



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:49 PM UTC

PDB ID : 1NPX / pdb_00001npx
Title : STRUCTURE OF NADH PEROXIDASE FROM STREPTOCOCCUS FAECALIS 10C1 REFINED AT 2.16 ANGSTROMS RESOLUTION
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Deposited on : 1991-08-02
Resolution : 2.16 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 ⓘ](#)) 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

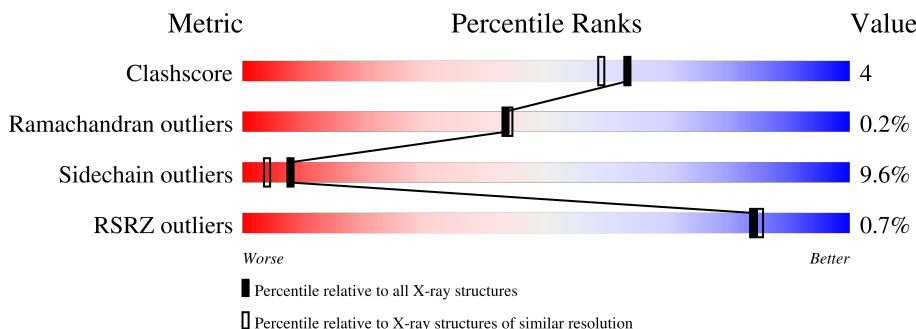
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.16 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))
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	447	78% 15% 6%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3915 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 NADH PEROXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	447	Total	C	N	O	S	0	0	0
			3493	2225	573	683	12			

- 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	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

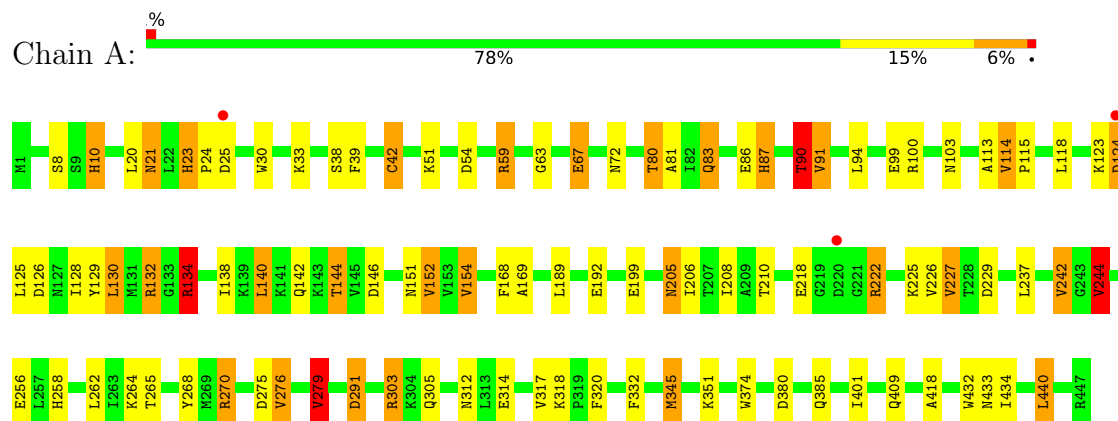
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	369	Total	O	0	0
			369	369		

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: NADH PEROXIDASE



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	77.20Å 134.50Å 145.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.00 – 2.16 7.00 – 2.16	Depositor EDS
% Data completeness (in resolution range)	(Not available) (7.00-2.16) 88.8 (7.00-2.16)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.177 , (Not available) 0.171 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	19.7	Xtriage
Anisotropy	0.195	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 70.1	EDS
L-test for twinning ¹	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3915	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Theoretical 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: OCS, 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	1.03	12/3547 (0.3%)	1.74	57/4807 (1.2%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	154	VAL	CA-CB	9.49	1.67	1.54
1	A	276	VAL	CA-CB	7.82	1.64	1.54
1	A	87	HIS	CD2-NE2	-6.52	1.30	1.37
1	A	258	HIS	CD2-NE2	-6.09	1.31	1.37
1	A	23	HIS	CD2-NE2	-6.04	1.31	1.37
1	A	90	THR	CA-CB	5.81	1.62	1.53
1	A	227	VAL	CA-CB	5.80	1.62	1.54
1	A	87	HIS	CG-ND1	-5.60	1.32	1.38
1	A	10	HIS	CD2-NE2	-5.18	1.32	1.37
1	A	23	HIS	CG-ND1	-5.12	1.32	1.38
1	A	152	VAL	CA-CB	5.11	1.61	1.54
1	A	401	ILE	CA-CB	5.07	1.61	1.54

