



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:17 PM UTC

PDB ID : 2B5O / pdb_00002b5o
Title : ferredoxin-NADP reductase
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Yeates, T.O.
Deposited on : 2005-09-29
Resolution : 2.50 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

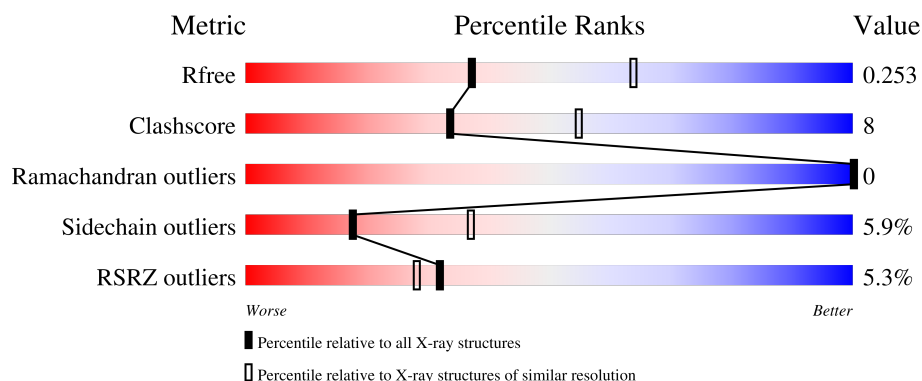
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	402	6% 61% 10% • 27%
1	B	402	6% 56% 14% • 27%

2 Entry composition [i](#)

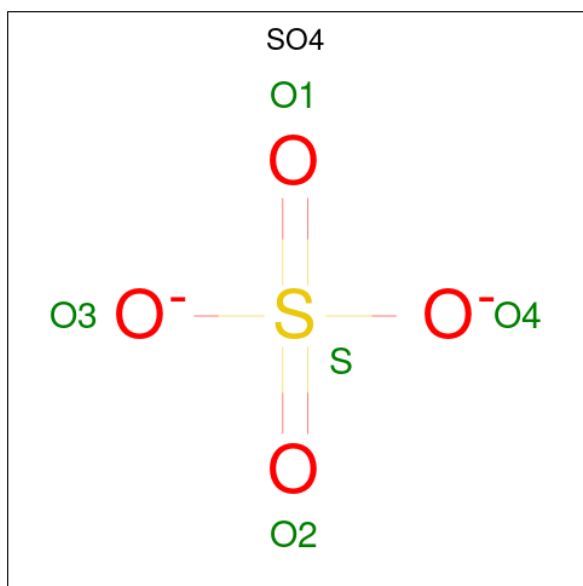
There are 4 unique types of molecules in this entry. The entry contains 4933 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Ferredoxin–NADP reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	0	0	0
			2340	1482	394	449	15			
1	B	292	Total	C	N	O	S	0	0	0
			2340	1482	394	449	15			

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



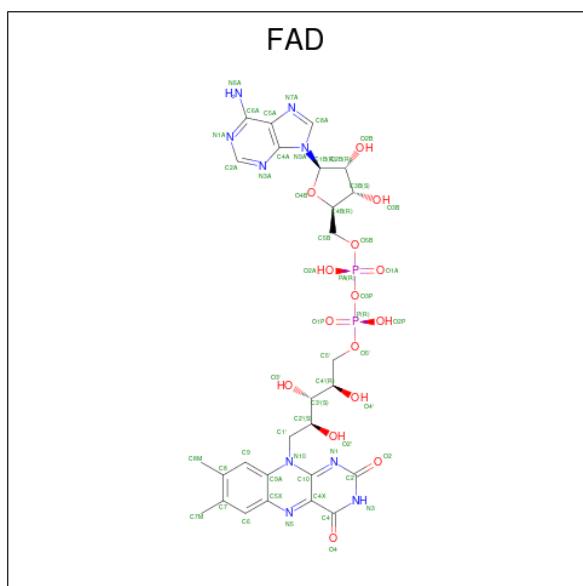
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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[illegible]

