



# wwPDB NMR Structure Validation Summary Report ⓘ

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PDB ID : 1R36  
Title : NMR-based structure of autoinhibited murine Ets-1 deltaN301  
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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  
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

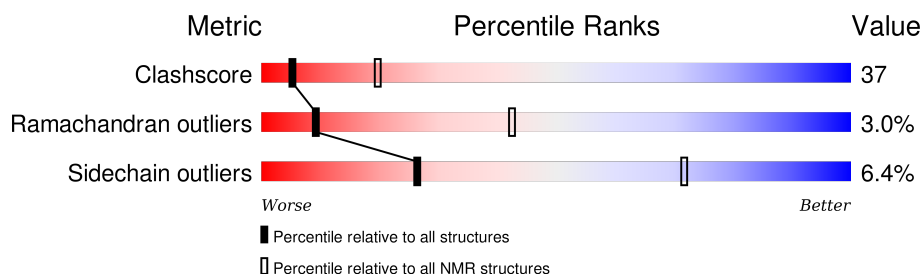
# 1 Overall quality at a glance

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  $\geq 3$ , 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	140	 41% 49% • 7%

## 2 Ensemble composition and analysis

This entry contains 25 models. Model 10 is the overall representative, medoid model (most similar to other models). The authors have identified model 12 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:302-A:378, A:384-A:436 (130)	0.19	10

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 4 single-model clusters were found.

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

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2291 atoms, of which 1143 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called C-ets-1 protein.

Mol	Chain	Residues	Atoms						Trace
1	A	140	Total	C	H	N	O	S	0
			2291	737	1143	198	210	3	

There is a discrepancy between the modelled and reference sequences:

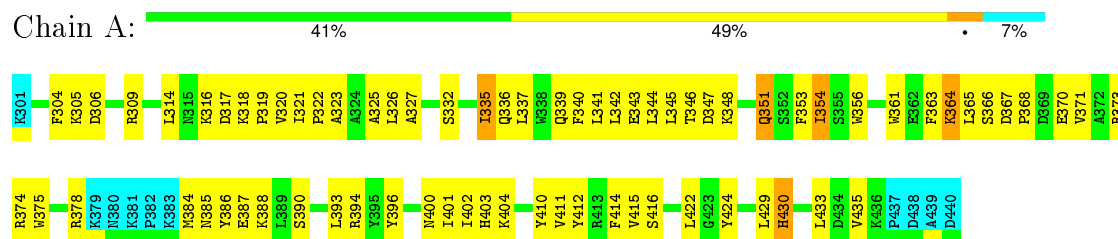
Chain	Residue	Modelled	Actual	Comment	Reference
A	416	SER	CYS	ENGINEERED	UNP P27577

## 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.

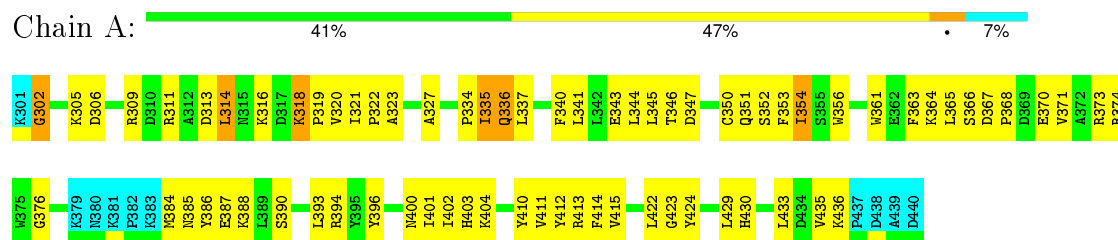
- Molecule 1: C-ets-1 protein



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

The representative model is number 10. Colouring as in section 4.1 above.

- Molecule 1: C-ets-1 protein



## 5 Refinement protocol and experimental data overview

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

Of the 50 calculated structures, 25 were deposited, based on the following criterion: *structures with the lowest energy*.

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

Software name	Classification	Version
Aria	refinement	1.2
CNS	refinement	1.1

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 5991
Number of chemical shift lists	1
Total number of shifts	1838
Number of shifts mapped to atoms	1838
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

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 (average) 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	0.46±0.04	1±1/1097 (0.1±0.1%)	0.51±0.01	0±0/1484 (0.0±0.0%)
All	All	0.47	24/27425 (0.1%)	0.51	0/37100 (0.0%)

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	0.0±0.2
All	All	0	1

5 of 7 unique bond outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	412	TYR	CE1-CZ	8.11	1.49	1.38	7	8
1	A	412	TYR	CE2-CZ	-7.47	1.28	1.38	7	9
1	A	424	TYR	CE1-CZ	-5.49	1.31	1.38	21	1
1	A	395	TYR	CE2-CZ	5.23	1.45	1.38	2	2
1	A	395	TYR	CE1-CZ	-5.07	1.31	1.38	2	2

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	412	TYR	Sidechain	1

## 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	1068	1056	1054	78±5
All	All	26700	26400	26350	1943

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 37.

5 of 313 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:320:VAL:HB	1:A:346:THR:HB	0.99	1.35	1	25
1:A:319:PRO:HG2	1:A:343:GLU:HB2	0.85	1.46	10	7
1:A:319:PRO:HB3	1:A:347:ASP:HB2	0.82	1.51	6	25
1:A:361:TRP:HB3	1:A:414:PHE:HB2	0.82	1.51	16	25
1:A:306:ASP:HA	1:A:309:ARG:HB3	0.81	1.51	6	15

## 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	130/140 (93%)	113±2 (87±2%)	13±2 (10±1%)	4±1 (3±1%)	9	42
All	All	3250/3500 (93%)	2837 (87%)	314 (10%)	99 (3%)	9	42

5 of 13 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	335	ILE	25
1	A	354	ILE	25
1	A	317	ASP	11

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Mol	Chain	Res	Type	Models (Total)
1	A	302	GLY	8
1	A	318	LYS	7

### 6.3.2 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 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	112/121 (93%)	105±2 (94±2%)	7±2 (6±2%)	26	72
All	All	2800/3025 (93%)	2621 (94%)	179 (6%)	26	72

5 of 34 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	430	HIS	20
1	A	351	GLN	18
1	A	385	ASN	14
1	A	336	GLN	14
1	A	332	SER	13

### 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

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

### 7.1 Chemical shift list 1

File name: BMRB entry 5991

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	1838
Number of shifts mapped to atoms	1838
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	12

#### 7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	140	$-0.33 \pm 0.08$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}_\beta$	130	$0.44 \pm 0.19$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}'$	127	$-0.28 \pm 0.09$	None needed ( $< 0.5$ ppm)
$^{15}\text{N}$	139	$0.66 \pm 0.23$	Should be applied

#### 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. 1542 atoms were assigned a chemical shift out of a possible 1677. 21 out of 21 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^1\text{H}$	$^{13}\text{C}$	$^{15}\text{N}$
Backbone	628/640 (98%)	255/255 (100%)	249/260 (96%)	124/125 (99%)
Sidechain	745/856 (87%)	459/505 (91%)	274/306 (90%)	12/45 (27%)

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	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Aromatic	169/181 (93%)	87/93 (94%)	78/80 (98%)	4/8 (50%)
Overall	1542/1677 (92%)	801/853 (94%)	601/646 (93%)	140/178 (79%)

#### 7.1.4 Statistically unusual chemical shifts [i](#)

The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	379	LYS	HD3	-0.15	2.75 – 0.45	-7.6
1	A	379	LYS	HD2	0.14	2.76 – 0.46	-6.4
1	A	426	PRO	HG3	-0.09	3.56 – 0.26	-6.1
1	A	339	GLN	CG	26.99	39.38 – 28.18	-6.1
1	A	379	LYS	CD	22.14	34.86 – 23.06	-5.8
1	A	351	GLN	CD	172.16	186.62 – 172.72	-5.4
1	A	336	GLN	CD	172.17	186.62 – 172.72	-5.4
1	A	426	PRO	HA	2.64	6.05 – 2.75	-5.3
1	A	339	GLN	CD	172.53	186.62 – 172.72	-5.1
1	A	345	LEU	HD11	-0.67	2.16 – -0.64	-5.1
1	A	345	LEU	HD13	-0.67	2.16 – -0.64	-5.1
1	A	345	LEU	HD12	-0.67	2.16 – -0.64	-5.1

#### 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:

