



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:33 PM UTC

PDB ID : 5O8P / pdb_00005o8p
Title : The crystal structure of DfoA bound to FAD, the desferrioxamine biosynthetic pathway cadaverine monooxygenase from the fire blight disease pathogen *Erwinia amylovora*
Authors : Salomone-Stagni, M.; Bartho, J.D.; Polsinelli, I.; Bellini, D.; Walsh, M.A.; Demitri, N.; Benini, S.
Deposited on : 2017-06-14
Resolution : 2.75 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references](#) ①) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

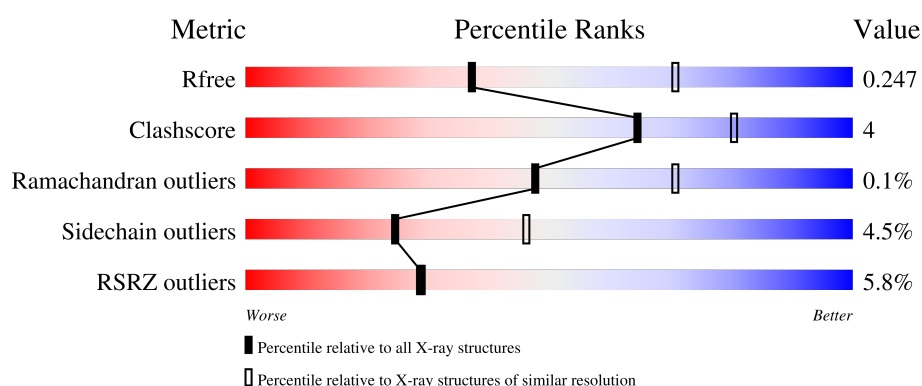
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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}	180053	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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 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 $\leq 5\%$. 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	433	10% 85% 12% ..
1	B	433	2% 89% 9% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7130 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 L-lysine 6-monooxygenase involved in desferrioxamine biosynthesis.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	427	Total	C	N	O	S	0	1	0
			3530	2254	604	658	14			
1	B	428	Total	C	N	O	S	0	0	0
			3526	2250	604	657	15			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP D4I246
A	-1	ALA	-	expression tag	UNP D4I246
A	0	MET	-	expression tag	UNP D4I246
A	1	ALA	-	expression tag	UNP D4I246
B	-2	GLY	-	expression tag	UNP D4I246
B	-1	ALA	-	expression tag	UNP D4I246
B	0	MET	-	expression tag	UNP D4I246
B	1	ALA	-	expression tag	UNP D4I246

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



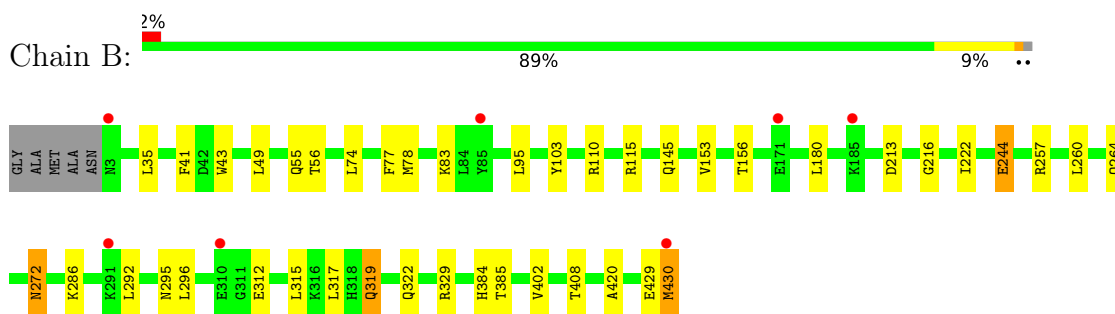
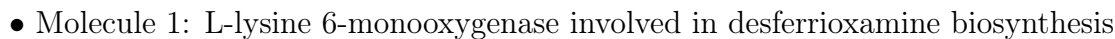
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	12	Total O 12 12	0	0
3	B	9	Total O 9 9	0	0

