



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 07:17 am BST

PDB ID : 1E1N
Title : Structure of adrenodoxin reductase at 2.4 Angstrom in crystal form A'
Authors : Ziegler, G.A.; Schulz, G.E.
Deposited on : 2000-05-09
Resolution : 2.40 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

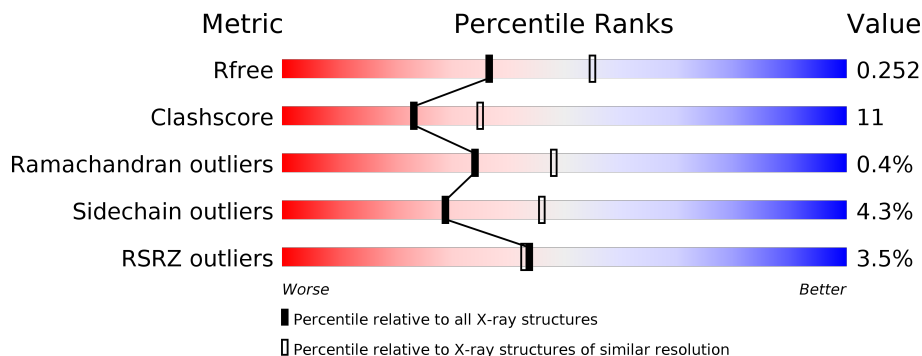
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	460	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3677 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 ADRENODOXIN REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	455	3505	2215	634	645	11	7	0	0

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



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

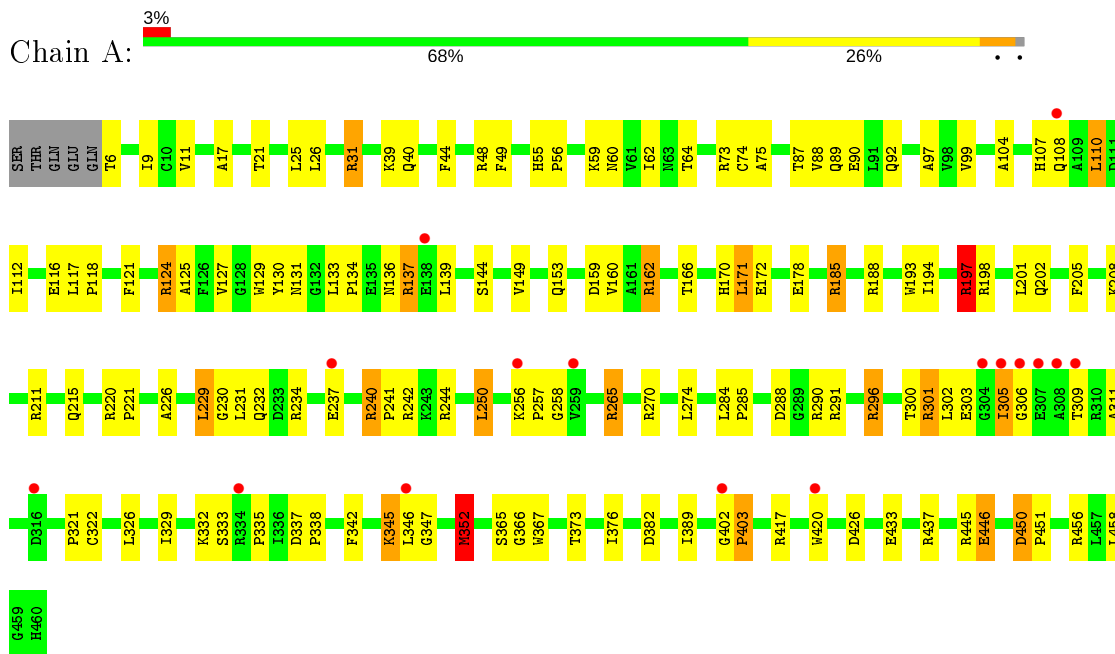
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	119	Total	O	0	0
			119	119		

