



wwPDB X-ray Structure Validation Summary Report i

Feb 27, 2014 – 07:51 AM GMT

PDB ID : 2RC5
Title : Refined structure of FNR from *Leptospira interrogans*
Authors : Nascimento, A.S.; Catalano-Dupuy, D.L.; Polikarpov, I.; Ceccarelli, E.A.
Deposited on : 2007-09-19
Resolution : 2.43 Å(reported)

This is a wwPDB validation summary 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>

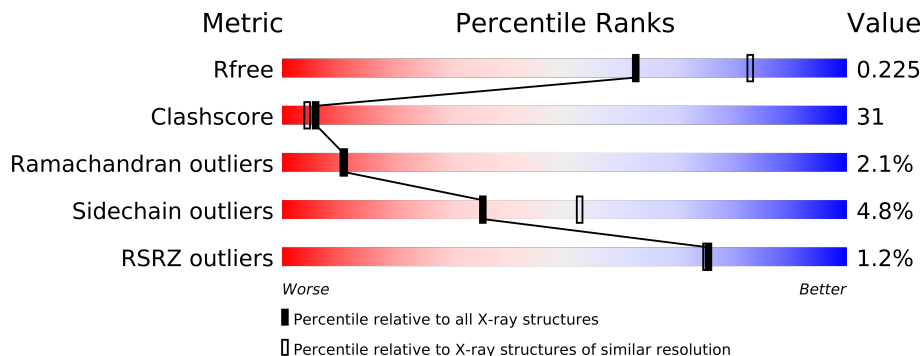
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.43 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	2989 (2.48-2.40)
Clashscore	79885	3698 (2.48-2.40)
Ramachandran outliers	78287	3639 (2.48-2.40)
Sidechain outliers	78261	3640 (2.48-2.40)
RSRZ outliers	66119	2993 (2.48-2.40)

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. 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	314	
1	B	314	
1	C	314	
1	D	314	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	SO4	D	315	-	X

2 Entry composition

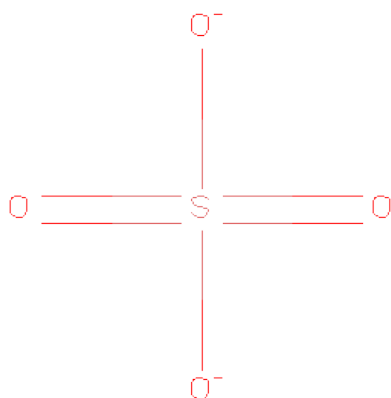
There are 5 unique types of molecules in this entry. The entry contains 10963 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Ferredoxin-NADP reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	306	Total	C	N	O	S	0	0	0
			2374	1516	397	449	12			
1	B	301	Total	C	N	O	S	0	0	0
			2348	1496	393	447	12			
1	C	304	Total	C	N	O	S	0	0	0
			2373	1513	397	451	12			
1	D	306	Total	C	N	O	S	0	0	0
			2382	1520	399	451	12			

- Molecule 2 is SULFATE ION (three-letter code: 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		

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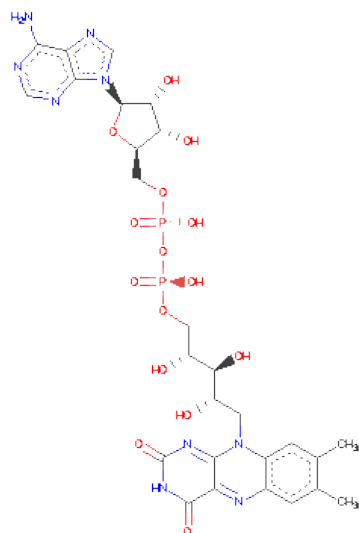
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

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

- Molecule 5 is water.

