



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 03:04 PM GMT

PDB ID : 4FGK
Title : Oxidized quinone reductase 2 in complex with chloroquine
Authors : Leung, K.K.; Shilton, B.H.
Deposited on : 2012-06-04
Resolution : 1.40 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

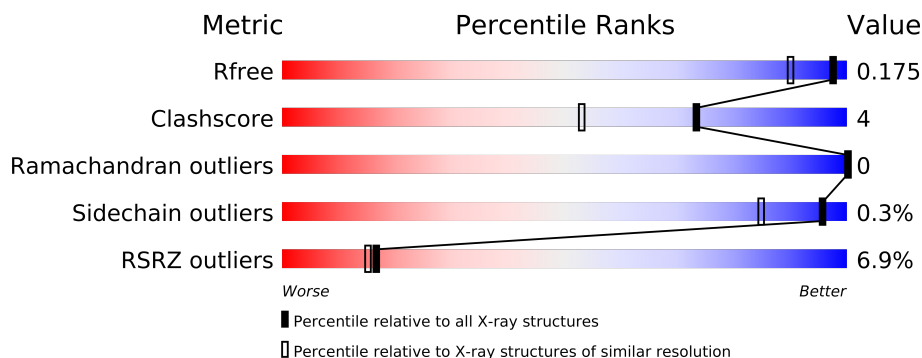
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.40 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}	66092	1097 (1.42-1.38)
Clashscore	79885	1246 (1.42-1.38)
Ramachandran outliers	78287	1206 (1.42-1.38)
Sidechain outliers	78261	1205 (1.42-1.38)
RSRZ outliers	66119	1097 (1.42-1.38)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	233	
1	B	233	

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 8091 atoms, of which 3622 are hydrogens and 0 are deuterium.

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.

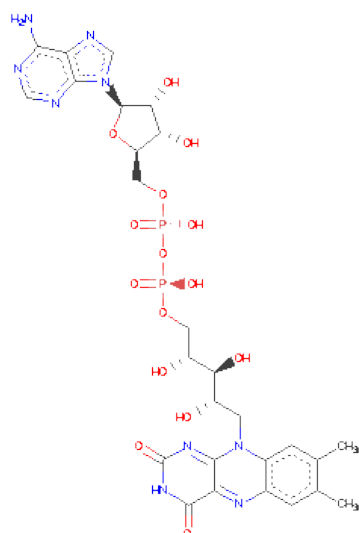
- Molecule 1 is a protein called Ribosyldihydronicotinamidedehydrogenase [quinone].

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	230	Total	C	H	N	O	S	0	14	0
			3677	1239	1750	320	359	9			
1	B	230	Total	C	H	N	O	S	0	5	0
			3641	1194	1784	310	344	9			

There are 6 discrepancies between the modelled and reference sequences:

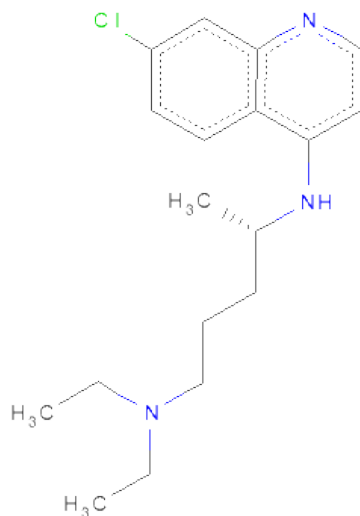
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP P16083
A	-1	ALA	-	EXPRESSION TAG	UNP P16083
A	46	PHE	LEU	SEE REMARK 999	UNP P16083
B	-2	GLY	-	EXPRESSION TAG	UNP P16083
B	-1	ALA	-	EXPRESSION TAG	UNP P16083
B	46	PHE	LEU	SEE REMARK 999	UNP P16083

