



Full wwPDB NMR Structure Validation Report ⓘ

Apr 26, 2016 – 07:19 PM BST

PDB ID : 2B88
Title : Structural basis for molecular recognition in an affibody:affibody complex
Authors : Lendel, C.; Dogan, J.; Hard, T.
Deposited on : 2005-10-06

This is a Full wwPDB NMR 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/NMRValidationReportHelp>

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:

Cyrange : Kirchner and Güntert (2011)
NmrClust : Kelley et al. (1996)
MolProbity : 4.02b-467
Mogul : unknown
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
ShiftChecker : rb-20027457
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027457

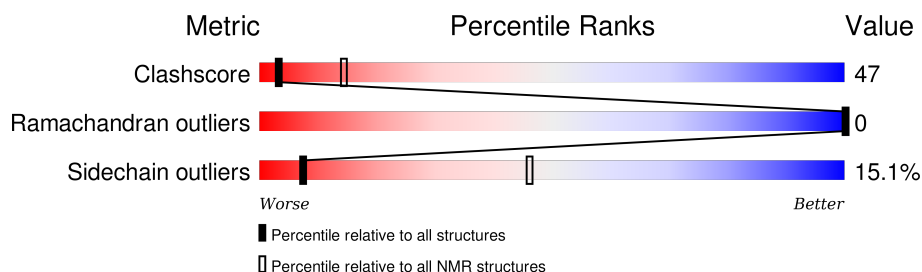
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 92%.

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)	NMR archive (#Entries)
Clashscore	114402	11133
Ramachandran outliers	111179	9975
Sidechain outliers	111093	9958

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and 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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	58	<div> <div></div> <div>28%</div> <div>45%</div> <div>7%</div> <div>21%</div> </div>

2 Ensemble composition and analysis

This entry contains 40 models. Model 34 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *closest to the average*.

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:11-A:56 (46)	0.12	34

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 6 clusters and 2 single-model clusters were found.

Cluster number	Models
1	1, 8, 9, 11, 12, 13, 19, 20, 22, 23, 27, 28, 33, 34, 35, 37, 39
2	2, 3, 4, 10, 14, 21
3	7, 17, 25, 29, 32
4	6, 16, 24, 38
5	5, 18, 31
6	15, 26, 40
Single-model clusters	30; 36

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 913 atoms, of which 454 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called ZTaq affibody.

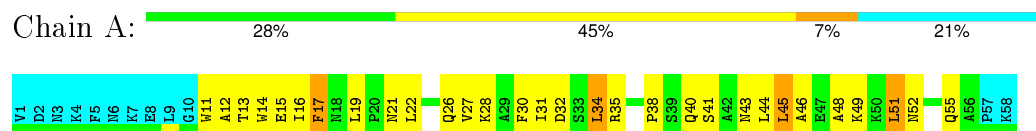
Mol	Chain	Residues	Atoms					Trace
1	A	58	Total	C	H	N	O	0
			913	291	454	79	89	

4 Residue-property plots

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA and DNA chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-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. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

