



# wwPDB NMR Structure Validation Summary Report ⓘ

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PDB ID : 1ZY3  
Title : Structural model of complex of Bcl-w protein with Bid BH3-peptide  
Authors : Denisov, A.Y.; Gehring, K.  
Deposited on : 2005-06-09

This is a wwPDB NMR Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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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
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	trunk28760
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

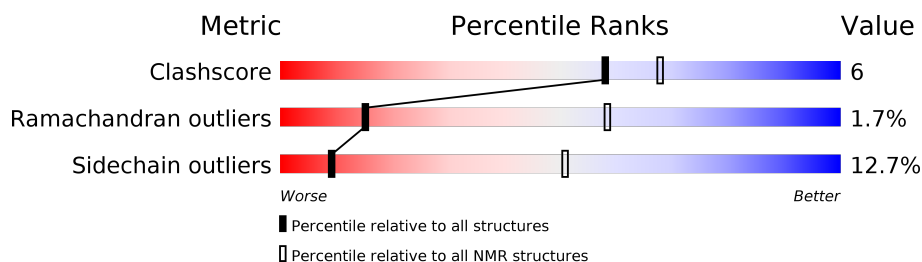
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment was not calculated.

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	125131	11601
Ramachandran outliers	121729	10391
Sidechain outliers	121581	10367

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  $\geq 3$ , 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	178	 63% 21% 10%
2	B	20	 60% 30% 10%

## 2 Ensemble composition and analysis

This entry contains 10 models. Model 8 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

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:1-A:152, B:201-B:218 (170)	0.64	8

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 2 clusters. No single-model clusters were found.

Cluster number	Models
1	3, 4, 5, 6, 7, 8, 9, 10
2	1, 2

### 3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2887 atoms, of which 1420 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Apoptosis regulator Bcl-W.

Mol	Chain	Residues	Atoms						Trace
1	A	170	Total	C	H	N	O	S	0
			2564	827	1253	235	244	5	

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	116	VAL	PRO	ENGINEERED	UNP Q92843
A	171	LEU	-	EXPRESSION TAG	UNP Q92843
A	172	GLU	-	EXPRESSION TAG	UNP Q92843
A	173	HIS	-	EXPRESSION TAG	UNP Q92843
A	174	HIS	-	EXPRESSION TAG	UNP Q92843
A	175	HIS	-	EXPRESSION TAG	UNP Q92843
A	176	HIS	-	EXPRESSION TAG	UNP Q92843
A	177	HIS	-	EXPRESSION TAG	UNP Q92843
A	178	HIS	-	EXPRESSION TAG	UNP Q92843

- Molecule 2 is a protein called BH3-peptide from BH3 interacting domain death agonist protein.

Mol	Chain	Residues	Atoms						Trace
2	B	20	Total	C	H	N	O	S	0
			323	95	167	31	29	1	

There is a discrepancy between the modelled and reference sequences:

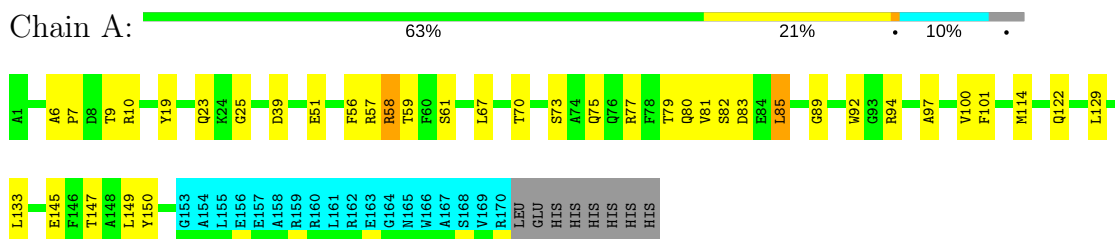
Chain	Residue	Modelled	Actual	Comment	Reference
B	203	LYS	ARG	ENGINEERED	UNP P55957

## 4 Residue-property plots [i](#)

### 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: Apoptosis regulator Bcl-W



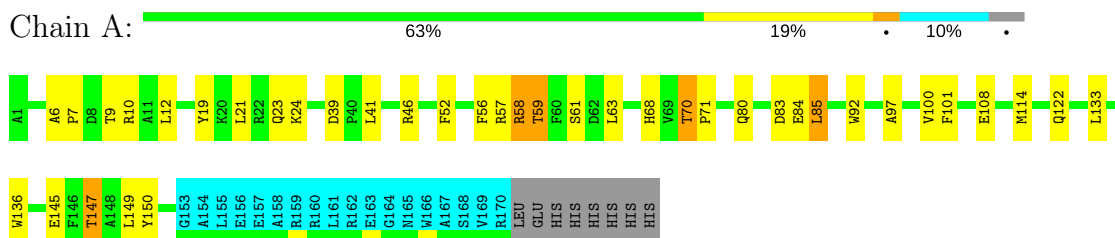
- Molecule 2: BH3-peptide from BH3 interacting domain death agonist protein



### 4.2 Residue scores for the representative (medoid) model from the NMR ensemble

The representative model is number 8. Colouring as in section 4.1 above.

- Molecule 1: Apoptosis regulator Bcl-W



- Molecule 2: BH3-peptide from BH3 interacting domain death agonist protein



I201	I202	K203	N204	I205	A206	R207	H208	L209	D214	R218	S219	I220
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## 5 Refinement protocol and experimental data overview

The models were refined using the following method: *Rigid body docking, semi-flexible simulated annealing, refinement using explicit water.*

Of the 100 calculated structures, 10 were deposited, based on the following criterion: *structures with the least restraint violations, structures with the lowest energy.*

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

Software name	Classification	Version
HADDOCK	refinement	1.2

No chemical shift data was provided. 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.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	1166	1110	1107	14±3
2	B	141	151	147	3±2
All	All	13070	12610	12540	147

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 6.

5 of 82 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:93:GLY:HA3	2:B:214:ASP:HA	0.88	1.42	4	3
1:A:70:THR:HG22	1:A:71:PRO:HD2	0.76	1.57	8	1
1:A:56:PHE:HA	2:B:216:MET:HG2	0.74	1.56	5	2
1:A:66:GLN:HG3	1:A:67:LEU:HD22	0.70	1.64	1	1
1:A:80:GLN:HG2	2:B:201:ILE:HD12	0.65	1.67	8	1

### 6.3 Torsion angles [i](#)

#### 6.3.1 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 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	151/178 (85%)	133±2 (88±1%)	16±2 (10±1%)	3±1 (2±1%)	14	57
2	B	17/20 (85%)	16±1 (94±4%)	1±1 (5±4%)	0±0 (1±2%)	20	66
All	All	1680/1980 (85%)	1488 (89%)	163 (10%)	29 (2%)	15	58

5 of 9 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	57	ARG	10
1	A	25	GLY	5
1	A	89	GLY	4
1	A	35	GLY	4
2	B	202	ILE	2

### 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	114/135 (84%)	100±2 (88±2%)	14±2 (12±2%)	9	51
2	B	15/17 (88%)	13±1 (85±7%)	2±1 (15±7%)	7	45
All	All	1290/1520 (85%)	1126 (87%)	164 (13%)	9	51

5 of 44 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	58	ARG	10
1	A	85	LEU	10
1	A	114	MET	10
2	B	218	ARG	9
1	A	122	GLN	9

### 6.3.3 RNA [i](#)

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

No chemical shift data were provided