



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 5MA8
Title : GFP-binding DARPin 3G124nc
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Deposited on : unknown
Resolution : 2.35 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20030345

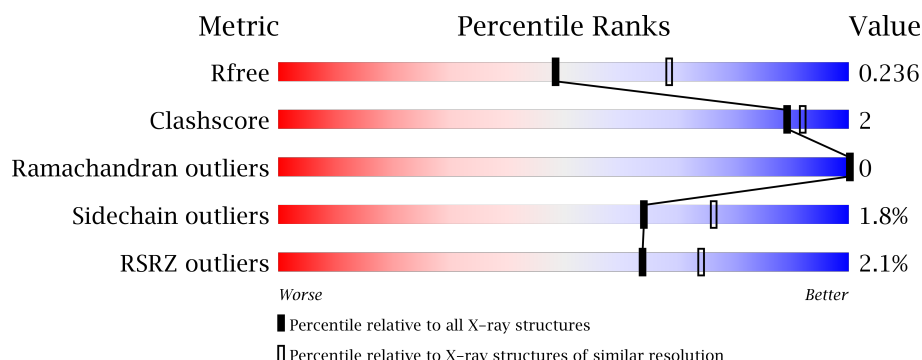
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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}	100719	1522 (2.38-2.34)
Clashscore	112137	1626 (2.38-2.34)
Ramachandran outliers	110173	1605 (2.38-2.34)
Sidechain outliers	110143	1606 (2.38-2.34)
RSRZ outliers	101464	1528 (2.38-2.34)

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	161	<div> <div>91%</div> <div>6% ..</div> </div>
1	C	161	<div> <div>%</div> <div>96% ..</div> </div>
2	B	243	<div> <div>2%</div> <div>87% 5% 7%</div> </div>
2	D	243	<div> <div>5%</div> <div>88% . 7%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11989 atoms, of which 5842 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 GA-binding protein subunit beta-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	157	Total	C	H	N	O	S	0	0	0
			2337	738	1163	207	228	1			
1	C	158	Total	C	H	N	O	S	0	0	0
			2345	740	1166	208	230	1			

- Molecule 2 is a protein called Green fluorescent protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	226	Total	C	H	N	O	S	0	0	0
			3570	1152	1760	305	347	6			
2	D	225	Total	C	H	N	O	S	0	0	0
			3556	1148	1753	304	345	6			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	GLY	-	expression tag	UNP P42212
B	-3	PRO	-	expression tag	UNP P42212
B	-2	GLY	-	expression tag	UNP P42212
B	-1	SER	-	expression tag	UNP P42212
B	0	MET	-	expression tag	UNP P42212
B	1	VAL	-	expression tag	UNP P42212
B	64	LEU	PHE	conflict	UNP P42212
B	66	CRO	SER	chromophore	UNP P42212
B	66	CRO	TYR	chromophore	UNP P42212
B	66	CRO	GLY	chromophore	UNP P42212
B	231	LEU	HIS	conflict	UNP P42212
B	239	GLN	-	expression tag	UNP P42212
B	240	ALA	-	expression tag	UNP P42212
D	-4	GLY	-	expression tag	UNP P42212
D	-3	PRO	-	expression tag	UNP P42212
D	-2	GLY	-	expression tag	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-1	SER	-	expression tag	UNP P42212
D	0	MET	-	expression tag	UNP P42212
D	1	VAL	-	expression tag	UNP P42212
D	64	LEU	PHE	conflict	UNP P42212
D	66	CRO	SER	chromophore	UNP P42212
D	66	CRO	TYR	chromophore	UNP P42212
D	66	CRO	GLY	chromophore	UNP P42212
D	231	LEU	HIS	conflict	UNP P42212
D	239	GLN	-	expression tag	UNP P42212
D	240	ALA	-	expression tag	UNP P42212

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	45	Total O 45 45	0	0
3	C	37	Total O 37 37	0	0
3	B	52	Total O 52 52	0	0
3	D	47	Total O 47 47	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: GA-binding protein subunit beta-1

Chain A:  91% 6% ..




- Molecule 1: GA-binding protein subunit beta-1

Chain C:  96% ..