- Molecule 1: ZTaq affibody

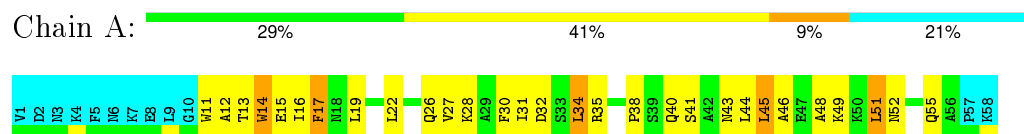


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

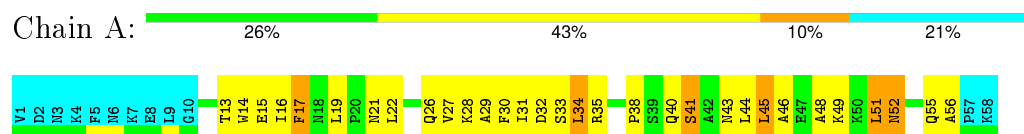
4.2.1 Score per residue for model 1

- Molecule 1: ZTaq affibody



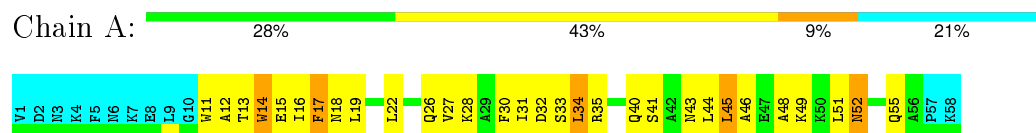
4.2.2 Score per residue for model 2

- Molecule 1: ZTaq affibody



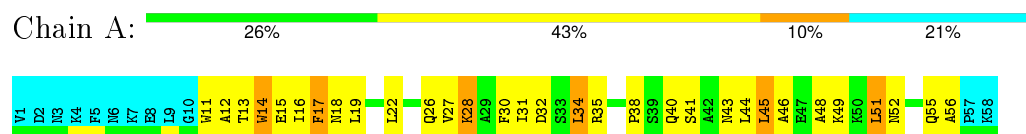
4.2.3 Score per residue for model 3

- Molecule 1: ZTaq affibody



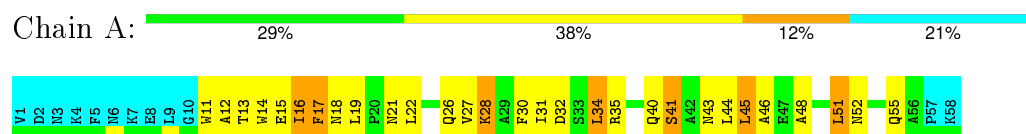
4.2.4 Score per residue for model 4

- Molecule 1: ZTaq affibody



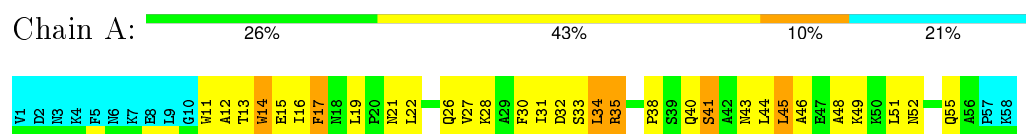
4.2.5 Score per residue for model 5

- Molecule 1: ZTaq affibody



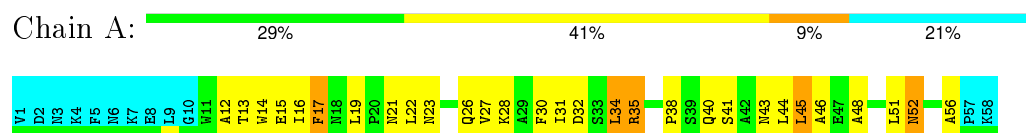
4.2.6 Score per residue for model 6

- Molecule 1: ZTaq affibody



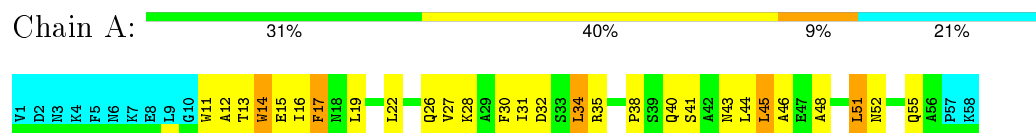
4.2.7 Score per residue for model 7

- Molecule 1: ZTaq affibody



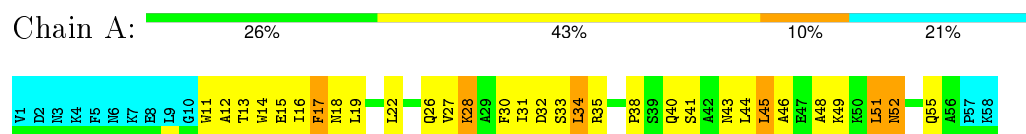
4.2.8 Score per residue for model 8

- Molecule 1: ZTaq affibody



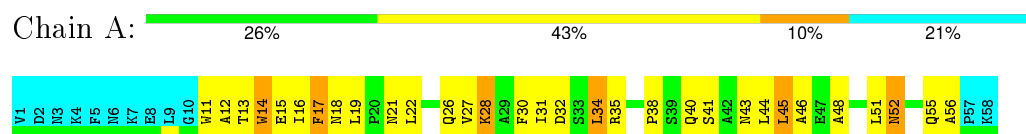
4.2.9 Score per residue for model 9

- Molecule 1: ZTaq affibody



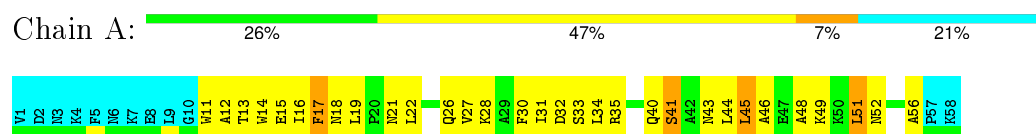
