



wwPDB NMR Structure Validation Summary Report ⓘ

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PDB ID : 2LVZ
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Title : Solution structure of a Eosinophil Cationic Protein-trisaccharide heparin mimetic complex
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<https://www.wwpdb.org/validation/2017/NMRValidationReportHelp>

with specific help available everywhere you see the ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 2022.3.0, CSD as543be (2022)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
wwPDB-RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
wwPDB-ShiftChecker : v1.2
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

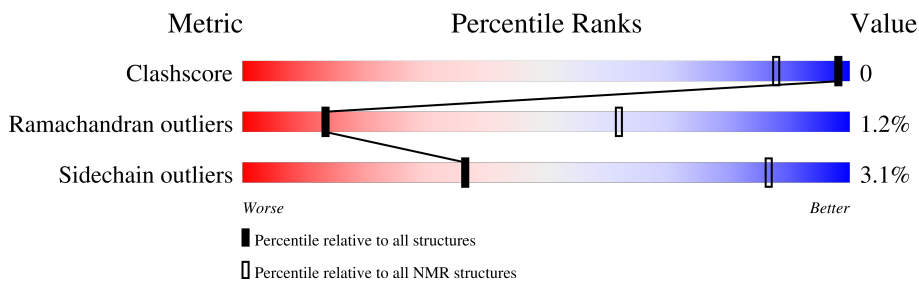
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 77%.

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	210492	14027
Ramachandran outliers	207382	12486
Sidechain outliers	206894	12463

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	133	 77% 22% .
2	B	3	 100%

2 Ensemble composition and analysis

This entry contains 20 models. Model 13 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:2-A:133 (132)	0.75	13

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 and 6 single-model clusters were found.

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

3 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 2259 atoms, of which 1107 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Eosinophil cationic protein.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	133	2169	677	1075	224	184	9	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	97	ARG	THR	conflict	UNP P12724

- Molecule 2 is an oligosaccharide called 2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-propan-2-yl 2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranoside.

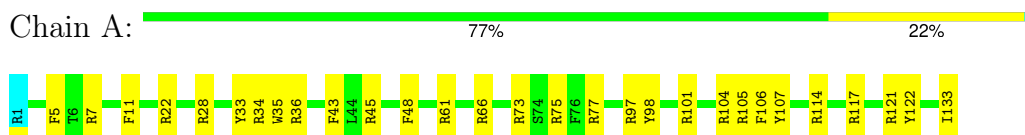
Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
2	B	3	90	21	32	2	30	5	0


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.

- Molecule 1: Eosinophil cationic protein



Chain B:  100%

LVZ1
LDS2
SGH3

5 Refinement protocol and experimental data overview

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

Of the 100 calculated structures, 20 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
Amber	structure solution	
CYANA	structure solution	
Amber	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)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	1481
Number of shifts mapped to atoms	1481
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	77%

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: IDS, SGN, LVZ

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	1.56±0.03	5±2/1113 (0.5± 0.2%)	2.17±0.09	42±5/1515 (2.7± 0.3%)
All	All	1.56	107/22260 (0.5%)	2.17	830/30300 (2.7%)

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	5.8±2.1
All	All	0	115

5 of 81 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	23	CYS	CB-SG	-7.98	1.68	1.82	19	1
1	A	63	PRO	N-CD	-7.97	1.36	1.47	12	2
1	A	61	ARG	CZ-NH1	-7.52	1.23	1.33	6	4
1	A	33	TYR	CG-CD1	7.24	1.48	1.39	17	1
1	A	22	ARG	CZ-NH1	-7.21	1.23	1.33	14	1

5 of 252 unique angle 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(°)	Ideal(°)	Models	
								Worst	Total
1	A	77	ARG	NE-CZ-NH1	24.71	132.65	120.30	5	10
1	A	34	ARG	NE-CZ-NH1	24.68	132.64	120.30	13	10
1	A	105	ARG	NE-CZ-NH1	22.62	131.61	120.30	17	14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	97	ARG	NE-CZ-NH1	21.13	130.86	120.30	2	12
1	A	117	ARG	NE-CZ-NH1	20.34	130.47	120.30	7	13

There are no chirality outliers.

