



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:20 AM UTC

PDB ID : 7VJA / pdb_00007vja
Title : class II photolyase MmCPDII semiquinone to fully reduced TR-SFX studies (10 ns time-point)
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Deposited on : 2021-09-28
Resolution : 2.15 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)

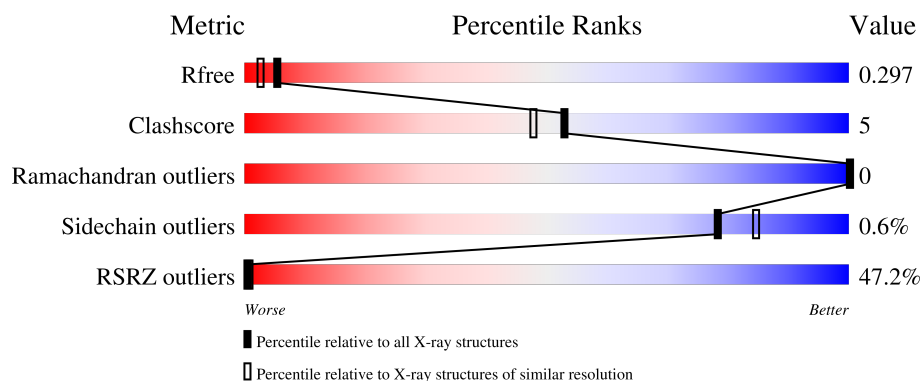
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
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	482	44% 81% 13% 6%

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

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3832 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 DNA photolyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	453	Total	C	N	O	S	0	7	0
			3574	2317	584	657	16			

There are 21 discrepancies between the modelled and reference sequences:

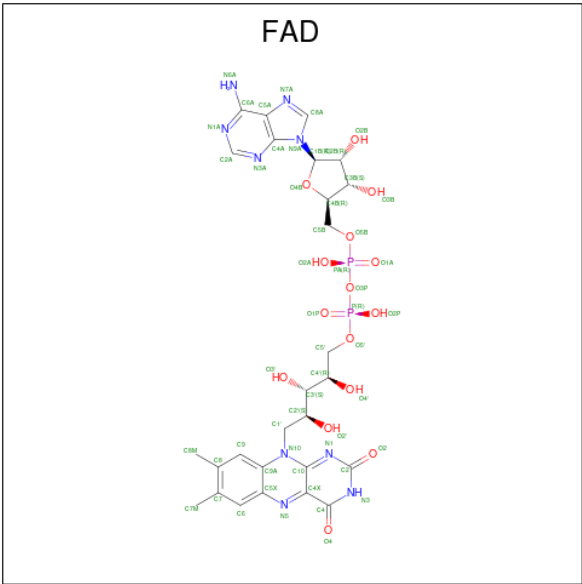
Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	MET	-	initiating methionine	UNP Q8PYK9
A	-16	GLY	-	expression tag	UNP Q8PYK9
A	-15	SER	-	expression tag	UNP Q8PYK9
A	-14	SER	-	expression tag	UNP Q8PYK9
A	-13	HIS	-	expression tag	UNP Q8PYK9
A	-12	HIS	-	expression tag	UNP Q8PYK9
A	-11	HIS	-	expression tag	UNP Q8PYK9
A	-10	HIS	-	expression tag	UNP Q8PYK9
A	-9	HIS	-	expression tag	UNP Q8PYK9
A	-8	HIS	-	expression tag	UNP Q8PYK9
A	-7	SER	-	expression tag	UNP Q8PYK9
A	-6	SER	-	expression tag	UNP Q8PYK9
A	-5	GLY	-	expression tag	UNP Q8PYK9
A	-4	LEU	-	expression tag	UNP Q8PYK9
A	-3	VAL	-	expression tag	UNP Q8PYK9
A	-2	PRO	-	expression tag	UNP Q8PYK9
A	-1	ARG	-	expression tag	UNP Q8PYK9
A	0	GLY	-	expression tag	UNP Q8PYK9
A	1	SER	-	expression tag	UNP Q8PYK9
A	2	HIS	-	expression tag	UNP Q8PYK9
A	377	THR	MET	engineered mutation	UNP Q8PYK9

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



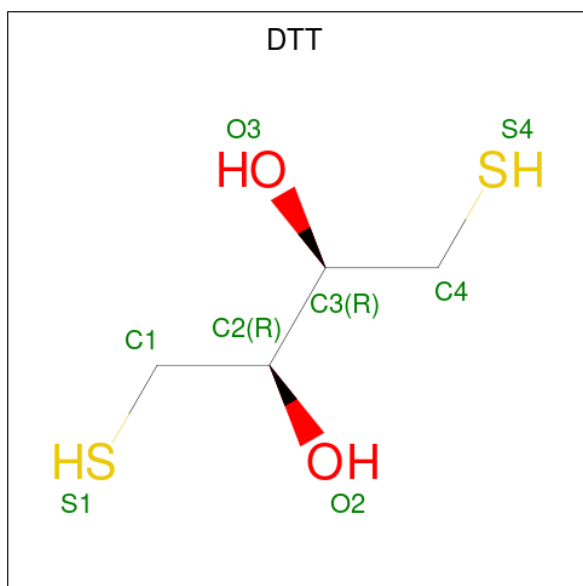
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is 2,3-DIHYDROXY-1,4-DITHIOBUTANE (CCD ID: DTT) (formula: $C_4H_{10}O_2S_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	S	0	0
			8	4	2	2		

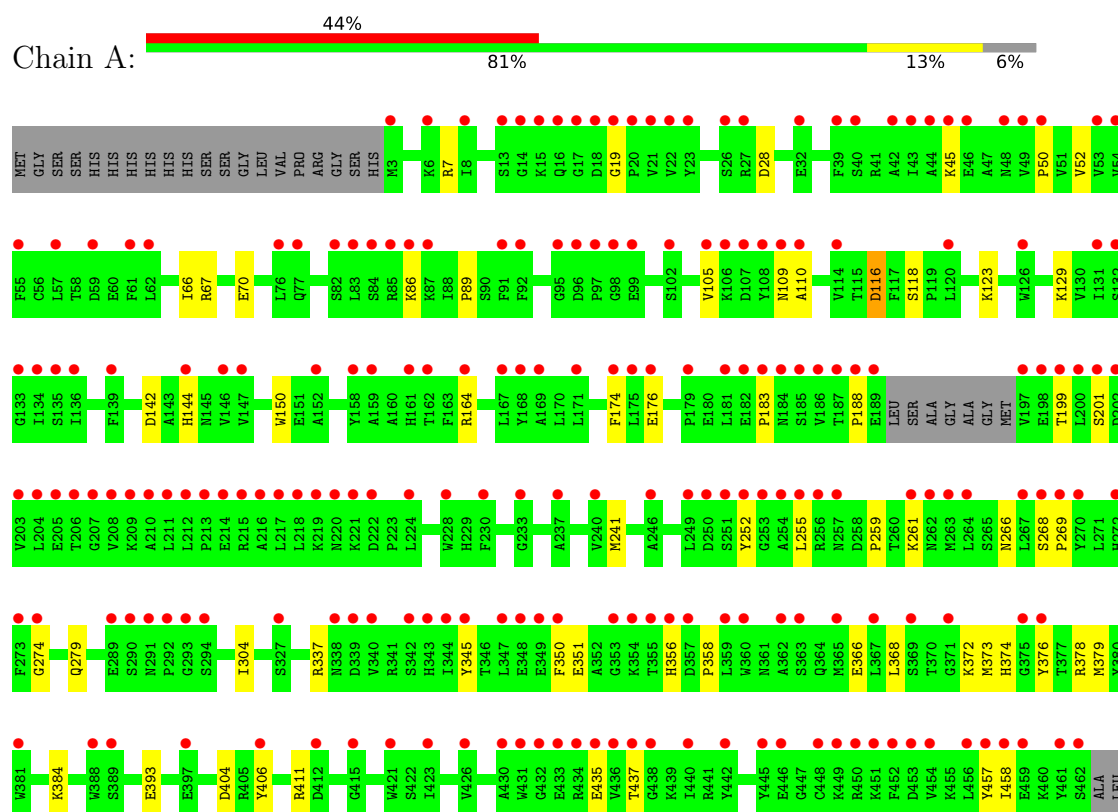
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	182	Total	O	0	0
			182	182		