i

- Molecule 1: L-lysine 6-monooxygenase involved in desferrioxamine biosynthesis



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	164.09Å 164.09Å 168.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	117.57 – 2.75 117.57 – 2.75	Depositor EDS
% Data completeness (in resolution range)	100.0 (117.57-2.75) 100.0 (117.57-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 2.73Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.200 , 0.224 0.227 , 0.247	Depositor DCC
R_{free} test set	3071 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	62.3	Xtriage
Anisotropy	0.402	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 26.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.022 for -h,l,k 0.003 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7130	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.21% 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: 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	0.67	0/3624	0.87	4/4916 (0.1%)
1	B	0.67	0/3619	0.85	1/4908 (0.0%)
All	All	0.67	0/7243	0.86	5/9824 (0.1%)

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	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	213	ASP	CB-CA-C	-5.90	100.99	110.79
1	A	370	ILE	N-CA-CB	-5.71	105.93	112.21
1	A	252	LEU	CA-C-N	-5.47	113.00	119.84
1	A	252	LEU	C-N-CA	-5.47	113.00	119.84
1	A	217	TYR	N-CA-C	5.17	117.80	110.50

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	216	GLY	Peptide

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Mol	Chain	Res	Type	Group
1	B	216	GLY	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	3530	0	3397	38	0
1	B	3526	0	3398	16	0
2	B	53	0	31	0	0
3	A	12	0	0	0	0
3	B	9	0	0	0	0
All	All	7130	0	6826	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 4.

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:251:SER:CB	1:A:253:PRO:HD3	2.02	0.89
1:A:251:SER:HB2	1:A:253:PRO:HD3	1.56	0.88
1:A:253:PRO:O	1:A:257:ARG:HD3	1.80	0.80
1:A:250:HIS:HA	1:A:257:ARG:HH22	1.51	0.76
1:A:354:ILE:HG23	1:A:370:ILE:HG22	1.73	0.70
1:A:252:LEU:H	1:A:257:ARG:NH2	1.89	0.69
1:B:244:GLU:OE1	1:B:420:ALA:HB3	1.95	0.67
1:A:389:VAL:CG1	1:A:395:MET:HG3	2.25	0.66
1:A:251:SER:C	1:A:253:PRO:HD3	2.20	0.65
1:A:251:SER:O	1:A:252:LEU:HB3	1.98	0.64
1:A:362:TYR:HB3	1:A:370:ILE:HD12	1.79	0.63
1:B:55:GLN:OE1	1:B:272:ASN:HB2	1.98	0.63
1:A:385:THR:HG23	1:A:386:HIS:HD2	1.67	0.58
1:A:251:SER:HB2	1:A:253:PRO:CD	2.32	0.58
1:A:286:LYS:HB3	1:A:292:LEU:HD21	1.87	0.56
1:A:364:VAL:HB	1:A:385:THR:HG21	1.91	0.53
1:A:49:LEU:HD23	1:A:180:LEU:HD21	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:286:LYS:HB3	1:B:292:LEU:HD21	1.91	0.52
1:A:252:LEU:N	1:A:253:PRO:HD3	2.24	0.52
1:A:222:ILE:HD13	1:A:317:LEU:HD11	1.90	0.52
1:A:380:ASN:O	1:A:381:ALA:HB3	2.09	0.52
1:A:354:ILE:HG23	1:A:370:ILE:CG2	2.40	0.51
1:B:319:GLN:HG3	1:B:322:GLN:HB3	1.93	0.51
1:B:49:LEU:HD23	1:B:180:LEU:HD21	1.92	0.51
1:A:387:GLY:O	1:A:390:THR:HG23	2.12	0.49
1:A:419:ILE:H	1:A:419:ILE:HD13	1.76	0.49
1:A:256:LYS:HA	1:A:259:GLU:HG2	1.93	0.49
1:A:245:TRP:CH2	1:A:249:PHE:CD1	3.01	0.49
1:B:74:LEU:HG	1:B:95:LEU:HD11	1.95	0.48
1:B:222:ILE:HD13	1:B:317:LEU:HD11	1.95	0.48
1:B:295:ASN:C	1:B:296:LEU:HD12	2.38	0.47
1:A:295:ASN:C	1:A:296:LEU:HD12	2.38	0.47
1:A:233[B]:TYR:CD1	1:A:233[B]:TYR:N	2.78	0.46
1:A:74:LEU:HG	1:A:95:LEU:HD11	1.98	0.46
1:A:250:HIS:HA	1:A:257:ARG:NH2	2.26	0.46
1:A:252:LEU:N	1:A:253:PRO:CD	2.80	0.44
1:B:260:LEU:HD12	1:B:264:GLN:NE2	2.33	0.44
1:B:286:LYS:CB	1:B:292:LEU:HD21	2.47	0.44
1:A:286:LYS:CB	1:A:292:LEU:HD21	2.48	0.43
1:A:364:VAL:HB	1:A:385:THR:CG2	2.48	0.43
1:B:384:HIS:CD2	1:B:385:THR:HG23	2.53	0.43
1:A:251:SER:OG	1:A:253:PRO:HD3	2.19	0.43
1:B:429:GLU:O	1:B:430:MET:HG2	2.18	0.43
1:B:78:MET:HG3	1:B:83:LYS:HG3	2.01	0.42
1:B:312:GLU:OE1	1:B:329:ARG:NH1	2.53	0.41
1:B:77:PHE:CE2	1:B:103:TYR:HA	2.56	0.41
1:A:47:MET:HE3	1:A:176:SER:HB2	2.03	0.41
1:A:259:GLU:O	1:A:262:ALA:O	2.38	0.41
1:A:252:LEU:O	1:A:252:LEU:HG	2.20	0.41
1:A:249:PHE:CZ	1:A:257:ARG:HG3	2.56	0.41
1:A:251:SER:CB	1:A:253:PRO:CD	2.85	0.41
1:B:41:PHE:CE2	1:B:43:TRP:HB2	2.56	0.41
1:A:267:LEU:HB3	1:A:388:PHE:CD2	2.56	0.40
1:A:78:MET:HG3	1:A:83:LYS:HG3	2.02	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	426/433 (98%)	406 (95%)	19 (4%)	1 (0%)	43	64
1	B	426/433 (98%)	410 (96%)	16 (4%)	0	100	100
All	All	852/866 (98%)	816 (96%)	35 (4%)	1 (0%)	48	71