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: ADRENODOXIN REDUCTASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.80Å 62.00Å 83.00Å 90.00° 107.10° 90.00°	Depositor
Resolution (Å)	29.00 – 2.40 28.87 – 2.35	Depositor EDS
% Data completeness (in resolution range)	77.6 (29.00-2.40) 74.2 (28.87-2.35)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 2.36Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.201 , 0.265 0.200 , 0.252	Depositor DCC
R_{free} test set	896 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	29.7	Xtrriage
Anisotropy	1.253	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 35.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3677	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.54% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.74	3/3584 (0.1%)	1.43	41/4871 (0.8%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	352	MET	CG-SD	17.10	2.25	1.81
1	A	345	LYS	CD-CE	-10.46	1.25	1.51
1	A	365	SER	CB-OG	-6.53	1.33	1.42

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	240	ARG	NE-CZ-NH1	11.67	126.13	120.30
1	A	417	ARG	CD-NE-CZ	10.61	138.45	123.60
1	A	185	ARG	CD-NE-CZ	10.09	137.72	123.60
1	A	291	ARG	NE-CZ-NH1	9.95	125.27	120.30
1	A	345	LYS	CD-CE-NZ	9.92	134.52	111.70
1	A	345	LYS	CG-CD-CE	9.46	140.29	111.90
1	A	73	ARG	NE-CZ-NH2	8.91	124.75	120.30
1	A	445	ARG	NE-CZ-NH2	-8.67	115.97	120.30
1	A	296	ARG	CD-NE-CZ	8.42	135.39	123.60
1	A	197	ARG	NE-CZ-NH1	8.38	124.49	120.30
1	A	382	ASP	CB-CG-OD1	-7.39	111.65	118.30
1	A	31	ARG	NE-CZ-NH2	-7.26	116.67	120.30
1	A	240	ARG	CD-NE-CZ	7.21	133.69	123.60
1	A	291	ARG	CD-NE-CZ	7.20	133.67	123.60
1	A	352	MET	CB-CG-SD	-7.09	91.13	112.40
1	A	130	TYR	CA-CB-CG	7.05	126.79	113.40
1	A	301	ARG	CA-CB-CG	7.02	128.84	113.40
1	A	162	ARG	NE-CZ-NH1	6.99	123.80	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	39	LYS	CA-CB-CG	6.90	128.57	113.40
1	A	389	ILE	CB-CA-C	-6.88	97.83	111.60
1	A	197	ARG	CD-NE-CZ	6.87	133.22	123.60
1	A	265	ARG	NE-CZ-NH1	6.54	123.57	120.30
1	A	270	ARG	CD-NE-CZ	6.49	132.69	123.60
1	A	75	ALA	N-CA-CB	6.45	119.13	110.10
1	A	445	ARG	NE-CZ-NH1	6.42	123.51	120.30
1	A	257	PRO	N-CA-C	6.31	128.50	112.10
1	A	403	PRO	N-CA-CB	-6.28	95.70	102.60
1	A	265	ARG	CD-NE-CZ	6.22	132.31	123.60
1	A	417	ARG	NE-CZ-NH1	6.14	123.37	120.30
1	A	456	ARG	NE-CZ-NH2	-5.91	117.34	120.30
1	A	124	ARG	NE-CZ-NH2	-5.91	117.35	120.30
1	A	240	ARG	NH1-CZ-NH2	-5.79	113.03	119.40
1	A	244	ARG	CB-CA-C	5.79	121.98	110.40
1	A	446	GLU	N-CA-CB	5.78	121.00	110.60
1	A	137	ARG	CA-CB-CG	5.74	126.02	113.40
1	A	73	ARG	NE-CZ-NH1	-5.69	117.45	120.30
1	A	450	ASP	CB-CG-OD1	5.67	123.41	118.30
1	A	208	LYS	CA-CB-CG	5.51	125.52	113.40
1	A	445	ARG	CD-NE-CZ	5.32	131.04	123.60
1	A	456	ARG	NE-CZ-NH1	5.23	122.92	120.30
1	A	456	ARG	CD-NE-CZ	5.08	130.72	123.60

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	3505	0	3535	80	1
2	A	53	0	31	3	0
3	A	119	0	0	16	1
All	All	3677	0	3566	80	1

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 11.