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $\text{C}_{27}\text{H}_{33}\text{N}_9\text{O}_{15}\text{P}_2$).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	56	Total 56	O 56	0	0
4	B	36	Total 36	O 36	0	0

- Molecule 1: Ferredoxin–NADP reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	121.98Å 121.98Å 74.73Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.14 – 2.50 47.14 – 2.50	Depositor EDS
% Data completeness (in resolution range)	96.5 (47.14-2.50) 96.3 (47.14-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.65 (at 2.51Å)	Xtriage
Refinement program	REFMAC refmac_5.2.0005	Depositor
R, R_{free}	0.178 , 0.235 0.216 , 0.253	Depositor DCC
R_{free} test set	2149 reflections (9.55%)	wwPDB-VP
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.035 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4933	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 3.20% 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, SO4

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	0.65	0/2396	1.16	2/3238 (0.1%)
1	B	0.58	0/2396	1.11	7/3238 (0.2%)
All	All	0.61	0/4792	1.14	9/6476 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	166	PRO	O-C-N	7.18	124.52	121.15
1	A	371	GLU	N-CA-C	6.46	119.31	111.82
1	B	205	GLU	CB-CA-C	-5.81	97.58	109.38
1	B	322	SER	N-CA-C	5.67	117.46	111.28
1	B	153	ASP	CB-CA-C	-5.58	101.73	110.94
1	B	237	GLY	CA-C-N	-5.54	114.26	119.85
1	B	237	GLY	C-N-CA	-5.54	114.26	119.85
1	A	303	ASP	CB-CA-C	-5.45	102.25	110.92
1	B	241	LYS	N-CA-C	-5.10	107.00	114.39

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	2340	0	2275	28	0
1	B	2340	0	2275	43	0
2	A	25	0	0	2	0
2	B	30	0	0	1	0
3	A	53	0	31	4	0
3	B	53	0	31	1	0
4	A	56	0	0	0	0
4	B	36	0	0	2	0
All	All	4933	0	4612	73	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 8.