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	252	LEU

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	379/381 (100%)	360 (95%)	19 (5%)	22	42
1	B	379/381 (100%)	364 (96%)	15 (4%)	28	50
All	All	758/762 (100%)	724 (96%)	34 (4%)	24	46

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	LEU
1	A	53	HIS
1	A	56	THR
1	A	110	ARG
1	A	153	VAL

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Mol	Chain	Res	Type
1	A	156	THR
1	A	174	VAL
1	A	185	LYS
1	A	251	SER
1	A	257	ARG
1	A	260	LEU
1	A	267	LEU
1	A	272	ASN
1	A	315	LEU
1	A	319	GLN
1	A	359	LYS
1	A	361	ARG
1	A	402	VAL
1	A	419	ILE
1	B	35	LEU
1	B	56	THR
1	B	110	ARG
1	B	115	ARG
1	B	145	GLN
1	B	153	VAL
1	B	156	THR
1	B	244	GLU
1	B	257	ARG
1	B	272	ASN
1	B	315	LEU
1	B	319	GLN
1	B	402	VAL
1	B	408	THR
1	B	430	MET

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

Mol	Chain	Res	Type
1	A	37	GLN
1	A	55	GLN
1	A	218	GLN
1	A	220	ASN
1	A	272	ASN
1	A	295	ASN
1	A	299	HIS
1	A	342	GLN
1	B	31	ASN

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Mol	Chain	Res	Type
1	B	128	ASN
1	B	218	GLN
1	B	220	ASN
1	B	264	GLN
1	B	295	ASN
1	B	299	HIS
1	B	373	HIS
1	B	427	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	B	501	-	58,58,58	1.72	10 (17%)	85,89,89	1.69	21 (24%)

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	B	501	-	-	9/34/50/50	0/6/6/6