4.2.10 Score per residue for model 10

- Molecule 1: ZTaq affibody



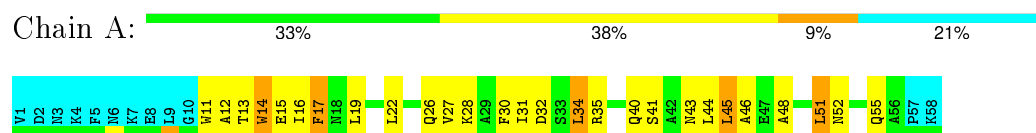
4.2.11 Score per residue for model 11

- Molecule 1: ZTaq affibody



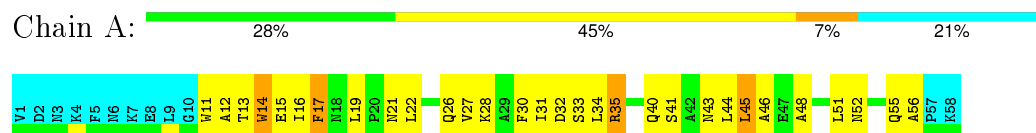
4.2.12 Score per residue for model 12

- Molecule 1: ZTaq affibody



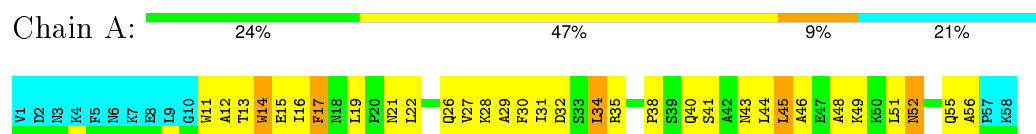
4.2.13 Score per residue for model 13

- Molecule 1: ZTaq affibody



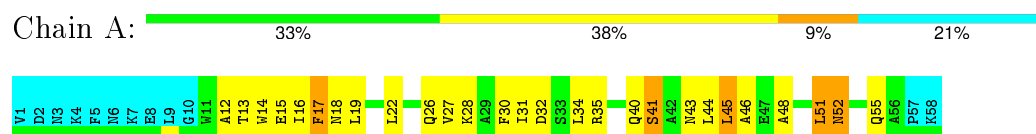
4.2.14 Score per residue for model 14

- Molecule 1: ZTaq affibody



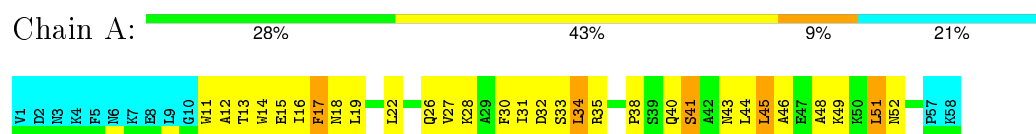
4.2.15 Score per residue for model 15

- Molecule 1: ZTaq affibody



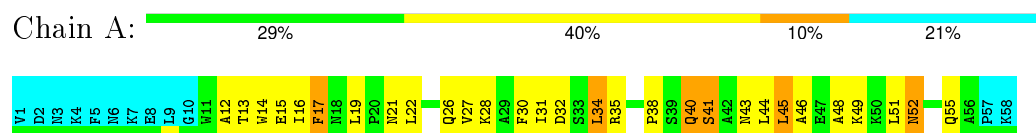
4.2.16 Score per residue for model 16

- Molecule 1: ZTaq affibody



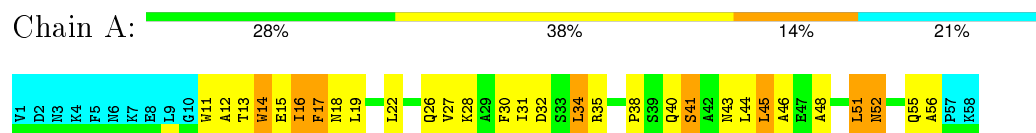
4.2.17 Score per residue for model 17

- Molecule 1: ZTaq affibody



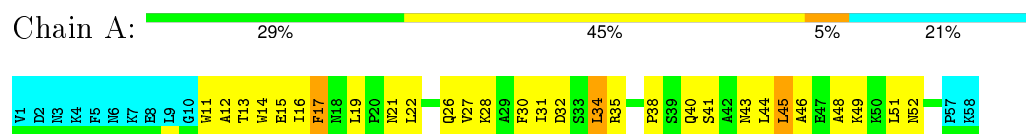
4.2.18 Score per residue for model 18

- Molecule 1: ZTaq affibody



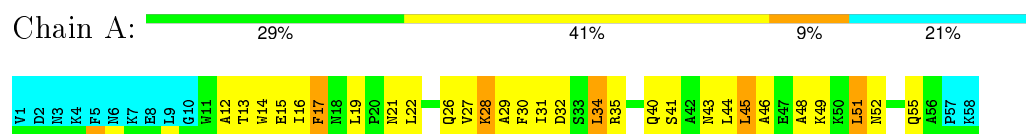
4.2.19 Score per residue for model 19

- Molecule 1: ZTaq affibody



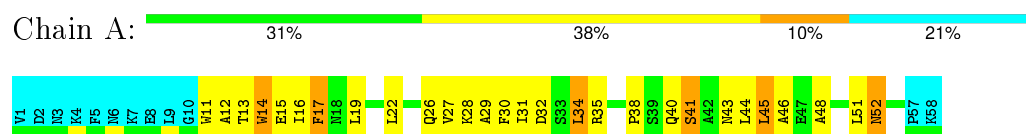
4.2.20 Score per residue for model 20

- Molecule 1: ZTaq affibody