All (73) 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:B:382:LEU:HD21	1:B:387:MET:SD	1.94	1.08
3:A:403:FAD:H3B	3:A:403:FAD:H8A	1.67	0.77
1:B:323:ARG:HH11	1:B:323:ARG:HG3	1.51	0.74
1:B:382:LEU:HD21	1:B:387:MET:CE	2.19	0.73
1:A:323:ARG:HG3	1:A:323:ARG:HH11	1.56	0.68
1:B:382:LEU:CD2	1:B:387:MET:SD	2.79	0.64
1:B:129:GLU:HG2	1:B:146:THR:HB	1.81	0.61
1:A:245:LEU:HD11	1:A:272:MET:HE2	1.84	0.60
3:A:403:FAD:H3B	3:A:403:FAD:C8A	2.32	0.59
2:A:410:SO4:O4	3:B:403:FAD:N6A	2.36	0.58
1:B:323:ARG:HG3	1:B:323:ARG:NH1	2.19	0.57
1:B:359:MET:HG3	1:B:365:MET:HE2	1.86	0.56
1:A:274:LYS:HB2	1:A:307:MET:HE2	1.87	0.56
1:A:327:THR:HG22	1:A:329:ASP:H	1.70	0.55
1:A:363:LYS:HA	1:A:401:VAL:HG13	1.87	0.55
1:A:203:GLN:NE2	1:A:221:THR:OG1	2.40	0.54
1:B:130:ASN:OD1	1:B:143:ARG:HD3	2.07	0.53
1:A:240:GLY:HA2	2:A:414:SO4:O4	2.09	0.53
3:A:403:FAD:C8A	3:A:403:FAD:C3B	2.87	0.53
1:B:366:GLN:HB3	1:B:367:PRO:HD3	1.92	0.52
1:A:117:TYR:H	1:A:237:GLY:HA3	1.73	0.52
1:A:245:LEU:HD11	1:A:272:MET:CE	2.40	0.51
1:B:274:LYS:HB2	1:B:307:MET:HE2	1.92	0.51
1:B:384:TRP:HB3	4:B:487:HOH:O	2.10	0.50
1:A:130:ASN:OD1	1:A:143:ARG:HD3	2.11	0.50
3:A:403:FAD:H8A	3:A:403:FAD:C3B	2.39	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:327:THR:HG22	1:B:329:ASP:H	1.77	0.49
1:A:129:GLU:HG2	1:A:146:THR:HB	1.93	0.49
1:A:323:ARG:HG3	1:A:323:ARG:NH1	2.26	0.49
1:A:366:GLN:HB3	1:A:367:PRO:HD3	1.94	0.49
1:B:122:PRO:HD3	4:B:456:HOH:O	2.13	0.49
1:B:129:GLU:HG2	1:B:146:THR:CB	2.43	0.49
1:B:117:TYR:H	1:B:237:GLY:HA3	1.77	0.49
1:B:384:TRP:CH2	1:B:388:ARG:HD3	2.47	0.48
1:A:367:PRO:HB2	1:A:368:PRO:HD3	1.95	0.48
1:B:171:LYS:HG3	1:B:172:ASN:ND2	2.29	0.48
1:B:203:GLN:NE2	1:B:221:THR:OG1	2.47	0.46
1:B:363:LYS:HA	1:B:401:VAL:HG13	1.97	0.46
1:B:176:HIS:ND1	2:B:405:SO4:O4	2.48	0.46
1:B:385:GLU:O	1:B:389:ARG:HG3	2.15	0.46
1:B:323:ARG:NH1	1:B:323:ARG:CG	2.79	0.46
1:A:170:ASP:HB3	1:A:174:LYS:HD2	1.97	0.46
1:A:171:LYS:HG3	1:A:172:ASN:ND2	2.30	0.46
1:A:213:GLU:HB3	1:B:293:PRO:HG3	1.97	0.46
1:A:385:GLU:O	1:A:389:ARG:HG3	2.16	0.46
1:A:268:PHE:HB3	1:A:272:MET:HE3	1.98	0.45
1:B:245:LEU:HD11	1:B:272:MET:HE2	1.98	0.45
1:A:170:ASP:HB3	1:A:174:LYS:CD	2.47	0.45
1:B:373:PHE:HB3	1:B:384:TRP:CD1	2.52	0.45
1:B:382:LEU:HD21	1:B:387:MET:HE1	1.97	0.45
1:A:327:THR:HG22	1:A:329:ASP:N	2.33	0.44
1:A:361:GLY:O	1:A:401:VAL:HA	2.17	0.44
1:B:205:GLU:CG	1:B:216:TYR:CE1	3.01	0.44
1:B:361:GLY:O	1:B:401:VAL:HA	2.18	0.44
1:A:367:PRO:HB2	1:A:368:PRO:CD	2.48	0.44
1:B:205:GLU:HG2	1:B:216:TYR:CE1	2.52	0.44
1:A:158:GLU:HG3	1:A:186:THR:HA	2.01	0.43
1:B:171:LYS:HG3	1:B:172:ASN:HD22	1.83	0.43
1:A:323:ARG:NH1	1:A:323:ARG:CG	2.82	0.43
1:B:268:PHE:HB3	1:B:272:MET:HE3	2.00	0.43
1:B:384:TRP:CZ2	1:B:388:ARG:HD3	2.54	0.43
1:A:329:ASP:N	1:A:329:ASP:OD1	2.52	0.42
1:B:347:PHE:O	1:B:351:GLN:HG3	2.19	0.42
1:B:216:TYR:O	1:B:217:GLY:C	2.62	0.42
1:B:208:ASP:HA	1:B:209:PRO:HD3	1.90	0.41
1:B:360:CYS:HA	1:B:400:GLU:O	2.20	0.41
1:A:293:PRO:HG3	1:B:213:GLU:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:HIS:HA	1:B:194:ASN:OD1	2.20	0.41
1:B:300:TYR:O	1:B:301:LYS:C	2.64	0.41
1:B:172:ASN:HB2	1:B:174:LYS:HE2	2.02	0.41
1:B:178:LEU:HD12	1:B:178:LEU:C	2.46	0.40
1:B:216:TYR:HB3	1:B:221:THR:OG1	2.21	0.40
1:A:205:GLU:OE1	1:B:205:GLU:HB3	2.22	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	290/402 (72%)	281 (97%)	9 (3%)	0	100	100
1	B	290/402 (72%)	281 (97%)	9 (3%)	0	100	100
All	All	580/804 (72%)	562 (97%)	18 (3%)	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	254/344 (74%)	239 (94%)	15 (6%)	18	37
1	B	254/344 (74%)	239 (94%)	15 (6%)	18	37