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: DNA photolyase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	70.55Å 70.55Å 245.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.65 – 2.15 31.65 – 2.15	Depositor EDS
% Data completeness (in resolution range)	100.0 (31.65-2.15) 99.9 (31.65-2.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.01 (at 1.83Å)	Xtriage
Refinement program	PHENIX (1.19_4092: ???)	Depositor
R, R_{free}	0.198 , 0.220 0.280 , 0.297	Depositor DCC
R_{free} test set	2485 reflections (2.20%)	wwPDB-VP
Wilson B-factor (Å ²)	5.9	Xtriage
Anisotropy	1.404	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 40.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	3832	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.88% 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: SO4, DTT, 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	0.07	0/3696	0.20	0/5025

There are no bond length outliers.

There are no bond angle outliers.

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	3574	0	3346	34	0
2	A	15	0	0	0	0
3	A	53	0	28	2	0
4	A	8	0	10	0	0
5	A	182	0	0	2	0
All	All	3832	0	3384	35	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 (35) 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:266:ASN:HA	1:A:372:LYS:HE3	1.75	0.68
1:A:118:SER:HB3	1:A:123:LYS:HG2	1.78	0.65
1:A:376:TYR:HA	1:A:379:MET:HE2	1.82	0.61
3:A:503:FAD:H8A	3:A:503:FAD:O5B	2.03	0.59
1:A:351:GLU:OE2	1:A:406:TYR:OH	2.21	0.58
1:A:105:VAL:HG13	1:A:110:ALA:HB3	1.83	0.58
1:A:373:MET:O	1:A:378:ARG:NH1	2.37	0.57
1:A:86:LYS:HD2	1:A:183:PRO:HG3	1.89	0.55
1:A:255:LEU:HD22	1:A:261:LYS:HG3	1.90	0.54
1:A:45:LYS:NZ	5:A:621:HOH:O	2.40	0.54
1:A:345:TYR:OH	1:A:356:HIS:ND1	2.37	0.52
1:A:358:PRO:HB2	1:A:458:ILE:HD13	1.91	0.52
1:A:199:THR:HG22	1:A:201:SER:H	1.75	0.51
1:A:52:VAL:HG12	1:A:89:PRO:HG2	1.93	0.51
1:A:252:TYR:OH	3:A:503:FAD:O1A	2.19	0.49
1:A:176:GLU:O	1:A:279:GLN:NE2	2.42	0.49
1:A:7:ARG:NH2	1:A:174:PHE:O	2.46	0.48
1:A:241:MET:HE2	1:A:241:MET:HB3	1.77	0.48
1:A:435:GLU:O	1:A:437:THR:N	2.45	0.47
1:A:28:ASP:HA	1:A:274:GLY:HA3	1.96	0.47
1:A:366:GLU:OE1	1:A:457:TYR:OH	2.23	0.47
1:A:268:SER:OG	1:A:269:PRO:HD3	2.16	0.45
1:A:404:ASP:HB3	1:A:411:ARG:HG3	1.98	0.45
1:A:259:PRO:HA	1:A:374:HIS:CD2	2.52	0.45
1:A:116:ASP:O	1:A:123:LYS:NZ	2.50	0.44
1:A:164:ARG:HA	1:A:304:ILE:HD12	1.99	0.44
1:A:345:TYR:HB2	1:A:350:PHE:CZ	2.53	0.43
1:A:19:GLY:HA3	1:A:109:ASN:O	2.17	0.43
1:A:337:ARG:NH2	1:A:393:GLU:OE2	2.35	0.43
1:A:50:PRO:HG2	1:A:188:PRO:HB3	2.01	0.42
1:A:129:LYS:NZ	5:A:617:HOH:O	2.38	0.42
1:A:67:ARG:NH2	1:A:368:LEU:O	2.54	0.41
1:A:142:ASP:HB2	1:A:150:TRP:CH2	2.56	0.41
1:A:384:LYS:HA	1:A:384:LYS:HD3	1.97	0.41
1:A:66:ILE:HG22	1:A:70:GLU:HG3	2.04	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	456/482 (95%)	448 (98%)	8 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	356/416 (86%)	354 (99%)	2 (1%)	78	84

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	116	ASP
1	A	144	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	403	ASN

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](#)

5 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	SO4	A	504	-	4,4,4	0.23	0	6,6,6	0.08	0
2	SO4	A	502	-	4,4,4	0.24	0	6,6,6	0.07	0
2	SO4	A	501	-	4,4,4	0.23	0	6,6,6	0.08	0
3	FAD	A	503	-	58,58,58	3.53	28 (48%)	85,89,89	2.05	23 (27%)
4	DTT	A	505	-	7,7,7	1.03	1 (14%)	4,8,8	0.95	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
3	FAD	A	503	-	-	3/34/50/50	0/6/6/6
4	DTT	A	505	-	-	4/8/8/8	-

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	503	FAD	PA-O3P	8.26	1.68	1.59
3	A	503	FAD	P-O3P	7.42	1.67	1.59
3	A	503	FAD	O2'-C2'	-7.41	1.27	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	503	FAD	O4'-C4'	-6.87	1.28	1.43
3	A	503	FAD	C3B-C4B	-6.79	1.35	1.53
3	A	503	FAD	C4'-C3'	-6.52	1.42	1.53
3	A	503	FAD	C9A-N10	-6.39	1.30	1.41
3	A	503	FAD	O4-C4	5.89	1.34	1.23
3	A	503	FAD	C2'-C3'	-5.83	1.43	1.53
3	A	503	FAD	C2B-C1B	-5.81	1.35	1.53
3	A	503	FAD	C8A-N9A	-5.73	1.27	1.37
3	A	503	FAD	C2B-C3B	-5.69	1.37	1.53
3	A	503	FAD	O2-C2	5.47	1.35	1.24
3	A	503	FAD	C6A-N6A	4.55	1.45	1.34
3	A	503	FAD	C4-N3	-4.20	1.31	1.38
3	A	503	FAD	O3'-C3'	-3.54	1.34	1.43
3	A	503	FAD	C4A-N9A	-3.50	1.30	1.37
3	A	503	FAD	C1B-N9A	-3.38	1.37	1.46
3	A	503	FAD	P-O5'	2.96	1.71	1.59
3	A	503	FAD	C5B-C4B	-2.87	1.42	1.51
3	A	503	FAD	C10-N10	-2.71	1.31	1.37
3	A	503	FAD	C2-N3	-2.68	1.33	1.39
3	A	503	FAD	PA-O2A	2.55	1.67	1.55
3	A	503	FAD	PA-O5B	2.53	1.69	1.59
4	A	505	DTT	O3-C3	-2.52	1.38	1.43
3	A	503	FAD	C8M-C8	2.47	1.55	1.51
3	A	503	FAD	P-O2P	2.30	1.66	1.55
3	A	503	FAD	C5A-C4A	2.24	1.43	1.39
3	A	503	FAD	C5X-N5	-2.07	1.35	1.39

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	503	FAD	C4A-N9A-C8A	8.17	114.31	105.74
3	A	503	FAD	C5A-C4A-N3A	-5.67	118.91	126.72
3	A	503	FAD	N3A-C4A-N9A	5.60	136.69	127.17
3	A	503	FAD	C4A-C5A-N7A	-4.57	105.35	110.58
3	A	503	FAD	N3A-C2A-N1A	-4.19	122.23	128.58
3	A	503	FAD	N9A-C8A-N7A	-3.74	108.63	113.94
3	A	503	FAD	C5X-N5-C4X	-3.66	112.17	118.09
3	A	503	FAD	C2A-N3A-C4A	3.42	120.18	111.83
3	A	503	FAD	C4X-C10-N1	-3.34	116.41	124.59
3	A	503	FAD	C4X-C10-N10	3.28	121.17	116.48
3	A	503	FAD	C4X-C4-N3	3.04	120.99	113.25
3	A	503	FAD	C5'-C4'-C3'	-3.03	106.51	112.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	503	FAD	C4-N3-C2	-2.99	120.33	125.64
3	A	503	FAD	C4B-O4B-C1B	-2.69	103.52	109.47
3	A	503	FAD	C5X-C9A-N10	2.48	120.21	117.97
3	A	503	FAD	C5A-N7A-C8A	2.45	107.30	103.45
3	A	503	FAD	C6A-C5A-N7A	2.44	136.80	132.09
3	A	503	FAD	O4-C4-C4X	-2.33	120.39	126.53
3	A	503	FAD	C4-C4X-N5	-2.23	115.13	118.21
3	A	503	FAD	C10-N1-C2	2.19	121.58	116.85
3	A	503	FAD	C9A-N10-C10	-2.09	117.56	120.75
3	A	503	FAD	C3B-C2B-C1B	2.07	105.38	101.46
3	A	503	FAD	C4A-N9A-C1B	-2.05	121.84	126.63