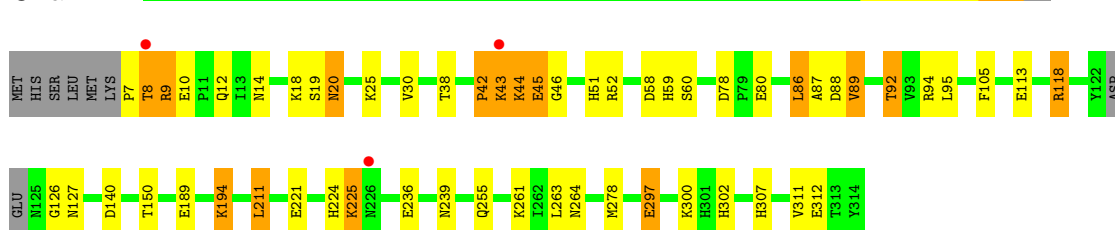
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	297	Total	O	0	0
			297	297		
5	B	307	Total	O	0	0
			307	307		
5	C	329	Total	O	0	0
			329	329		
5	D	305	Total	O	0	0
			305	305		

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.

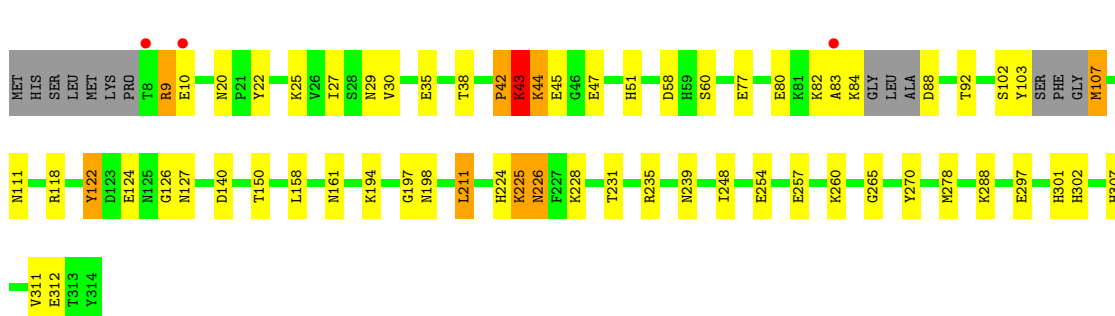
- Molecule 1: Ferredoxin-NADP reductase

Chain A:



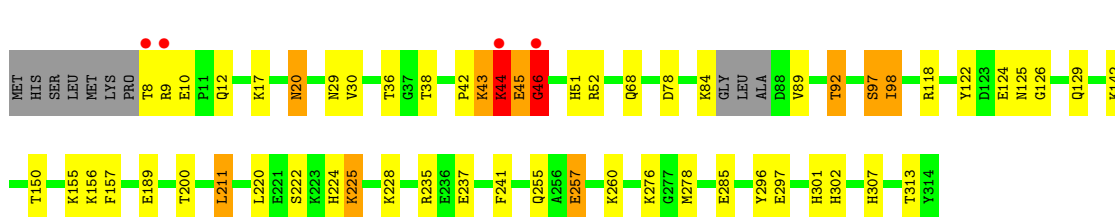
- Molecule 1: Ferredoxin-NADP reductase

Chain B:



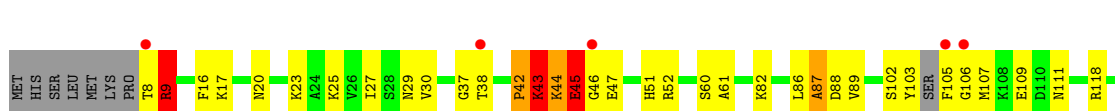
- Molecule 1: Ferredoxin-NADP reductase

Chain C:



- Molecule 1: Ferredoxin-NADP reductase

Chain D:



Y122	D123	E124	N125	G126	N127	I128	D140	T150	N161	T162	D163	F164	S165	E185	E189	H190	G197	L211	E221	H224	K225	N226	K238	R253	K261	R269	N278	E279	K280	K288	N292	Y296	E297	H301	H307	Y311	Y314
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	65.11Å 111.83Å 89.91Å 90.00° 92.76° 90.00°	Depositor
Resolution (Å)	44.90 – 2.43 44.90 – 2.43	Depositor EDS
% Data completeness (in resolution range)	98.5 (44.90-2.43) 98.5 (44.90-2.43)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.93 (at 2.42Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.225 , 0.255 0.197 , 0.225	Depositor DCC
R_{free} test set	2412 reflections (5.06%)	DCC
Wilson B-factor (Å ²)	24.5	Xtriage
Anisotropy	0.266	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 19.0	EDS
Estimated twinning fraction	0.065 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 47690 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10963	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.78% 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, SO4, 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	0.53	0/2425	0.64	1/3265 (0.0%)
1	B	0.51	0/2396	0.77	3/3224 (0.1%)
1	C	0.56	1/2423 (0.0%)	0.66	1/3260 (0.0%)
1	D	0.54	0/2432	0.65	0/3272
All	All	0.54	1/9676 (0.0%)	0.68	5/13021 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	2
1	C	0	3
1	D	0	2
All	All	0	10

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	97	SER	C-O	-5.43	1.13	1.23

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	43	LYS	N-CA-CB	-23.21	68.83	110.60
1	C	211	LEU	CA-CB-CG	6.93	131.24	115.30
1	A	211	LEU	CA-CB-CG	5.83	128.70	115.30
1	B	211	LEU	CA-CB-CG	5.33	127.55	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	107	MET	C-N-CA	5.28	134.89	121.70

There are no chirality outliers.

5 of 10 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	42	PRO	Peptide
1	A	43	LYS	Peptide
1	A	45	GLU	Peptide
1	B	42	PRO	Peptide
1	B	43	LYS	Peptide

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	2374	0	0	68	1
1	B	2348	0	0	66	4
1	C	2373	0	5	83	1
1	D	2382	0	0	80	3
2	A	10	0	0	9	4
2	B	5	0	0	4	1
2	C	15	0	0	17	3
2	D	5	0	0	5	1
3	A	1	0	0	0	0
4	A	53	0	0	1	0
4	B	53	0	0	0	0
4	C	53	0	0	3	0
4	D	53	0	0	0	0
5	A	297	0	0	45	0
5	B	307	0	0	49	1
5	C	329	0	0	54	1
5	D	305	0	0	59	2
All	All	10963	0	5	303	11

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 31.

The worst 5 of 303 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:17:LYS:CD	5:C:1011:HOH:O	1.84	1.25
1:B:82:LYS:CE	5:B:556:HOH:O	1.83	1.23
1:D:124:GLU:CB	5:D:584:HOH:O	1.84	1.22
1:C:42:PRO:O	1:C:45:GLU:HB2	1.09	1.22
1:C:224:HIS:NE2	2:C:317:SO4:O3	1.72	1.21

The worst 5 of 11 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:297:GLU:OE2	2:B:315:SO4:O3[2_555]	1.49	0.71
1:B:297:GLU:OE1	2:A:316:SO4:O1[2_545]	1.80	0.40
1:B:301:HIS:ND1	2:A:316:SO4:O2[2_545]	1.86	0.34
2:C:317:SO4:O4	5:D:584:HOH:O[1_655]	1.87	0.33
1:D:297:GLU:OE1	2:C:316:SO4:O3[2_454]	1.97	0.23

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	302/314 (96%)	281 (93%)	14 (5%)	7 (2%)	10	9
1	B	295/314 (94%)	284 (96%)	9 (3%)	2 (1%)	30	41
1	C	300/314 (96%)	284 (95%)	8 (3%)	8 (3%)	8	6
1	D	302/314 (96%)	280 (93%)	14 (5%)	8 (3%)	8	7
All	All	1199/1256 (96%)	1129 (94%)	45 (4%)	25 (2%)	11	11

5 of 25 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	8	THR
1	C	46	GLY

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Mol	Chain	Res	Type
1	C	89	VAL
1	D	124	GLU
1	A	9	ARG

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	254/266 (96%)	241 (95%)	13 (5%)	33	49
1	B	253/266 (95%)	240 (95%)	13 (5%)	33	49
1	C	256/266 (96%)	247 (96%)	9 (4%)	48	68
1	D	255/266 (96%)	241 (94%)	14 (6%)	30	45
All	All	1018/1064 (96%)	969 (95%)	49 (5%)	35	53

5 of 49 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	226	ASN
1	C	30	VAL
1	D	211	LEU
1	B	278	MET
1	C	92	THR