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	508/688 (74%)	478 (94%)	30 (6%)	18	37

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	139	SER
1	A	141	THR
1	A	151	GLU
1	A	153	ASP
1	A	154	LEU
1	A	174	LYS
1	A	204	LEU
1	A	303	ASP
1	A	327	THR
1	A	329	ASP
1	A	354	ASN
1	A	386	GLU
1	A	391	MET
1	A	395	HIS
1	A	401	VAL
1	B	139	SER
1	B	141	THR
1	B	151	GLU
1	B	153	ASP
1	B	154	LEU
1	B	174	LYS
1	B	203	GLN
1	B	204	LEU
1	B	303	ASP
1	B	329	ASP
1	B	337	SER
1	B	354	ASN
1	B	391	MET
1	B	395	HIS
1	B	401	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	172	ASN
1	A	203	GLN

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Mol	Chain	Res	Type
1	A	276	GLN
1	B	172	ASN
1	B	203	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](#)

13 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	$\# Z > 2$
3	FAD	B	403	-	58,58,58	1.21	8 (13%)	85,89,89	1.66	12 (14%)
2	SO4	B	413	-	4,4,4	0.25	0	6,6,6	0.32	0
2	SO4	A	408	-	4,4,4	0.19	0	6,6,6	0.81	0
2	SO4	B	406	-	4,4,4	0.28	0	6,6,6	0.38	0
2	SO4	A	407	-	4,4,4	0.26	0	6,6,6	0.92	0
2	SO4	A	414	-	4,4,4	0.26	0	6,6,6	0.46	0
3	FAD	A	403	-	58,58,58	1.12	6 (10%)	85,89,89	1.52	12 (14%)
2	SO4	B	405	-	4,4,4	0.32	0	6,6,6	1.04	1 (16%)
2	SO4	A	410	-	4,4,4	0.32	0	6,6,6	0.46	0
2	SO4	B	412	-	4,4,4	0.26	0	6,6,6	0.14	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	B	411	-	4,4,4	0.23	0	6,6,6	0.40	0
2	SO4	A	409	-	4,4,4	0.23	0	6,6,6	0.83	0
2	SO4	B	415	-	4,4,4	0.23	0	6,6,6	0.57	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	403	-	-	2/34/50/50	0/6/6/6
3	FAD	B	403	-	-	2/34/50/50	0/6/6/6

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	403	FAD	C8A-N7A	3.31	1.38	1.31
3	B	403	FAD	C4X-N5	3.16	1.37	1.30
3	A	403	FAD	C4X-N5	3.11	1.37	1.30
3	B	403	FAD	C2A-N1A	2.77	1.38	1.33
3	A	403	FAD	C2A-N3A	2.75	1.38	1.33
3	A	403	FAD	C2A-N1A	2.74	1.38	1.33
3	A	403	FAD	C5A-N7A	-2.60	1.34	1.39
3	A	403	FAD	C10-N1	2.57	1.38	1.33
3	B	403	FAD	PA-O3P	2.54	1.62	1.59
3	B	403	FAD	C10-N1	2.35	1.38	1.33
3	B	403	FAD	P-O3P	2.22	1.61	1.59
3	B	403	FAD	C2A-N3A	2.05	1.37	1.33
3	B	403	FAD	C4X-C10	-2.04	1.38	1.44
3	A	403	FAD	C9A-N10	-2.03	1.37	1.41

