

Full wwPDB NMR Structure Validation Report (i)

Jun 6, 2023 – 01:38 pm BST

PDB ID	:	5NDA
BMRB ID	:	34114
Title	:	NMR Structural Characterisation of Pharmaceutically Relevant Proteins Ob-
		tained Through a Novel Recombinant Production: The Case of The Pulmonary
		Surfactant Polypeptide C Analogue rSP-C33Leu.
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Deposited on	:	2017-03-08

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 https://www.wwpdb.org/validation/2017/NMRValidationReportHelp with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

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

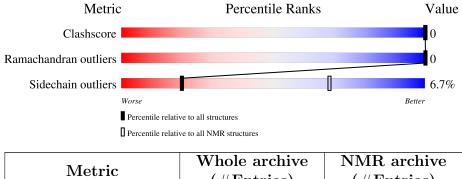
wwPDB-RCI	:	4.02b-467 20191225.v01 (using entries in the PDB archive December 25th 2019) $v_1n_11_5_13_A$ (Berjanski et al., 2005)
		Wang et al. (2010)
wwPDB-ShiftChecker BMRB Restraints Analysis		
Ideal geometry (proteins)		
Ideal geometry (DNA, RNA)	:	0
Validation Pipeline (wwPDB-VP)	:	2.33

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 is 68%.

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	(#Entries)	(#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	33	61%	•	36%	



2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 1 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:5-A:25 (21)	0.31	1		

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 4 clusters. No single-model clusters were found.

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



3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 574 atoms, of which 321 are hydrogens and 0 are deuteriums.

• Molecule 1 is a protein called rSP-C33Leu -RECOMBINANT PULMONARY SURFACTA NT-ASSOCIATED POLYPEPTIDE C ANALOGUE-.

Mol	Chain	Residues	Atoms				Trace	
1	Δ	33	Total	С	Η	Ν	0	0
	A		574	178	321	40	35	0



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.

 \bullet Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-



4.2 Scores per residue for each member of the ensemble

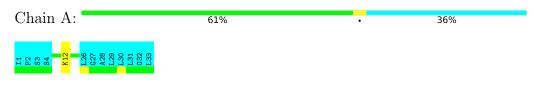
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1 (medoid)

 \bullet Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-



4.2.2 Score per residue for model 2





4.2.3 Score per residue for model 3

• Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-



4.2.4 Score per residue for model 4

 \bullet Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-

Chain A:	61%	·	36%
11 13 24 25 24 24 25 12 26 12 26 12 26 12 26 27 12 30 27 12 30 23 23 23 23 23 23 23 23 23 24 25 26 27 26 27 26 26 26 26 26 26 26 26 26 26 26 26 26			

4.2.5 Score per residue for model 5

 \bullet Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-

Chain A:		55%	9%	36%
11 P2 S3 H7 H7 H7	K11 K12 C27 C27 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C27 C28 C28 C28 C28 C28 C28 C28 C28 C28 C28			

4.2.6 Score per residue for model 6

• Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-



4.2.7 Score per residue for model 7



Chain A: 58% 6% 36%

4.2.8 Score per residue for model 8

• Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-

Chain A:	61%	•	36%
11 8 23 8 35 8 4 8 4 8 4 8 2 1 2 6 1 2 0 1 2 0 1			

4.2.9 Score per residue for model 9

• Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-

Chain A:	61%	·	36%
11 82 84 84 84 126 128 128 128 123 133 133 133 133 133			

4.2.10 Score per residue for model 10

• Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-



4.2.11 Score per residue for model 11





4.2.12 Score per residue for model 12

• Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-



4.2.13 Score per residue for model 13

 \bullet Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-

Chain A:	58%	6%	36%
11 83 84 84 84 84 128 128 128 128 128 128 128 128 128 128			

4.2.14 Score per residue for model 14

 \bullet Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-

Chain A:	61%	·	36%	
11 P2 S4 K12 K12 K12 C31 L30 L30 L31 L31 L33 L33 L33 L33 L33				

4.2.15 Score per residue for model 15

• Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-



4.2.16 Score per residue for model 16



36%

Chain A: 61%



4.2.17 Score per residue for model 17

• Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-

Chain A:	61%	·	36%
11 82 83 83 84 84 84 82 828 129 129 133 133 133			

4.2.18 Score per residue for model 18

• Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-

Chain A:	58%	6%	36%
11 82 83 84 84 84 84 84 84 84 84 84 84 84 84 84			

4.2.19 Score per residue for model 19

 \bullet Molecule 1: rSP-C33Leu -RECOMBINANT PULMONARY SURFACTANT-ASSOCIATED POLYPEPTIDE C ANALOGUE-



4.2.20 Score per residue for model 20





5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *distance geometry*.