- Molecule 2: Green fluorescent protein

Chain B:  87% 5% 7% 2%



- Molecule 2: Green fluorescent protein

Chain D:  88% 5% 7% 1%



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	62.14Å 62.14Å 213.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.94 – 2.35 46.78 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.9 (43.94-2.35) 99.9 (46.78-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 2.34Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.189 , 0.236 0.190 , 0.236	Depositor DCC
R_{free} test set	1657 reflections (4.95%)	DCC
Wilson B-factor (Å ²)	42.4	Xtriage
Anisotropy	0.283	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 40.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.069 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11989	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CRO

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	0.24	0/1192	0.43	0/1623
1	C	0.25	0/1197	0.45	0/1629
2	B	0.26	0/1828	0.47	0/2470
2	D	0.26	0/1821	0.48	1/2460 (0.0%)
All	All	0.25	0/6038	0.46	1/8182 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	194	LEU	CA-CB-CG	5.01	126.83	115.30

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	1174	1163	1163	6	0
1	C	1179	1166	1166	1	0
2	B	1810	1760	1759	11	0
2	D	1803	1753	1752	9	0
3	A	45	0	0	1	0
3	B	52	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	37	0	0	1	0
3	D	47	0	0	5	0
All	All	6147	5842	5840	26	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 2.

All (26) 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)
2:D:126:LYS:NZ	3:D:301:HOH:O	1.94	1.00
1:A:69:TYR:OH	3:A:201:HOH:O	2.03	0.75
2:B:36:ASP:OD2	3:B:301:HOH:O	2.05	0.74
2:D:141:LEU:O	3:D:302:HOH:O	2.05	0.73
2:B:111:GLU:OE1	3:B:302:HOH:O	2.10	0.69
1:C:160:ASP:OD2	3:C:201:HOH:O	2.10	0.68
1:A:147:LYS:NZ	1:A:155:ASP:OD2	2.26	0.68
2:B:163:VAL:HB	2:B:183:GLN:HG2	1.81	0.63
2:B:21:ASP:OD2	3:B:303:HOH:O	2.16	0.61
1:A:56:GLN:OE1	1:A:57:ARG:NH1	2.37	0.57
2:B:163:VAL:HB	2:B:183:GLN:CG	2.38	0.54
2:D:122:ARG:NH1	3:D:306:HOH:O	2.37	0.52
1:A:23:ARG:NH2	2:B:34:GLU:OE2	2.44	0.51
2:D:85:LYS:NZ	3:D:310:HOH:O	2.44	0.50
2:B:96:ARG:HG2	2:B:183:GLN:CB	2.43	0.49
2:D:86:SER:HB3	2:D:194:LEU:CD2	2.42	0.49
2:B:139:HIS:ND1	3:B:311:HOH:O	2.35	0.48
2:D:3:LYS:N	3:D:312:HOH:O	2.48	0.47
2:B:96:ARG:HG2	2:B:183:GLN:HB2	1.98	0.45
2:D:145:TYR:HB2	2:D:207:LEU:CD1	2.48	0.43
2:B:136:ILE:N	2:B:136:ILE:CD1	2.82	0.43
1:A:91:GLY:HA2	1:A:128:ILE:HD12	2.02	0.42
2:D:207:LEU:HD12	2:D:207:LEU:N	2.35	0.41
2:B:66:CRO:N2	2:B:66:CRO:HD2	2.35	0.41
2:D:66:CRO:N2	2:D:66:CRO:HD2	2.35	0.41
1:A:148:THR:HG23	1:A:149:PRO:HD2	2.03	0.41

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	155/161 (96%)	155 (100%)	0	0	100	100
1	C	156/161 (97%)	156 (100%)	0	0	100	100
2	B	221/243 (91%)	217 (98%)	4 (2%)	0	100	100
2	D	220/243 (90%)	217 (99%)	3 (1%)	0	100	100
All	All	752/808 (93%)	745 (99%)	7 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	117/119 (98%)	115 (98%)	2 (2%)	66	77
1	C	118/119 (99%)	115 (98%)	3 (2%)	53	66
2	B	197/210 (94%)	192 (98%)	5 (2%)	53	66
2	D	196/210 (93%)	195 (100%)	1 (0%)	91	95
All	All	628/658 (95%)	617 (98%)	11 (2%)	64	76

