



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 08:08 am BST

PDB ID : 5MXU
Title : Structure of the Y503F mutant of vanillyl alcohol oxidase
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Deposited on : 2017-01-24
Resolution : 2.80 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

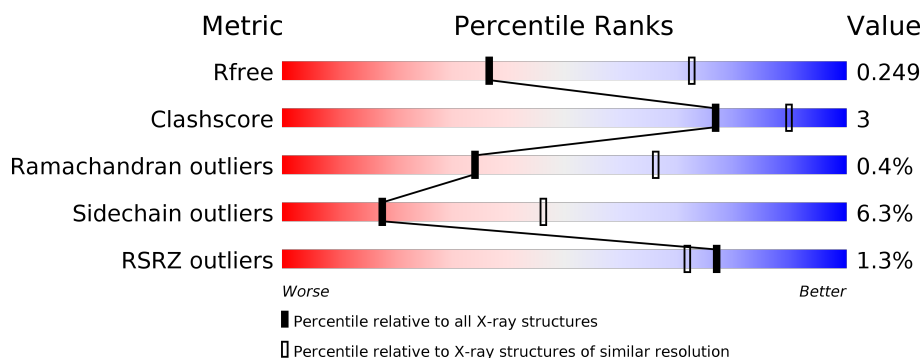
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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	560	<div> <div>2%</div> <div> <div></div> <div>87%</div> <div>10%</div> <div>..</div> </div> </div>
1	B	560	<div> <div>%</div> <div> <div></div> <div>85%</div> <div>13%</div> <div>..</div> </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9056 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 Vanillyl-alcohol oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	553	Total	C	N	O	S	6	0	0
			4345	2789	740	793	23			
1	B	555	Total	C	N	O	S	6	0	0
			4361	2797	743	798	23			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	503	PHE	TYR	engineered mutation	UNP P56216
B	503	PHE	TYR	engineered mutation	UNP P56216

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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



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

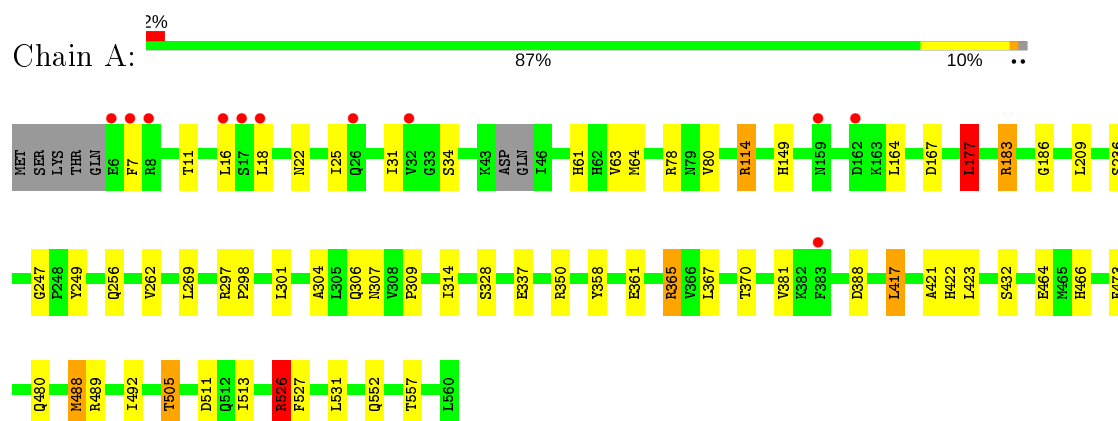
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	101	Total	O	0	0
			101	101		
4	B	131	Total	O	0	0
			131	131		