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	FAD	C9A-C5X	6.34	1.51	1.41
2	B	501	FAD	C5A-C4A	4.72	1.47	1.39
2	B	501	FAD	C8-C7	3.61	1.49	1.40
2	B	501	FAD	PA-O3P	3.54	1.63	1.59
2	B	501	FAD	P-O3P	3.51	1.63	1.59
2	B	501	FAD	C5A-C6A	3.00	1.49	1.41
2	B	501	FAD	C8A-N7A	2.62	1.36	1.31
2	B	501	FAD	C10-N10	2.53	1.42	1.37
2	B	501	FAD	C4X-N5	2.21	1.35	1.30
2	B	501	FAD	C4-N3	-2.06	1.35	1.38

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	FAD	C5A-C4A-N3A	-5.45	119.21	126.72
2	B	501	FAD	N3A-C4A-N9A	4.19	134.30	127.17
2	B	501	FAD	C2A-N3A-C4A	3.91	121.38	111.83
2	B	501	FAD	N3A-C2A-N1A	-3.84	122.78	128.58
2	B	501	FAD	C4A-C5A-N7A	-3.80	106.24	110.58
2	B	501	FAD	O2P-P-O3P	3.08	115.59	107.27
2	B	501	FAD	C4A-N9A-C8A	3.04	108.93	105.74
2	B	501	FAD	C5A-N7A-C8A	2.85	107.93	103.45
2	B	501	FAD	C4X-C10-N1	-2.79	117.75	124.59
2	B	501	FAD	C6A-C5A-N7A	2.71	137.32	132.09
2	B	501	FAD	C4-C4X-N5	2.70	121.94	118.21
2	B	501	FAD	O4B-C1B-N9A	2.62	113.12	108.09
2	B	501	FAD	C4X-C10-N10	2.60	120.21	116.48
2	B	501	FAD	C10-N1-C2	2.54	122.35	116.85
2	B	501	FAD	N9A-C8A-N7A	-2.49	110.41	113.94
2	B	501	FAD	O4-C4-C4X	-2.39	120.23	126.53
2	B	501	FAD	C4X-C4-N3	2.27	119.03	113.25
2	B	501	FAD	O2-C2-N1	-2.24	118.09	121.80
2	B	501	FAD	C9A-N10-C10	-2.21	117.38	120.75
2	B	501	FAD	C4-N3-C2	-2.08	121.94	125.64
2	B	501	FAD	C2A-N1A-C6A	2.04	122.08	118.73

There are no chirality outliers.