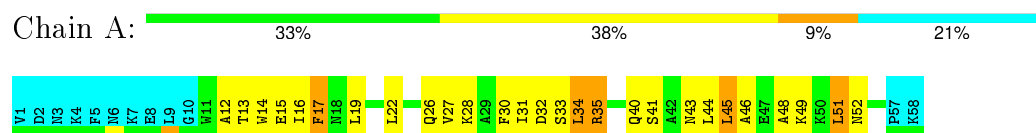
4.2.21 Score per residue for model 21

- Molecule 1: ZTaq affibody



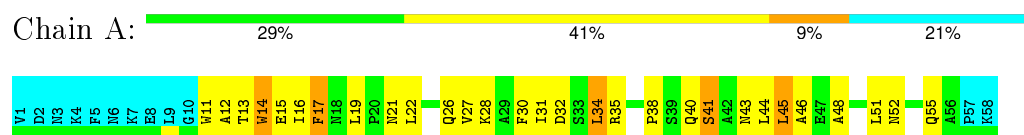
4.2.22 Score per residue for model 22

- Molecule 1: ZTaq affibody



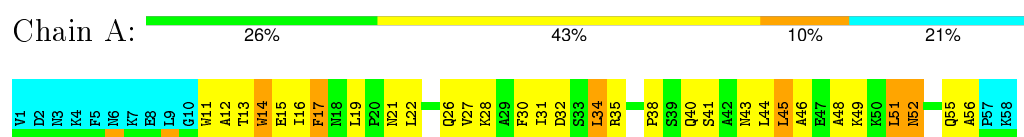
4.2.23 Score per residue for model 23

- Molecule 1: ZTaq affibody



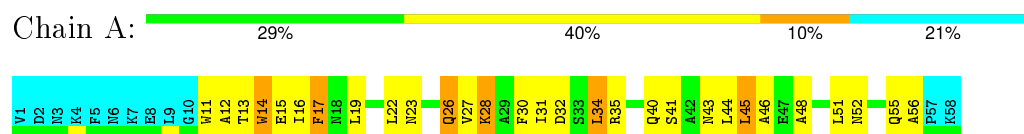
4.2.24 Score per residue for model 24

- Molecule 1: ZTaq affibody



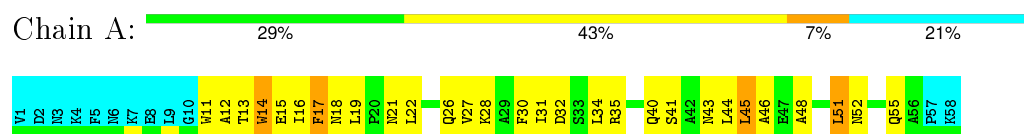
4.2.25 Score per residue for model 25

- Molecule 1: ZTaq affibody



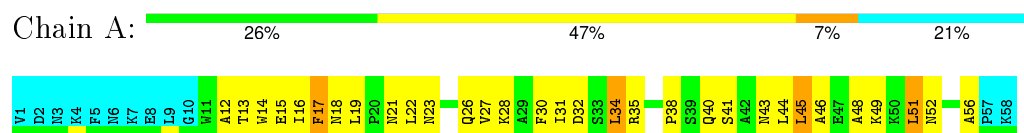
4.2.26 Score per residue for model 26

- Molecule 1: ZTaq affibody



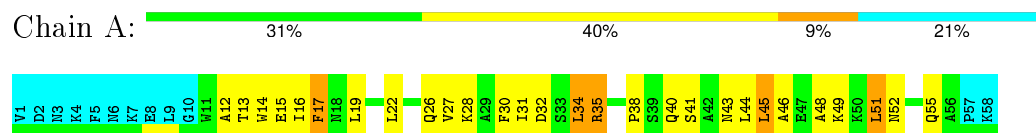
4.2.27 Score per residue for model 27

- Molecule 1: ZTaq affibody



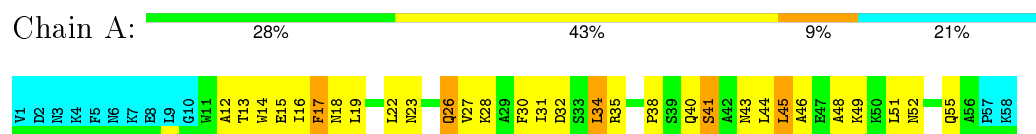
4.2.28 Score per residue for model 28

- Molecule 1: ZTaq affibody



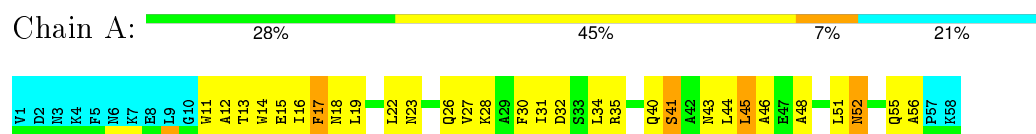
4.2.29 Score per residue for model 29

- Molecule 1: ZTaq affibody



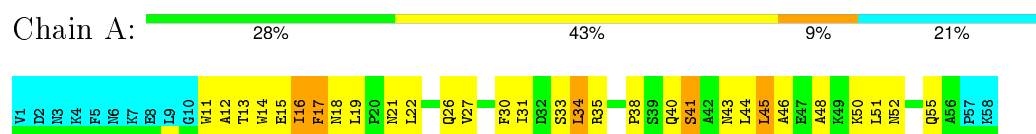
4.2.30 Score per residue for model 30

- Molecule 1: ZTaq affibody



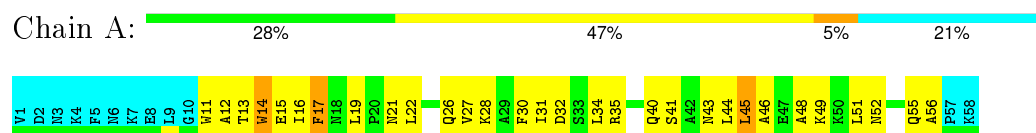