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	403	FAD	N3A-C2A-N1A	-5.63	120.06	128.58
3	B	403	FAD	C4A-N9A-C1B	-4.92	115.12	126.63
3	A	403	FAD	N3A-C2A-N1A	-4.58	121.64	128.58
3	B	403	FAD	C1B-N9A-C8A	4.46	136.99	127.09
3	B	403	FAD	C2B-C1B-N9A	-4.26	102.72	113.30
3	A	403	FAD	C2B-C1B-N9A	-3.91	103.60	113.30
3	B	403	FAD	N9A-C8A-N7A	-3.50	108.97	113.94
3	A	403	FAD	C6A-C5A-C4A	3.50	121.95	117.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	403	FAD	C4X-C10-N10	3.48	121.46	116.48
3	A	403	FAD	O2A-PA-O3P	3.30	116.19	107.27
3	A	403	FAD	C5A-C4A-N3A	-3.08	122.47	126.72
3	B	403	FAD	O2A-PA-O3P	2.90	115.10	107.27
3	B	403	FAD	C9A-C5X-N5	-2.73	119.56	122.45
3	B	403	FAD	O4-C4-C4X	-2.62	119.61	126.53
3	A	403	FAD	C5A-C4A-N9A	2.59	108.63	105.81
3	B	403	FAD	O3'-C3'-C2'	-2.47	103.32	108.93
3	A	403	FAD	C4-N3-C2	-2.40	121.39	125.64
3	B	403	FAD	C10-C4X-N5	-2.32	120.08	124.81
3	A	403	FAD	C10-C4X-N5	-2.29	120.12	124.81
3	A	403	FAD	O4B-C1B-N9A	2.24	112.39	108.09
2	B	405	SO4	O3-S-O1	-2.24	97.87	109.56
3	A	403	FAD	C2A-N1A-C6A	2.06	122.11	118.73
3	B	403	FAD	C2A-N1A-C6A	2.06	122.11	118.73
3	B	403	FAD	C4X-C10-N10	2.04	119.41	116.48
3	A	403	FAD	O3P-PA-O1A	-2.01	104.65	110.70

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	403	FAD	PA-O3P-P-O1P
3	B	403	FAD	C4B-C5B-O5B-PA
3	A	403	FAD	PA-O3P-P-O2P
3	B	403	FAD	PA-O3P-P-O2P