All (80) 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:234:ARG:HB3	3:A:2064:HOH:O	1.77	0.85
1:A:59:LYS:HE2	3:A:2116:HOH:O	1.80	0.80
1:A:345:LYS:HG3	1:A:346:LEU:HD23	1.68	0.76
1:A:178:GLU:HG3	3:A:2112:HOH:O	1.88	0.74
1:A:258:GLY:HA3	3:A:2069:HOH:O	1.90	0.71
1:A:284:LEU:HD21	1:A:296:ARG:HB2	1.79	0.64
1:A:144:SER:HA	1:A:188:ARG:HG3	1.79	0.62
1:A:87:THR:OG1	1:A:90:GLU:HG3	1.99	0.62
1:A:88:VAL:O	1:A:92:GLN:HG3	1.99	0.62
1:A:166:THR:HG21	1:A:171:LEU:HD13	1.80	0.61
1:A:9:ILE:HG23	1:A:97:ALA:HB3	1.82	0.61
1:A:198:ARG:NH1	1:A:202:GLN:HG2	2.17	0.59
1:A:232:GLN:HB2	1:A:250:LEU:HD11	1.85	0.59
1:A:159:ASP:HA	1:A:162:ARG:HH12	1.69	0.58
1:A:300:THR:HB	1:A:311:ALA:HB1	1.83	0.58
1:A:134:PRO:HA	3:A:2040:HOH:O	2.04	0.57
1:A:160:VAL:HG21	1:A:326:LEU:HD13	1.87	0.57
1:A:107:HIS:HB3	3:A:2035:HOH:O	2.05	0.57
1:A:221:PRO:HD3	3:A:2072:HOH:O	2.03	0.57
1:A:108:GLN:HG2	1:A:332:LYS:HZ2	1.69	0.56
1:A:402:GLY:HA2	3:A:2099:HOH:O	2.06	0.56
1:A:116:GLU:HG2	1:A:117:LEU:N	2.20	0.55
1:A:211:ARG:HG3	1:A:215:GLN:HE21	1.71	0.55
1:A:302:LEU:HD22	1:A:309:THR:HB	1.88	0.55
1:A:433:GLU:O	1:A:437:ARG:HG2	2.07	0.55
1:A:159:ASP:HA	1:A:162:ARG:NH1	2.22	0.54
1:A:226:ALA:HA	1:A:229:LEU:HD22	1.88	0.54
1:A:127:VAL:O	1:A:131:ASN:HB2	2.10	0.52
1:A:121:PHE:CD2	1:A:139:LEU:HD11	2.45	0.52
1:A:205:PHE:HZ	1:A:274:LEU:HD21	1.75	0.52
1:A:17:ALA:HB2	1:A:366:GLY:HA3	1.92	0.51
1:A:153:GLN:HG2	1:A:194:ILE:HG22	1.93	0.51
1:A:220:ARG:HD3	1:A:265:ARG:O	2.11	0.51
1:A:197:ARG:HH11	1:A:197:ARG:HG3	1.76	0.50
1:A:301:ARG:NH1	1:A:303:GLU:OE2	2.45	0.49
1:A:376:ILE:HG13	2:A:801:FAD:N1	2.28	0.48
1:A:367:TRP:CZ2	2:A:801:FAD:H3'	2.49	0.48
1:A:129:TRP:HB2	1:A:139:LEU:HD23	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:GLU:HB3	3:A:2063:HOH:O	2.14	0.47
1:A:321:PRO:O	1:A:322:CYS:HB3	2.14	0.47
1:A:332:LYS:HZ2	1:A:332:LYS:HB2	1.79	0.47
1:A:125:ALA:HB1	1:A:136:ASN:ND2	2.29	0.47
1:A:108:GLN:HG2	1:A:332:LYS:NZ	2.29	0.47
1:A:44:PHE:CE2	1:A:62:ILE:HG23	2.49	0.47
1:A:201:LEU:O	1:A:242:ARG:NH2	2.48	0.47
1:A:333:SER:HB2	1:A:347:GLY:O	2.16	0.46
1:A:26:LEU:HD11	1:A:74:CYS:HB2	1.98	0.46
1:A:124:ARG:HH12	1:A:367:TRP:HZ3	1.64	0.45
1:A:240:ARG:HB3	1:A:241:PRO:HD3	1.98	0.45
1:A:241:PRO:HB2	3:A:2065:HOH:O	2.16	0.45
1:A:450:ASP:HA	1:A:451:PRO:HD3	1.79	0.45
1:A:48:ARG:HB2	3:A:2051:HOH:O	2.17	0.45
1:A:256:LYS:HG2	3:A:2068:HOH:O	2.17	0.44
1:A:337:ASP:CG	1:A:338:PRO:HD2	2.38	0.44
1:A:178:GLU:HB2	1:A:446:GLU:OE1	2.18	0.44
1:A:17:ALA:HB3	3:A:2004:HOH:O	2.16	0.43
1:A:230:GLY:O	1:A:234:ARG:HG3	2.18	0.43
1:A:335:PRO:HG3	1:A:342:PHE:CD2	2.53	0.43
1:A:426:ASP:HB3	1:A:458:LEU:HD11	2.00	0.43
1:A:172:GLU:OE1	1:A:185:ARG:NH2	2.51	0.43
1:A:110:LEU:HB3	1:A:112:ILE:HG12	2.00	0.43
1:A:11:VAL:HG12	1:A:99:VAL:HB	2.01	0.43
1:A:376:ILE:HG13	2:A:801:FAD:C2	2.49	0.43
1:A:305:ILE:HG22	1:A:306:GLY:N	2.34	0.42
1:A:104:ALA:HA	1:A:333:SER:HA	2.00	0.42
1:A:117:LEU:HA	1:A:118:PRO:HD3	1.93	0.42
1:A:197:ARG:HH11	1:A:197:ARG:CG	2.33	0.42
1:A:40:GLN:NE2	1:A:40:GLN:HA	2.34	0.42
1:A:21:THR:O	1:A:25:LEU:HG	2.19	0.42
1:A:124:ARG:NH1	1:A:367:TRP:HZ3	2.17	0.42
1:A:284:LEU:HA	1:A:285:PRO:HD3	1.85	0.42
1:A:49:PHE:CE1	1:A:133:LEU:HA	2.55	0.41
1:A:89:GLN:HB2	3:A:2082:HOH:O	2.20	0.41
1:A:197:ARG:CZ	3:A:2058:HOH:O	2.67	0.41
1:A:337:ASP:OD1	1:A:338:PRO:HD2	2.21	0.41
1:A:149:VAL:HG22	1:A:193:TRP:HB2	2.03	0.41
1:A:55:HIS:N	1:A:56:PRO:CD	2.83	0.41
1:A:194:ILE:HB	1:A:274:LEU:HD12	2.03	0.40
1:A:288:ASP:OD2	1:A:290:ARG:HG3	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:ARG:HB2	3:A:2065:HOH:O	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:MET:CE	3:A:2103:HOH:O[2_545]	1.34	0.86

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	453/460 (98%)	428 (94%)	23 (5%)	2 (0%)	34 48

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	403	PRO
1	A	329	ILE

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	374/379 (99%)	358 (96%)	16 (4%)	29 46

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	THR
1	A	31	ARG
1	A	60	ASN
1	A	64	THR
1	A	110	LEU
1	A	137	ARG
1	A	170	HIS
1	A	171	LEU
1	A	197	ARG
1	A	229	LEU
1	A	231	LEU
1	A	250	LEU
1	A	305	ILE
1	A	352	MET
1	A	373	THR
1	A	420	TRP

Some 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	GLN
1	A	28	HIS
1	A	40	GLN
1	A	136	ASN
1	A	215	GLN
1	A	282	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 carbohydrates in this entry.

5.6 Ligand geometry

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	801	-	51,58,58	1.73	8 (15%)	60,89,89	2.08	18 (30%)

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	801	-	-	10/30/50/50	0/6/6/6

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	FAD	C4X-C10	4.70	1.43	1.38
2	A	801	FAD	O5'-C5'	4.15	1.60	1.44
2	A	801	FAD	O4B-C1B	4.05	1.46	1.41
2	A	801	FAD	PA-O2A	-3.90	1.37	1.55
2	A	801	FAD	P-O2P	-3.43	1.39	1.55
2	A	801	FAD	C2-N1	-3.05	1.32	1.38
2	A	801	FAD	C5'-C4'	2.91	1.55	1.51
2	A	801	FAD	C2-N3	2.29	1.42	1.38

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	FAD	C4-N3-C2	8.54	122.35	115.14
2	A	801	FAD	C4X-C4-N3	-5.21	116.31	123.43
2	A	801	FAD	O4B-C1B-C2B	-4.23	100.75	106.93
2	A	801	FAD	O4'-C4'-C3'	3.42	117.42	109.10
2	A	801	FAD	O2'-C2'-C3'	3.20	116.88	109.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	FAD	C10-C4X-N5	2.99	123.33	121.26
2	A	801	FAD	O2'-C2'-C1'	-2.93	102.54	109.59
2	A	801	FAD	O3'-C3'-C2'	-2.90	101.81	108.81
2	A	801	FAD	O2A-PA-O1A	2.61	125.17	112.24
2	A	801	FAD	C5A-C6A-N1A	-2.39	114.93	120.35
2	A	801	FAD	C5A-C6A-N6A	2.30	123.84	120.35
2	A	801	FAD	O5'-P-O1P	-2.27	100.18	109.07
2	A	801	FAD	O5B-PA-O1A	-2.25	100.29	109.07
2	A	801	FAD	C7M-C7-C8	-2.13	116.37	120.74
2	A	801	FAD	C4A-C5A-N7A	2.07	111.55	109.40
2	A	801	FAD	O3'-C3'-C4'	-2.04	103.87	108.81
2	A	801	FAD	C5'-C4'-C3'	-2.01	108.32	112.20
2	A	801	FAD	C1'-N10-C9A	2.01	119.87	118.29

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	FAD	C5B-O5B-PA-O3P
2	A	801	FAD	O3'-C3'-C4'-O4'
2	A	801	FAD	O3'-C3'-C4'-C5'
2	A	801	FAD	O4B-C4B-C5B-O5B
2	A	801	FAD	C3B-C4B-C5B-O5B
2	A	801	FAD	C2'-C3'-C4'-O4'
2	A	801	FAD	PA-O3P-P-O5'
2	A	801	FAD	C5B-O5B-PA-O1A
2	A	801	FAD	C5B-O5B-PA-O2A
2	A	801	FAD	P-O3P-PA-O2A

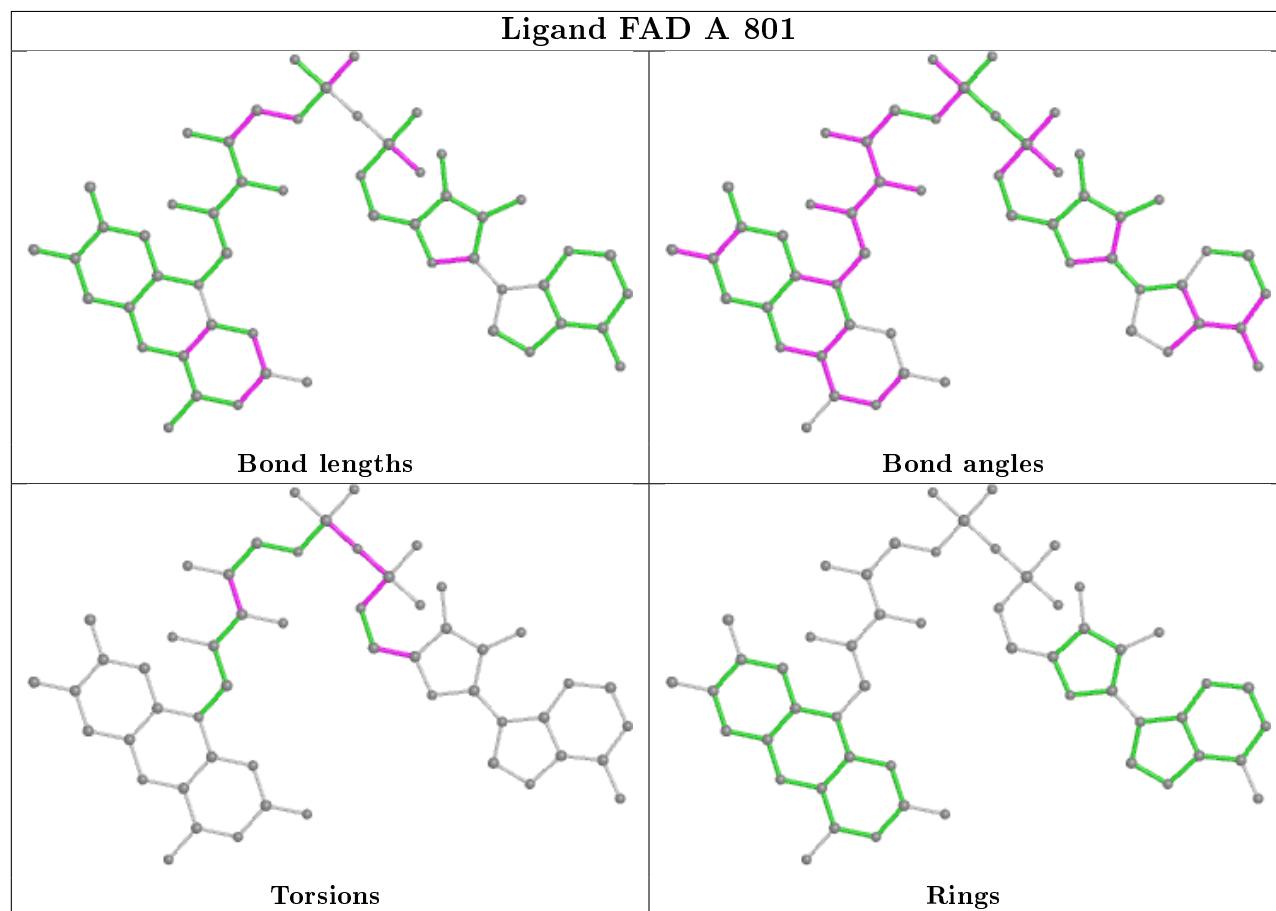
There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	FAD	3	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	455/460 (98%)	-0.19	16 (3%) 44 43	15, 40, 73, 104	3 (0%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	305	ILE	4.5
1	A	309	THR	3.8
1	A	307	GLU	3.8
1	A	256	LYS	3.4
1	A	346	LEU	3.3
1	A	420	TRP	3.3
1	A	402	GLY	3.3
1	A	306	GLY	3.2
1	A	108	GLN	2.6
1	A	308	ALA	2.5
1	A	259	VAL	2.2
1	A	237	GLU	2.2
1	A	334	ARG	2.1
1	A	304	GLY	2.1
1	A	138	GLU	2.0
1	A	316	ASP	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](#)

