



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 09:23 AM GMT

PDB ID : 4JW3
Title : Selection of specific protein binders for pre-defined targets from an optimized library of artificial helicoidal repeat proteins (alphaRep)
Authors : Guellouz, A.; Valerio-Lepiniec, M.; Urvoas, A.; Chevrel, A.; Graille, M.; Fourati-Kammoun, Z.; Desmadril, M.; van Tilbeurgh, H.; Minard, P.
Deposited on : 2013-03-27
Resolution : 2.60 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>

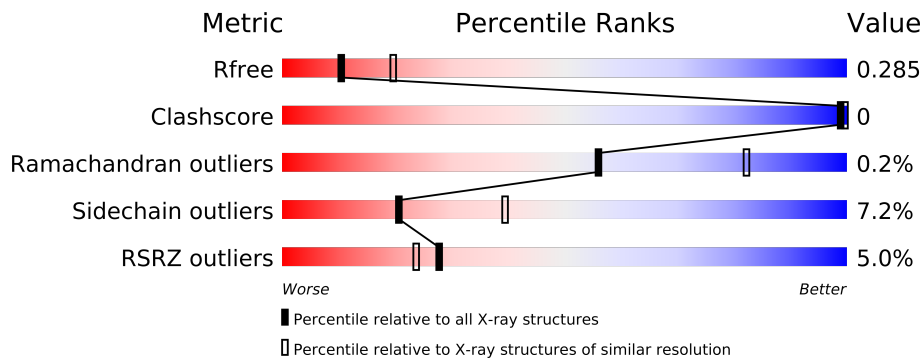
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.60 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	1718 (2.60-2.60)
Clashscore	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)
RSRZ outliers	66119	1718 (2.60-2.60)

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	120	
1	B	120	
2	C	139	
2	D	139	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3095 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 Neocarzinostatin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	111	Total	C	N	O	S	0	0	0
			785	486	135	162	2			
1	B	112	Total	C	N	O	S	0	0	0
			790	489	136	163	2			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	33	TRP	ASP	ENGINEERED MUTATION	UNP P0A3S0
A	35	ALA	GLY	ENGINEERED MUTATION	UNP P0A3S0
A	37	TRP	CYS	ENGINEERED MUTATION	UNP P0A3S0
A	39	ARG	TRP	ENGINEERED MUTATION	UNP P0A3S0
A	45	TRP	LEU	ENGINEERED MUTATION	UNP P0A3S0
A	47	TYR	CYS	ENGINEERED MUTATION	UNP P0A3S0
A	52	ASN	PHE	ENGINEERED MUTATION	UNP P0A3S0
A	113	ALA	-	EXPRESSION TAG	UNP P0A3S0
A	114	ALA	-	EXPRESSION TAG	UNP P0A3S0
A	115	HIS	-	EXPRESSION TAG	UNP P0A3S0
A	116	HIS	-	EXPRESSION TAG	UNP P0A3S0
A	117	HIS	-	EXPRESSION TAG	UNP P0A3S0
A	118	HIS	-	EXPRESSION TAG	UNP P0A3S0
A	119	HIS	-	EXPRESSION TAG	UNP P0A3S0
A	120	HIS	-	EXPRESSION TAG	UNP P0A3S0
B	33	TRP	ASP	ENGINEERED MUTATION	UNP P0A3S0
B	35	ALA	GLY	ENGINEERED MUTATION	UNP P0A3S0
B	37	TRP	CYS	ENGINEERED MUTATION	UNP P0A3S0
B	39	ARG	TRP	ENGINEERED MUTATION	UNP P0A3S0
B	45	TRP	LEU	ENGINEERED MUTATION	UNP P0A3S0
B	47	TYR	CYS	ENGINEERED MUTATION	UNP P0A3S0
B	52	ASN	PHE	ENGINEERED MUTATION	UNP P0A3S0
B	113	ALA	-	EXPRESSION TAG	UNP P0A3S0
B	114	ALA	-	EXPRESSION TAG	UNP P0A3S0
B	115	HIS	-	EXPRESSION TAG	UNP P0A3S0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	116	HIS	-	EXPRESSION TAG	UNP P0A3S0
B	117	HIS	-	EXPRESSION TAG	UNP P0A3S0
B	118	HIS	-	EXPRESSION TAG	UNP P0A3S0
B	119	HIS	-	EXPRESSION TAG	UNP P0A3S0
B	120	HIS	-	EXPRESSION TAG	UNP P0A3S0

- Molecule 2 is a protein called Alpha-helical artificial proteins.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	97	Total	C	N	O	S	0	0	0
			762	480	137	144	1			
2	D	96	Total	C	N	O	S	0	0	0
			752	474	134	143	1			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	O	0	0
			1	1		
3	B	2	Total	O	0	0
			2	2		
3	C	1	Total	O	0	0
			1	1		
3	D	2	Total	O	0	0
			2	2		

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.