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	132	ARG	NE-CZ-NH2	-13.47	107.08	119.20
1	A	132	ARG	NE-CZ-NH1	11.59	133.09	121.50
1	A	21	ASN	CA-CB-CG	11.28	123.88	112.60
1	A	205	ASN	CA-CB-CG	10.28	122.88	112.60
1	A	317	VAL	N-CA-CB	-9.35	101.92	112.21
1	A	124	ASP	N-CA-C	9.19	124.42	107.60
1	A	103	ASN	OD1-CG-ND2	-8.80	113.80	122.60
1	A	114	VAL	N-CA-CB	-8.66	99.09	111.21
1	A	80	THR	N-CA-CB	-8.42	97.70	110.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	303	ARG	NE-CZ-NH1	7.66	129.16	121.50
1	A	244	VAL	N-CA-CB	-7.42	99.29	111.45
1	A	279	VAL	N-CA-CB	-7.14	97.35	112.00
1	A	38	SER	N-CA-C	7.14	121.25	111.54
1	A	144	THR	N-CA-CB	-7.11	99.27	110.22
1	A	83	GLN	OE1-CD-NE2	-7.05	115.55	122.60
1	A	205	ASN	N-CA-C	6.82	125.33	110.80
1	A	291	ASP	CA-CB-CG	-6.81	105.79	112.60
1	A	433	ASN	CA-CB-CG	6.72	119.33	112.60
1	A	125	LEU	N-CA-C	-6.58	101.84	110.53
1	A	72	ASN	CA-CB-CG	-6.56	106.04	112.60
1	A	226	VAL	N-CA-C	-6.44	99.09	108.11
1	A	24	PRO	CA-C-N	6.43	130.08	120.31
1	A	24	PRO	C-N-CA	6.43	130.08	120.31
1	A	385	GLN	OE1-CD-NE2	-6.19	116.41	122.60
1	A	33	LYS	CB-CG-CD	-6.18	97.07	111.30
1	A	59	ARG	NE-CZ-NH1	6.14	127.64	121.50
1	A	169	ALA	N-CA-C	-6.10	104.55	111.14
1	A	434	ILE	N-CA-CB	-5.98	103.00	110.47
1	A	374	TRP	CB-CG-CD1	-5.96	117.95	126.90
1	A	168	PHE	CA-CB-CG	-5.88	107.92	113.80
1	A	320	PHE	CA-C-N	5.50	125.12	119.56
1	A	320	PHE	C-N-CA	5.50	125.12	119.56
1	A	409	GLN	OE1-CD-NE2	-5.48	117.12	122.60
1	A	67	GLU	CA-CB-CG	-5.38	103.34	114.10
1	A	199	GLU	CA-CB-CG	5.35	124.80	114.10
1	A	409	GLN	O-C-N	-5.32	116.48	122.12
1	A	229	ASP	CA-CB-CG	5.29	117.89	112.60
1	A	134	ARG	NE-CZ-NH2	-5.28	114.45	119.20
1	A	258	HIS	CB-CG-CD2	-5.23	124.40	131.20
1	A	132	ARG	CB-CG-CD	5.23	123.32	111.30
1	A	128	ILE	N-CA-C	-5.22	99.43	106.85
1	A	332	PHE	CA-CB-CG	-5.20	108.60	113.80
1	A	146	ASP	CA-CB-CG	5.18	117.78	112.60
1	A	432	TRP	CG-CD2-CE3	5.16	139.06	133.90
1	A	10	HIS	N-CA-C	5.13	117.54	111.33
1	A	59	ARG	NE-CZ-NH2	-5.12	114.59	119.20
1	A	54	ASP	CA-CB-CG	5.11	117.71	112.60
1	A	303	ARG	CB-CG-CD	5.11	123.04	111.30
1	A	151	ASN	CA-CB-CG	-5.10	107.50	112.60
1	A	380	ASP	CA-C-N	5.09	124.70	119.05
1	A	380	ASP	C-N-CA	5.09	124.70	119.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	124	ASP	CA-CB-CG	-5.07	107.53	112.60
1	A	242	VAL	N-CA-CB	-5.05	101.05	111.76
1	A	434	ILE	N-CA-C	5.05	116.38	110.62
1	A	25	ASP	N-CA-C	5.02	117.64	111.82
1	A	80	THR	CA-CB-OG1	-5.01	102.08	109.60
1	A	206	ILE	N-CA-C	-5.00	100.55	107.80

