



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 3V74
Title : crystal structure of FBF-2 in complex with gld-1 FBEa13 RNA
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Deposited on : 2011-12-20
Resolution : 2.30 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

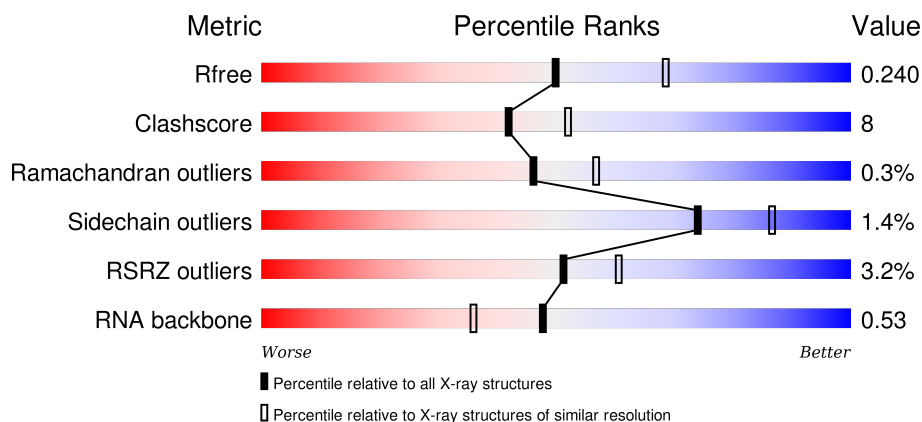
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)
RNA backbone	2183	1011 (2.84-1.76)

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. 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	413	 3% 79% 16% • 5%
2	B	13	 69% 8% 8% 15%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3689 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 Fem-3 mRNA-binding factor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	394	Total	C	N	O	S	0	0	0
			3139	1976	548	588	27			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	163	GLY	-	EXPRESSION TAG	UNP Q09312

- Molecule 2 is a RNA chain called RNA (5'-R(*UP*CP*AP*UP*GP*UP*GP*CP*CP*AP*UP*AP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	11	Total	C	N	O	P	0	0	0
			232	104	40	77	11			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	279	Total	O	0	0
			279	279		
3	B	39	Total	O	0	0
			39	39		

- Molecule 1: Fem-3 mRNA-binding factor 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	98.47Å 98.47Å 107.98Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	24.01 – 2.30 24.01 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (24.01-2.30) 99.8 (24.01-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.51 (at 2.31Å)	Xtriage
Refinement program	cns1.1 and phenix	Depositor
R, R_{free}	0.177 , 0.227 0.196 , 0.240	Depositor DCC
R_{free} test set	1361 reflections (5.16%)	DCC
Wilson B-factor (Å ²)	34.6	Xtriage
Anisotropy	0.056	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.1	EDS
Estimated twinning fraction	0.056 for h,-h-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 26479 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3689	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.77% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.45	0/3184	0.56	0/4285
2	B	0.83	0/258	1.41	2/399 (0.5%)
All	All	0.49	0/3442	0.68	2/4684 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	B	3	A	O4'-C1'-N9	8.95	115.36	108.20
2	B	3	A	C5'-C4'-O4'	6.71	117.16	109.10

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	3139	0	3175	56	0
2	B	232	0	119	0	0
3	A	279	0	0	10	0
3	B	39	0	0	0	0
All	All	3689	0	3294	56	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 8.

All (56) 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:411:ASN:HD22	1:A:414:ALA:H	1.13	0.94
1:A:225:GLU:OE2	1:A:230:ARG:NH1	2.17	0.78
1:A:547:HIS:HB3	3:A:844:HOH:O	1.84	0.77
1:A:174:ASP:HB3	1:A:176:ASN:H	1.55	0.71
1:A:422:VAL:O	1:A:431:ARG:HD3	1.91	0.70
1:A:519:GLN:HE21	1:A:521:LYS:H	1.37	0.69
1:A:251:ILE:HG12	1:A:269:LEU:HD21	1.76	0.67
1:A:354:ARG:HB2	1:A:404:ARG:HH21	1.60	0.67
1:A:354:ARG:HD3	1:A:404:ARG:NH2	2.10	0.66
1:A:266:GLN:HG2	1:A:301:LEU:HG	1.78	0.65
1:A:266:GLN:HE22	1:A:298:ASP:H	1.44	0.64
1:A:354:ARG:HD3	1:A:404:ARG:HH22	1.64	0.63
1:A:354:ARG:HB2	1:A:404:ARG:NH2	2.16	0.60
1:A:519:GLN:NE2	1:A:521:LYS:H	2.00	0.60
1:A:318:ILE:HG23	1:A:324:ALA:HB1	1.84	0.59
1:A:420:HIS:HD2	3:A:664:HOH:O	1.84	0.59
1:A:404:ARG:O	1:A:408:LEU:HD22	2.03	0.58
1:A:358:SER:HA	1:A:411:ASN:ND2	2.20	0.56
1:A:533:ASP:HB2	3:A:726:HOH:O	2.07	0.55
1:A:440:MET:O	1:A:441:ARG:HB2	2.08	0.53
1:A:411:ASN:ND2	1:A:414:ALA:H	1.95	0.52
1:A:411:ASN:ND2	1:A:413:TYR:H	2.08	0.52
1:A:181:SER:O	1:A:182:ARG:HB2	2.10	0.50
1:A:230:ARG:HH11	1:A:230:ARG:HG2	1.76	0.50
1:A:230:ARG:HH11	1:A:230:ARG:CG	2.25	0.49
1:A:358:SER:HA	1:A:411:ASN:HD21	1.77	0.49
1:A:529:ILE:HG13	3:A:622:HOH:O	2.13	0.48
1:A:527:HIS:HE1	3:A:627:HOH:O	1.96	0.47
1:A:420:HIS:HE1	3:A:624:HOH:O	1.98	0.46
1:A:388:ASN:HB3	1:A:392:ARG:NH1	2.32	0.45
1:A:351:GLU:OE2	1:A:354:ARG:NH2	2.50	0.44
1:A:354:ARG:HH11	1:A:354:ARG:HG2	1.82	0.44
1:A:179:MET:HE2	1:A:179:MET:HB3	1.61	0.44
1:A:232:ASP:HB2	3:A:757:HOH:O	2.17	0.44
1:A:519:GLN:HE21	1:A:520:THR:N	2.15	0.44
1:A:179:MET:HE3	1:A:206:PHE:CG	2.53	0.43
1:A:429:VAL:HG12	3:A:792:HOH:O	2.18	0.43
1:A:469:ALA:HA	1:A:534:TRP:CZ2	2.54	0.43
1:A:301:LEU:HA	1:A:301:LEU:HD13	1.90	0.43
1:A:547:HIS:HE1	3:A:755:HOH:O	2.01	0.43
1:A:494:MET:HE3	1:A:495:PHE:CE2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:GLN:HE21	1:A:330:LYS:NZ	2.17	0.42
1:A:283:ASP:HB3	1:A:286:ALA:HB3	2.01	0.42
1:A:303:CYS:O	1:A:307:GLN:HG2	2.20	0.42
1:A:514:VAL:HG21	1:A:535:LEU:HD23	2.02	0.42
1:A:354:ARG:CB	1:A:404:ARG:HH21	2.31	0.42
1:A:338:LYS:HD3	3:A:681:HOH:O	2.20	0.41
1:A:549:LEU:O	1:A:555:GLY:HA3	2.21	0.41
1:A:391:GLU:O	1:A:395:GLN:HG3	2.21	0.41
1:A:310:PRO:C	1:A:311:ARG:HD3	2.41	0.41
1:A:519:GLN:HE21	1:A:520:THR:H	1.69	0.41
1:A:179:MET:HE3	1:A:206:PHE:CD1	2.56	0.41
1:A:343:ILE:O	1:A:347:VAL:HG23	2.21	0.41
1:A:512:ASP:OD1	1:A:517:ARG:NH1	2.51	0.40
1:A:356:ILE:O	1:A:362:GLY:HA3	2.21	0.40
1:A:264:LYS:H	1:A:264:LYS:HG2	1.64	0.40

There are no symmetry-related clashes.

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	390/413 (94%)	385 (99%)	4 (1%)	1 (0%)	46 57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	488	ASP

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	353/370 (95%)	348 (99%)	5 (1%)	74 86

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	204	CYS
1	A	264	LYS
1	A	304	LYS
1	A	408	LEU
1	A	529	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	205	GLN
1	A	266	GLN
1	A	276	GLN
1	A	295	GLN
1	A	325	ASN
1	A	329	GLN
1	A	379	ASN
1	A	388	ASN
1	A	411	ASN
1	A	420	HIS
1	A	519	GLN
1	A	527	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	10/13 (76%)	2 (20%)	1 (10%)

All (2) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	3	A
2	B	12	A

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	3	A

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 [i](#)

There are no ligands in this entry.

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	394/413 (95%)	-0.28	13 (3%) 50 59	21, 35, 65, 103	0
2	B	11/13 (84%)	-0.59	0 100 100	37, 38, 53, 70	0
All	All	405/426 (95%)	-0.28	13 (3%) 51 60	21, 35, 65, 103	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	564	ASN	4.1
1	A	182	ARG	4.1
1	A	566	ARG	3.4
1	A	480	ILE	3.4
1	A	175	SER	3.3
1	A	503	VAL	3.2
1	A	479	TYR	2.9
1	A	176	ASN	2.9
1	A	263	THR	2.9
1	A	567	SER	2.7
1	A	180	ARG	2.3
1	A	365	VAL	2.1
1	A	366	VAL	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

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.