



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 6, 2021 – 05:05 am GMT

PDB ID : 7AAW  
Title : Thioredoxin Reductase from *Bacillus cereus*  
Authors : Shoor, M.; Gudim, I.; Hersleth, H.-P.; Hammerstad, M.  
Deposited on : 2020-09-04  
Resolution : 2.25 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.4 (270009), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.23.2  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0267  
CCP4 : 7.1.010 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.23.2

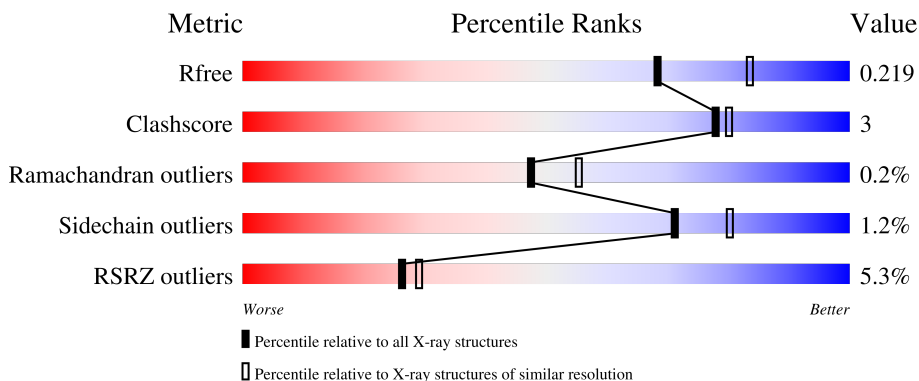
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

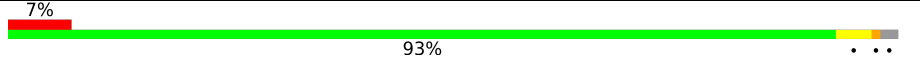
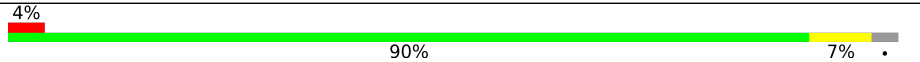
The reported resolution of this entry is 2.25 Å.

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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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  $\geq 3$ , 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	321	 7% 93%
1	B	321	 4% 90% 7%

## 2 Entry composition i

There are 7 unique types of molecules in this entry. The entry contains 5262 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 Thioredoxin reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	315	Total 2430	C 1536	N 409	O 476	S 9	0	2	0
1	B	311	Total 2416	C 1530	N 407	O 470	S 9	0	4	0

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:  $C_{27}H_{33}N_9O_{15}P_2$ ) (labeled as "Ligand of Interest" by depositor).



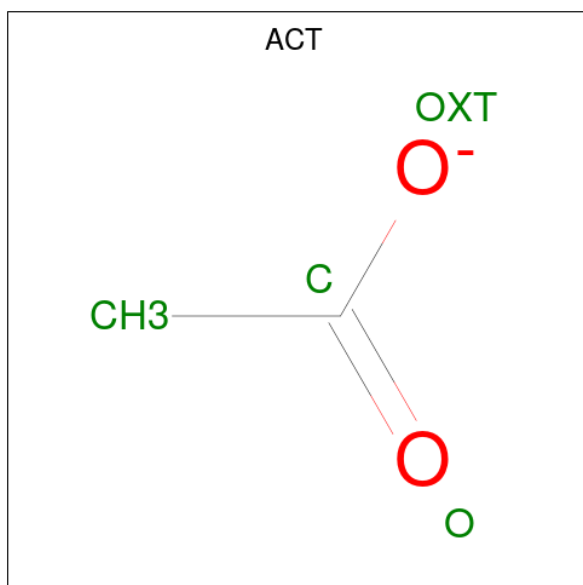
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total 53	C 27	N 9	O 15	P 2	0	0
2	B	1	Total 53	C 27	N 9	O 15	P 2	0	0

- Molecule 3 is alpha-D-glucopyranose (three-letter code: GLC) (formula:  $C_6H_{12}O_6$ ).



