



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 16, 2017 – 12:34 AM EDT

PDB ID : 5WED  
Title : Structure of bacterial type II NADH dehydrogenase from *Caldalkalibacillus thermarum* at 2.15Å resolution  
Authors : Nakatani, Y.; Aragao, D.; Cook, G.M.  
Deposited on : unknown  
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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030345  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20030345

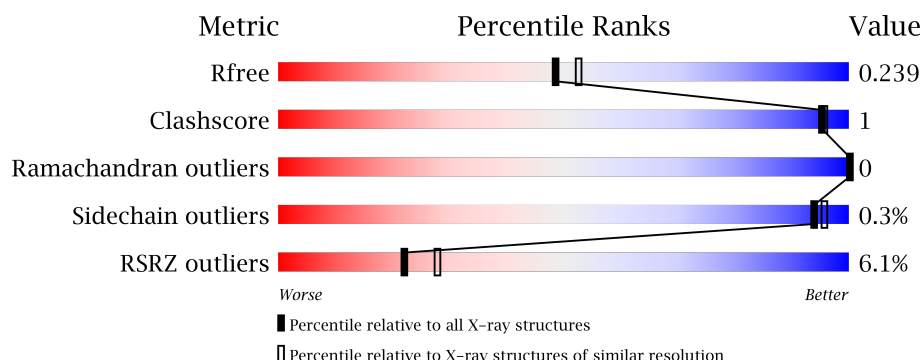
# 1 Overall quality at a glance

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))
$R_{free}$	100719	1170 (2.16-2.16)
Clashscore	112137	1278 (2.16-2.16)
Ramachandran outliers	110173	1256 (2.16-2.16)
Sidechain outliers	110143	1255 (2.16-2.16)
RSRZ outliers	101464	1175 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	405	<div> <div>6%</div> <div>94%</div> <div>• •</div> </div>
1	B	405	<div> <div>4%</div> <div>93%</div> <div>• •</div> </div>
1	C	405	<div> <div>7%</div> <div>93%</div> <div>• •</div> </div>
1	D	405	<div> <div>7%</div> <div>94%</div> <div>• •</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12435 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 FAD-dependent pyridine nucleotide-disulfide oxidoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	394	Total	C	N	O	S	0	0	0
			2948	1882	509	549	8			
1	A	394	Total	C	N	O	S	0	0	0
			2946	1880	509	549	8			
1	C	394	Total	C	N	O	S	0	0	0
			2933	1876	496	552	9			
1	D	394	Total	C	N	O	S	0	0	0
			2894	1855	506	525	8			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	400	HIS	-	expression tag	UNP F5L3B8
B	401	HIS	-	expression tag	UNP F5L3B8
B	402	HIS	-	expression tag	UNP F5L3B8
B	403	HIS	-	expression tag	UNP F5L3B8
B	404	HIS	-	expression tag	UNP F5L3B8
B	405	HIS	-	expression tag	UNP F5L3B8
A	400	HIS	-	expression tag	UNP F5L3B8
A	401	HIS	-	expression tag	UNP F5L3B8
A	402	HIS	-	expression tag	UNP F5L3B8
A	403	HIS	-	expression tag	UNP F5L3B8
A	404	HIS	-	expression tag	UNP F5L3B8
A	405	HIS	-	expression tag	UNP F5L3B8
C	400	HIS	-	expression tag	UNP F5L3B8
C	401	HIS	-	expression tag	UNP F5L3B8
C	402	HIS	-	expression tag	UNP F5L3B8
C	403	HIS	-	expression tag	UNP F5L3B8
C	404	HIS	-	expression tag	UNP F5L3B8
C	405	HIS	-	expression tag	UNP F5L3B8
D	400	HIS	-	expression tag	UNP F5L3B8
D	401	HIS	-	expression tag	UNP F5L3B8
D	402	HIS	-	expression tag	UNP F5L3B8

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Chain	Residue	Modelled	Actual	Comment	Reference
D	403	HIS	-	expression tag	UNP F5L3B8
D	404	HIS	-	expression tag	UNP F5L3B8
D	405	HIS	-	expression tag	UNP F5L3B8

- # FAD

- Molecule 3 is water.