- 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
2	A	1	Total	C	H	N	O	P	
			84	27	31	9	15	2	0
2	B	1	Total	C	H	N	O	P	
			84	27	31	9	15	2	0

- Molecule 3 is (4S)-N 4 -(7-CHLOROQUINOLIN-4-YL)-N 1 ,N 1 -DIETHYLPENTANE-1,4-DIAMINE (three-letter code: OTX) (formula: C₁₈H₂₆ClN₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N			
			44	36	2	6		0	1
3	A	1	Total	C	Cl	H	N		
			48	18	1	26	3	0	0

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Zn		
			1	1	0	0
4	A	1	Total	Zn		
			1	1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	221	Total	O		
			221	221	0	0

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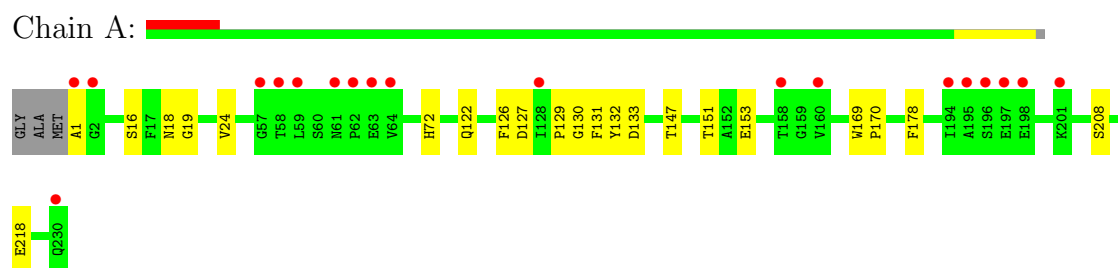
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	290	Total 290	O 290	0	0

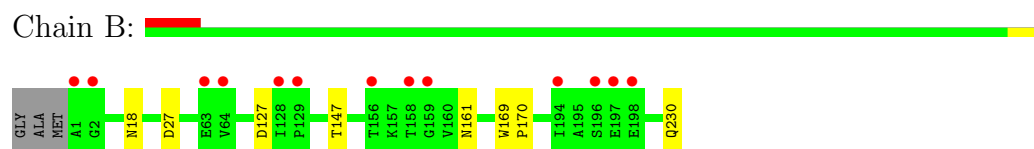
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Ribosyldihydronicotinamidedehydrogenase [quinone]



- Molecule 1: Ribosyldihydronicotinamidedehydrogenase [quinone]



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	56.37Å 83.11Å 106.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.67 – 1.40 32.67 – 1.40	Depositor EDS
% Data completeness (in resolution range)	99.5 (32.67-1.40) 99.5 (32.67-1.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.08 (at 1.40Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.133 , 0.174 0.133 , 0.175	Depositor DCC
R_{free} test set	1998 reflections (2.03%)	DCC
Wilson B-factor (Å ²)	13.0	Xtriage
Anisotropy	0.586	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.46 , 43.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 98466 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8091	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.03% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, FAD, OTX

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	0/1980	0.59	0/2684
1	B	0.36	0/1907	0.57	0/2588
All	All	0.37	0/3887	0.58	0/5272

