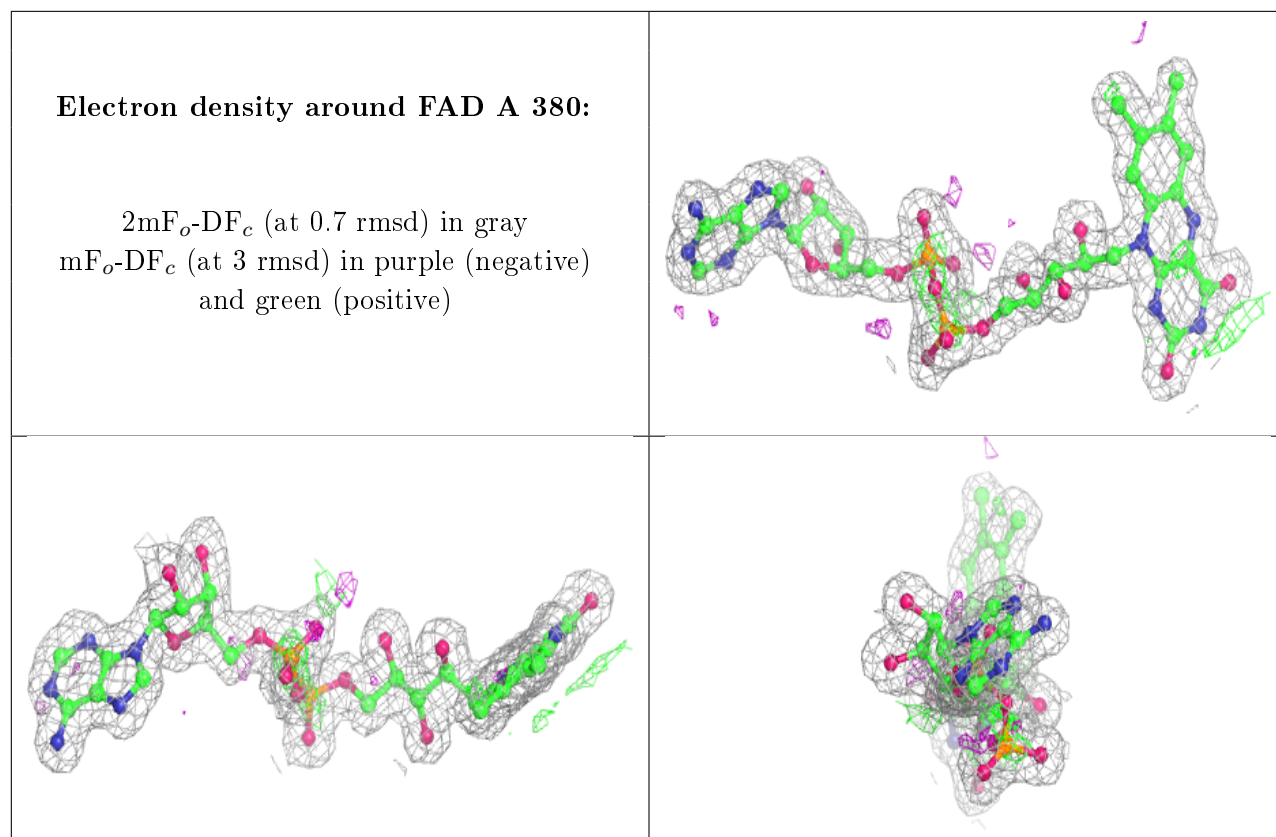
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	B	382	6/6	0.84	0.25	28,37,45,45	0
4	GOL	A	382	6/6	0.85	0.13	22,39,44,51	0
3	BME	B	381	4/4	0.91	0.18	41,44,46,51	0
4	GOL	A	384	6/6	0.93	0.15	22,36,42,43	0
4	GOL	A	383	6/6	0.94	0.26	28,37,39,41	0
4	GOL	B	383	6/6	0.94	0.13	27,34,39,39	0
2	FAD	B	380	53/53	0.96	0.10	15,21,31,33	0
2	FAD	A	380	53/53	0.97	0.10	8,14,17,19	0
3	BME	A	381	4/4	0.99	0.07	16,16,25,27	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 380:

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.