Some 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 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 12 ligands modelled in this entry, 1 is monoatomic - leaving 11 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	SO4	A	315	-	4,4,4	1.16	0	6,6,6	0.44	0
2	SO4	A	316	-	4,4,4	1.48	1 (25%)	6,6,6	0.87	0
4	FAD	A	415	-	58,58,58	1.05	3 (5%)	85,89,89	1.74	11 (12%)
2	SO4	B	315	-	4,4,4	0.99	0	6,6,6	0.45	0
4	FAD	B	415	-	58,58,58	0.99	2 (3%)	85,89,89	1.89	11 (12%)
2	SO4	C	315	-	4,4,4	1.05	0	6,6,6	0.63	0
2	SO4	C	316	-	4,4,4	1.45	0	6,6,6	0.45	0
2	SO4	C	317	-	4,4,4	0.74	0	6,6,6	0.88	0
4	FAD	C	415	-	58,58,58	2.30	30 (51%)	85,89,89	1.91	16 (18%)
2	SO4	D	315	-	4,4,4	1.20	0	6,6,6	0.56	0
4	FAD	D	415	-	58,58,58	1.03	3 (5%)	85,89,89	1.88	11 (12%)

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	SO4	A	315	-	-	0/0/0/0	0/0/0/0
2	SO4	A	316	-	-	0/0/0/0	0/0/0/0
4	FAD	A	415	-	-	0/34/50/50	0/1/6/6
2	SO4	B	315	-	-	0/0/0/0	0/0/0/0
4	FAD	B	415	-	-	0/34/50/50	0/1/6/6
2	SO4	C	315	-	-	0/0/0/0	0/0/0/0
2	SO4	C	316	-	-	0/0/0/0	0/0/0/0
2	SO4	C	317	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	FAD	C	415	-	-	0/34/50/50	0/1/6/6
2	SO4	D	315	-	-	0/0/0/0	0/0/0/0
4	FAD	D	415	-	-	0/34/50/50	0/1/6/6

The worst 5 of 39 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	415	FAD	C6-C5X	-5.21	1.35	1.41
4	C	415	FAD	C4-C4X	-4.18	1.34	1.41
4	C	415	FAD	C4A-N9A	-3.60	1.32	1.37
4	C	415	FAD	O4-C4	-3.55	1.17	1.24
4	C	415	FAD	O4B-C4B	-3.45	1.36	1.45

The worst 5 of 49 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	415	FAD	N3A-C2A-N1A	-11.80	118.85	128.71
4	D	415	FAD	N3A-C2A-N1A	-10.59	119.86	128.71
4	A	415	FAD	N3A-C2A-N1A	-10.51	119.92	128.71
4	C	415	FAD	N3A-C2A-N1A	-8.58	121.54	128.71
4	D	415	FAD	O4B-C1B-N9A	6.65	114.62	108.44

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	306/314 (97%)	-0.19	3 (0%) 79 79	12, 24, 47, 61	1 (0%)
1	B	301/314 (95%)	-0.13	3 (0%) 79 79	12, 23, 51, 74	2 (0%)
1	C	304/314 (96%)	-0.23	4 (1%) 74 73	12, 22, 43, 75	5 (1%)
1	D	306/314 (97%)	-0.09	5 (1%) 68 67	13, 23, 56, 72	3 (0%)
All	All	1217/1256 (96%)	-0.16	15 (1%) 75 75	12, 23, 50, 75	11 (0%)

The worst 5 of 15 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	8	THR	8.2
1	D	105	PHE	8.2
1	A	8	THR	5.4
1	D	106	GLY	4.3
1	D	8	THR	4.2

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(\AA^2)	Q<0.9
2	SO4	D	315	5/5	0.20	2.29	22,23,23,23	0
2	SO4	C	315	5/5	0.20	1.51	20,20,20,20	0
4	FAD	C	415	53/53	0.17	1.20	44,45,46,46	0
4	FAD	B	415	53/53	0.16	0.89	27,32,33,34	0
4	FAD	D	415	53/53	0.16	0.38	24,25,26,26	0
2	SO4	A	315	5/5	0.15	0.18	21,21,21,22	0
2	SO4	C	316	5/5	0.19	0.06	19,19,19,19	0
2	SO4	A	316	5/5	0.14	0.01	18,18,25,25	0
4	FAD	A	415	53/53	0.14	-0.12	24,26,27,29	0
2	SO4	B	315	5/5	0.13	-1.20	24,24,25,25	0
2	SO4	C	317	5/5	0.12	-1.21	28,29,29,29	0
3	ZN	A	317	1/1	0.04	-11.31	43,43,43,43	0

6.5 Other polymers ⓘ

There are no such residues in this entry.