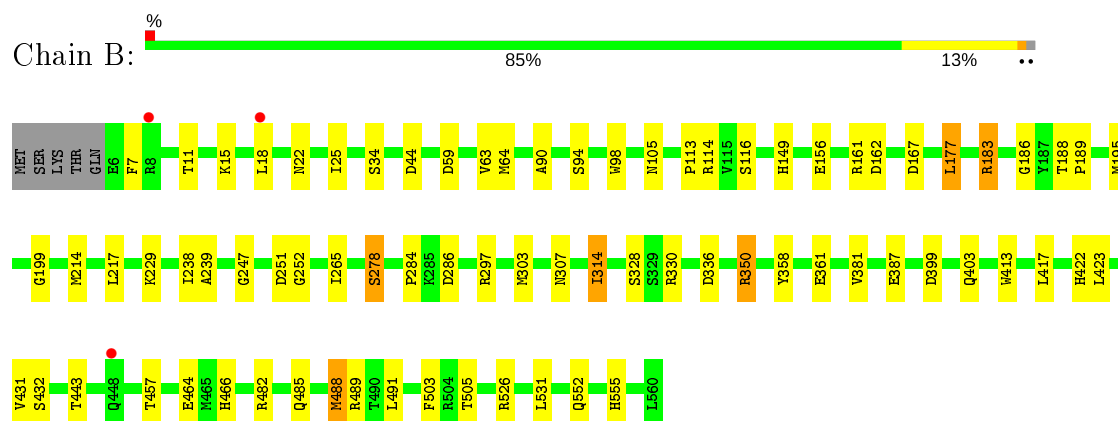
3 Residue-property plots [i](#)

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: Vanillyl-alcohol oxidase



- Molecule 1: Vanillyl-alcohol oxidase



4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	138.01Å 138.01Å 132.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	54.94 – 2.80 54.94 – 2.80	Depositor EDS
% Data completeness (in resolution range)	90.7 (54.94-2.80) 90.7 (54.94-2.80)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.54 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.175 , 0.249 0.182 , 0.249	Depositor DCC
R_{free} test set	1353 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	53.5	Xtriage
Anisotropy	0.006	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 53.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.005 for l,-k,h 0.021 for -l,-k,-h 0.010 for -h,-l,-k 0.004 for -h,l,k 0.024 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9056	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, FAD

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.69	2/4464 (0.0%)	0.77	4/6071 (0.1%)
1	B	0.68	2/4481 (0.0%)	0.77	5/6096 (0.1%)
All	All	0.69	4/8945 (0.0%)	0.77	9/12167 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	328	SER	CB-OG	-21.67	1.14	1.42
1	A	328	SER	CB-OG	-21.20	1.14	1.42
1	A	149	HIS	CB-CG	-18.90	1.16	1.50
1	B	149	HIS	CB-CG	-17.35	1.18	1.50

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	526	ARG	NE-CZ-NH1	6.21	123.41	120.30
1	A	183	ARG	NE-CZ-NH1	5.80	123.20	120.30
1	B	328	SER	CA-CB-OG	-5.63	95.99	111.20
1	A	183	ARG	NE-CZ-NH2	-5.51	117.54	120.30
1	B	526	ARG	NE-CZ-NH1	5.43	123.02	120.30
1	B	183	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	B	183	ARG	NE-CZ-NH2	-5.26	117.67	120.30
1	B	59	ASP	CB-CG-OD2	-5.21	113.61	118.30
1	A	177	LEU	CA-CB-CG	5.13	127.10	115.30

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	4345	0	4246	27	0
1	B	4361	0	4257	28	0
2	A	53	0	29	1	0
2	B	53	0	29	2	0
3	B	12	0	16	0	0
4	A	101	0	0	0	0
4	B	131	0	0	2	0
All	All	9056	0	8577	52	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 3.