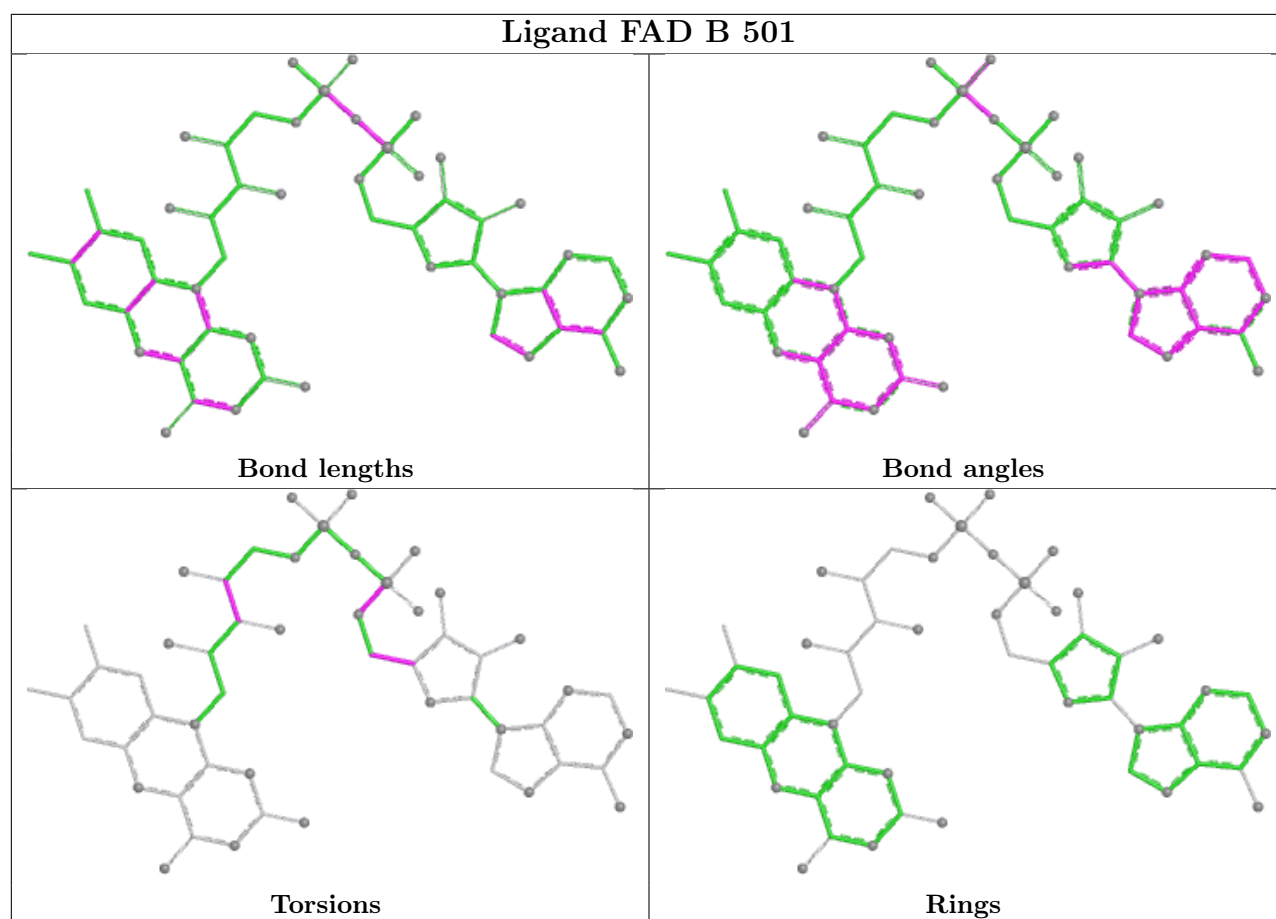
All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	501	FAD	C5B-O5B-PA-O1A
2	B	501	FAD	C5B-O5B-PA-O2A
2	B	501	FAD	C5B-O5B-PA-O3P
2	B	501	FAD	C2'-C3'-C4'-O4'
2	B	501	FAD	C2'-C3'-C4'-C5'
2	B	501	FAD	O3'-C3'-C4'-O4'
2	B	501	FAD	O4B-C4B-C5B-O5B
2	B	501	FAD	C3B-C4B-C5B-O5B
2	B	501	FAD	O3'-C3'-C4'-C5'

There are no ring outliers.

No monomer is involved in short contacts.

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

6.1 Protein, DNA and RNA chains

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	427/433 (98%)	0.48	43 (10%) 12 11	22, 39, 98, 136	1 (0%)
1	B	428/433 (98%)	0.18	7 (1%) 70 69	29, 42, 63, 87	0
All	All	855/866 (98%)	0.33	50 (5%) 29 29	22, 41, 78, 136	1 (0%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	43	TRP	6.7
1	A	252	LEU	6.6
1	A	233[A]	TYR	6.4
1	A	44	HIS	5.6
1	A	156	THR	5.5
1	A	383	LEU	5.2
1	A	45	THR	4.8
1	A	255	ALA	4.7
1	A	384	HIS	4.7
1	A	391	PRO	4.7
1	A	268	TYR	4.6
1	A	269	LYS	4.5
1	A	385	THR	4.4
1	A	263	SER	4.2
1	A	253	PRO	4.0
1	B	3	ASN	4.0
1	A	388	PHE	3.9
1	A	254	ALA	3.9
1	A	48	MET	3.8
1	A	389	VAL	3.8
1	A	46	GLY	3.7
1	A	381	ALA	3.6
1	B	430	MET	3.4
1	A	340	HIS	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	390	THR	3.3
1	A	47	MET	3.1
1	A	265	LYS	3.0
1	A	291	LYS	2.9
1	A	250	HIS	2.9
1	B	185	LYS	2.8
1	B	171	GLU	2.7
1	A	256	LYS	2.7
1	A	157	GLY	2.6
1	A	259	GLU	2.6
1	A	341	TYR	2.6
1	A	251	SER	2.6
1	A	262	ALA	2.6
1	B	85	TYR	2.6
1	A	260	LEU	2.5
1	A	258	ASP	2.4
1	A	386	HIS	2.4
1	A	360	ASP	2.3
1	A	55	GLN	2.3
1	A	419	ILE	2.3
1	B	310	GLU	2.3
1	A	352	GLN	2.3
1	A	197	SER	2.2
1	A	357	ASP	2.2
1	B	291	LYS	2.1
1	A	387	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](#)

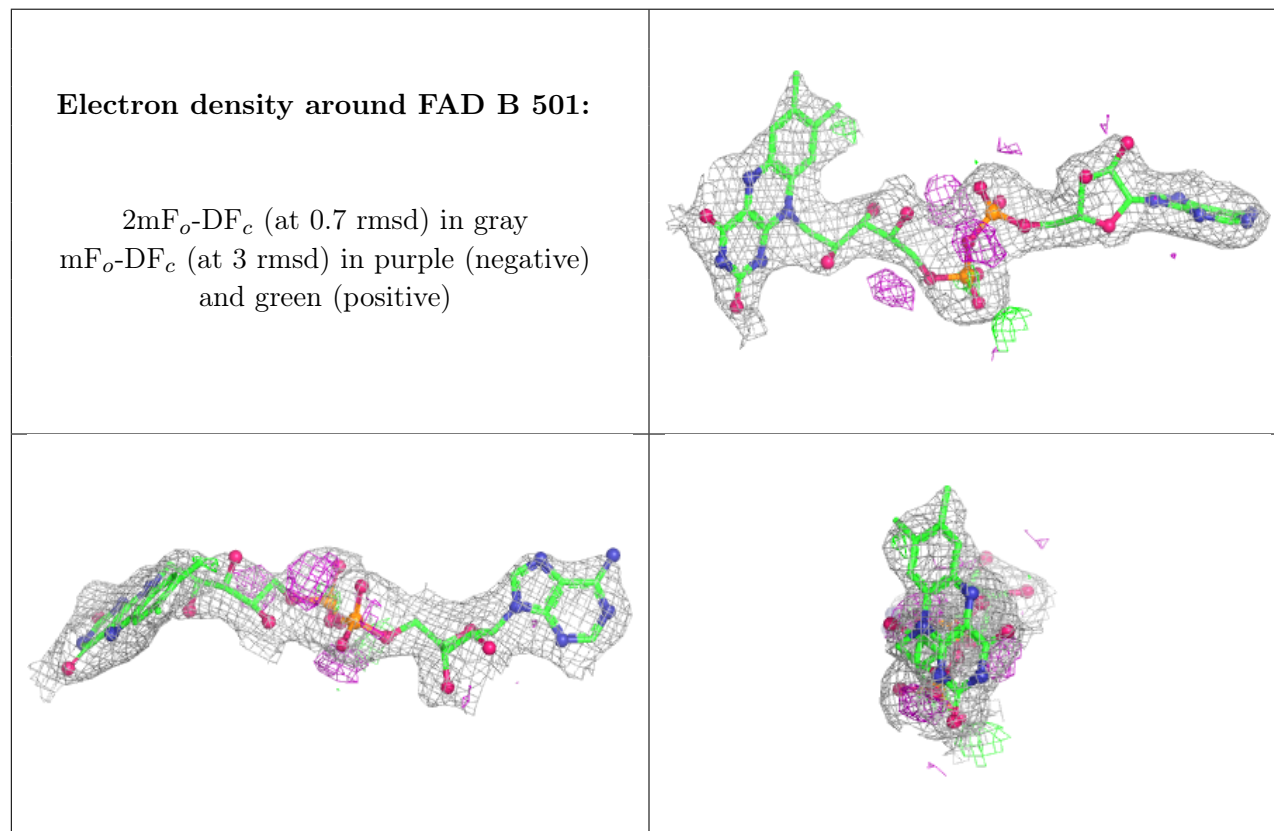
There are no oligosaccharides in this entry.

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($\text{\AA}^2$)	Q<0.9
2	FAD	B	501	53/53	0.91	0.16	68,92,105,113	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 ⓘ

There are no such residues in this entry.