There are no chirality outliers.

All (7) torsion outliers are listed below:

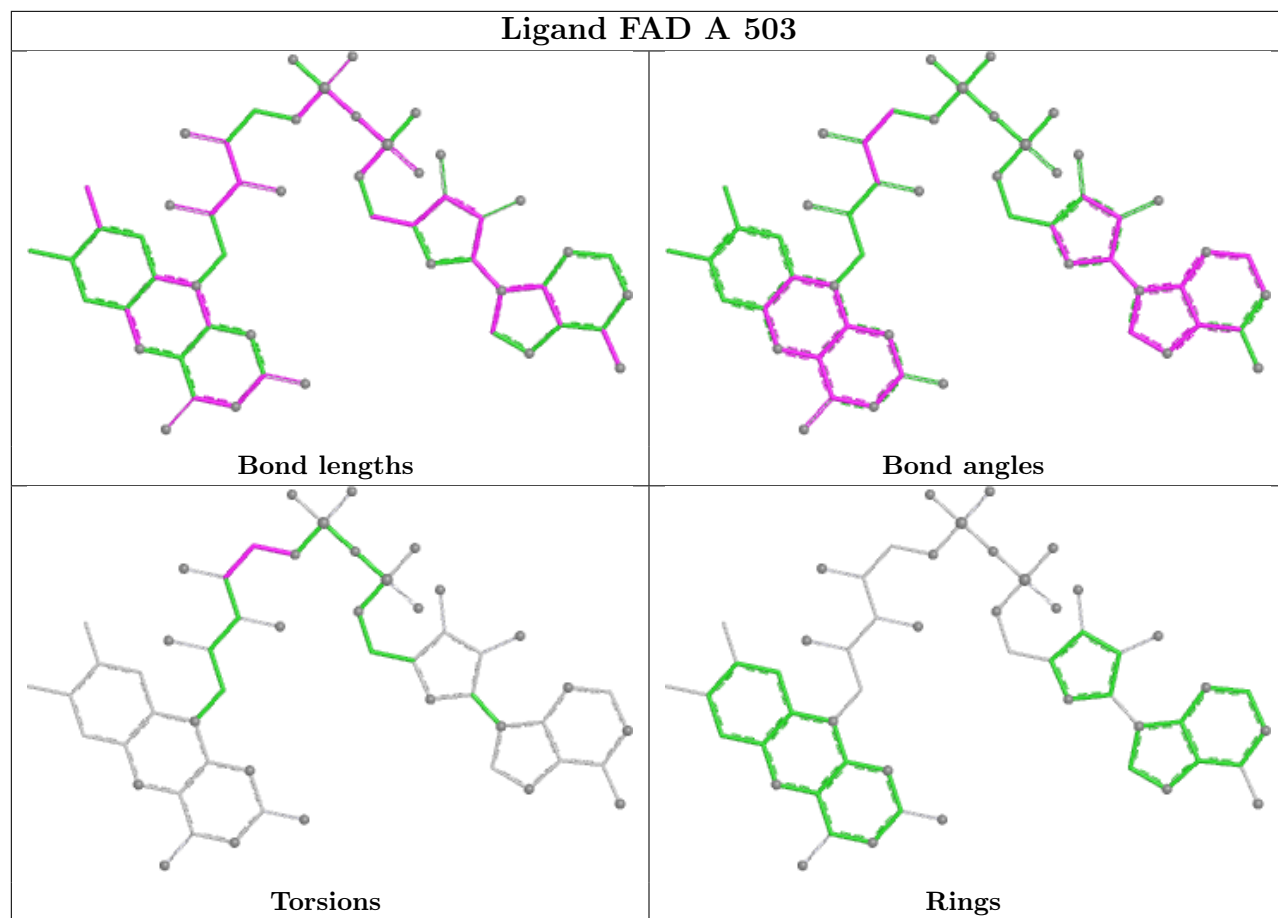
Mol	Chain	Res	Type	Atoms
3	A	503	FAD	O4'-C4'-C5'-O5'
4	A	505	DTT	C1-C2-C3-O3
4	A	505	DTT	C1-C2-C3-C4
4	A	505	DTT	O2-C2-C3-O3
4	A	505	DTT	O2-C2-C3-C4
3	A	503	FAD	C3'-C4'-C5'-O5'
3	A	503	FAD	C4'-C5'-O5'-P