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	3493	0	3489	28	0
2	A	53	0	31	1	0
3	A	369	0	0	3	0
All	All	3915	0	3520	28	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 (28) 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:268:TYR:HB2	1:A:270:ARG:HD2	1.74	0.68
1:A:418:ALA:HB1	1:A:440:LEU:HD13	1.77	0.67
1:A:23:HIS:HE1	3:A:573:HOH:O	1.82	0.61
1:A:312:ASN:HD21	1:A:318:LYS:H	1.51	0.58
1:A:83:GLN:HE22	1:A:90:THR:HG22	1.67	0.58
1:A:87:HIS:HE1	1:A:275:ASP:OD2	1.89	0.56
1:A:91:VAL:HG22	1:A:100:ARG:HG2	1.86	0.56
1:A:279:VAL:HG22	1:A:305:GLN:HB3	1.88	0.56
1:A:81:ALA:HB3	1:A:90:THR:HG23	1.87	0.55
1:A:115:PRO:HB2	1:A:130:LEU:HG	1.89	0.54
1:A:345:MET:HG2	3:A:758:HOH:O	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:ARG:NH2	1:A:244:VAL:HG21	2.26	0.50
1:A:39:PHE:CD1	1:A:134:ARG:HG2	2.48	0.48
1:A:218:GLU:HB2	1:A:225:LYS:HB2	1.96	0.47
1:A:208:ILE:HG22	1:A:210:THR:HG23	1.95	0.47
1:A:8:SER:HA	1:A:30:TRP:CZ2	2.50	0.46
1:A:126:ASP:HB2	1:A:222:ARG:HB3	1.98	0.46
1:A:129:TYR:HD2	1:A:140:LEU:HD13	1.81	0.45
1:A:10:HIS:ND1	1:A:303:ARG:HD2	2.32	0.45
1:A:63:GLY:O	1:A:67:GLU:HG3	2.17	0.45
1:A:312:ASN:HD21	1:A:318:LYS:N	2.13	0.44
1:A:91:VAL:CG2	1:A:100:ARG:HG2	2.49	0.43
1:A:42:OCS:OD2	2:A:449:FAD:C10	2.67	0.43
1:A:113:ALA:HB1	1:A:244:VAL:HG13	2.02	0.42
1:A:256:GLU:HB3	1:A:264:LYS:HE2	2.02	0.42
1:A:138:ILE:O	1:A:142:GLN:HG2	2.21	0.40
1:A:265:THR:HA	1:A:270:ARG:O	2.21	0.40
1:A:210:THR:HG22	3:A:812:HOH:O	2.20	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	444/447 (99%)	433 (98%)	10 (2%)	1 (0%)	43 44

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	123	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	375/375 (100%)	339 (90%)	36 (10%)	8 4

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LEU
1	A	21	ASN
1	A	51	LYS
1	A	59	ARG
1	A	80	THR
1	A	86	GLU
1	A	90	THR
1	A	91	VAL
1	A	94	LEU
1	A	99	GLU
1	A	114	VAL
1	A	118	LEU
1	A	124	ASP
1	A	130	LEU
1	A	134	ARG
1	A	140	LEU
1	A	144	THR
1	A	152	VAL
1	A	154	VAL
1	A	189	LEU
1	A	192	GLU
1	A	205	ASN
1	A	222	ARG
1	A	227	VAL
1	A	237	LEU
1	A	242	VAL
1	A	244	VAL
1	A	262	LEU
1	A	270	ARG
1	A	276	VAL

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Mol	Chain	Res	Type
1	A	279	VAL
1	A	291	ASP
1	A	314	GLU
1	A	345	MET
1	A	351	LYS
1	A	440	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	HIS
1	A	87	HIS
1	A	247	ASN
1	A	295	ASN
1	A	312	ASN
1	A	445	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue 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	$\# Z > 2$
1	OCS	A	42	1	6,8,9	1.72	2 (33%)	7,11,13	4.33	5 (71%)

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
1	OCS	A	42	1	-	0/4/7/9	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	42	OCS	OD2-SG	-2.75	1.36	1.47
1	A	42	OCS	CB-CA	2.50	1.55	1.53

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	42	OCS	OD3-SG-CB	7.52	117.98	106.76
1	A	42	OCS	OD1-SG-CB	5.63	115.16	106.76
1	A	42	OCS	OD2-SG-OD1	-3.75	102.02	111.40
1	A	42	OCS	OD2-SG-OD3	-3.42	102.84	111.40
1	A	42	OCS	CA-CB-SG	-3.18	108.88	113.61

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	42	OCS	1	0

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	A	449	-	58,58,58	1.08	2 (3%)	85,89,89	1.14	9 (10%)

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	449	-	-	1/34/50/50	0/6/6/6

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	449	FAD	PA-O5B	-2.55	1.49	1.59
2	A	449	FAD	C5A-C4A	2.29	1.43	1.39

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	449	FAD	C5A-C4A-N3A	-3.36	122.09	126.72
2	A	449	FAD	O2A-PA-O3P	2.65	114.43	107.27
2	A	449	FAD	C4A-C5A-N7A	-2.60	107.61	110.58
2	A	449	FAD	C6A-C5A-C4A	2.56	120.67	117.18
2	A	449	FAD	C4-N3-C2	-2.51	121.18	125.64
2	A	449	FAD	O2-C2-N1	-2.44	117.75	121.80
2	A	449	FAD	N3A-C4A-N9A	2.27	131.03	127.17
2	A	449	FAD	C5A-N7A-C8A	2.03	106.63	103.45
2	A	449	FAD	O4-C4-C4X	-2.01	121.24	126.53

There are no chirality outliers.

All (1) torsion outliers are listed below:

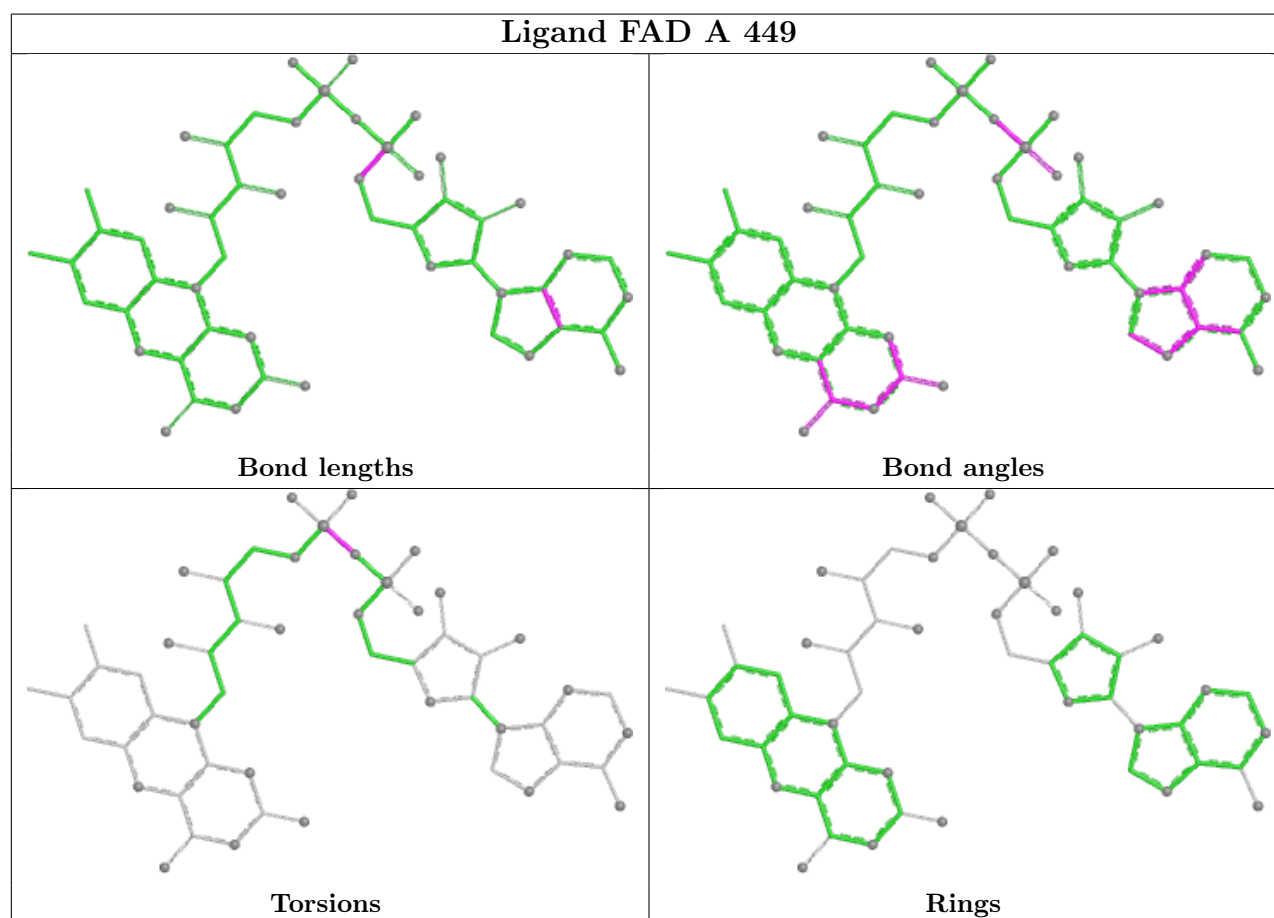
Mol	Chain	Res	Type	Atoms
2	A	449	FAD	PA-O3P-P-O5'

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	449	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.



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	446/447 (99%)	-0.70	3 (0%) 84 85	6, 16, 35, 57	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	124	ASP	3.8
1	A	25	ASP	2.4
1	A	220	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	OCS	A	42	9/10	0.99	0.05	10,11,13,15	0

6.3 Carbohydrates [i](#)

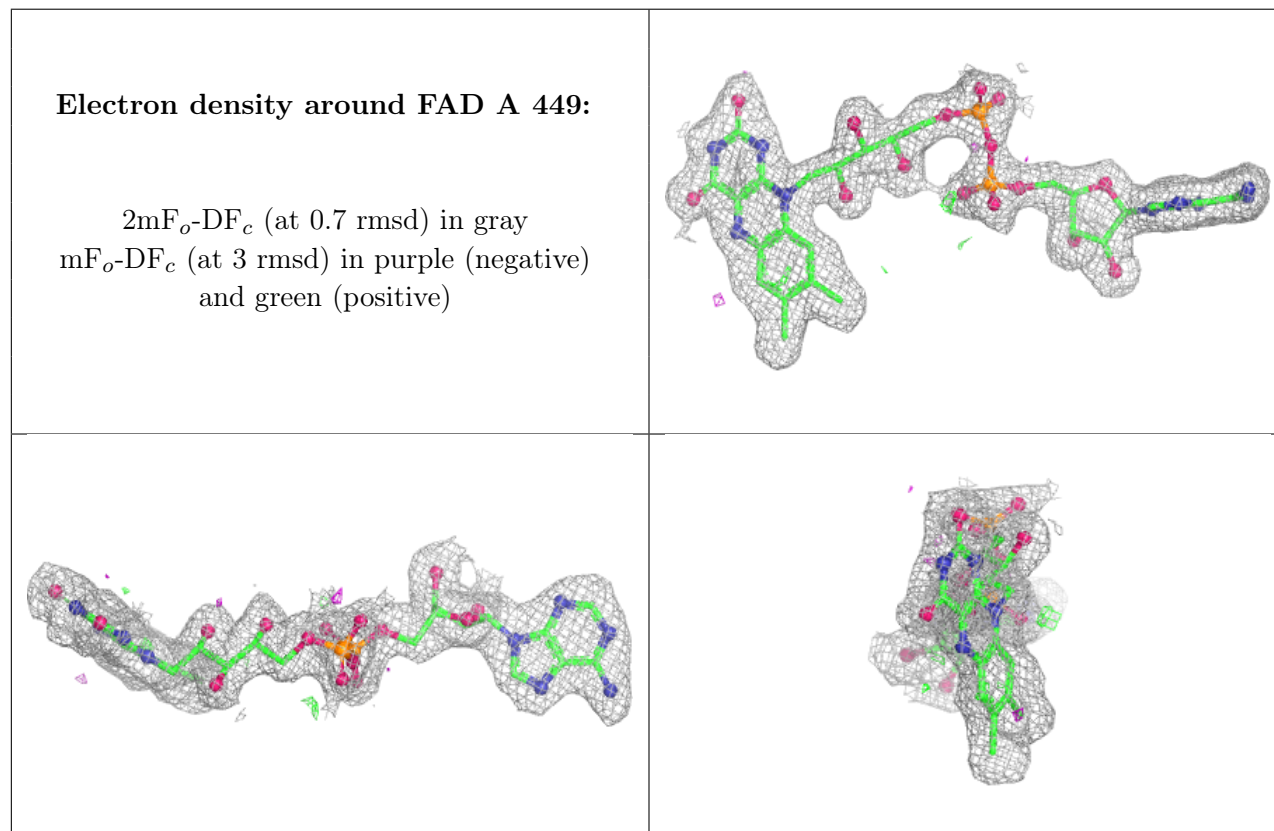
There are no oligosaccharides in this entry.

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	A	449	53/53	0.99	0.04	3,6,11,11	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.