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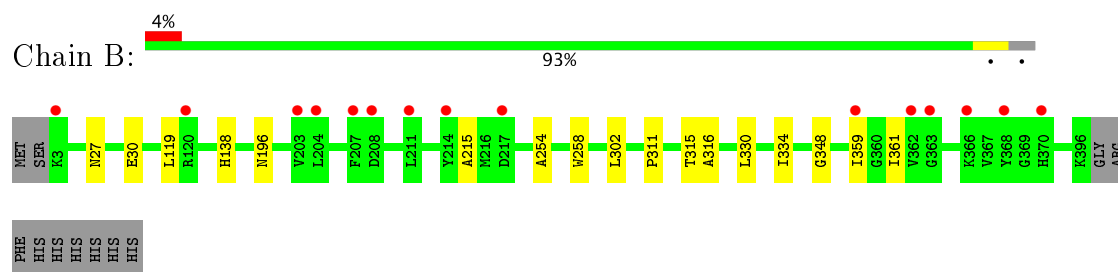
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	99	Total	O	0	0
			99	99		

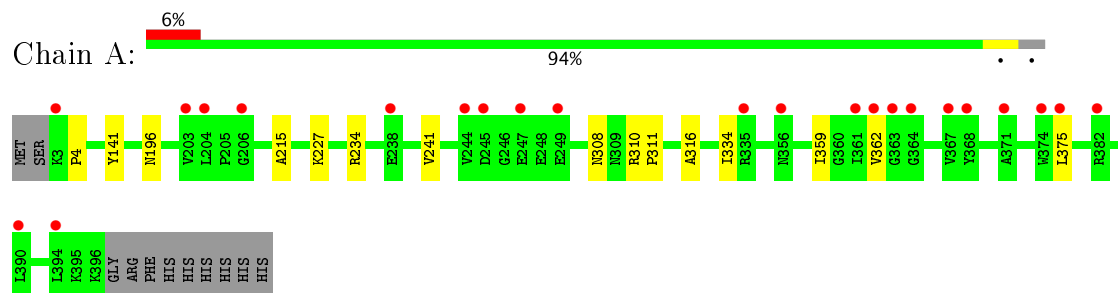
### 3 Residue-property plots [i](#)

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.

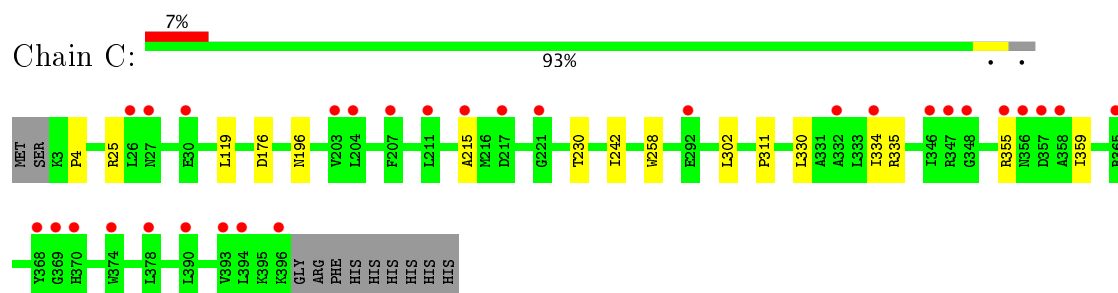
- Molecule 1: FAD-dependent pyridine nucleotide-disulfide oxidoreductase



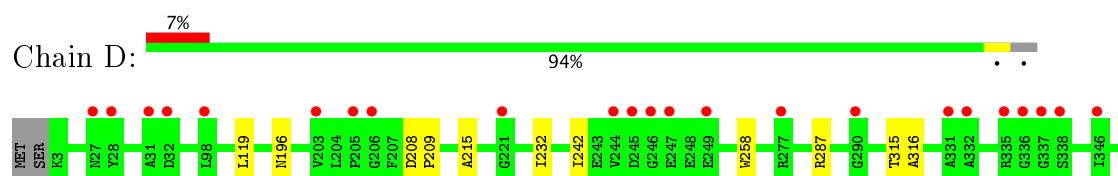
- Molecule 1: FAD-dependent pyridine nucleotide-disulfide oxidoreductase