All (52) 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:183:ARG:NH2	1:B:247:GLY:O	2.09	0.84
1:B:167:ASP:OD1	1:B:186:GLY:HA3	1.89	0.72
1:B:161:ARG:HD2	4:B:708:HOH:O	1.92	0.70
1:A:505:THR:HG21	1:A:513:ILE:CD1	2.27	0.65
1:A:361:GLU:OE2	1:A:365:ARG:NH1	2.31	0.64
1:B:64:MET:SD	1:B:485:GLN:OE1	2.56	0.63
1:B:105:ASN:HB2	2:B:601:FAD:O1A	1.99	0.62
1:B:297:ARG:HG3	1:B:431:VAL:HB	1.83	0.58
1:B:314:ILE:HD13	1:B:314:ILE:O	2.04	0.58
1:B:423:LEU:HD21	1:B:488:MET:HG3	1.87	0.57
1:A:247:GLY:O	1:B:183:ARG:NH2	2.37	0.56
1:B:63:VAL:HG11	1:B:423:LEU:HD22	1.87	0.56
1:A:31:ILE:O	1:A:78:ARG:NH1	2.38	0.56
1:A:505:THR:HG21	1:A:513:ILE:HD12	1.86	0.56
1:A:61:HIS:HB3	1:A:421:ALA:HB1	1.87	0.55
1:B:278:SER:OG	1:B:399:ASP:OD1	2.24	0.55
1:A:63:VAL:HG11	1:A:423:LEU:HD22	1.90	0.54
1:B:413:TRP:CZ2	2:B:601:FAD:HM72	2.45	0.51
1:A:167:ASP:OD1	1:A:186:GLY:HA3	2.11	0.51
1:A:177:LEU:HD21	1:A:262:VAL:HG12	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7:PHE:O	1:B:18:LEU:HD11	2.13	0.48
1:A:297:ARG:HB3	1:A:298:PRO:HD3	1.95	0.48
1:B:443:THR:HA	1:B:491:LEU:HD21	1.97	0.47
1:B:482:ARG:NH1	4:B:702:HOH:O	2.47	0.47
1:B:177:LEU:HD22	1:B:265:ILE:HG22	1.95	0.47
1:A:417:LEU:CD2	1:A:480:GLN:HE22	2.28	0.47
1:A:18:LEU:O	1:A:22:ASN:ND2	2.38	0.46
1:A:7:PHE:HA	1:A:22:ASN:OD1	2.16	0.46
2:A:600:FAD:O5B	2:A:600:FAD:H8A	2.16	0.46
1:B:214:MET:HB2	1:B:239:ALA:HA	1.97	0.46
1:B:98:TRP:CD2	1:B:113:PRO:HA	2.51	0.45
1:A:306:GLN:HE22	1:A:358:TYR:H	1.65	0.45
1:A:531:LEU:HD22	1:B:531:LEU:HD22	1.99	0.44
1:B:307:ASN:HB3	1:B:358:TYR:CE1	2.53	0.43
1:B:90:ALA:O	1:B:94:SER:N	2.50	0.43
1:B:188:THR:HB	1:B:189:PRO:HD2	2.01	0.42
1:B:22:ASN:HA	1:B:25:ILE:HG22	2.01	0.42
1:A:417:LEU:HD21	1:A:480:GLN:HE22	1.85	0.42
1:B:238:ILE:HG22	1:B:238:ILE:O	2.19	0.42
1:A:249:TYR:O	1:B:252:GLY:HA3	2.19	0.42
1:A:63:VAL:HG12	1:A:473:PHE:CZ	2.55	0.42
1:B:286:ASP:HB2	1:B:350:ARG:HD3	2.02	0.42
1:A:527:PHE:CE2	1:A:531:LEU:HD11	2.55	0.42
1:A:114:ARG:NH1	1:A:511:ASP:OD2	2.53	0.42
1:A:164:LEU:HD13	1:A:269:LEU:HD12	2.01	0.41
1:A:526:ARG:HD3	1:A:526:ARG:HA	1.91	0.41
1:A:80:VAL:HG11	1:A:209:LEU:HD11	2.02	0.41
1:B:156:GLU:HA	1:B:156:GLU:OE1	2.21	0.41
1:A:304:ALA:HA	1:A:367:LEU:HD22	2.03	0.41
1:A:488:MET:CE	1:A:492:ILE:HD11	2.50	0.40
1:B:399:ASP:O	1:B:403:GLN:HG2	2.22	0.40
1:A:307:ASN:O	1:A:309:PRO:HD3	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

of similar resolution.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	549/560 (98%)	516 (94%)	31 (6%)	2 (0%)	34	66
1	B	553/560 (99%)	529 (96%)	22 (4%)	2 (0%)	34	66
All	All	1102/1120 (98%)	1045 (95%)	53 (5%)	4 (0%)	34	66

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	236	SER
1	B	284	PRO
1	A	256	GLN
1	B	199	GLY

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	466/482 (97%)	440 (94%)	26 (6%)	21	51
1	B	468/482 (97%)	435 (93%)	33 (7%)	14	39
All	All	934/964 (97%)	875 (94%)	59 (6%)	18	46

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	THR
1	A	16	LEU
1	A	25	ILE
1	A	34	SER
1	A	64	MET
1	A	114	ARG
1	A	177	LEU
1	A	301	LEU

