

wwPDB NMR Structure Validation Summary Report (i)

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PDB ID	:	2ADR
Title	:	ADR1 DNA-BINDING DOMAIN FROM SACCHAROMYCES CERE-
		VISIAE, NMR, 25 STRUCTURES
Authors	:	Bowers, P.M.; Kleivt, R.E.
Deposited on	:	1998-03-20

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

We welcome your comments at *validation@mail.wwpdb.org* A user guide is available at https://www.wwpdb.org/validation/2017/NMRValidationReportHelp with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

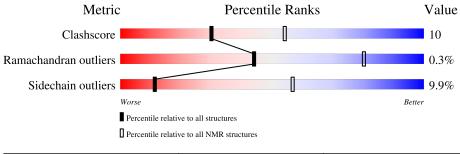
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.23.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

1 Overall quality at a glance (i)

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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f NMR} \ { m archive} \ (\#{ m Entries})$
Clashscore	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

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 <=5%

Mol	Chain	Length	Quality of chain			
1	А	60	48%	27%	••	22%



2 Ensemble composition and analysis (i)

This entry contains 25 models. Model 3 is the overall representative, medoid model (most similar to other models).

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:104-A:126 (23)	0.25	20		
2	A:132-A:155 (24)	0.20	3		

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. No single-model clusters were found.

Cluster number	Models
1	2, 11, 12, 17, 20, 23
2	3, 9, 10, 14, 18, 19
3	4, 5, 8, 22
4	6, 7, 15, 24
5	1, 16, 25
6	13, 21



3 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 989 atoms, of which 488 are hydrogens and 0 are deuteriums.

• Molecule 1 is a protein called ADR1.

Mol	Chain	Residues	Atoms				Trace		
1	٨	60	Total	С	Η	Ν	0	S	0
	A	60	987	306	488	105	84	4	U

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	140	ALA	CYS	engineered mutation	UNP P07248

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

Μ	ol	Chain	Residues	Aton	ns
2		А	2	Total 2	Zn 2

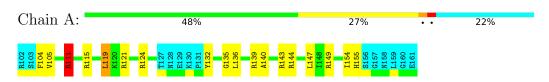


4 Residue-property plots (i)

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide 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: ADR1



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

The representative model is number 3. Colouring as in section 4.1 above.

• Molecule 1: ADR1

Chain A:	47%	23%	7%• 22%	
R102 F104 V105 R111 R111 R115	L119 R121 R121 R124 R124 F128 F129 F129 F132 C135 C135 C135 C135 C135 R139 R139 R139	R143 R144 R144 R144 I147 R149 R149 R153 H155 S155 S155 S155	615 N158 6160 E161	



5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *DISTANCE GEOMETRY, SIMULATED ANNEALING*.

Of the 50 calculated structures, 25 were deposited, based on the following criterion: LOWEST ENERGY TERM.

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

Software name	Classification	Version
X-PLOR	refinement	3.1
NMRPIPES	structure solution	
X-PLOR	structure solution	

No chemical shift data was provided.



6 Model quality (i)

6.1 Standard geometry (i)

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

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	А	$0.0{\pm}0.0$	$8.0{\pm}0.0$
All	All	0	200

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

5 of 8 unique planar outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Group	Models (Total)
1	А	111	ARG	Sidechain	25
1	А	115	ARG	Sidechain	25
1	А	121	ARG	Sidechain	25
1	А	124	ARG	Sidechain	25
1	А	139	ARG	Sidechain	25

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	А	402	397	397	8±1
All	All	10100	9925	9925	200

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 10.

Atom-1	Atom-2	Clash(Å)	$\operatorname{Distance}(\operatorname{\AA})$	Models	
Atom-1	Atom-2	Clash(A)		Worst	Total
1:A:136:LEU:HD12	1:A:155:HIS:CD2	0.78	2.14	16	18
1:A:108:VAL:HG21	1:A:126:HIS:HB3	0.71	1.60	1	3
1:A:105:VAL:HG23	1:A:111:ARG:O	0.70	1.86	20	24
1:A:147:LEU:O	1:A:147:LEU:HD23	0.69	1.87	9	22
1:A:132:TYR:O	1:A:147:LEU:HD12	0.68	1.89	13	11

5 of 40 unique clashes are listed below, sorted by their clash magnitude.

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	А	47/60~(78%)	$41\pm2~(87\pm4\%)$	$6\pm2~(13\pm4\%)$	0±0 (0±1%)	44 80
All	All	1175/1500 (78%)	1017 (87%)	154 (13%)	4 (0%)	44 80

All 4 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	А	111	ARG	1
1	А	114	ALA	1
1	А	134	CYS	1
1	А	135	GLY	1

6.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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 side chain conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	А	42/53~(79%)	$38\pm2(90\pm4\%)$	$4\pm2~(10\pm4\%)$	11 57

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1050/1325~(79%)	946 (90%)	104 (10%)	11 57

5 of 20 unique residues with a non-rotameric side chain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	А	119	LEU	14
1	А	111	ARG	14
1	А	149	ARG	12
1	А	152	GLN	8
1	А	123	TYR	8

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 monosaccharides in this entry.

6.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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 (i)

No chemical shift data were provided