Of the 100 calculated structures, 20 were deposited, based on the following criterion: target function.

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

Software name Classification		Version
TALOS	geometry optimization	
CYANA	structure calculation	3.97
OPAL	refinement	3.97

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)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	367
Number of shifts mapped to atoms	367
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	68%



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
All	All	3460	4540	4540	-

The all-atom clash score is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clash score for this structure is -.

There are no clashes.

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	Perce	ntiles
1	А	21/33~(64%)	21±0 (100±0%)	0±0 (0±0%)	0±0 (0±0%)	100	100
All	All	420/660~(64%)	420 (100%)	0 (0%)	0 (0%)	100	100

There are no Ramachandran outliers.



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	Perce	entiles
1	А	21/30~(70%)	20 ± 1 (93 $\pm3\%$)	$1 \pm 1 (7 \pm 3\%)$	20	68
All	All	420/600 (70%)	392 (93%)	28~(7%)	20	68

All 4 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	А	12	LYS	20
1	А	10	ARG	6
1	А	7	HIS	1
1	А	9	LYS	1

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)

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 (i)

The completeness of assignment taking into account all chemical shift lists is 68% for the well-defined parts and 68% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: D_1200003777.bmrb

7.1.1 Bookkeeping (i)

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	367
Number of shifts mapped to atoms	367
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

7.1.2 Chemical shift referencing (i)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	$\textbf{Correction} \pm \textbf{precision}, \textit{ppm}$	Suggested action
$^{13}C_{\alpha}$	31	0.92 ± 0.14	Should be checked
$^{13}C_{\beta}$	5		None (insufficient data)
$^{13}C'$	0		None (insufficient data)
¹⁵ N	30	0.69 ± 0.19	Should be applied

7.1.3 Completeness of resonance assignments (i)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 68%, i.e. 252 atoms were assigned a chemical shift out of a possible 371. 0 out of 14 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^{1}\mathrm{H}$	$^{13}\mathrm{C}$	$^{15}\mathbf{N}$
Backbone	81/103~(79%)	41/41~(100%)	20/42~(48%)	20/20~(100%)
Sidechain	169/261~(65%)	162/177~(92%)	6/79~(8%)	1/5~(20%)

Continued on next page...



	Total	$^{1}\mathbf{H}$	$^{13}\mathrm{C}$	$^{15}\mathbf{N}$
Aromatic	2/7~(29%)	2/4~(50%)	0/2~(0%)	0/1~(0%)
Overall	252/371~(68%)	205/222~(92%)	26/123~(21%)	21/26~(81%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 68%, i.e. 361 atoms were assigned a chemical shift out of a possible 528. 0 out of 19 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^{1}\mathrm{H}$	$^{13}\mathrm{C}$	$^{15}\mathbf{N}$
Backbone	126/163~(77%)	65/66~(98%)	31/66~(47%)	30/31~(97%)
Sidechain	233/358~(65%)	222/244~(91%)	10/109~(9%)	1/5~(20%)
Aromatic	2/7~(29%)	2/4~(50%)	0/2~(0%)	0/1~(0%)
Overall	361/528~(68%)	289/314~(92%)	41/177~(23%)	31/37~(84%)

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.

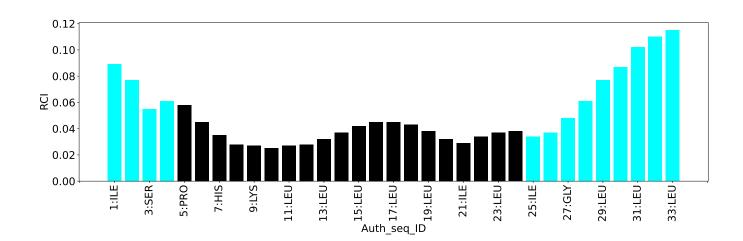
List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	А	10	ARG	NE	116.57	76.53 - 92.65	19.8

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. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:







8 NMR restraints analysis (i)

8.1 Conformationally restricting restraints (i)

The following table provides the summary of experimentally observed NMR restraints in different categories. Restraints are classified into different categories based on the sequence separation of the atoms involved.

Description	Value
Total distance restraints	259
Intra-residue (i-j =0)	157
Sequential (i-j =1)	55
Medium range ($ i-j >1$ and $ i-j <5$)	47
Long range $(i-j \ge 5)$	0
Inter-chain	0
Hydrogen bond restraints	0
Disulfide bond restraints	0
Total dihedral-angle restraints	0
Number of unmapped restraints	0
Number of restraints per residue	7.8
Number of long range restraints per residue ¹	0.0

¹Long range hydrogen bonds and disulfide bonds are counted as long range restraints while calculating the number of long range restraints per residue

8.2 Residual restraint violations (i)

This section provides the overview of the restraint violations analysis. The violations are binned as small, medium and large violations based on its absolute value. Average number of violations per model is calculated by dividing the total number of violations in each bin by the size of the ensemble.

8.2.1 Average number of distance violations per model (i)

Distance violations less than 0.1 Å are not included in the calculation. There are no distance violations

8.2.2 Average number of dihedral-angle violations per model (i)

Dihedral-angle violations less than 1° are not included in the calculation. There are no dihedral-angle violations



9 Distance violation analysis (i)

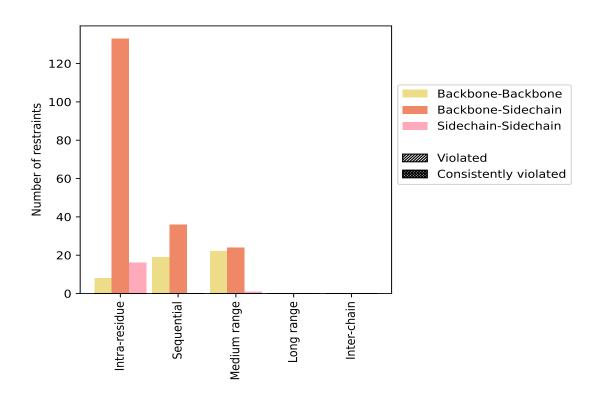
9.1 Summary of distance violations (i)

The following table shows the summary of distance violations in different restraint categories based on the sequence separation of the atoms involved. Each category is further sub-divided into three sub-categories based on the atoms involved. Violations less than 0.1 Å are not included in the statistics.

Destroints torms	Count	071	$_1$ Violated ³			Consistently Violated ⁴		
Restraints type	Count	$\%^1$	Count	$\%^2$	$\%^1$	Count	$\%^2$	$\%^1$
Intra-residue (i-j =0)	157	60.6	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	8	3.1	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	133	51.4	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	16	6.2	0	0.0	0.0	0	0.0	0.0
Sequential (i-j =1)	55	21.2	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	19	7.3	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	36	13.9	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Medium range ($ i-j > 1 \& i-j < 5$)	47	18.1	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	22	8.5	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	24	9.3	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	1	0.4	0	0.0	0.0	0	0.0	0.0
Long range $(i-j \ge 5)$	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Inter-chain	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Hydrogen bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Disulfide bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Total	259	100.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	49	18.9	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	193	74.5	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	17	6.6	0	0.0	0.0	0	0.0	0.0

 1 percentage calculated with respect to the total number of distance restraints, 2 percentage calculated with respect to the number of restraints in a particular restraint category, 3 violated in at least one model, 4 violated in all the models





9.1.1 Bar chart : Distribution of distance restraints and violations (i)

Violated and consistently violated restraints are shown using different hatch patterns in their respective categories. The hydrogen bonds and disulfied bonds are counted in their appropriate category on the x-axis

9.2 Distance violation statistics for each model (i)

No violations found

9.3 Distance violation statistics for the ensemble (i)

No violations found

9.4 Most violated distance restraints in the ensemble (i)

No violations found

9.5 All violated distance restraints (i)

No violations found



10 Dihedral-angle violation analysis (i)

Dihedral angle analysis failed due to data error in the dihedral angle restraints, possibly missing target value

