



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 01:07 PM GMT

PDB ID : 3D2Q
Title : Crystal structure of MBNL1 tandem zinc finger 3 and 4 domain
Authors : Teplova, M.; Patel, D.J.
Deposited on : 2008-05-08
Resolution : 1.50 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>

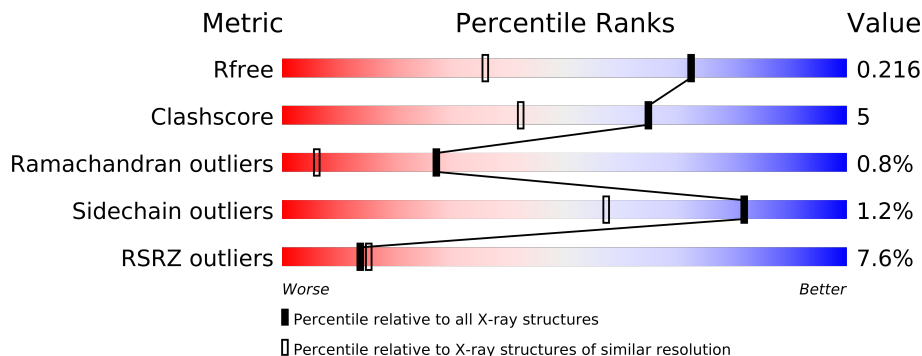
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.50 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	1513 (1.50-1.50)
Clashscore	79885	1768 (1.50-1.50)
Ramachandran outliers	78287	1720 (1.50-1.50)
Sidechain outliers	78261	1718 (1.50-1.50)
RSRZ outliers	66119	1514 (1.50-1.50)

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	70	
1	B	70	
1	C	70	
1	D	70	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2476 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 Muscleblind-like protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	66	Total	C	N	O	S	0	0	0
			532	319	103	102	8			
1	B	67	Total	C	N	O	S	0	0	0
			545	326	107	104	8			
1	C	66	Total	C	N	O	S	0	0	0
			532	319	103	102	8			
1	D	69	Total	C	N	O	S	0	0	0
			557	334	109	106	8			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	177	SER	-	EXPRESSION TAG	UNP Q9NR56
B	177	SER	-	EXPRESSION TAG	UNP Q9NR56
C	177	SER	-	EXPRESSION TAG	UNP Q9NR56
D	177	SER	-	EXPRESSION TAG	UNP Q9NR56

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Zn	0	0
			2	2		
2	A	2	Total	Zn	0	0
			2	2		
2	D	2	Total	Zn	0	0
			2	2		
2	C	2	Total	Zn	0	0
			2	2		

- Molecule 3 is water.

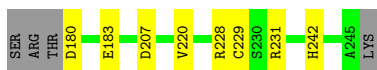
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	102	Total 102	O 102	0	0
3	B	78	Total 78	O 78	0	0
3	C	58	Total 58	O 58	0	0
3	D	64	Total 64	O 64	0	0

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: Muscblind-like protein 1

Chain A: 



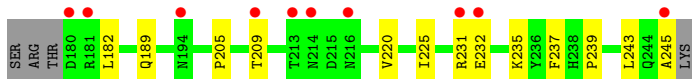
- Molecule 1: Muscblind-like protein 1

Chain B: 



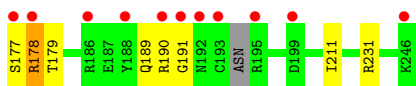
- Molecule 1: Muscblind-like protein 1

Chain C: 



- Molecule 1: Muscblind-like protein 1

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	29.17Å 80.46Å 55.36Å 90.00° 101.76° 90.00°	Depositor
Resolution (Å)	20.00 – 1.50 32.30 – 1.50	Depositor EDS
% Data completeness (in resolution range)	97.6 (20.00-1.50) 97.6 (32.30-1.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.93 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.197 , 0.217 0.196 , 0.216	Depositor DCC
R_{free} test set	1980 reflections (5.36%)	DCC
Wilson B-factor (Å ²)	11.7	Xtriage
Anisotropy	0.280	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 39.2	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 38922 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2476	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 47.84 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 9.3748e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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	0.47	0/542	0.59	0/729
1	B	0.44	0/555	0.68	0/746
1	C	0.43	0/542	0.63	0/729
1	D	0.43	0/566	0.61	0/758
All	All	0.44	0/2205	0.63	0/2962

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	532	0	490	7	0
1	B	545	0	505	4	1
1	C	532	0	490	9	1
1	D	557	0	521	8	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	102	0	0	1	0
3	B	78	0	0	3	0
3	C	58	0	0	2	0
3	D	64	0	0	3	0
All	All	2476	0	2006	22	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 5.

All (22) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:207:ASP:OD2	1:D:231:ARG:HD2	1.71	0.90
1:C:220:VAL:HG11	1:C:225:ILE:HD11	1.67	0.74
1:D:189:GLN:NE2	3:D:413:HOH:O	2.28	0.65
1:D:211:ILE:HG21	3:D:413:HOH:O	1.98	0.64
1:C:235:LYS:HG2	3:C:345:HOH:O	2.00	0.62
1:A:231:ARG:O	1:B:178:ARG:HD2	2.00	0.62
1:C:245:ALA:CB	1:D:231:ARG:HE	2.14	0.60
1:C:220:VAL:CG1	1:C:225:ILE:HD11	2.33	0.58
1:C:209:THR:HG21	3:C:320:HOH:O	2.05	0.56
1:B:178:ARG:NH1	3:B:389:HOH:O	2.41	0.52
1:A:229:CYS:SG	3:B:389:HOH:O	2.60	0.49
1:A:183:GLU:HG3	3:A:382:HOH:O	2.13	0.48
1:C:232:GLU:HA	1:D:177:SER:HB3	1.94	0.48
1:A:228:ARG:HD3	1:B:212:ASP:CG	2.34	0.47
1:C:205:PRO:HB2	1:C:209:THR:OG1	2.15	0.46
1:C:231:ARG:O	1:D:177:SER:HB3	2.16	0.46
1:B:226:LYS:HG3	1:B:226:LYS:O	2.15	0.46
1:D:211:ILE:HD13	3:D:413:HOH:O	2.16	0.45
1:A:180:ASP:HB3	1:A:220:VAL:HB	2.02	0.41
1:D:178:ARG:HG3	1:D:179:THR:HG23	2.03	0.41
1:A:242:HIS:HA	3:B:441:HOH:O	2.22	0.40
1:C:237:PHE:O	1:C:239:PRO:HD3	2.22	0.40

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:B:231:ARG:NH2	1:C:189:GLN:O[1_656]	2.16	0.04

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	64/70 (91%)	64 (100%)	0	0	100	100
1	B	65/70 (93%)	64 (98%)	1 (2%)	0	100	100
1	C	64/70 (91%)	63 (98%)	1 (2%)	0	100	100
1	D	65/70 (93%)	57 (88%)	6 (9%)	2 (3%)	7	0
All	All	258/280 (92%)	248 (96%)	8 (3%)	2 (1%)	27	6

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	178	ARG
1	D	191	GLY

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	59/63 (94%)	59 (100%)	0	100	100
1	B	61/63 (97%)	61 (100%)	0	100	100
1	C	59/63 (94%)	57 (97%)	2 (3%)	49	13
1	D	62/63 (98%)	61 (98%)	1 (2%)	75	44
All	All	241/252 (96%)	238 (99%)	3 (1%)	82	57

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	182	LEU

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Mol	Chain	Res	Type
1	C	243	LEU
1	D	190	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	244	GLN
1	D	214	ASN

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 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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	66/70 (94%)	-0.01	0	100 100	6, 10, 16, 19	0
1	B	67/70 (95%)	0.12	0	100 100	7, 12, 20, 22	0
1	C	66/70 (94%)	0.71	10 (15%)	3 3	9, 16, 27, 30	0
1	D	69/70 (98%)	0.82	11 (15%)	3 2	8, 14, 30, 37	0
All	All	268/280 (95%)	0.42	21 (7%)	14 14	6, 13, 27, 37	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	191	GLY	6.6
1	D	192	ASN	5.6
1	D	177	SER	5.6
1	D	190	ARG	4.8
1	D	199	ASP	3.8
1	D	195	ARG	3.8
1	C	180	ASP	3.4
1	D	186	ARG	3.4
1	D	246	LYS	3.1
1	D	178	ARG	3.1
1	C	181	ARG	3.0
1	C	216	ASN	3.0
1	D	193	CYS	2.8
1	C	194	ASN	2.7
1	C	214	ASN	2.7
1	C	232	GLU	2.7
1	D	188	TYR	2.5
1	C	209	THR	2.2
1	C	213	THR	2.1
1	C	245	ALA	2.1
1	C	231	ARG	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
2	ZN	D	303	1/1	0.07	-0.73	20,20,20,20	0
2	ZN	C	303	1/1	0.05	-1.32	16,16,16,16	0
2	ZN	D	304	1/1	0.03	-1.35	12,12,12,12	0
2	ZN	B	303	1/1	0.03	-2.88	15,15,15,15	0
2	ZN	A	303	1/1	0.04	-3.76	12,12,12,12	0
2	ZN	C	304	1/1	0.03	-4.41	15,15,15,15	0
2	ZN	A	304	1/1	0.02	-5.19	12,12,12,12	0
2	ZN	B	304	1/1	0.02	-9.65	12,12,12,12	0

6.5 Other polymers

There are no such residues in this entry.