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Mol	Chain	Res	Type
1	A	314	ILE
1	A	337	GLU
1	A	350	ARG
1	A	365	ARG
1	A	370	THR
1	A	381	VAL
1	A	388	ASP
1	A	417	LEU
1	A	422	HIS
1	A	432	SER
1	A	464	GLU
1	A	466	HIS
1	A	488	MET
1	A	489	ARG
1	A	505	THR
1	A	526	ARG
1	A	552	GLN
1	A	557	THR
1	B	11	THR
1	B	15	LYS
1	B	34	SER
1	B	44	ASP
1	B	114	ARG
1	B	116	SER
1	B	162	ASP
1	B	177	LEU
1	B	195	MET
1	B	217	LEU
1	B	229	LYS
1	B	251	ASP
1	B	278	SER
1	B	303	MET
1	B	314	ILE
1	B	330	ARG
1	B	336	ASP
1	B	350	ARG
1	B	361	GLU
1	B	381	VAL
1	B	387	GLU
1	B	417	LEU
1	B	422	HIS
1	B	432	SER

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Mol	Chain	Res	Type
1	B	457	THR
1	B	464	GLU
1	B	466	HIS
1	B	488	MET
1	B	489	ARG
1	B	503	PHE
1	B	505	THR
1	B	552	GLN
1	B	555	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

4 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$
2	FAD	A	600	1	51,58,58	2.05	8 (15%)	60,89,89	2.07	9 (15%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	B	602	-	5,5,5	0.53	0	5,5,5	0.41	0
2	FAD	B	601	1	51,58,58	2.06	7 (13%)	60,89,89	2.16	11 (18%)
3	GOL	B	603	-	5,5,5	0.39	0	5,5,5	0.53	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	A	600	1	-	10/30/50/50	0/6/6/6
3	GOL	B	602	-	-	4/4/4/4	-
2	FAD	B	601	1	-	9/30/50/50	0/6/6/6
3	GOL	B	603	-	-	0/4/4/4	-

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	600	FAD	C4X-C10	10.92	1.49	1.38
2	B	601	FAD	C4X-C10	10.92	1.49	1.38
2	A	600	FAD	C9A-C5X	4.44	1.51	1.42
2	B	601	FAD	C9A-C5X	3.91	1.50	1.42
2	A	600	FAD	C8-C7	3.72	1.50	1.40
2	B	601	FAD	C4-C4X	3.59	1.47	1.41
2	B	601	FAD	C8-C7	3.28	1.49	1.40
2	B	601	FAD	C9A-N10	3.13	1.42	1.38
2	A	600	FAD	C4-C4X	2.96	1.46	1.41
2	A	600	FAD	C9A-N10	2.90	1.42	1.38
2	B	601	FAD	C2B-C1B	-2.82	1.49	1.53
2	B	601	FAD	C5A-C4A	2.36	1.47	1.40
2	A	600	FAD	C10-N1	2.30	1.36	1.33
2	A	600	FAD	C5A-C4A	2.22	1.46	1.40
2	A	600	FAD	C2A-N3A	2.04	1.35	1.32

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	FAD	C4-N3-C2	9.08	122.81	115.14
2	A	600	FAD	C4-N3-C2	8.76	122.53	115.14
2	A	600	FAD	C1'-N10-C9A	7.06	123.85	118.29
2	B	601	FAD	C1'-N10-C9A	6.61	123.49	118.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	600	FAD	C4-C4X-C10	-4.83	116.75	119.95
2	B	601	FAD	C4-C4X-C10	-4.30	117.11	119.95
2	B	601	FAD	C4X-C4-N3	-4.27	117.59	123.43
2	B	601	FAD	N3A-C2A-N1A	-3.86	122.64	128.68
2	B	601	FAD	C9A-N10-C10	-3.70	117.06	121.91
2	B	601	FAD	P-O3P-PA	-3.65	120.30	132.83
2	A	600	FAD	N3A-C2A-N1A	-3.51	123.19	128.68
2	A	600	FAD	C4X-C4-N3	-3.49	118.65	123.43
2	A	600	FAD	C9A-N10-C10	-3.31	117.58	121.91
2	B	601	FAD	C4X-N5-C5X	3.08	119.85	116.77
2	A	600	FAD	C10-C4X-N5	2.98	123.32	121.26
2	B	601	FAD	C5X-C9A-N10	2.89	119.81	117.72
2	B	601	FAD	O2'-C2'-C3'	-2.23	103.68	109.10
2	B	601	FAD	C2A-N1A-C6A	2.10	122.35	118.75
2	A	600	FAD	N6A-C6A-N1A	2.05	122.83	118.57
2	A	600	FAD	C2A-N1A-C6A	2.00	122.18	118.75