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	503	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 ⓘ

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	453/482 (93%)	2.11	214 (47%) 0 1	18, 39, 63, 96	10 (2%)

All (214) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	197	VAL	8.0
1	A	158	TYR	6.2
1	A	251	SER	5.6
1	A	220	ASN	5.2
1	A	430	ALA	5.2
1	A	233	GLY	5.2
1	A	291	ASN	5.0
1	A	203	VAL	4.8
1	A	255	LEU	4.8
1	A	202	ASP	4.7
1	A	221	LYS	4.6
1	A	18	ASP	4.5
1	A	49	VAL	4.5
1	A	198	GLU	4.5
1	A	290	SER	4.5
1	A	218	LEU	4.5
1	A	186	VAL	4.4
1	A	262	ASN	4.3
1	A	174	PHE	4.3
1	A	342	SER	4.3
1	A	458	ILE	4.3
1	A	199	THR	4.3
1	A	188	PRO	4.3
1	A	108	TYR	4.3
1	A	344	ILE	4.2
1	A	249	LEU	4.2
1	A	189	GLU	4.2

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Mol	Chain	Res	Type	RSRZ
1	A	84	SER	4.2
1	A	456	LEU	4.1
1	A	212	LEU	4.1
1	A	132	SER	4.1
1	A	3	MET	4.0
1	A	432	GLY	4.0
1	A	362	ALA	3.9
1	A	462	SER	3.9
1	A	16	GLN	3.9
1	A	256	ARG	3.9
1	A	184	ASN	3.9
1	A	187	THR	3.9
1	A	228	TRP	3.9
1	A	431	TRP	3.9
1	A	219	LYS	3.9
1	A	350	PHE	3.8
1	A	168	TYR	3.8
1	A	292	PRO	3.8
1	A	365[A]	MET	3.7
1	A	46	GLU	3.7
1	A	389	SER	3.7
1	A	204	LEU	3.6
1	A	211	LEU	3.6
1	A	50	PRO	3.6
1	A	454	VAL	3.6
1	A	107	ASP	3.6
1	A	345	TYR	3.6
1	A	161	HIS	3.6
1	A	8	ILE	3.6
1	A	213	PRO	3.6
1	A	99	GLU	3.6
1	A	450	ARG	3.6
1	A	421	TRP	3.5
1	A	48	ASN	3.5
1	A	267	LEU	3.4
1	A	206	THR	3.4
1	A	442	TYR	3.4
1	A	215	ARG	3.4
1	A	167	LEU	3.4
1	A	440	ILE	3.3
1	A	340	VAL	3.3
1	A	445	TYR	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	457	TYR	3.3
1	A	91	PHE	3.2
1	A	343	HIS	3.2
1	A	214	GLU	3.2
1	A	327	SER	3.2
1	A	355	THR	3.2
1	A	13	SER	3.2
1	A	338	ASN	3.2
1	A	208	VAL	3.1
1	A	216	ALA	3.1
1	A	254	ALA	3.1
1	A	109	ASN	3.1
1	A	95	GLY	3.1
1	A	347	LEU	3.1
1	A	126[A]	TRP	3.0
1	A	207	GLY	3.0
1	A	61	PHE	3.0
1	A	273	PHE	3.0
1	A	96	ASP	3.0
1	A	83	LEU	3.0
1	A	436	VAL	3.0
1	A	87	LYS	3.0
1	A	263	MET	3.0
1	A	217	LEU	3.0
1	A	182	GLU	3.0
1	A	435	GLU	3.0
1	A	134	ILE	3.0
1	A	86	LYS	3.0
1	A	201	SER	2.9
1	A	210	ALA	2.9
1	A	356	HIS	2.9
1	A	85	ARG	2.9
1	A	433	GLU	2.9
1	A	252	TYR	2.9
1	A	381	TRP	2.9
1	A	39	PHE	2.9
1	A	102	SER	2.9
1	A	388	TRP	2.9
1	A	434	ARG	2.8
1	A	54	VAL	2.8
1	A	240	VAL	2.8
1	A	437	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	294	SER	2.8
1	A	222	ASP	2.8
1	A	77	GLN	2.8
1	A	375	GLY	2.8
1	A	453	ASP	2.8
1	A	6	LYS	2.8
1	A	261	LYS	2.8
1	A	269	PRO	2.8
1	A	22	VAL	2.8
1	A	423	ILE	2.8
1	A	176	GLU	2.7
1	A	446	GLU	2.7
1	A	438	GLY	2.7
1	A	45	LYS	2.7
1	A	164	ARG	2.7
1	A	257	ASN	2.7
1	A	136	ILE	2.7
1	A	253	GLY	2.7
1	A	293	GLY	2.7
1	A	159	ALA	2.7
1	A	246	ALA	2.7
1	A	14	GLY	2.7
1	A	17	GLY	2.7
1	A	459	GLU	2.6
1	A	367	LEU	2.6
1	A	146	VAL	2.6
1	A	42	ALA	2.6
1	A	230	PHE	2.6
1	A	357	ASP	2.6
1	A	105	VAL	2.6
1	A	451	LYS	2.6
1	A	82	SER	2.6
1	A	139	PHE	2.5
1	A	21	VAL	2.5
1	A	185	SER	2.5
1	A	57	LEU	2.5
1	A	20	PRO	2.5
1	A	369	SER	2.5
1	A	171	LEU	2.5
1	A	250[A]	ASP	2.5
1	A	106	LYS	2.5
1	A	449	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	23	TYR	2.4
1	A	92	PHE	2.4
1	A	461	TYR	2.4
1	A	169	ALA	2.4
1	A	264	LEU	2.4
1	A	133	GLY	2.4
1	A	426	VAL	2.4
1	A	26	SER	2.4
1	A	363	SER	2.4
1	A	183	PRO	2.4
1	A	270	TYR	2.4
1	A	237	ALA	2.4
1	A	120	LEU	2.4
1	A	181	LEU	2.4
1	A	452	PHE	2.4
1	A	205	GLU	2.3
1	A	19	GLY	2.3
1	A	415	GLY	2.3
1	A	272	HIS	2.3
1	A	359	LEU	2.3
1	A	144	HIS	2.3
1	A	110	ALA	2.3
1	A	76	LEU	2.3
1	A	53	VAL	2.3
1	A	147	VAL	2.3
1	A	354	LYS	2.3
1	A	175	LEU	2.3
1	A	32	GLU	2.2
1	A	339	ASP	2.2
1	A	371	GLY	2.2
1	A	289	GLU	2.2
1	A	348	GLU	2.2
1	A	268	SER	2.2
1	A	152	ALA	2.2
1	A	448	CYS	2.2
1	A	59	ASP	2.2
1	A	274	GLY	2.2
1	A	397	GLU	2.2
1	A	40	SER	2.2
1	A	131	ILE	2.1
1	A	200	LEU	2.1
1	A	224	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	376	TYR	2.1
1	A	97	PRO	2.1
1	A	360	TRP	2.1
1	A	353	GLY	2.1
1	A	15	LYS	2.1
1	A	135	SER	2.1
1	A	406	TYR	2.1
1	A	43	ILE	2.1
1	A	98	GLY	2.1
1	A	209	LYS	2.1
1	A	44	ALA	2.1
1	A	27	ARG	2.1
1	A	162	THR	2.1
1	A	55	PHE	2.1
1	A	114	VAL	2.1
1	A	349	GLU	2.0
1	A	412	ASP	2.0
1	A	62	LEU	2.0
1	A	179	PRO	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.

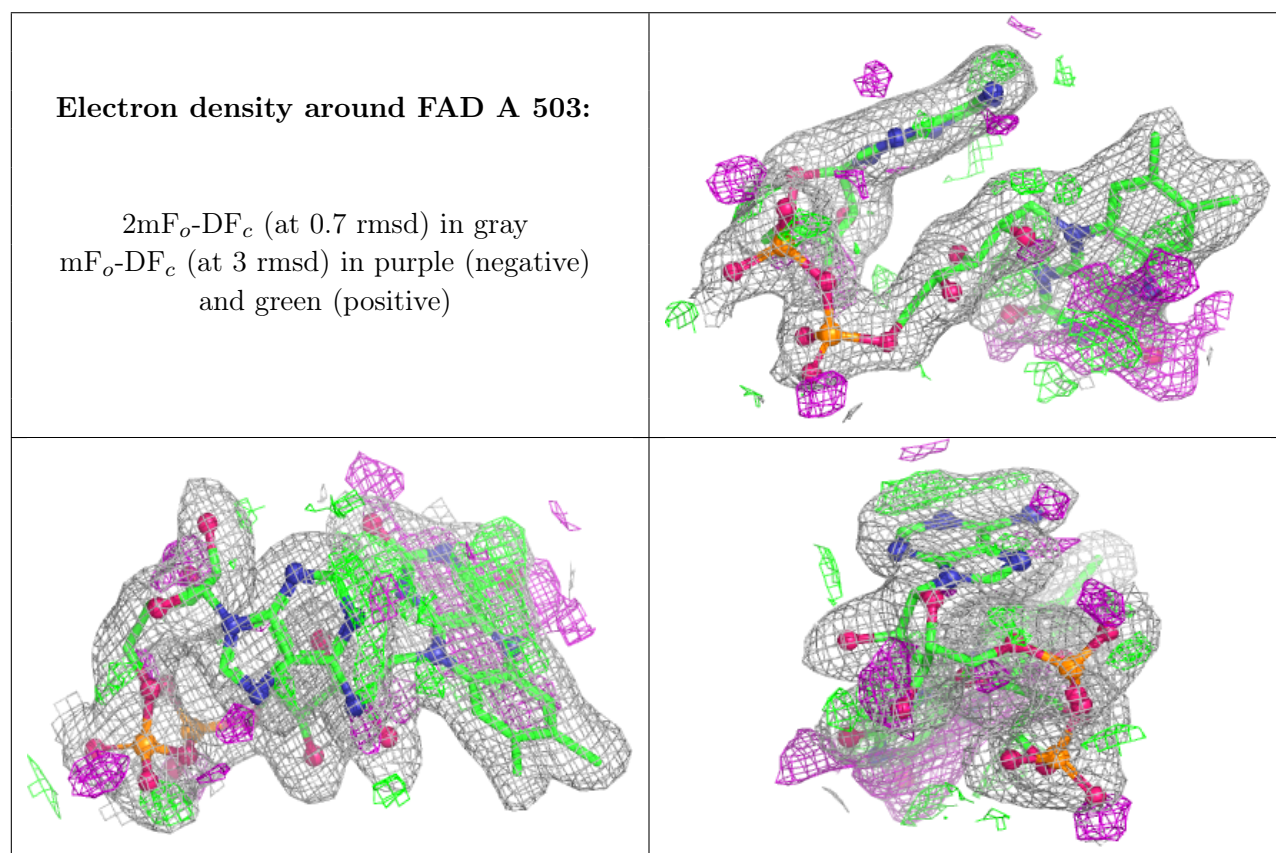
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	A	504	5/5	0.41	0.17	87,90,93,98	0
4	DTT	A	505	8/8	0.66	0.28	44,45,47,63	4
2	SO4	A	502	5/5	0.75	0.15	63,66,68,69	0
3	FAD	A	503	53/53	0.87	0.15	23,25,26,28	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	A	501	5/5	0.95	0.12	35,37,41,43	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.