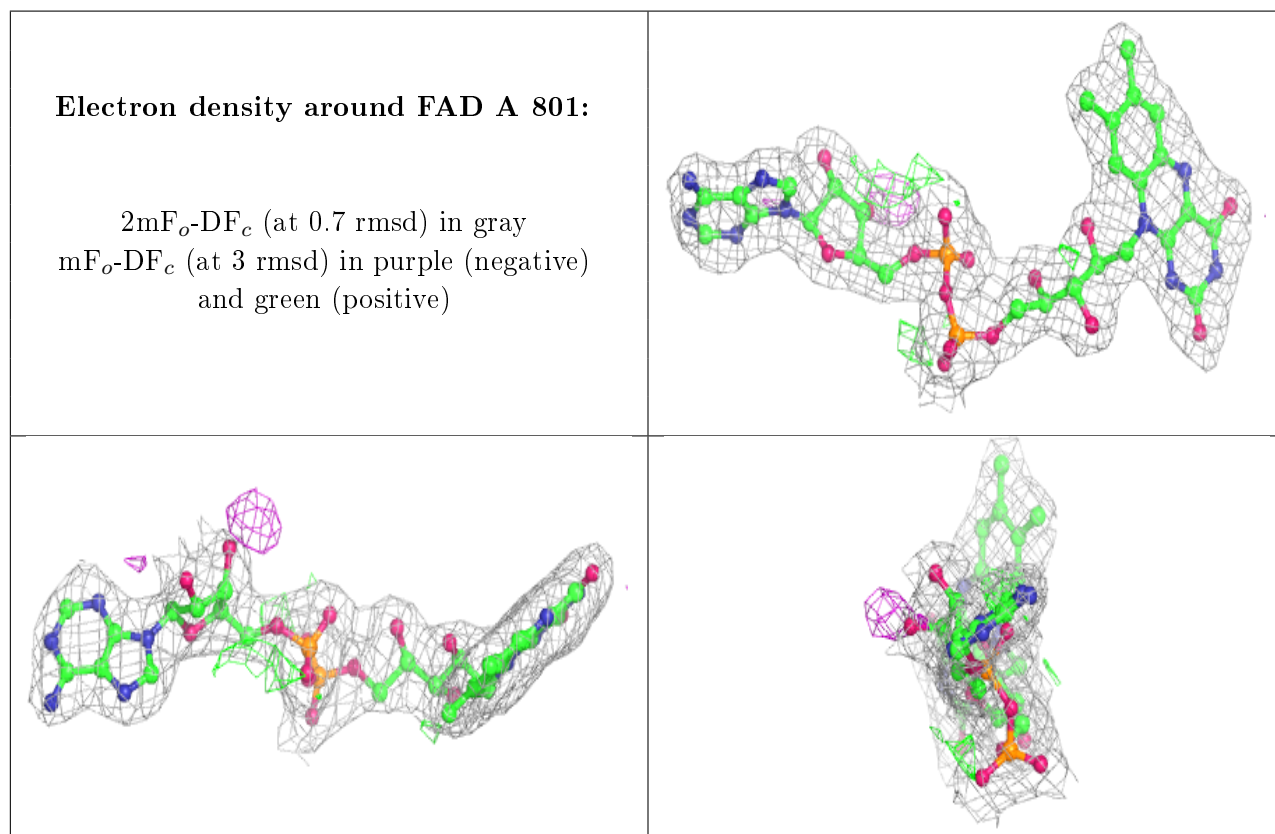
There are no carbohydrates in this entry.

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(Å ²)	Q<0.9
2	FAD	A	801	53/53	0.97	0.15	16,26,30,32	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.