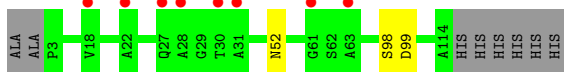
- Molecule 1: Neocarzinostatin

Chain A: 



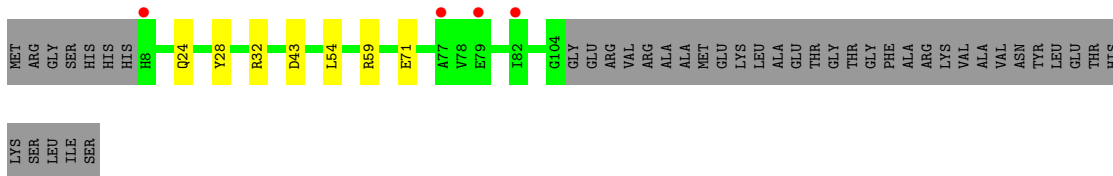
- Molecule 1: Neocarzinostatin

Chain B: 



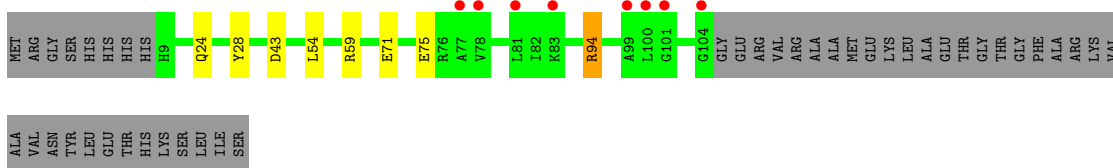
- Molecule 2: Alpha-helical artificial proteins

Chain C: 



- Molecule 2: Alpha-helical artificial proteins

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	118.21Å 59.21Å 65.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.77 – 2.60 39.40 – 2.60	Depositor EDS
% Data completeness (in resolution range)	95.5 (33.77-2.60) 96.0 (39.40-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 2.61Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.224 , 0.264 0.232 , 0.285	Depositor DCC
R_{free} test set	704 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	60.0	Xtriage
Anisotropy	0.731	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 52.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 14076 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3095	wwPDB-VP
Average B, all atoms (Å ²)	76.0	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 31.64 % 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 1.0708e-03. 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

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	0.48	0/802	0.71	0/1101
1	B	0.44	0/807	0.68	0/1107
2	C	0.56	0/775	0.70	0/1044
2	D	0.54	0/764	0.70	0/1029
All	All	0.51	0/3148	0.70	0/4281

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	785	0	0	1	0
1	B	790	0	0	0	0
2	C	762	0	0	0	0
2	D	752	0	0	1	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
3	C	1	0	0	0	0
3	D	2	0	0	0	0
All	All	3095	0	0	1	0

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 0.

All (1) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:81:THR:OG1	2:D:94:ARG:NH2	2.42	0.52

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	109/120 (91%)	105 (96%)	3 (3%)	1 (1%)	25	49
1	B	110/120 (92%)	106 (96%)	4 (4%)	0	100	100
2	C	95/139 (68%)	93 (98%)	2 (2%)	0	100	100
2	D	94/139 (68%)	92 (98%)	2 (2%)	0	100	100
All	All	408/518 (79%)	396 (97%)	11 (3%)	1 (0%)	56	82

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	103	ASN

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	79/85 (93%)	75 (95%)	4 (5%)	33	60
1	B	79/85 (93%)	76 (96%)	3 (4%)	44	74

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	C	75/108 (69%)	68 (91%)	7 (9%)	13	24
2	D	74/108 (68%)	66 (89%)	8 (11%)	9	17
All	All	307/386 (80%)	285 (93%)	22 (7%)	21	39

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	SER
1	A	52	ASN
1	A	82	ARG
1	A	98	SER
1	B	52	ASN
1	B	98	SER
1	B	99	ASP
2	C	24	GLN
2	C	28	TYR
2	C	32	ARG
2	C	43	ASP
2	C	54	LEU
2	C	59	ARG
2	C	71	GLU
2	D	24	GLN
2	D	28	TYR
2	D	43	ASP
2	D	54	LEU
2	D	59	ARG
2	D	71	GLU
2	D	75	GLU
2	D	94	ARG

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 ⓘ

There are no ligands in this entry.

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	111/120 (92%)	0.21	1 (0%) 81 82	45, 69, 88, 97	0
1	B	112/120 (93%)	0.63	8 (7%) 16 12	44, 92, 120, 143	0
2	C	97/139 (69%)	0.28	4 (4%) 35 32	51, 72, 108, 111	0
2	D	96/139 (69%)	0.34	8 (8%) 11 8	51, 74, 103, 122	0
All	All	416/518 (80%)	0.37	21 (5%) 28 24	44, 75, 109, 143	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	77	ALA	3.7
2	C	77	ALA	3.4
1	B	28	ALA	3.3
2	C	82	ILE	3.1
1	B	61	GLY	3.0
1	B	31	ALA	3.0
1	A	63	ALA	2.9
2	D	101	GLY	2.7
1	B	63	ALA	2.7
2	D	100	LEU	2.7
2	D	104	GLY	2.5
1	B	27	GLN	2.5
1	B	30	THR	2.5
2	D	78	VAL	2.4
2	D	81	LEU	2.4
2	C	79	GLU	2.4
1	B	22	ALA	2.2
2	D	99	ALA	2.2
2	C	8	HIS	2.0
1	B	18	VAL	2.0
2	D	83	LYS	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 ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.