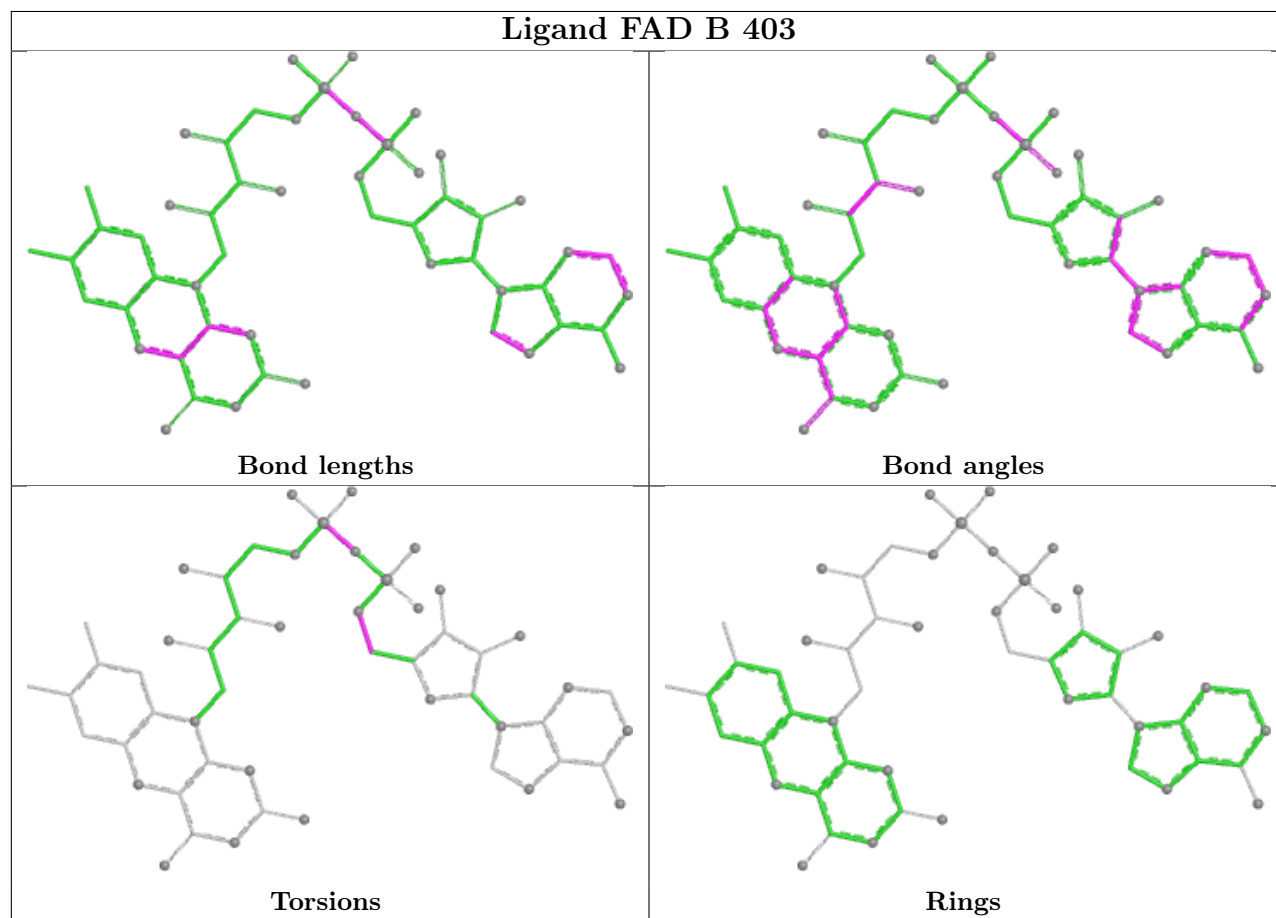
There are no ring outliers.