- Molecule 1: FAD-dependent pyridine nucleotide-disulfide oxidoreductase



- Molecule 1: FAD-dependent pyridine nucleotide-disulfide oxidoreductase



R355	
I359	
Y368	
W374	
L394	
K396	
GLY	
ARG	
PHE	
HIS	
HIS	
HIS	
HIS	
HIS	
HIS	

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	72.80 Å 113.57 Å 129.78 Å 90.00° 91.02° 90.00°	Depositor
Resolution (Å)	48.87 – 2.15 48.87 – 2.15	Depositor EDS
% Data completeness (in resolution range)	91.7 (48.87-2.15) 91.8 (48.87-2.15)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.47 (at 2.16 Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.208 , 0.238 0.208 , 0.239	Depositor DCC
$R_{free}$ test set	5247 reflections (4.99%)	DCC
Wilson B-factor (Å <sup>2</sup> )	41.4	Xtriage
Anisotropy	0.106	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 45.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.022 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	12435	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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

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	0.22	0/3004	0.41	0/4088
1	B	0.22	0/3005	0.42	0/4089
1	C	0.22	0/2989	0.41	0/4068
1	D	0.21	0/2952	0.40	0/4022
All	All	0.22	0/11950	0.41	0/16267

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	2946	0	2899	9	0
1	B	2948	0	2914	10	0
1	C	2933	0	2891	8	0
1	D	2894	0	2844	7	0
2	A	53	0	31	1	0
2	B	53	0	31	2	0
2	C	53	0	31	0	0
2	D	53	0	31	2	0
3	A	122	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	154	0	0	0	0
3	C	127	0	0	0	0
3	D	99	0	0	1	0
All	All	12435	0	11672	33	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 1.

All (33) 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:215:ALA:HB2	1:A:359:ILE:HD13	1.84	0.58
1:B:302:LEU:HD11	1:B:311:PRO:HB3	1.86	0.57
1:D:215:ALA:HB2	1:D:359:ILE:HD13	1.86	0.57
1:B:215:ALA:HB2	1:B:359:ILE:HD13	1.85	0.57
1:A:227:LYS:NZ	3:A:702:HOH:O	2.37	0.57
1:D:232:ILE:HD13	1:D:242:ILE:HG22	1.87	0.56
1:D:316:ALA:HB2	2:D:601:FAD:H2'	1.93	0.51
1:C:215:ALA:HB2	1:C:359:ILE:HD13	1.93	0.51
1:C:4:PRO:HG3	1:C:334:ILE:HD12	1.93	0.51
1:B:27:ASN:N	1:B:30:GLU:OE2	2.35	0.49
1:C:230:THR:HG21	1:C:242:ILE:HD12	1.94	0.48
1:C:330:LEU:O	1:C:334:ILE:HG12	2.13	0.48
1:A:234:ARG:HG2	1:A:241:VAL:HB	1.98	0.46
1:C:25:ARG:O	1:C:335:ARG:NH2	2.48	0.46
1:A:316:ALA:HB2	2:A:601:FAD:H2'	1.97	0.46
1:C:302:LEU:HD11	1:C:311:PRO:HB3	1.97	0.45
1:B:330:LEU:O	1:B:334:ILE:HG12	2.17	0.45
1:D:119:LEU:HD21	1:D:258:TRP:CE3	2.53	0.44
1:B:315:THR:HB	2:B:601:FAD:O2	2.17	0.44
1:C:176:ASP:OD2	1:C:355:ARG:N	2.42	0.44
1:B:316:ALA:HB2	2:B:601:FAD:H2'	1.99	0.44
1:B:119:LEU:HD21	1:B:258:TRP:CE3	2.52	0.43
1:A:4:PRO:HG3	1:A:334:ILE:HD12	1.99	0.43
1:A:310:ARG:HA	1:A:311:PRO:HD3	1.92	0.42
1:B:348:GLY:HA3	1:B:361:ILE:O	2.18	0.42
1:C:119:LEU:HD21	1:C:258:TRP:CE3	2.55	0.42
1:A:362:VAL:HG21	1:A:375:LEU:HD13	2.01	0.41
1:D:287:ARG:NH1	3:D:703:HOH:O	2.41	0.41
1:A:308:ASN:OD1	1:A:310:ARG:HG2	2.21	0.41
1:D:315:THR:HB	2:D:601:FAD:O2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:138:HIS:HE1	1:B:254:ALA:O	2.04	0.41
1:D:208:ASP:HA	1:D:209:PRO:HD3	1.96	0.41
1:B:138:HIS:HD2	1:A:141:TYR:OH	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	392/405 (97%)	383 (98%)	9 (2%)	0	100	100
1	B	392/405 (97%)	382 (97%)	10 (3%)	0	100	100
1	C	392/405 (97%)	382 (97%)	10 (3%)	0	100	100
1	D	392/405 (97%)	384 (98%)	8 (2%)	0	100	100
All	All	1568/1620 (97%)	1531 (98%)	37 (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	297/329 (90%)	296 (100%)	1 (0%)	94	96
1	B	298/329 (91%)	297 (100%)	1 (0%)	94	96

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	297/329 (90%)	296 (100%)	1 (0%)	94	96
1	D	283/329 (86%)	282 (100%)	1 (0%)	93	95
All	All	1175/1316 (89%)	1171 (100%)	4 (0%)	94	96

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	196	ASN
1	A	196	ASN
1	C	196	ASN
1	D	196	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	89	GLN
1	B	138	HIS
1	B	196	ASN
1	B	213	ASN
1	A	39	ASN
1	A	89	GLN
1	A	196	ASN
1	A	213	ASN
1	C	89	GLN
1	C	196	ASN
1	C	213	ASN
1	D	122	HIS
1	D	196	ASN
1	D	213	ASN
1	D	317	GLN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry

4 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	FAD	A	601	-	51,58,58	4.16	18 (35%)	54,89,89	2.64	10 (18%)
2	FAD	B	601	-	51,58,58	4.15	18 (35%)	54,89,89	2.67	11 (20%)
2	FAD	C	601	-	51,58,58	4.17	18 (35%)	54,89,89	2.66	11 (20%)
2	FAD	D	601	-	51,58,58	4.16	18 (35%)	54,89,89	2.66	10 (18%)

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	601	-	-	0/28/50/50	0/6/6/6
2	FAD	B	601	-	-	0/28/50/50	0/6/6/6
2	FAD	C	601	-	-	0/28/50/50	0/6/6/6
2	FAD	D	601	-	-	0/28/50/50	0/6/6/6

All (72) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	601	FAD	C2B-C1B	-15.42	1.29	1.53
2	C	601	FAD	C2B-C1B	-15.41	1.29	1.53
2	A	601	FAD	C2B-C1B	-15.26	1.29	1.53
2	B	601	FAD	C2B-C1B	-15.16	1.29	1.53
2	D	601	FAD	O4B-C4B	-6.97	1.29	1.45
2	C	601	FAD	O4B-C4B	-6.90	1.29	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	FAD	O4B-C4B	-6.86	1.29	1.45
2	B	601	FAD	O4B-C4B	-6.83	1.29	1.45
2	B	601	FAD	O4-C4	-3.68	1.15	1.24
2	A	601	FAD	O4-C4	-3.67	1.15	1.24
2	D	601	FAD	O4-C4	-3.64	1.15	1.24
2	C	601	FAD	O4-C4	-3.64	1.15	1.24
2	D	601	FAD	O4'-C4'	-2.68	1.37	1.43
2	A	601	FAD	O4'-C4'	-2.61	1.37	1.43
2	B	601	FAD	O4'-C4'	-2.61	1.37	1.43
2	C	601	FAD	O4'-C4'	-2.58	1.37	1.43
2	B	601	FAD	C7M-C7	2.74	1.56	1.51
2	D	601	FAD	C7M-C7	2.75	1.56	1.51
2	A	601	FAD	C7M-C7	2.78	1.56	1.51
2	C	601	FAD	C7M-C7	2.80	1.56	1.51
2	C	601	FAD	C6A-N6A	2.88	1.45	1.34
2	D	601	FAD	C6A-N6A	2.89	1.45	1.34
2	B	601	FAD	C6A-N6A	2.92	1.46	1.34
2	A	601	FAD	C6A-N6A	2.93	1.46	1.34
2	A	601	FAD	C2A-N3A	2.97	1.37	1.32
2	B	601	FAD	C2A-N3A	3.03	1.37	1.32
2	C	601	FAD	C2A-N3A	3.05	1.37	1.32
2	B	601	FAD	C8M-C8	3.06	1.57	1.51
2	D	601	FAD	C8M-C8	3.06	1.57	1.51
2	D	601	FAD	C2A-N3A	3.09	1.37	1.32
2	C	601	FAD	C8M-C8	3.10	1.57	1.51
2	A	601	FAD	C8M-C8	3.18	1.57	1.51
2	D	601	FAD	C2-N3	4.13	1.46	1.38
2	B	601	FAD	C2-N3	4.19	1.46	1.38
2	C	601	FAD	C2-N3	4.22	1.46	1.38
2	A	601	FAD	C2-N3	4.22	1.46	1.38
2	A	601	FAD	C2-N1	4.28	1.46	1.38
2	D	601	FAD	C4-C4X	4.34	1.49	1.41
2	B	601	FAD	C4-C4X	4.34	1.49	1.41
2	C	601	FAD	C2-N1	4.40	1.46	1.38
2	B	601	FAD	C2-N1	4.40	1.46	1.38
2	A	601	FAD	C4-C4X	4.41	1.49	1.41
2	C	601	FAD	C4-C4X	4.42	1.49	1.41
2	D	601	FAD	C2-N1	4.45	1.47	1.38
2	C	601	FAD	O2B-C2B	4.48	1.53	1.43
2	D	601	FAD	O2B-C2B	4.48	1.53	1.43
2	A	601	FAD	O2B-C2B	4.49	1.53	1.43
2	B	601	FAD	O2B-C2B	4.55	1.53	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	FAD	C4-N3	5.38	1.42	1.33
2	A	601	FAD	C9A-N10	5.39	1.45	1.38
2	D	601	FAD	C4-N3	5.52	1.43	1.33
2	C	601	FAD	C9A-N10	5.55	1.46	1.38
2	D	601	FAD	C9A-N10	5.56	1.46	1.38
2	C	601	FAD	C4-N3	5.58	1.43	1.33
2	B	601	FAD	C9A-N10	5.59	1.46	1.38
2	A	601	FAD	C4-N3	5.59	1.43	1.33
2	B	601	FAD	C4X-N5	6.43	1.42	1.33
2	D	601	FAD	C4X-N5	6.48	1.42	1.33
2	A	601	FAD	C4X-N5	6.69	1.43	1.33
2	C	601	FAD	C4X-N5	6.70	1.43	1.33
2	A	601	FAD	C10-N1	7.13	1.43	1.33
2	D	601	FAD	C10-N1	7.16	1.43	1.33
2	B	601	FAD	C10-N1	7.18	1.43	1.33
2	C	601	FAD	C10-N1	7.24	1.43	1.33
2	B	601	FAD	C5X-N5	8.75	1.48	1.35
2	C	601	FAD	C5X-N5	8.76	1.48	1.35
2	D	601	FAD	C5X-N5	8.78	1.48	1.35
2	A	601	FAD	C5X-N5	8.99	1.49	1.35
2	D	601	FAD	O4B-C1B	14.00	1.60	1.41
2	A	601	FAD	O4B-C1B	14.03	1.60	1.41
2	C	601	FAD	O4B-C1B	14.04	1.60	1.41
2	B	601	FAD	O4B-C1B	14.24	1.61	1.41