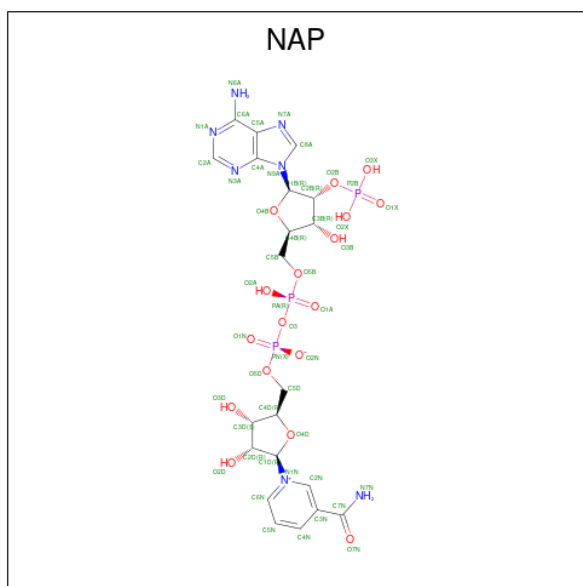
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 12 6 6	0	0
3	A	1	Total C O 12 6 6	0	0
3	A	1	Total C O 12 6 6	0	0
3	B	1	Total C O 12 6 6	0	0

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula:  $C_2H_3O_2^-$ ).



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

- Molecule 5 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula:  $C_{21}H_{28}N_7O_{17}P_3$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Na	0	0
			1	1		

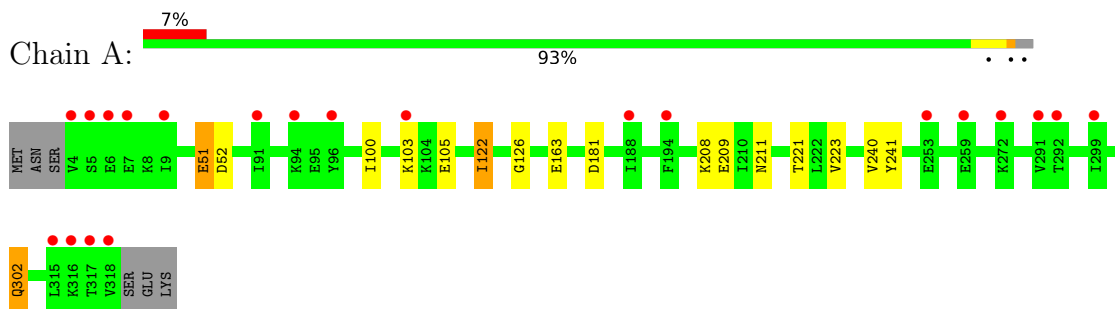
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	107	Total	O	0	0
			107	107		
7	B	102	Total	O	0	0
			102	102		

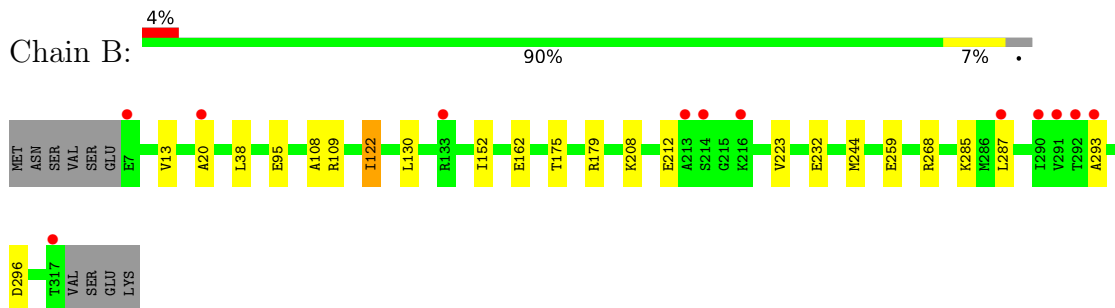
### 3 Residue-property plots i

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: Thioredoxin reductase



- Molecule 1: Thioredoxin reductase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	94.98Å 94.98Å 214.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.72 – 2.25 46.72 – 2.25	Depositor EDS
% Data completeness (in resolution range)	98.8 (46.72-2.25) 98.8 (46.72-2.25)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.83 (at 2.24Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660, REFMAC 5.8.0253	Depositor
R, $R_{free}$	0.178 , 0.219 0.179 , 0.219	Depositor DCC
$R_{free}$ test set	2289 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	45.8	Xtrriage
Anisotropy	0.214	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	(Not available) , (Not available)	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5262	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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, GLC, NA, NAP, ACT