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1927	1750	0	25	1
1	B	1857	1784	0	5	0
2	A	53	31	0	0	0
2	B	53	31	0	0	0
3	A	66	26	0	3	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	221	0	0	5	1
5	B	290	0	0	4	0
All	All	4469	3622	0	32	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (32) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:127[B]:ASP:OD2	5:A:621:HOH:O	1.95	0.85
3:A:304:OTX:CL	5:B:689:HOH:O	2.35	0.80
1:A:72:HIS:CG	1:A:122[B]:GLN:HG3	2.23	0.74
1:A:72:HIS:NE2	5:A:621:HOH:O	2.22	0.72
1:A:131[B]:PHE:CB	1:A:178:PHE:CZ	2.80	0.65
1:B:27:ASP:OD1	5:B:554:HOH:O	2.15	0.63
1:A:1[A]:ALA:N	5:A:619:HOH:O	2.32	0.62
1:A:169:TRP:HB3	1:A:170:PRO:HD3	1.84	0.59
1:A:127[B]:ASP:N	1:A:130[B]:GLY:O	2.25	0.59
1:A:122[B]:GLN:HB3	1:A:126:PHE:CZ	2.40	0.56
1:A:72:HIS:HB2	1:A:122[B]:GLN:CG	2.36	0.55
1:A:24:VAL:HG12	1:A:208:SER:HB2	1.89	0.55
1:A:127[B]:ASP:HB3	1:A:129[B]:PRO:O	2.08	0.54
1:A:131[B]:PHE:CD2	1:A:178:PHE:CE1	2.96	0.53
1:A:131[B]:PHE:CG	1:A:178:PHE:CZ	2.96	0.53
3:A:304:OTX:H61	5:B:689:HOH:O	2.09	0.53
1:A:16:SER:HG	1:A:19:GLY:H	1.57	0.52
1:B:18:ASN:HD21	1:B:147:THR:HG1	1.58	0.51
1:A:132[B]:TYR:CD1	1:A:178:PHE:HA	2.46	0.51
1:A:18:ASN:HD21	1:A:147:THR:HG1	1.57	0.51
1:A:132[B]:TYR:OH	1:B:161[B]:ASN:OD1	2.30	0.49
1:A:122[B]:GLN:HG2	5:A:473:HOH:O	2.12	0.48
1:A:151:THR:HG1	1:A:153:GLU:HG2	1.80	0.47
1:B:169:TRP:HB3	1:B:170:PRO:HD3	1.96	0.46
1:A:24:VAL:CG1	1:A:208:SER:HB2	2.47	0.44
1:A:72:HIS:CB	1:A:122[B]:GLN:HG3	2.48	0.43
1:A:72:HIS:HB2	1:A:122[B]:GLN:HG3	2.00	0.43
1:A:133[B]:ASP:OD1	5:A:584:HOH:O	2.21	0.42
1:A:132[B]:TYR:CZ	1:A:178:PHE:HD1	2.38	0.41
3:A:304:OTX:H122	3:A:304:OTX:H162	1.86	0.41
1:A:132[B]:TYR:CE1	1:A:178:PHE:HD1	2.39	0.41
1:B:127:ASP:OD1	5:B:451:HOH:O	2.22	0.41

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	Distance(Å)	Clash(Å)
1:A:218:GLU:OE1	5:A:621:HOH:O[4_445]	2.01	0.19

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	241/233 (103%)	232 (96%)	9 (4%)	0	100	100
1	B	233/233 (100%)	224 (96%)	9 (4%)	0	100	100
All	All	474/466 (102%)	456 (96%)	18 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	204/195 (105%)	204 (100%)	0	100	100
1	B	198/195 (102%)	197 (100%)	1 (0%)	94	80
All	All	402/390 (103%)	401 (100%)	1 (0%)	96	85

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	230	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	ASN
1	A	77	GLN
1	A	172	GLN
1	A	212	GLN
1	B	138	GLN

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Mol	Chain	Res	Type
1	B	172	GLN
1	B	212	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 for Mogul analysis.

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	301	-	58,58,58	1.64	13 (22%)	85,89,89	1.85	11 (12%)
3	OTX	A	302[A]	-	23,23,23	0.96	2 (8%)	30,30,30	1.49	4 (13%)
3	OTX	A	302[B]	-	23,23,23	1.07	2 (8%)	30,30,30	1.32	4 (13%)
3	OTX	A	304	-	23,23,23	1.08	2 (8%)	30,30,30	1.37	4 (13%)
2	FAD	B	302	-	58,58,58	1.64	13 (22%)	85,89,89	1.73	12 (14%)

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	301	-	-	0/34/50/50	0/1/6/6
3	OTX	A	302[A]	-	-	0/14/14/14	0/0/2/2
3	OTX	A	302[B]	-	-	0/14/14/14	0/0/2/2
3	OTX	A	304	-	-	0/14/14/14	0/0/2/2
2	FAD	B	302	-	-	0/34/50/50	0/1/6/6

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	302	FAD	C10-N1	4.87	1.44	1.35
2	A	301	FAD	C10-N1	4.66	1.44	1.35
2	B	302	FAD	C4-C4X	4.07	1.47	1.41
2	A	301	FAD	C4-C4X	4.00	1.47	1.41
2	A	301	FAD	C2B-C3B	-3.66	1.43	1.53
2	B	302	FAD	C2B-C3B	-3.48	1.43	1.53
3	A	304	OTX	C3-N2	3.45	1.46	1.37
2	B	302	FAD	C2-N3	3.19	1.43	1.37
2	A	301	FAD	C1'-N10	3.17	1.51	1.48
3	A	302[B]	OTX	C3-N2	3.11	1.46	1.37
3	A	302[A]	OTX	C3-N2	2.98	1.45	1.37
2	B	302	FAD	C2-N1	2.89	1.43	1.35
2	A	301	FAD	C2-N1	2.87	1.43	1.35
2	A	301	FAD	C2B-C1B	-2.84	1.49	1.53
2	A	301	FAD	O4'-C4'	-2.78	1.37	1.43
3	A	302[B]	OTX	C9-N1	-2.76	1.32	1.37
2	B	302	FAD	O4'-C4'	-2.69	1.37	1.43
2	B	302	FAD	C4-N3	2.63	1.41	1.37
2	A	301	FAD	C2-N3	2.58	1.42	1.37
2	A	301	FAD	C1'-C2'	2.56	1.54	1.51
2	A	301	FAD	O3B-C3B	-2.55	1.36	1.43
3	A	304	OTX	C9-N1	-2.53	1.33	1.37
3	A	302[A]	OTX	C9-N1	-2.47	1.33	1.37
2	B	302	FAD	C1'-C2'	2.45	1.53	1.51
2	B	302	FAD	C1'-N10	2.42	1.50	1.48
2	B	302	FAD	O3B-C3B	-2.42	1.37	1.43
2	A	301	FAD	C6A-N6A	2.37	1.42	1.35
2	B	302	FAD	C9A-C5X	-2.35	1.37	1.42
2	B	302	FAD	C6A-N6A	2.27	1.42	1.35
2	A	301	FAD	C9A-C5X	-2.25	1.37	1.42
2	B	302	FAD	C4A-N9A	-2.14	1.34	1.37
2	A	301	FAD	C4A-N9A	-2.12	1.34	1.37