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	600	FAD	N10-C1'-C2'-C3'
2	A	600	FAD	C3'-C4'-C5'-O5'
2	A	600	FAD	O4'-C4'-C5'-O5'
3	B	602	GOL	O1-C1-C2-C3
3	B	602	GOL	C1-C2-C3-O3
3	B	602	GOL	O2-C2-C3-O3
2	B	601	FAD	C3'-C4'-C5'-O5'
2	B	601	FAD	O4'-C4'-C5'-O5'
2	A	600	FAD	C2'-C3'-C4'-O4'
3	B	602	GOL	O1-C1-C2-O2
2	B	601	FAD	PA-O3P-P-O1P
2	B	601	FAD	C2'-C3'-C4'-C5'
2	B	601	FAD	C2'-C3'-C4'-O4'
2	B	601	FAD	C4'-C5'-O5'-P
2	A	600	FAD	C2'-C3'-C4'-C5'
2	B	601	FAD	C5B-O5B-PA-O1A
2	A	600	FAD	PA-O3P-P-O2P
2	A	600	FAD	C4'-C5'-O5'-P
2	A	600	FAD	O3'-C3'-C4'-O4'
2	B	601	FAD	O3'-C3'-C4'-O4'
2	A	600	FAD	PA-O3P-P-O1P

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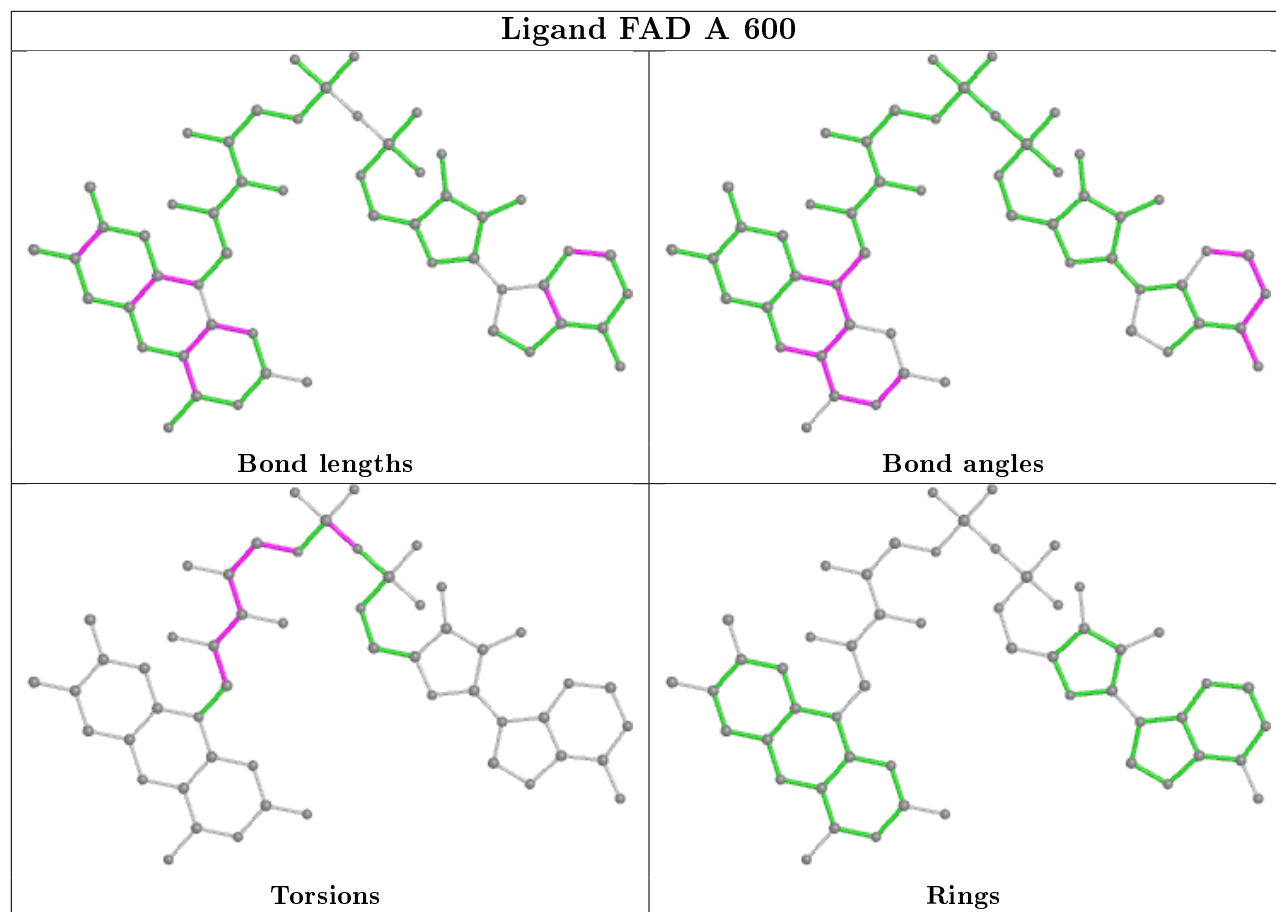
Mol	Chain	Res	Type	Atoms
2	B	601	FAD	PA-O3P-P-O2P
2	A	600	FAD	C1'-C2'-C3'-O3'