4.2.31 Score per residue for model 31

- Molecule 1: ZTaq affibody



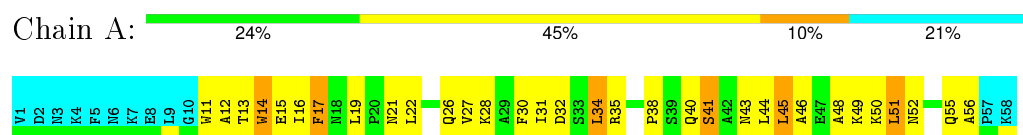
4.2.32 Score per residue for model 32

- Molecule 1: ZTaq affibody



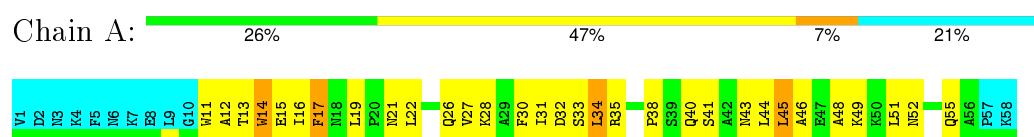
4.2.33 Score per residue for model 33

- Molecule 1: ZTaq affibody



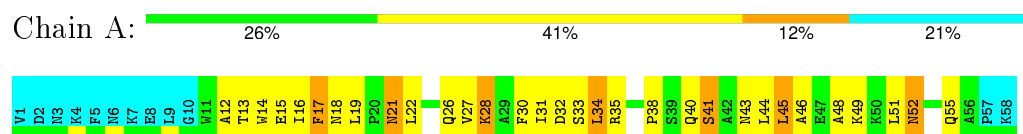
4.2.34 Score per residue for model 34 (medoid)

- Molecule 1: ZTaq affibody



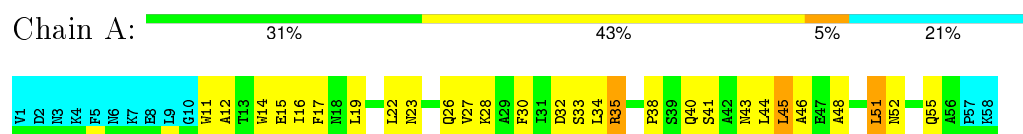
4.2.35 Score per residue for model 35

- Molecule 1: ZTaq affibody



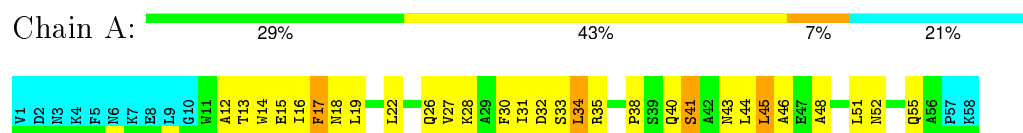
4.2.36 Score per residue for model 36

- Molecule 1: ZTaq affibody



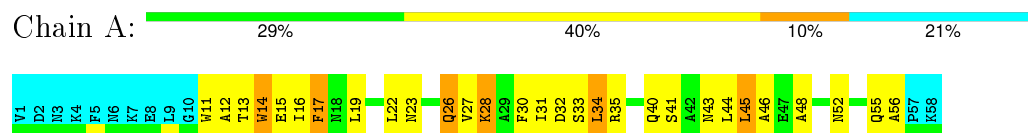
4.2.37 Score per residue for model 37

- Molecule 1: ZTaq affibody



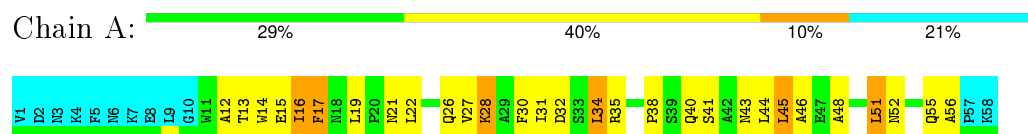
4.2.38 Score per residue for model 38

- Molecule 1: ZTaq affibody



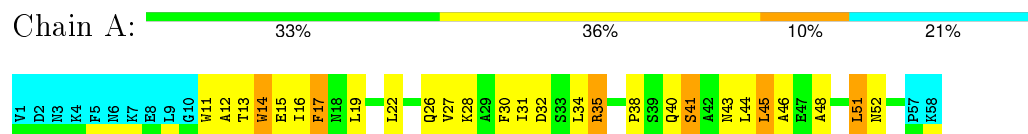
4.2.39 Score per residue for model 39

- Molecule 1: ZTaq affibody



4.2.40 Score per residue for model 40

- Molecule 1: ZTaq affibody



5 Refinement protocol and experimental data overview

The models were refined using the following method: *simulated annealing*.