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	0.37	0/2467	0.54	0/3323
1	B	0.39	0/2453	0.56	0/3304
All	All	0.38	0/4920	0.55	0/6627

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 [i](#)

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	2430	0	2433	13	0
1	B	2416	0	2426	15	0
2	A	53	0	31	0	0
2	B	53	0	31	0	0
3	A	36	0	36	2	0
3	B	12	0	12	0	0
4	A	4	0	3	0	0
5	B	48	0	25	2	0
6	B	1	0	0	0	0
7	A	107	0	0	1	0
7	B	102	0	0	1	0
All	All	5262	0	4997	27	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 3.

All (27) 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:130:LEU:HD21	1:B:212:GLU:HG2	1.80	0.62
1:B:285:LYS:NZ	1:B:296:ASP:OD2	2.30	0.60
1:A:209:GLU:HG2	1:A:221:THR:HB	1.88	0.55
1:B:268[B]:ARG:NH2	7:B:505:HOH:O	2.42	0.52
1:B:162:GLU:OE1	5:B:402:NAP:H4N	2.14	0.48
1:A:208:LYS:HB2	1:A:223:VAL:HG23	1.94	0.48
1:B:259:GLU:H	1:B:259:GLU:CD	2.17	0.47
1:A:181:ASP:O	7:A:501:HOH:O	2.20	0.47
1:A:208:LYS:HB2	1:A:223:VAL:CG2	2.45	0.47
1:B:13:VAL:HG23	1:B:108:ALA:HB2	1.97	0.46
1:A:100:ILE:HG12	1:A:105:GLU:HG2	1.97	0.46
1:B:208:LYS:HB2	1:B:223:VAL:HG23	1.98	0.46
1:A:163:GLU:HG3	1:A:241:TYR:CZ	2.52	0.45
1:B:130:LEU:HD11	1:B:212:GLU:HG2	1.98	0.45
1:B:152:ILE:HD13	1:B:175:THR:HB	1.98	0.44
1:B:20:ALA:HB2	1:B:293:ALA:HB1	2.00	0.44
1:A:126:GLY:CA	3:A:404:GLC:H5	2.48	0.44
1:A:211:ASN:OD1	3:A:404:GLC:O2	2.37	0.43
1:B:122:ILE:H	1:B:122:ILE:HG12	1.61	0.43
1:A:103:LYS:HD3	1:A:103:LYS:HA	1.82	0.43
1:A:122:ILE:HD11	1:A:240:VAL:HG22	2.00	0.43
1:A:163:GLU:HG3	1:A:241:TYR:CE1	2.54	0.42
1:B:179:ARG:HH21	5:B:402:NAP:P2B	2.43	0.42
1:B:13:VAL:CG2	1:B:108:ALA:HB2	2.50	0.42
1:B:95:GLU:O	1:B:109:ARG:NH2	2.49	0.41
1:A:51:GLU:HB3	1:A:52:ASP:H	1.60	0.41
1:A:302[B]:GLN:HG3	1:B:287:LEU:HD22	2.03	0.41

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	315/321 (98%)	310 (98%)	4 (1%)	1 (0%)	41	46
1	B	313/321 (98%)	307 (98%)	6 (2%)	0	100	100
All	All	628/642 (98%)	617 (98%)	10 (2%)	1 (0%)	47	55

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	51	GLU

### 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	255/259 (98%)	252 (99%)	3 (1%)	71	80
1	B	253/259 (98%)	249 (98%)	4 (2%)	62	73
All	All	508/518 (98%)	501 (99%)	7 (1%)	71	76

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	122	ILE
1	A	302[A]	GLN
1	A	302[B]	GLN
1	B	38	LEU
1	B	122	ILE
1	B	232	GLU
1	B	244	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates

There are no monosaccharides in this entry.

## 5.6 Ligand geometry

Of 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

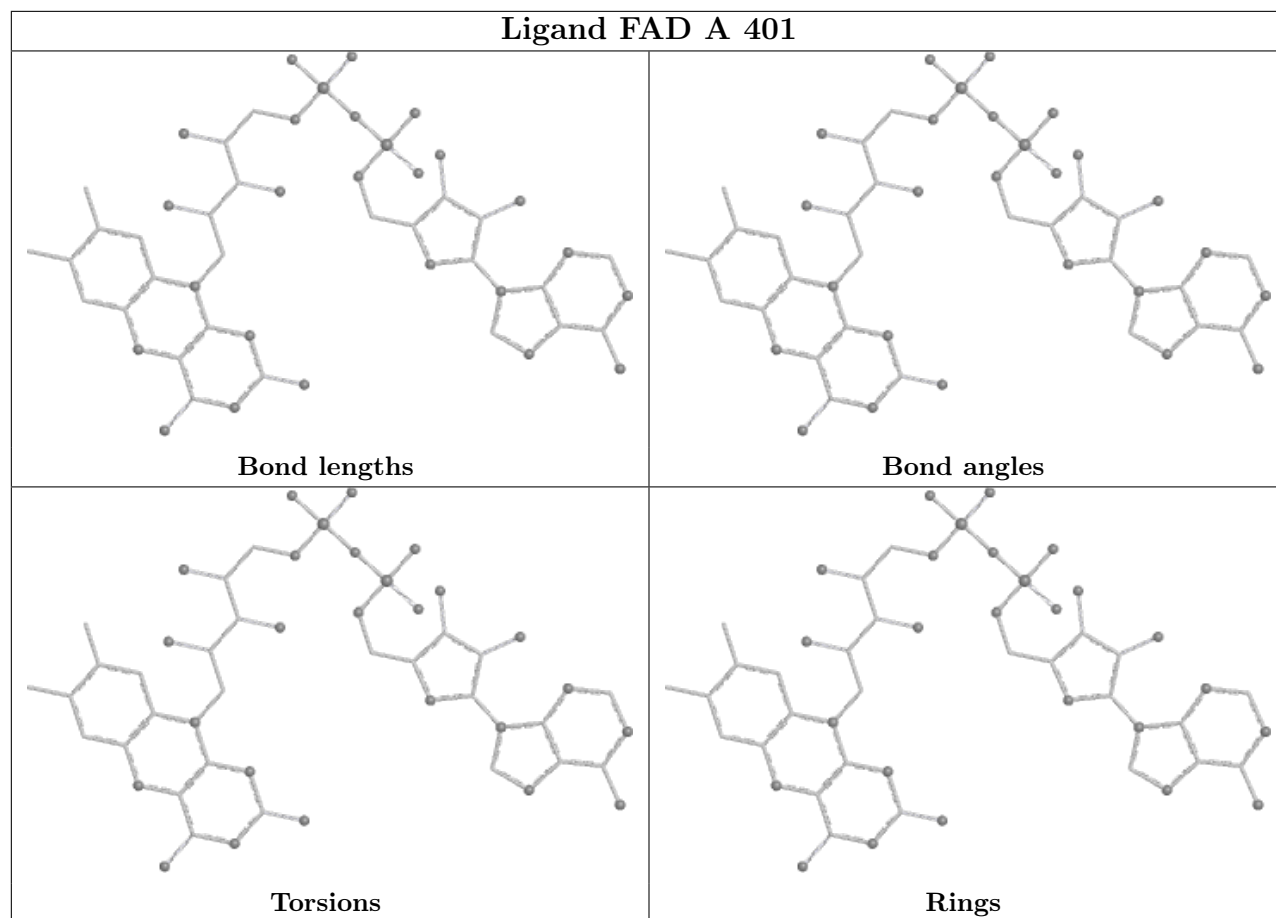
There are no chirality outliers.

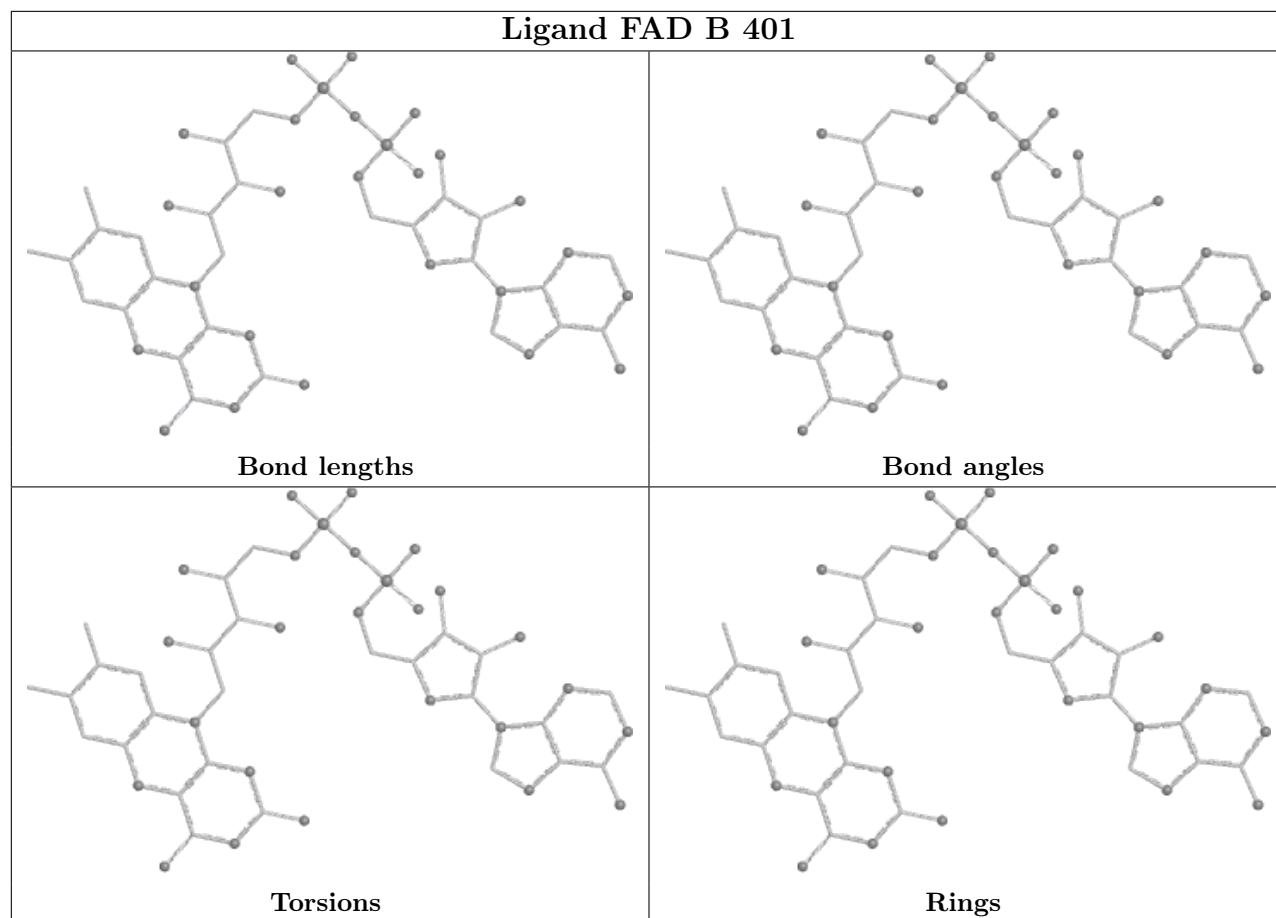
There are no torsion outliers.

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	315/321 (98%)	0.26	21 (6%) 17 19	31, 44, 72, 101	0
1	B	311/321 (96%)	0.19	12 (3%) 39 42	29, 44, 68, 90	0
All	All	626/642 (97%)	0.23	33 (5%) 26 29	29, 44, 71, 101	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	318	VAL	6.0
1	A	315	LEU	5.7
1	A	5	SER	4.8
1	A	103	LYS	3.9
1	A	4	VAL	3.5
1	B	213	ALA	3.3
1	A	188	ILE	2.9
1	B	291	VAL	2.8
1	B	317	THR	2.7
1	A	96	TYR	2.6
1	A	94	LYS	2.6
1	A	7	GLU	2.5
1	A	316	LYS	2.5
1	A	6	GLU	2.4
1	B	287	LEU	2.3
1	A	253	GLU	2.3
1	A	259	GLU	2.2
1	B	292	THR	2.2
1	A	299	ILE	2.2
1	B	290	ILE	2.2
1	B	133	ARG	2.2
1	B	214	SER	2.2
1	A	272	LYS	2.2
1	B	293	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	291	VAL	2.1
1	A	317	THR	2.1
1	A	292	THR	2.1
1	B	7	GLU	2.1
1	A	91	ILE	2.1
1	B	216	LYS	2.1
1	A	194	PHE	2.1
1	A	9	ILE	2.0
1	B	20	ALA	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 monosaccharides in this entry.

## 6.4 Ligands [\(i\)](#)

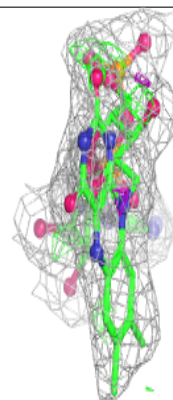
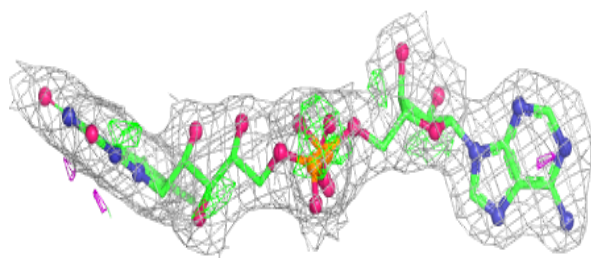
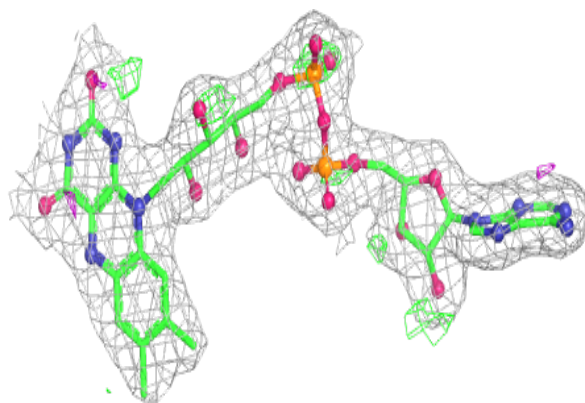
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
3	GLC	A	404	12/12	0.69	0.26	40,50,55,60	12
3	GLC	A	402	12/12	0.81	0.26	48,68,75,76	0
3	GLC	A	403	12/12	0.87	0.23	48,67,72,84	0
3	GLC	B	403	12/12	0.90	0.24	59,65,68,70	0
4	ACT	A	405	4/4	0.90	0.17	56,62,63,68	0
2	FAD	B	401	53/53	0.97	0.17	28,36,43,47	0
2	FAD	A	401	53/53	0.97	0.15	30,39,45,47	0
5	NAP	B	402	48/48	0.97	0.11	41,51,58,61	0
6	NA	B	404	1/1	0.97	0.13	47,47,47,47	0

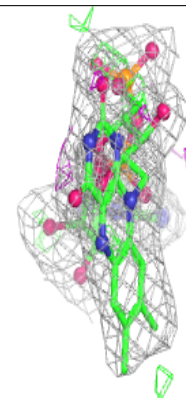
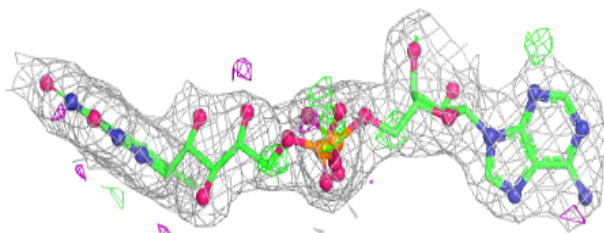
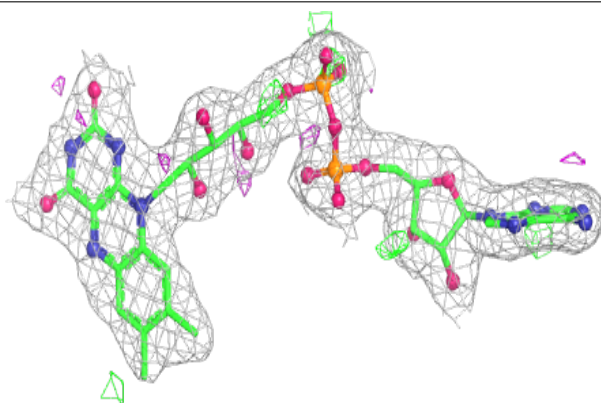
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around FAD B 401:**

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

**Electron density around FAD A 401:**

$2mF_o-DF_c$  (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.