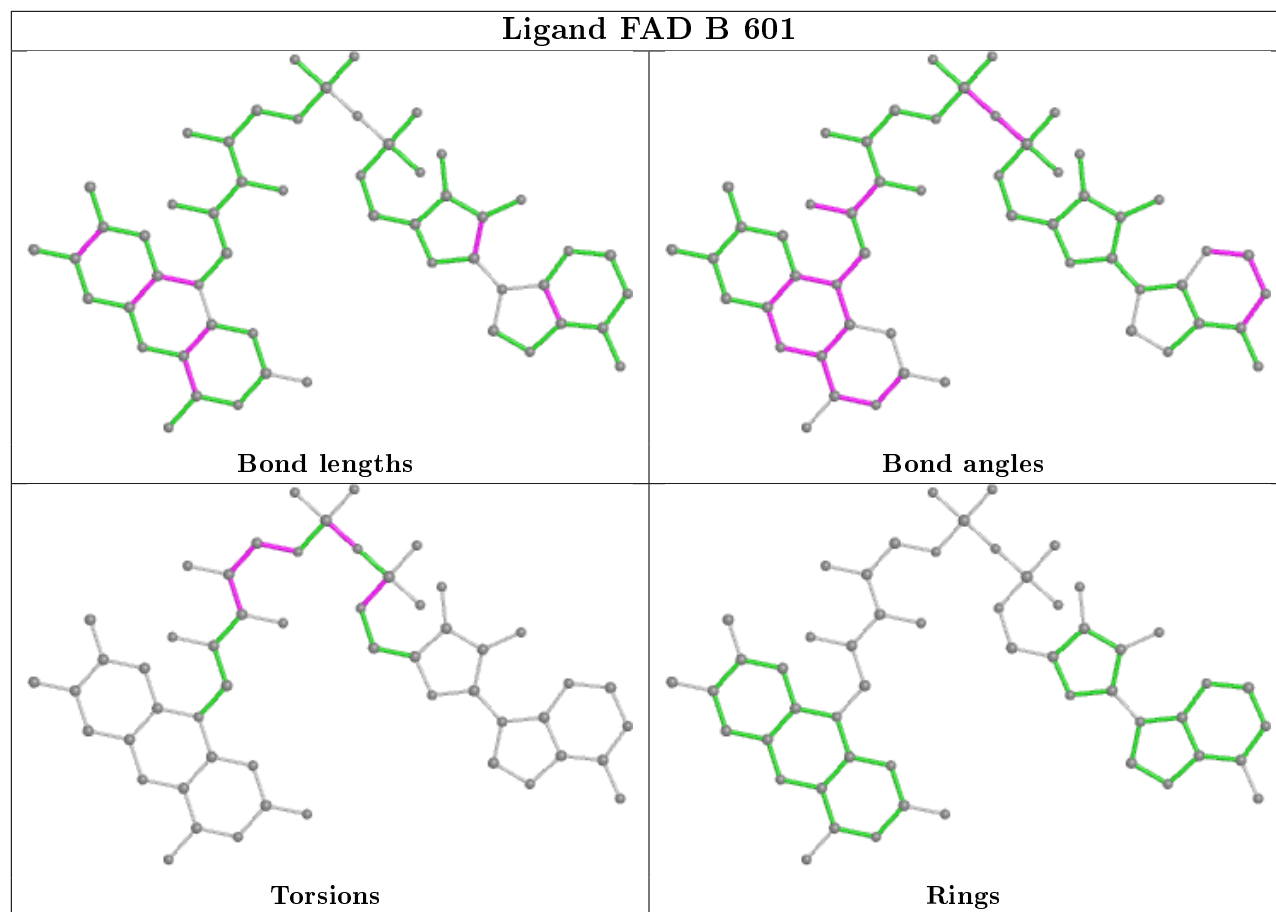
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	600	FAD	1	0
2	B	601	FAD	2	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	553/560 (98%)	-0.13	11 (1%) 65 56	34, 55, 77, 103	2 (0%)
1	B	555/560 (99%)	-0.11	3 (0%) 91 88	32, 49, 71, 89	2 (0%)
All	All	1108/1120 (98%)	-0.12	14 (1%) 77 72	32, 52, 74, 103	4 (0%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	6	GLU	4.2
1	A	159	ASN	3.7
1	A	26	GLN	3.3
1	A	8	ARG	3.1
1	A	17	SER	2.7
1	B	18	LEU	2.5
1	B	448	GLN	2.5
1	A	16	LEU	2.4
1	A	32	VAL	2.4
1	A	7	PHE	2.4
1	A	162	ASP	2.3
1	B	8	ARG	2.2
1	A	383	PHE	2.1
1	A	18	LEU	2.1

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](#)

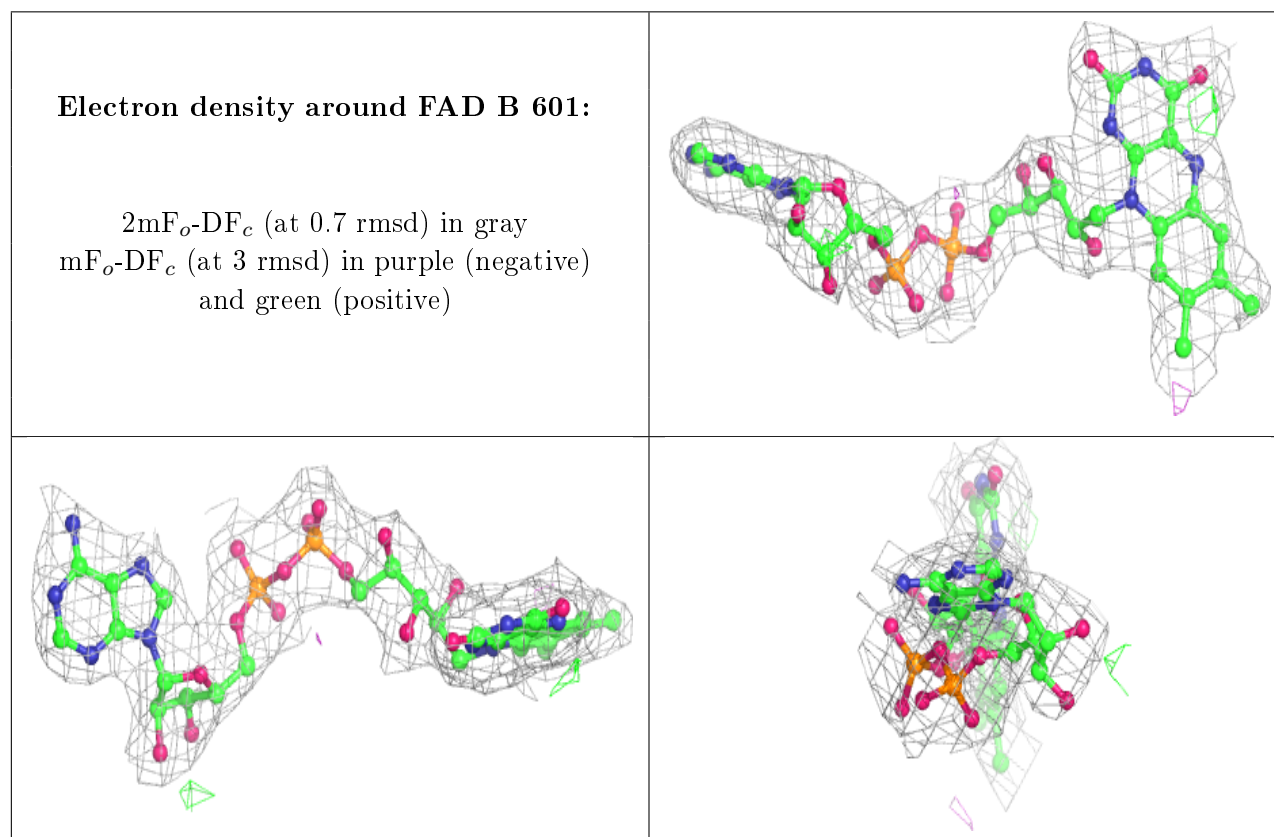
There are no carbohydrates in this entry.

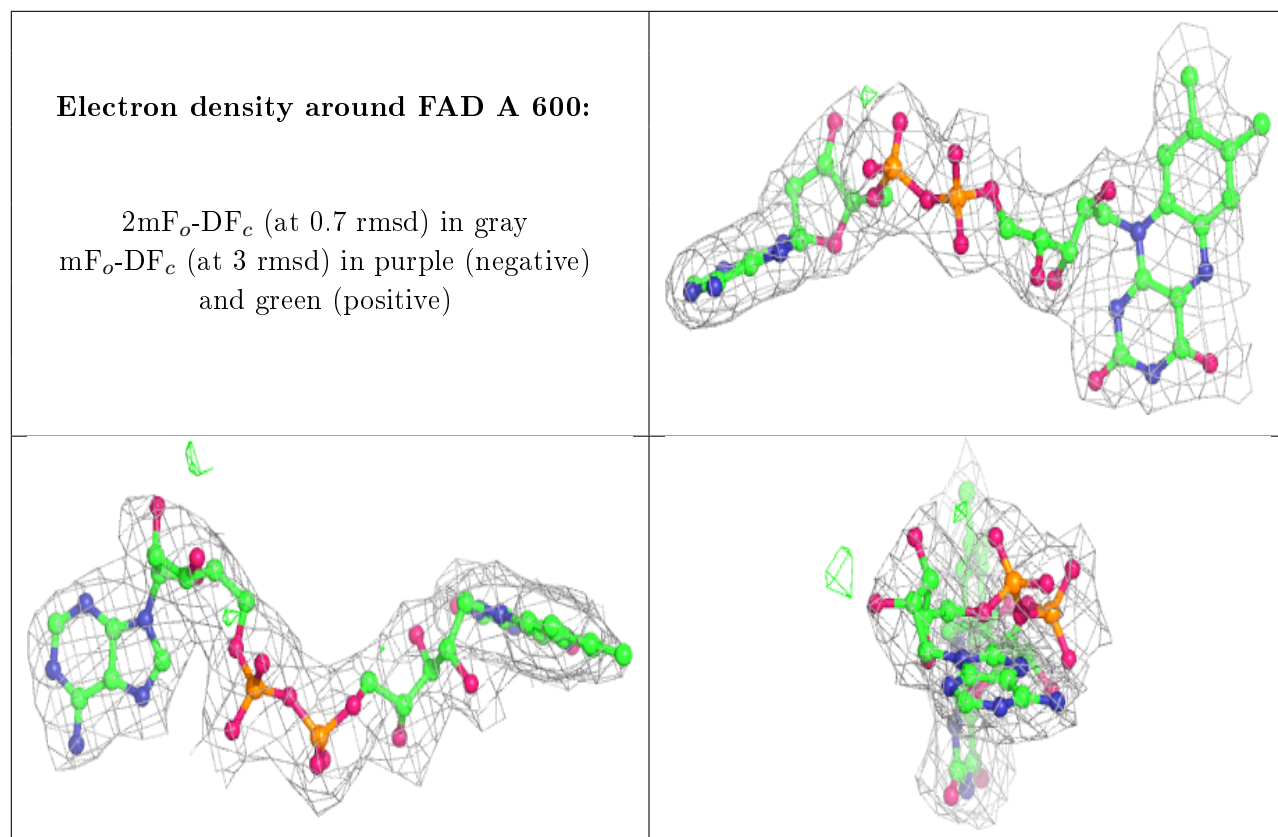
6.4 Ligands [i](#)

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(\AA^2)	Q<0.9
3	GOL	B	602	6/6	0.90	0.23	66,71,80,80	0
3	GOL	B	603	6/6	0.91	0.19	53,59,61,62	0
2	FAD	B	601	53/53	0.98	0.20	32,37,41,42	0
2	FAD	A	600	53/53	0.98	0.13	37,50,56,58	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.