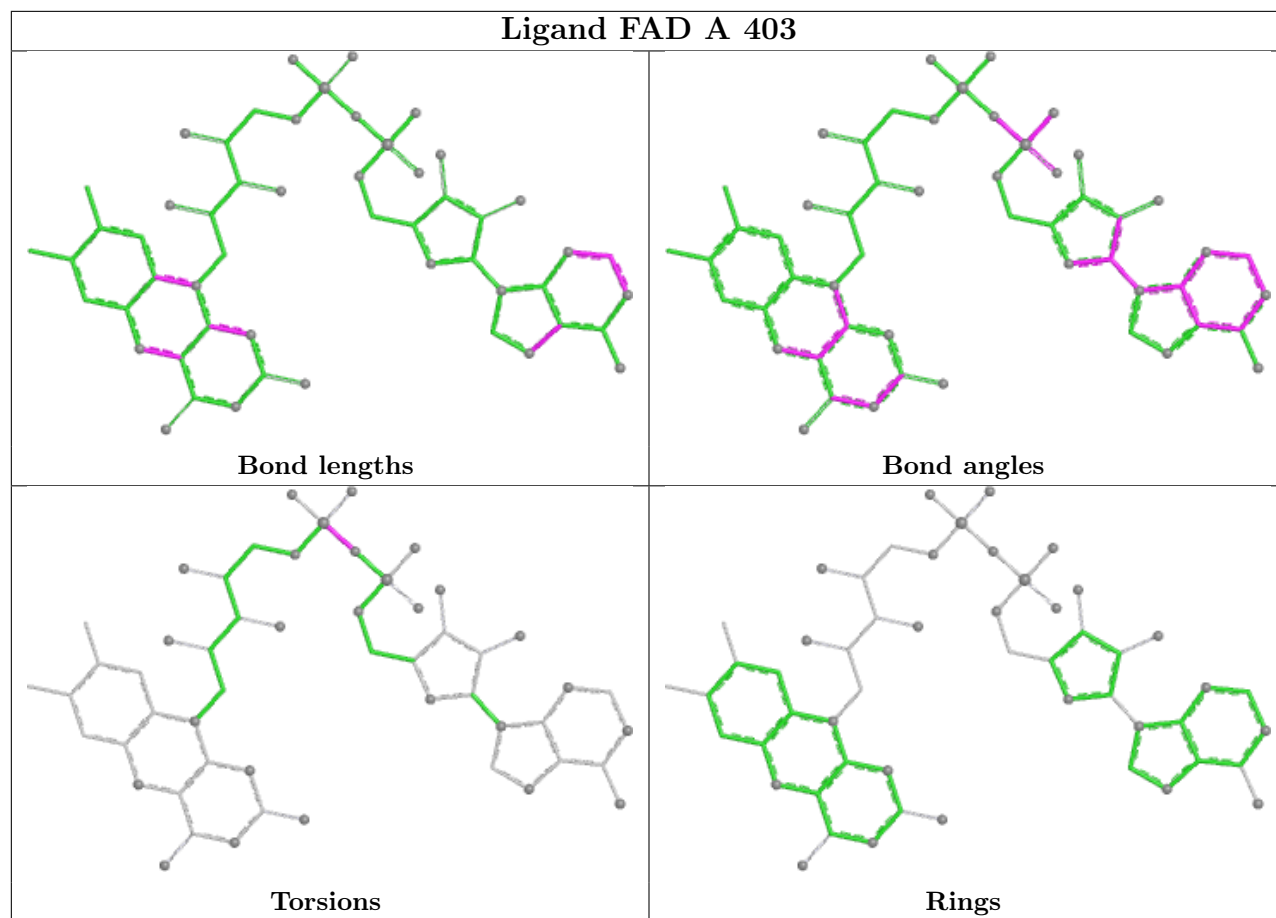
5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	403	FAD	1	0
2	A	414	SO4	1	0
3	A	403	FAD	4	0
2	B	405	SO4	1	0
2	A	410	SO4	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

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	292/402 (72%)	0.08	6 (2%) 63 59	24, 37, 65, 77	0
1	B	292/402 (72%)	0.62	25 (8%) 16 14	23, 38, 66, 81	0
All	All	584/804 (72%)	0.35	31 (5%) 32 28	23, 38, 66, 81	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	232	ASP	4.0
1	B	382	LEU	3.8
1	B	380	ARG	3.7
1	B	386	GLU	3.7
1	B	384	TRP	3.6
1	B	381	GLY	3.3
1	B	378	GLU	3.2
1	B	383	ASN	3.1
1	B	389	ARG	3.0
1	B	171	LYS	2.9
1	B	351	GLN	2.9
1	B	111	SER	2.7
1	A	170	ASP	2.6
1	A	174	LYS	2.5
1	B	385	GLU	2.5
1	A	119	PRO	2.4
1	B	278	GLU	2.2
1	B	112	VAL	2.2
1	B	119	PRO	2.2
1	B	393	LYS	2.2
1	B	328	ALA	2.1
1	A	382	LEU	2.1
1	A	279	ASP	2.1
1	B	149	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	153	ASP	2.1
1	B	170	ASP	2.1
1	B	329	ASP	2.1
1	B	114	VAL	2.1
1	B	370	ASP	2.1
1	B	115	ASN	2.0
1	B	237	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.