5 of 44 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	A	107	TYR	Sidechain	9
1	A	5	PHE	Sidechain,Peptide	8
1	A	22	ARG	Sidechain	5
1	A	11	PHE	Sidechain	4
1	A	34	ARG	Sidechain,Mainchain	4

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	1083	1060	1055	1±1
2	B	58	32	12	0±0
All	All	22820	21840	21347	13

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 0.

5 of 12 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:107:TYR:HB3	1:A:132:THR:HG22	0.53	1.79	13	1
1:A:80:LEU:HD12	1:A:80:LEU:C	0.51	2.25	19	1
1:A:128:HIS:CD2	1:A:129:LEU:H	0.47	2.28	17	1
1:A:60:ILE:HG13	1:A:71:CYS:SG	0.46	2.49	20	1
1:A:23:CYS:SG	1:A:100:ASP:HA	0.44	2.52	1	2

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	131/133 (98%)	113±4 (87±3%)	16±3 (12±3%)	2±1 (1±1%)	14	62
All	All	2620/2660 (98%)	2267 (87%)	322 (12%)	31 (1%)	14	62

5 of 21 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	86	ILE	3
1	A	24	THR	3
1	A	102	PRO	3
1	A	88	PRO	2
1	A	89	GLY	2

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	121/122 (99%)	117±2 (97±1%)	4±2 (3±1%)	37	86
All	All	2420/2440 (99%)	2344 (97%)	76 (3%)	37	86

5 of 28 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	133	ILE	20
1	A	35	TRP	9
1	A	30	ILE	6
1	A	101	ARG	5
1	A	80	LEU	3

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](#)

3 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length is the number of standard deviations the observed value is removed from the expected value. A bond length with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

Mol	Type	Chain	Res	Link	Bond lengths		
					Counts	RMSZ	#Z>2
2	LVZ	B	1	2	23,23,23	1.89±0.30	6±1 (25±5%)
2	IDS	B	2	2	16,16,17	1.89±0.34	4±1 (25±9%)
2	SGN	B	3	2	19,19,20	1.90±0.45	5±2 (25±8%)

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of angles that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond angle is the number of standard deviations the observed value is removed from the expected value. A bond angle with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
2	LVZ	B	1	2	28,35,35	1.88±0.28	6±2 (22±7%)
2	IDS	B	2	2	16,24,26	1.69±0.37	3±1 (20±9%)
2	SGN	B	3	2	23,29,31	2.05±0.33	7±2 (28±7%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means

no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	LVZ	B	1	2	-	0±0,15,35,35	0±0,1,1,1
2	IDS	B	2	2	-	0±0,9,26,29	0±0,1,1,1
2	SGN	B	3	2	-	0±0,11,28,31	0±0,1,1,1

5 of 43 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
2	B	3	SGN	S1-N2	8.36	1.70	1.59	1	17
2	B	1	LVZ	S19-N9	8.16	1.69	1.59	3	12
2	B	3	SGN	O1S-S1	7.12	1.50	1.42	4	7
2	B	2	IDS	O5-C1	6.76	1.55	1.43	8	12
2	B	2	IDS	C1-C2	6.43	1.62	1.51	10	9