Of the 100 calculated structures, 40 were deposited, based on the following criterion: *structures with the least restraint violations, structures with the lowest energy, good ramachandran plots*.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
X-PLOR	structure solution	
X-PLOR	refinement	

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	BMRB entry 6804
Number of chemical shift lists	1
Total number of shifts	722
Number of shifts mapped to atoms	722
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	92%

No validations of the models with respect to experimental NMR restraints is performed at this time.

6 Model quality [i](#)

6.1 Standard geometry [i](#)

There are no covalent bond-length or bond-angle outliers.

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	Chirality	Planarity
1	A	0.0±0.0	1.0±0.0
All	All	0	40

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

All unique planar outliers are listed below.

Mol	Chain	Res	Type	Group	Models (Total)
1	A	35	ARG	Sidechain	40

6.2 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 each 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 averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	361	352	352	34±3
All	All	14440	14080	14080	1342

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 47.

All unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:17:PHE:CD1	1:A:27:VAL:HG11	0.99	1.93	24	39
1:A:26:GLN:NE2	1:A:51:LEU:HD11	0.86	1.85	14	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:34:LEU:HD12	1:A:44:LEU:HD12	0.81	1.52	15	6
1:A:19:LEU:HD13	1:A:52:ASN:HB3	0.80	1.50	7	12
1:A:15:GLU:HB3	1:A:45:LEU:HD11	0.79	1.54	17	40
1:A:30:PHE:CZ	1:A:51:LEU:HD12	0.78	2.14	13	5
1:A:12:ALA:HB3	1:A:34:LEU:HD21	0.77	1.57	38	24
1:A:17:PHE:CG	1:A:27:VAL:HG11	0.75	2.16	30	38
1:A:17:PHE:CE1	1:A:27:VAL:HG11	0.73	2.19	25	39
1:A:12:ALA:HB2	1:A:41:SER:HB2	0.70	1.63	32	27
1:A:19:LEU:HD13	1:A:52:ASN:HB2	0.70	1.61	38	26
1:A:12:ALA:CB	1:A:34:LEU:HD21	0.68	2.18	38	29
1:A:27:VAL:HG12	1:A:31:ILE:CD1	0.67	2.20	10	18
1:A:12:ALA:HB2	1:A:41:SER:OG	0.67	1.89	40	6
1:A:34:LEU:HD22	1:A:44:LEU:HD12	0.67	1.65	29	25
1:A:22:LEU:HD13	1:A:27:VAL:HG22	0.67	1.67	14	38
1:A:16:ILE:O	1:A:22:LEU:HD12	0.65	1.91	14	39
1:A:27:VAL:HG12	1:A:31:ILE:HD12	0.64	1.70	40	18
1:A:33:SER:OG	1:A:44:LEU:HD22	0.63	1.94	22	2
1:A:34:LEU:HD22	1:A:44:LEU:CD1	0.63	2.24	12	24
1:A:16:ILE:HD13	1:A:30:PHE:HB3	0.63	1.70	36	35
1:A:26:GLN:NE2	1:A:51:LEU:HD12	0.63	2.08	35	28
1:A:45:LEU:HD12	1:A:49:LYS:NZ	0.62	2.09	34	3
1:A:22:LEU:CD2	1:A:51:LEU:HD12	0.62	2.24	14	1
1:A:34:LEU:CD1	1:A:44:LEU:HD12	0.62	2.24	32	6
1:A:19:LEU:HD21	1:A:49:LYS:NZ	0.61	2.10	3	8
1:A:19:LEU:HD12	1:A:22:LEU:HD11	0.61	1.73	31	38
1:A:26:GLN:CD	1:A:51:LEU:HD12	0.60	2.16	19	23
1:A:13:THR:HG23	1:A:31:ILE:CD1	0.59	2.27	20	18
1:A:22:LEU:CD1	1:A:27:VAL:HG22	0.59	2.28	20	39
1:A:17:PHE:CD1	1:A:27:VAL:CG1	0.58	2.87	36	1
1:A:34:LEU:HD11	1:A:41:SER:HB2	0.58	1.76	29	14
1:A:16:ILE:HD12	1:A:30:PHE:HB3	0.56	1.77	31	3
1:A:43:ASN:O	1:A:46:ALA:HB3	0.56	2.00	23	40
1:A:13:THR:HA	1:A:31:ILE:HD12	0.56	1.78	13	21
1:A:34:LEU:HD11	1:A:41:SER:HB3	0.55	1.78	2	12
1:A:33:SER:CB	1:A:44:LEU:HD22	0.55	2.32	36	11
1:A:12:ALA:HB2	1:A:41:SER:HB3	0.55	1.78	23	1
1:A:26:GLN:NE2	1:A:55:GLN:NE2	0.54	2.55	4	24
1:A:15:GLU:CB	1:A:45:LEU:HD11	0.54	2.32	35	15
1:A:16:ILE:HG21	1:A:30:PHE:HB2	0.54	1.79	36	39
1:A:11:TRP:O	1:A:14:TRP:CE3	0.54	2.61	18	20