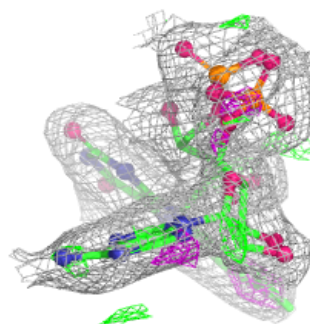
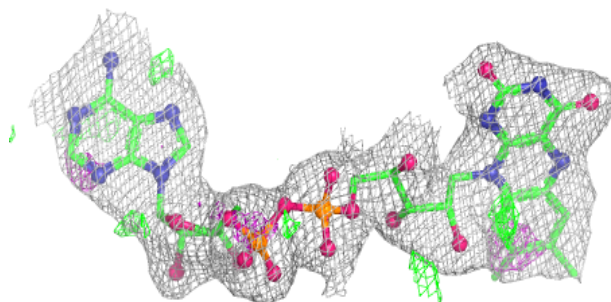
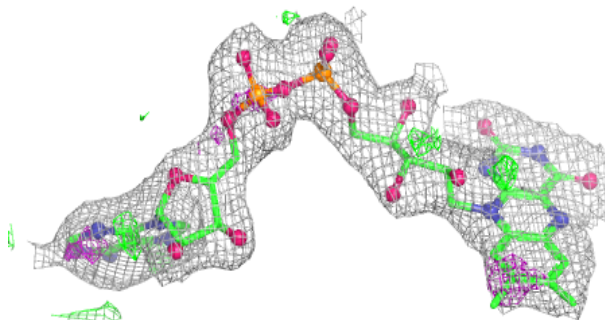
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SO4	B	415	5/5	0.62	0.13	83,88,98,105	0
2	SO4	B	412	5/5	0.64	0.13	92,93,104,105	0
2	SO4	B	411	5/5	0.71	0.15	101,101,113,116	0
2	SO4	B	413	5/5	0.81	0.15	77,86,92,96	0
2	SO4	A	410	5/5	0.82	0.17	65,67,94,98	0
2	SO4	B	405	5/5	0.84	0.14	61,65,77,79	0
2	SO4	A	409	5/5	0.86	0.15	59,63,75,78	0
2	SO4	A	414	5/5	0.90	0.11	68,69,73,83	0
2	SO4	A	407	5/5	0.90	0.10	66,75,76,83	0
2	SO4	B	406	5/5	0.92	0.09	45,56,72,72	0
3	FAD	B	403	53/53	0.94	0.08	18,32,43,49	0
3	FAD	A	403	53/53	0.95	0.08	21,30,53,58	0
2	SO4	A	408	5/5	0.97	0.08	42,45,48,51	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.

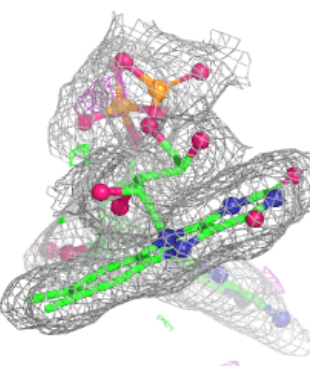
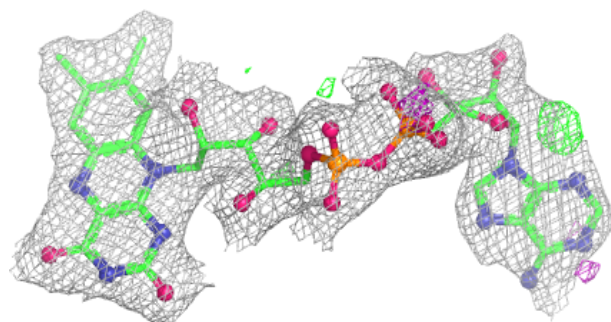
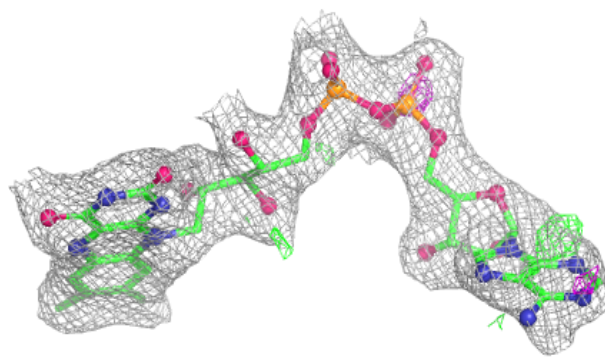
Electron density around FAD B 403:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around FAD A 403:

$2mF_o-DF_c$ (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.