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	ASP
1	A	148	THR

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Mol	Chain	Res	Type
1	C	63	VAL
1	C	68	LYS
1	C	112	ILE
2	B	3	LYS
2	B	30	SER
2	B	80	GLN
2	B	136	ILE
2	B	222	GLU
2	D	163	VAL

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 molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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	CRO	B	66	2	23,23,24	1.89	5 (21%)	27,32,34	3.06	8 (29%)
2	CRO	D	66	2	23,23,24	1.85	5 (21%)	27,32,34	2.76	8 (29%)

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	CRO	B	66	2	-	0/12/31/32	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	CRO	D	66	2	-	0/12/31/32	0/2/2/2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	66	CRO	CA3-N3	-2.01	1.43	1.47
2	B	66	CRO	C2-N3	2.21	1.45	1.39
2	D	66	CRO	CG2-CB2	2.76	1.52	1.46
2	B	66	CRO	CG2-CB2	2.77	1.52	1.46
2	D	66	CRO	C1-N3	3.20	1.43	1.37
2	B	66	CRO	CA2-C2	3.31	1.51	1.48
2	B	66	CRO	C1-N3	3.41	1.43	1.37
2	D	66	CRO	CA2-C2	3.59	1.52	1.48
2	D	66	CRO	C1-N2	5.45	1.40	1.32
2	B	66	CRO	C1-N2	5.70	1.41	1.32

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	66	CRO	O2-C2-CA2	-8.95	126.15	130.97
2	D	66	CRO	O2-C2-CA2	-6.79	127.31	130.97
2	B	66	CRO	CG2-CB2-CA2	-4.46	125.01	130.19
2	D	66	CRO	CG2-CB2-CA2	-4.19	125.32	130.19
2	B	66	CRO	C2-CA2-N2	-3.55	106.33	108.93
2	D	66	CRO	CA1-C1-N3	-3.48	120.58	124.75
2	D	66	CRO	C2-CA2-N2	-3.41	106.43	108.93
2	B	66	CRO	CA1-C1-N3	-3.24	120.86	124.75
2	D	66	CRO	O3-C3-CA3	-2.69	117.50	126.38
2	B	66	CRO	O3-C3-CA3	-2.07	119.54	126.38
2	D	66	CRO	CA1-C1-N2	2.72	127.69	123.89
2	B	66	CRO	CA1-C1-N2	2.75	127.74	123.89
2	D	66	CRO	CA2-N2-C1	3.81	108.74	105.75
2	B	66	CRO	CA2-N2-C1	4.37	109.18	105.75
2	D	66	CRO	CA2-C2-N3	8.79	107.22	103.30
2	B	66	CRO	CA2-C2-N3	9.35	107.47	103.30

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	66	CRO	1	0
2	D	66	CRO	1	0

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	64:LEU	C	66:CRO	N1	1.10

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	157/161 (97%)	0.23	0 100 100	32, 45, 67, 85	0
1	C	158/161 (98%)	0.21	1 (0%) 89 94	32, 48, 70, 85	0
2	B	225/243 (92%)	0.30	4 (1%) 69 78	31, 54, 79, 102	0
2	D	224/243 (92%)	0.46	11 (4%) 30 42	32, 63, 93, 116	0
All	All	764/808 (94%)	0.31	16 (2%) 64 74	31, 53, 85, 116	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	157	GLN	3.9
2	D	229	ILE	3.7
2	D	193	VAL	3.1
2	D	125	LEU	2.8
2	D	214	LYS	2.6
2	D	162	LYS	2.5
2	B	192	PRO	2.5
1	C	11	GLY	2.5
2	D	26	LYS	2.4
2	D	194	LEU	2.4
2	B	25	HIS	2.4
2	D	93	VAL	2.3
2	D	195	LEU	2.2
2	B	52	LYS	2.1
2	B	230	THR	2.1
2	D	80	GLN	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	CRO	D	66	22/23	0.91	0.12	-	20,30,36,39	0
2	CRO	B	66	22/23	0.95	0.11	-	21,28,34,36	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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.