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	601	FAD	N3A-C2A-N1A	-10.39	119.81	128.86
2	A	601	FAD	N3A-C2A-N1A	-10.09	120.07	128.86
2	B	601	FAD	N3A-C2A-N1A	-10.07	120.09	128.86
2	D	601	FAD	N3A-C2A-N1A	-10.05	120.11	128.86
2	B	601	FAD	C7M-C7-C6	-6.71	103.53	120.34
2	A	601	FAD	C7M-C7-C6	-6.67	103.63	120.34
2	D	601	FAD	C7M-C7-C6	-6.64	103.70	120.34
2	C	601	FAD	C7M-C7-C6	-6.47	104.13	120.34
2	B	601	FAD	N6A-C6A-N1A	-6.25	106.38	118.77
2	D	601	FAD	N6A-C6A-N1A	-6.14	106.59	118.77
2	A	601	FAD	N6A-C6A-N1A	-6.10	106.67	118.77
2	C	601	FAD	N6A-C6A-N1A	-6.02	106.83	118.77
2	C	601	FAD	C8M-C8-C7	-3.02	114.37	120.72
2	D	601	FAD	C8M-C8-C7	-3.01	114.41	120.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	601	FAD	C4X-C4-N3	-2.82	119.47	123.48
2	B	601	FAD	C4X-C4-N3	-2.81	119.48	123.48
2	B	601	FAD	C8M-C8-C7	-2.79	114.87	120.72
2	A	601	FAD	C4X-C4-N3	-2.75	119.56	123.48
2	C	601	FAD	C4X-C4-N3	-2.70	119.64	123.48
2	A	601	FAD	C8M-C8-C7	-2.62	115.22	120.72
2	C	601	FAD	C4'-C3'-C2'	-2.24	108.58	113.41
2	B	601	FAD	C4'-C3'-C2'	-2.15	108.77	113.41
2	B	601	FAD	C4X-N5-C5X	3.07	120.01	116.76
2	C	601	FAD	C4X-N5-C5X	3.08	120.01	116.76
2	D	601	FAD	C4X-N5-C5X	3.16	120.09	116.76
2	A	601	FAD	C4X-N5-C5X	3.18	120.12	116.76
2	D	601	FAD	C5X-C9A-N10	3.41	120.19	117.66
2	B	601	FAD	C5X-C9A-N10	3.44	120.21	117.66
2	A	601	FAD	C5X-C9A-N10	3.53	120.28	117.66
2	C	601	FAD	C5X-C9A-N10	3.67	120.38	117.66
2	D	601	FAD	C4-N3-C2	5.34	119.83	115.16
2	C	601	FAD	C4-N3-C2	5.35	119.84	115.16
2	B	601	FAD	C4-N3-C2	5.46	119.93	115.16
2	A	601	FAD	C4-N3-C2	5.49	119.96	115.16
2	C	601	FAD	C5A-C6A-N6A	6.28	133.28	120.47
2	A	601	FAD	C5A-C6A-N6A	6.33	133.38	120.47
2	D	601	FAD	C5A-C6A-N6A	6.40	133.51	120.47
2	B	601	FAD	C5A-C6A-N6A	6.51	133.74	120.47
2	C	601	FAD	C7M-C7-C8	7.24	135.91	120.72
2	A	601	FAD	C7M-C7-C8	7.39	136.23	120.72
2	D	601	FAD	C7M-C7-C8	7.40	136.25	120.72
2	B	601	FAD	C7M-C7-C8	7.58	136.63	120.72

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	FAD	1	0
2	B	601	FAD	2	0
2	D	601	FAD	2	0



## 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	394/405 (97%)	0.33	23 (5%)	24 30	29, 51, 83, 114	0
1	B	394/405 (97%)	0.30	15 (3%)	41 48	29, 47, 79, 97	0
1	C	394/405 (97%)	0.43	30 (7%)	15 19	33, 50, 84, 112	0
1	D	394/405 (97%)	0.45	28 (7%)	17 22	34, 57, 92, 113	0
All	All	1576/1620 (97%)	0.38	96 (6%)	22 28	29, 51, 86, 114	0

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	374	TRP	6.2
1	A	390	LEU	5.4
1	D	336	GLY	5.3
1	A	374	TRP	5.0
1	C	394	LEU	4.7
1	D	332	ALA	4.6
1	C	365	ARG	4.6
1	D	205	PRO	4.4
1	B	362	VAL	4.4
1	A	368	TYR	4.4
1	D	368	TYR	4.3
1	D	32	ASP	4.3
1	D	98	LEU	4.2
1	B	368	TYR	4.2
1	C	356	ASN	4.2
1	D	394	LEU	4.2
1	B	370	HIS	3.9
1	D	203	VAL	3.8
1	A	203	VAL	3.6
1	A	363	GLY	3.6
1	B	363	GLY	3.5

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Mol	Chain	Res	Type	RSRZ
1	C	26	LEU	3.4
1	D	346	ILE	3.4
1	A	244	VAL	3.2
1	C	368	TYR	3.1
1	C	355	ARG	3.1
1	C	221	GLY	3.0
1	C	215	ALA	3.0
1	A	371	ALA	3.0
1	A	394	LEU	2.9
1	D	244	VAL	2.9
1	C	347	ARG	2.8
1	C	390	LEU	2.8
1	A	245	ASP	2.8
1	B	204	LEU	2.8
1	D	246	GLY	2.7
1	D	290	GLY	2.7
1	C	332	ALA	2.7
1	D	31	ALA	2.7
1	A	206	GLY	2.7
1	B	207	PHE	2.7
1	A	356	ASN	2.6
1	A	249	GLU	2.6
1	D	331	ALA	2.6
1	C	378	LEU	2.6
1	C	346	ILE	2.6
1	D	27	ASN	2.6
1	D	245	ASP	2.5
1	D	359	ILE	2.5
1	B	211	LEU	2.5
1	A	204	LEU	2.5
1	A	361	ILE	2.5
1	D	28	TYR	2.5
1	A	375	LEU	2.4
1	A	362	VAL	2.4
1	C	292	GLU	2.4
1	B	120	ARG	2.4
1	C	207	PHE	2.3
1	C	358	ALA	2.3
1	C	396	LYS	2.3
1	C	203	VAL	2.3
1	B	3	LYS	2.3
1	D	337	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	370	HIS	2.3
1	B	208	ASP	2.2
1	C	30	GLU	2.2
1	D	206	GLY	2.2
1	C	348	GLY	2.2
1	D	249	GLU	2.2
1	B	359	ILE	2.2
1	C	334	ILE	2.2
1	C	393	VAL	2.2
1	D	277	ARG	2.2
1	D	221	GLY	2.2
1	A	335	ARG	2.2
1	D	335	ARG	2.2
1	B	203	VAL	2.2
1	B	214	TYR	2.2
1	A	364	GLY	2.2
1	A	367	VAL	2.1
1	D	374	TRP	2.1
1	D	338	SER	2.1
1	A	247	GLU	2.1
1	A	3	LYS	2.1
1	D	247	GLU	2.1
1	C	217	ASP	2.1
1	C	357	ASP	2.1
1	B	366	LYS	2.1
1	C	204	LEU	2.1
1	C	27	ASN	2.1
1	A	382	ARG	2.1
1	D	355	ARG	2.1
1	B	217	ASP	2.0
1	C	211	LEU	2.0
1	C	369	GLY	2.0
1	A	238	GLU	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	FAD	B	601	53/53	0.94	0.15	0.86	31,36,47,59	0
2	FAD	A	601	53/53	0.97	0.15	0.51	31,38,43,76	0
2	FAD	D	601	53/53	0.95	0.13	0.39	30,45,51,58	0
2	FAD	C	601	53/53	0.96	0.14	-0.00	32,39,49,57	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.