5 of 56 unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

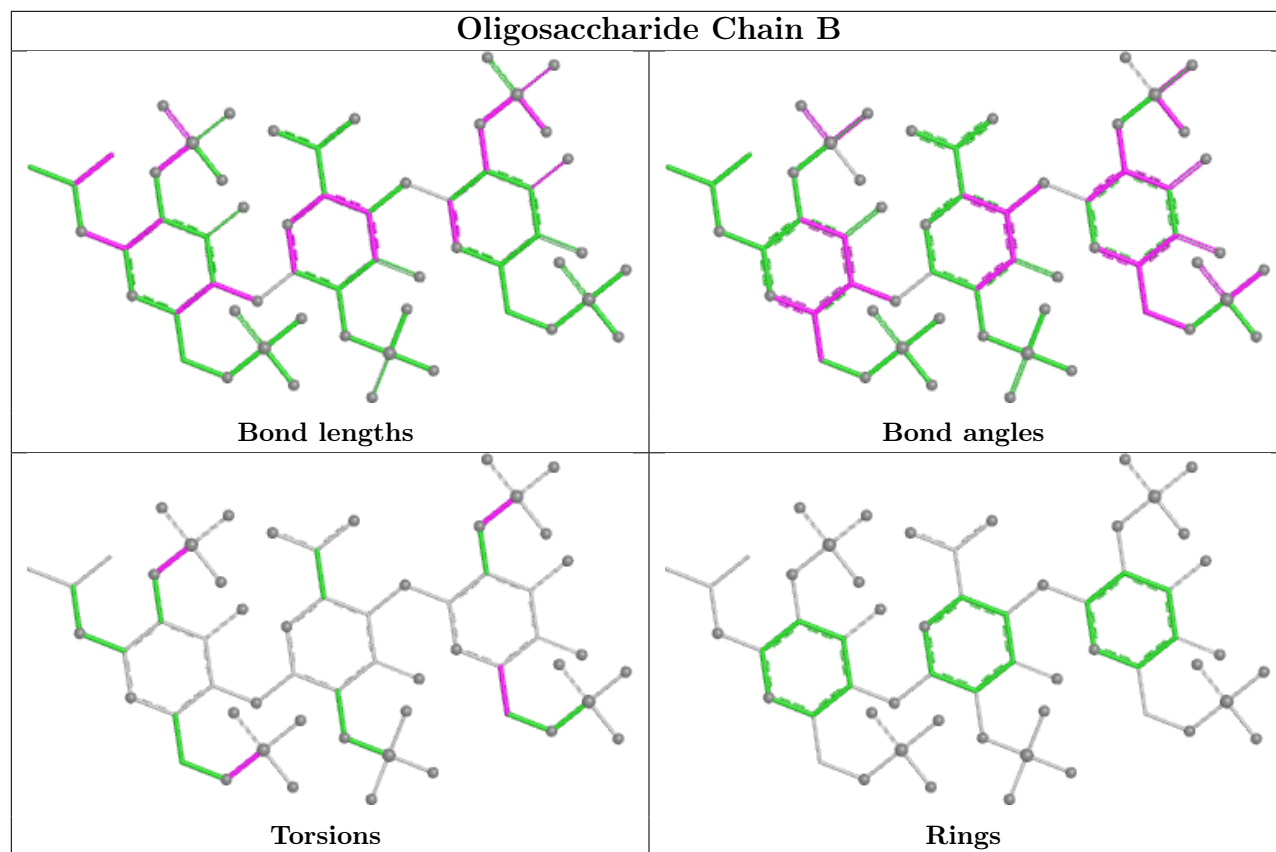
Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
2	B	3	SGN	C1-O5-C5	8.98	124.22	112.19	10	16
2	B	1	LVZ	O3S-S19-N9	7.85	95.75	108.88	4	14
2	B	2	IDS	O2-C2-C3	7.49	117.42	106.95	10	9
2	B	2	IDS	C4-C3-C2	6.52	122.06	110.23	10	12
2	B	1	LVZ	O2S-S19-O3S	6.45	106.01	120.36	6	20

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



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 77% for the well-defined parts and 77% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_1*

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	1481
Number of shifts mapped to atoms	1481
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	8

7.1.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	129	-0.00 ± 0.15	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	129	-0.09 ± 0.10	None needed (< 0.5 ppm)
$^{13}\text{C}'$	0	—	None (insufficient data)
^{15}N	120	1.10 ± 0.53	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 77%, i.e. 1470 atoms were assigned a chemical shift out of a possible 1898. 0 out of 16 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	503/639 (79%)	255/255 (100%)	128/264 (48%)	120/120 (100%)
Sidechain	888/1099 (81%)	599/707 (85%)	263/316 (83%)	26/76 (34%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	79/160 (49%)	62/78 (79%)	15/70 (21%)	2/12 (17%)
Overall	1470/1898 (77%)	916/1040 (88%)	406/650 (62%)	148/208 (71%)

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	A	117	ARG	H	0.30	5.25 – 11.22	-13.3
1	A	12	ALA	CB	38.40	10.19 – 27.75	11.1
1	A	14	GLN	HG3	0.58	0.91 – 3.68	-6.2
1	A	126	PRO	HB3	0.07	0.25 – 3.76	-5.5
1	A	14	GLN	HB2	0.68	0.80 – 3.29	-5.5
1	A	74	SER	HB3	2.44	2.49 – 5.20	-5.2
1	A	27	MET	HB3	0.28	0.33 – 3.66	-5.2
1	A	77	ARG	HG3	0.15	0.15 – 2.94	-5.0

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:

