



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 6FQR / pdb_00006fqr
Title : Crystal structure of IMP3 RRM12 in complex with RNA (CCCC)
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Deposited on : 2018-02-14
Resolution : 2.10 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

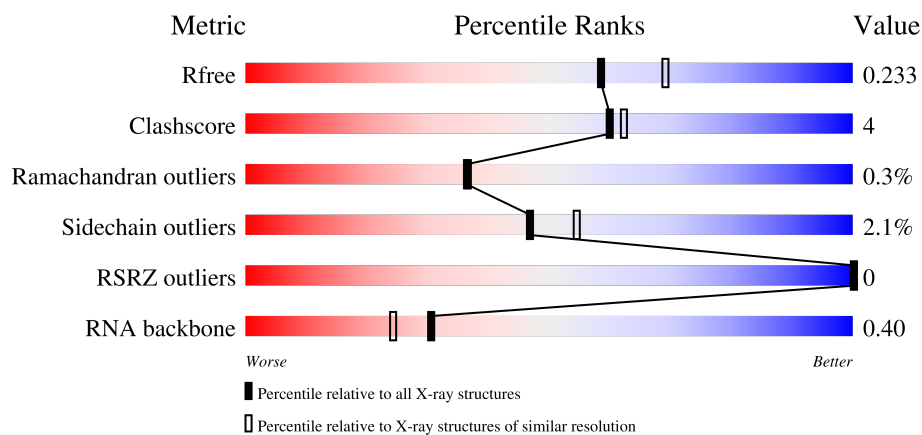
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)
RNA backbone	3983	1006 (2.42-1.78)

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 ≥ 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	170	 82% 11% 7%
1	B	170	 81% 13% 6%
2	C	4	 50% 25% 25%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2888 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 Insulin-like growth factor 2 mRNA-binding protein 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	158	Total	C	N	O	S	0	1	0
			1254	797	211	243	3			
1	B	160	Total	C	N	O	S	0	2	0
			1279	813	217	245	4			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP O00425
A	-1	GLY	-	expression tag	UNP O00425
A	0	SER	-	expression tag	UNP O00425
A	162	HIS	-	expression tag	UNP O00425
A	163	HIS	-	expression tag	UNP O00425
A	164	HIS	-	expression tag	UNP O00425
A	165	HIS	-	expression tag	UNP O00425
A	166	HIS	-	expression tag	UNP O00425
A	167	HIS	-	expression tag	UNP O00425
B	-2	GLY	-	expression tag	UNP O00425
B	-1	GLY	-	expression tag	UNP O00425
B	0	SER	-	expression tag	UNP O00425
B	162	HIS	-	expression tag	UNP O00425
B	163	HIS	-	expression tag	UNP O00425
B	164	HIS	-	expression tag	UNP O00425
B	165	HIS	-	expression tag	UNP O00425
B	166	HIS	-	expression tag	UNP O00425
B	167	HIS	-	expression tag	UNP O00425

- Molecule 2 is a RNA chain called RNA CCCC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	3	Total	C	N	O	P	0	0	0
			57	27	9	19	2			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	160	Total 160	O 160	0	0
3	B	129	Total 129	O 129	0	0
3	C	9	Total 9	O 9	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	30.06Å 41.19Å 72.27Å 92.14° 100.40° 108.65°	Depositor
Resolution (Å)	38.82 – 2.10 38.82 – 2.10	Depositor EDS
% Data completeness (in resolution range)	95.3 (38.82-2.10) 95.3 (38.82-2.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.167 , 0.230 0.169 , 0.233	Depositor DCC
R_{free} test set	895 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	27.2	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2888	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.43% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.30	0/1279	0.49	0/1736
1	B	0.28	0/1304	0.47	0/1767
2	C	0.30	0/62	0.54	0/94
All	All	0.29	0/2645	0.48	0/3597

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

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	1254	0	1255	11	0
1	B	1279	0	1289	11	1
2	C	57	0	35	1	0
3	A	160	0	0	3	1
3	B	129	0	0	4	1
3	C	9	0	0	0	0
All	All	2888	0	2579	22	3

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 4.

All (22) 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:110:CYS:SG	3:A:356:HOH:O	2.54	0.66
1:B:150:LYS:NZ	3:B:204:HOH:O	2.42	0.53
1:A:156:ASP:O	1:A:157:GLU:HG2	2.10	0.52
1:A:28:PRO:HD2	1:A:49:TRP:CE3	2.45	0.52
1:B:136:LEU:HD12	1:B:151:VAL:HG12	1.94	0.48
1:B:145:GLU:HG2	2:C:3:C:H2'	1.95	0.48
1:A:4:LEU:HD21	1:A:50:ALA:HB1	1.94	0.48
1:B:3:LYS:HB3	3:B:306:HOH:O	2.14	0.47
1:A:104:TYR:CE1	1:A:138:LYS:HG3	2.51	0.46
1:A:93:GLN:NE2	3:A:216:HOH:O	2.47	0.45
1:B:60:LYS:NZ	3:B:205:HOH:O	2.42	0.45
1:B:129:LYS:HE2	1:B:129:LYS:HB3	1.54	0.44
1:B:2:ASN:HA	1:B:50:ALA:HB2	1.99	0.43
1:B:89:PRO:HG3	1:B:147:PHE:HB2	2.00	0.43
1:B:28:PRO:HD2	1:B:49:TRP:CE3	2.54	0.43
1:B:82:LYS:NZ	3:B:214:HOH:O	2.50	0.43
1:B:117:SER:OG	1:B:119:THR:OG1	2.36	0.42
1:A:9:LEU:O	1:A:38:GLY:HA2	2.20	0.42
1:A:23:LYS:HB2	1:A:23:LYS:HE2	1.90	0.41
1:A:65:GLY:O	3:A:202:HOH:O	2.22	0.41
1:A:80:ILE:HG22	1:A:81:ARG:HG3	2.01	0.41
1:A:45:PRO:HG2	1:A:49:TRP:CD1	2.55	0.41

All (3) 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	Interatomic distance (Å)	Clash overlap (Å)
3:A:328:HOH:O	3:A:330:HOH:O[1_655]	2.14	0.06
1:B:90:PRO:O	1:B:133:ARG:NH2[1_455]	2.17	0.03
3:B:211:HOH:O	3:B:297:HOH:O[1_455]	2.17	0.03

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	157/170 (92%)	154 (98%)	3 (2%)	0	100	100
1	B	160/170 (94%)	156 (98%)	3 (2%)	1 (1%)	21	18
All	All	317/340 (93%)	310 (98%)	6 (2%)	1 (0%)	36	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	80	ILE

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	141/148 (95%)	140 (99%)	1 (1%)	76	83
1	B	143/148 (97%)	138 (96%)	5 (4%)	32	35
All	All	284/296 (96%)	278 (98%)	6 (2%)	47	54

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	LYS
1	B	11	GLU
1	B	27	ILE
1	B	37	THR
1	B	48	SER
1	B	150	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	93	GLN
1	B	134	GLN
1	B	143	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	2/4 (50%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

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 oligosaccharides 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	158/170 (92%)	-0.41	0 100 100	13, 25, 44, 61	1 (0%)
1	B	160/170 (94%)	-0.30	0 100 100	13, 28, 46, 56	2 (1%)
2	C	3/4 (75%)	-0.56	0 100 100	24, 24, 38, 44	0
All	All	321/344 (93%)	-0.35	0 100 100	13, 26, 45, 61	3 (0%)

There are no RSRZ outliers to report.

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.