1:A:12:ALA:O	1:A:16:ILE:CG1	0.53	2.57	18	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:21:ASN:ND2	1:A:56:ALA:HB2	0.53	2.18	13	7
1:A:40:GLN:O	1:A:44:LEU:HG	0.53	2.04	14	40
1:A:40:GLN:O	1:A:43:ASN:N	0.53	2.42	29	40
1:A:19:LEU:HD13	1:A:52:ASN:CB	0.53	2.34	36	31
1:A:45:LEU:HD12	1:A:49:LYS:HE2	0.53	1.80	16	7
1:A:16:ILE:HG23	1:A:48:ALA:CB	0.52	2.35	20	31
1:A:26:GLN:NE2	1:A:51:LEU:HD22	0.52	2.19	30	5
1:A:34:LEU:HD11	1:A:41:SER:CB	0.52	2.35	39	9
1:A:28:LYS:O	1:A:32:ASP:N	0.52	2.43	15	39
1:A:51:LEU:HD12	1:A:55:GLN:OE1	0.51	2.04	6	1
1:A:22:LEU:CB	1:A:27:VAL:HG22	0.51	2.36	36	1
1:A:22:LEU:HB3	1:A:27:VAL:HG22	0.51	1.82	36	1
1:A:12:ALA:O	1:A:16:ILE:HG13	0.50	2.06	39	4
1:A:13:THR:HG23	1:A:31:ILE:HD13	0.49	1.83	30	13
1:A:33:SER:OG	1:A:44:LEU:HD13	0.49	2.07	22	2
1:A:45:LEU:O	1:A:48:ALA:HB3	0.49	2.07	30	31
1:A:22:LEU:HB3	1:A:27:VAL:CG2	0.48	2.38	36	6
1:A:30:PHE:CE2	1:A:48:ALA:HA	0.48	2.43	30	8
1:A:26:GLN:NE2	1:A:55:GLN:HE21	0.48	2.06	9	13
1:A:23:ASN:H	1:A:26:GLN:NE2	0.48	2.07	25	3
1:A:16:ILE:HD12	1:A:30:PHE:CB	0.47	2.37	31	3
1:A:23:ASN:N	1:A:26:GLN:NE2	0.47	2.62	25	3
1:A:17:PHE:CD2	1:A:27:VAL:HG11	0.47	2.45	31	2
1:A:12:ALA:HB2	1:A:41:SER:CB	0.47	2.38	32	3
1:A:51:LEU:O	1:A:51:LEU:HD13	0.47	2.10	14	1
1:A:45:LEU:HD12	1:A:49:LYS:HE3	0.46	1.86	11	2
1:A:14:TRP:C	1:A:14:TRP:CE3	0.46	2.89	22	11
1:A:26:GLN:O	1:A:30:PHE:CD2	0.46	2.69	36	10
1:A:51:LEU:HD13	1:A:51:LEU:C	0.46	2.32	14	1
1:A:14:TRP:CE3	1:A:14:TRP:C	0.46	2.89	15	9
1:A:17:PHE:CE1	1:A:27:VAL:HG12	0.46	2.45	36	1
1:A:26:GLN:O	1:A:29:ALA:HB3	0.45	2.10	14	4
1:A:34:LEU:HD12	1:A:44:LEU:CD1	0.45	2.39	40	1
1:A:41:SER:HA	1:A:44:LEU:HD12	0.45	1.87	10	3
1:A:19:LEU:HD21	1:A:49:LYS:HZ2	0.45	1.72	24	7
1:A:27:VAL:HG12	1:A:31:ILE:HD11	0.45	1.89	5	7
1:A:26:GLN:NE2	1:A:55:GLN:HE22	0.45	2.09	25	3
1:A:52:ASN:O	1:A:56:ALA:N	0.45	2.50	7	10
1:A:14:TRP:CH2	1:A:15:GLU:CD	0.45	2.90	39	3
1:A:30:PHE:CE2	1:A:51:LEU:HD12	0.44	2.47	3	5
1:A:16:ILE:HD13	1:A:30:PHE:CB	0.44	2.42	36	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:17:PHE:HD1	1:A:27:VAL:CG1	0.43	2.26	36	1
1:A:41:SER:O	1:A:45:LEU:HD23	0.43	2.13	14	17
1:A:21:ASN:N	1:A:52:ASN:OD1	0.43	2.51	20	9
1:A:23:ASN:O	1:A:27:VAL:HG23	0.43	2.13	30	2
1:A:22:LEU:HD23	1:A:51:LEU:HD12	0.43	1.89	14	1
1:A:21:ASN:N	1:A:21:ASN:OD1	0.43	2.51	35	1
1:A:34:LEU:HD13	1:A:44:LEU:HD12	0.43	1.89	24	1
1:A:31:ILE:O	1:A:34:LEU:HB2	0.43	2.14	25	3
1:A:33:SER:HB2	1:A:44:LEU:HD13	0.43	1.91	11	1
1:A:17:PHE:CD1	1:A:27:VAL:HG21	0.43	2.48	30	1
1:A:17:PHE:CE1	1:A:27:VAL:CG1	0.43	3.01	16	13
1:A:33:SER:HB3	1:A:44:LEU:HD22	0.42	1.90	36	1
1:A:31:ILE:N	1:A:31:ILE:HD13	0.42	2.28	23	10
1:A:31:ILE:HD13	1:A:31:ILE:N	0.42	2.28	2	6
1:A:21:ASN:HB3	1:A:55:GLN:CB	0.42	2.44	24	3
1:A:26:GLN:CD	1:A:51:LEU:HD22	0.42	2.34	3	1
1:A:22:LEU:HD23	1:A:55:GLN:HE22	0.42	1.74	6	1
1:A:51:LEU:CD1	1:A:55:GLN:NE2	0.42	2.83	14	1
1:A:11:TRP:HA	1:A:14:TRP:CD1	0.42	2.50	19	6
1:A:12:ALA:HB3	1:A:34:LEU:CD2	0.42	2.44	30	1
1:A:16:ILE:HG23	1:A:48:ALA:HB1	0.41	1.90	20	1
1:A:40:GLN:CG	1:A:44:LEU:HD11	0.41	2.44	17	1
1:A:23:ASN:H	1:A:26:GLN:HE21	0.41	1.59	7	1
1:A:22:LEU:HB3	1:A:27:VAL:HG23	0.41	1.93	9	6
1:A:34:LEU:HD13	1:A:34:LEU:HA	0.40	1.78	20	1
1:A:11:TRP:O	1:A:14:TRP:CD1	0.40	2.75	31	2
1:A:21:ASN:O	1:A:55:GLN:NE2	0.40	2.54	23	2
1:A:21:ASN:HD22	1:A:52:ASN:ND2	0.40	2.15	10	1
1:A:34:LEU:HA	1:A:34:LEU:HD13	0.40	1.77	24	1
1:A:32:ASP:HA	1:A:35:ARG:NE	0.40	2.31	6	1
1:A:26:GLN:NE2	1:A:55:GLN:OE1	0.40	2.55	25	2
1:A:19:LEU:HA	1:A:19:LEU:HD23	0.40	1.74	3	1