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	FAD	N3A-C2A-N1A	-9.71	120.59	128.71
2	B	302	FAD	N3A-C2A-N1A	-8.32	121.75	128.71
2	A	301	FAD	O4B-C1B-N9A	5.21	113.29	108.44
2	B	302	FAD	C2-N1-C10	4.91	119.93	114.98
2	A	301	FAD	C2-N1-C10	4.84	119.85	114.98
2	B	302	FAD	N3A-C4A-N9A	4.17	132.96	125.43
2	B	302	FAD	C4B-O4B-C1B	-4.11	105.29	109.75
3	A	302[A]	0TX	C8-C9-N1	4.07	121.56	117.86
3	A	302[A]	0TX	C1-N1-C9	3.94	122.49	116.94
2	A	301	FAD	C4X-C10-N10	-3.79	118.62	120.51
2	A	301	FAD	C4B-O4B-C1B	-3.63	105.80	109.75
2	A	301	FAD	N3A-C4A-N9A	3.53	131.81	125.43
2	A	301	FAD	C1B-N9A-C4A	-3.49	120.60	126.64
3	A	304	0TX	C8-C9-N1	3.29	120.85	117.86
2	B	302	FAD	C4X-C10-N10	-3.28	118.87	120.51
2	B	302	FAD	C5X-C9A-N10	3.22	119.98	116.80
2	A	301	FAD	C8A-N9A-C4A	3.17	109.31	106.90
3	A	304	0TX	C1-N1-C9	3.12	121.34	116.94
2	A	301	FAD	C5X-C9A-N10	3.03	119.79	116.80
2	B	302	FAD	C8A-N9A-C4A	2.98	109.18	106.90
3	A	302[B]	0TX	C1-N1-C9	2.98	121.14	116.94
2	B	302	FAD	C1B-N9A-C4A	-2.93	121.57	126.64
3	A	304	0TX	C3-C4-C9	2.89	119.48	118.00
2	B	302	FAD	C4X-N5-C5X	2.87	119.92	116.69
3	A	302[A]	0TX	C4-C9-N1	-2.84	119.94	122.91
3	A	302[B]	0TX	C8-C9-N1	2.81	120.42	117.86
3	A	304	0TX	C4-C9-N1	-2.69	120.10	122.91
3	A	302[A]	0TX	C11-C12-C13	-2.43	106.47	113.71
2	B	302	FAD	C5A-C4A-N3A	-2.35	120.59	125.70
2	B	302	FAD	C1'-N10-C9A	2.30	121.11	118.87
3	A	302[B]	0TX	C4-C9-N1	-2.24	120.56	122.91
2	A	301	FAD	C4A-C5A-N7A	-2.19	107.64	109.52
2	A	301	FAD	N7A-C8A-N9A	-2.14	108.30	114.36
3	A	302[B]	0TX	C12-C11-C10	-2.06	108.88	113.93
2	B	302	FAD	N7A-C8A-N9A	-2.00	108.70	114.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	230/233 (98%)	0.16	19 (8%)	11 11	6, 19, 39, 51	0
1	B	230/233 (98%)	-0.20	13 (5%)	23 22	7, 13, 34, 49	0
All	All	460/466 (98%)	-0.02	32 (6%)	17 15	6, 16, 36, 51	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	64	VAL	6.7
1	A	198	GLU	4.6
1	B	1	ALA	4.5
1	A	194	ILE	4.3
1	B	128	ILE	4.0
1	B	158[A]	THR	3.6
1	A	197	GLU	3.6
1	A	63	GLU	3.6
1	A	128[A]	ILE	3.5
1	B	194	ILE	3.3
1	B	129	PRO	3.2
1	B	2	GLY	3.1
1	B	198	GLU	3.1
1	B	159[A]	GLY	3.1
1	A	62	PRO	3.1
1	B	197	GLU	3.0
1	A	158	THR	2.9
1	A	196	SER	2.9
1	A	230	GLN	2.8
1	A	201	LYS	2.7
1	A	57	GLY	2.7
1	A	58	THR	2.5
1	B	64	VAL	2.4
1	A	1[A]	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	156	THR	2.3
1	A	59	LEU	2.3
1	A	160	VAL	2.2
1	A	195	ALA	2.2
1	A	61	ASN	2.1
1	B	196	SER	2.1
1	B	63	GLU	2.1
1	A	2[A]	GLY	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 ⓘ

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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	0TX	A	302[A]	22/22	0.19	1.92	8,14,25,27	22
3	0TX	A	304	22/22	0.19	1.65	26,47,57,57	26
3	0TX	A	302[B]	22/22	0.19	1.35	9,22,33,34	22
2	FAD	A	301	53/53	0.08	-0.13	8,15,31,38	0
2	FAD	B	302	53/53	0.07	-0.27	10,15,31,39	0
4	ZN	A	303	1/1	0.04	-1.07	22,22,22,22	0
4	ZN	B	301	1/1	0.02	-2.22	11,11,11,11	0

6.5 Other polymers ⓘ

There are no such residues in this entry.