6.3 Torsion angles ⓘ

6.3.1 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 PDB entries followed by that with respect to all NMR entries. 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	46/58 (79%)	44±1 (95±1%)	2±1 (5±1%)	0±0 (0±0%)	100	100
All	All	1840/2320 (79%)	1746 (95%)	94 (5%)	0 (0%)	100	100

There are no Ramachandran outliers.

6.3.2 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 PDB entries followed by that with respect to all NMR entries. 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	38/49 (78%)	32±1 (85±4%)	6±1 (15±4%)	7	46
All	All	1520/1960 (78%)	1291 (85%)	229 (15%)	7	46

All 17 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	45	LEU	40
1	A	17	PHE	39
1	A	34	LEU	32
1	A	51	LEU	24
1	A	14	TRP	20
1	A	41	SER	17
1	A	18	ASN	16
1	A	52	ASN	13
1	A	28	LYS	9
1	A	35	ARG	5
1	A	16	ILE	4
1	A	26	GLN	3
1	A	21	ASN	2
1	A	50	LYS	2
1	A	23	ASN	1
1	A	40	GLN	1
1	A	49	LYS	1

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

The completeness of assignment taking into account all chemical shift lists is 92% for the well-defined parts and 91% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 6804

Chemical shift list name: *assigned_chem_shift_list_1*

7.1.1 Bookkeeping

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	722
Number of shifts mapped to atoms	722
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	58	-0.21 ± 0.19	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	56	0.21 ± 0.06	None needed (< 0.5 ppm)
$^{13}\text{C}'$	57	-0.20 ± 0.12	None needed (< 0.5 ppm)
^{15}N	53	-0.40 ± 0.32	None needed (< 0.5 ppm)

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 92%, i.e. 515 atoms were assigned a chemical shift out of a possible 561. 8 out of 8 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	226/226 (100%)	90/90 (100%)	92/92 (100%)	44/44 (100%)
Sidechain	267/293 (91%)	165/170 (97%)	93/109 (85%)	9/14 (64%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	22/42 (52%)	20/22 (91%)	0/18 (0%)	2/2 (100%)
Overall	515/561 (92%)	275/282 (98%)	185/219 (84%)	55/60 (92%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 91%, i.e. 654 atoms were assigned a chemical shift out of a possible 719. 10 out of 10 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	279/284 (98%)	111/113 (98%)	115/116 (99%)	53/55 (96%)
Sidechain	351/384 (91%)	219/224 (98%)	121/141 (86%)	11/19 (58%)
Aromatic	24/51 (47%)	22/27 (81%)	0/22 (0%)	2/2 (100%)
Overall	654/719 (91%)	352/364 (97%)	236/279 (85%)	66/76 (87%)

7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

7.1.5 Random Coil Index (RCI) plots [i](#)

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition.

Random coil index (RCI) for chain A:

