



wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 12, 2024 – 02:47 PM EST

PDB ID : 3DSL
Title : The Three-dimensional Structure of Bothropasin, the Main Hemorrhagic Factor from Bothrops jararaca venom.
Authors : Muniz, J.R.C.; Ambrosio, A.; Selistre-de-Araujo, H.S.; Oliva, G.; Garratt, R.C.; Souza, D.H.F.
Deposited on : 2008-07-13
Resolution : 2.70 Å(reported)

This is a wwPDB X-ray 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/XrayValidationReportHelp>

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)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

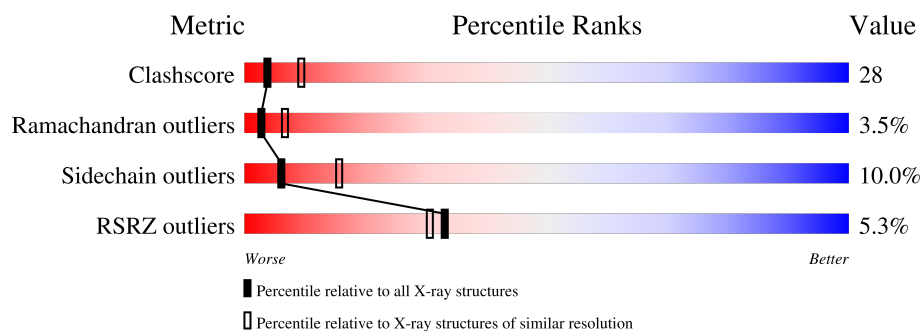
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

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)	Similar resolution (#Entries, resolution range(Å))
Clashscore	180529	3684 (2.70-2.70)
Ramachandran outliers	177936	3633 (2.70-2.70)
Sidechain outliers	177891	3633 (2.70-2.70)
RSRZ outliers	164620	3333 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	419	
1	B	419	

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 6675 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Zinc metalloproteinase-disintegrin bothropasin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	416	Total	C	N	O	S	0	0	0
			3219	1998	543	631	47			
1	B	412	Total	C	N	O	S	0	0	0
			3190	1981	538	624	47			

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Ca	0	0
			3	3		

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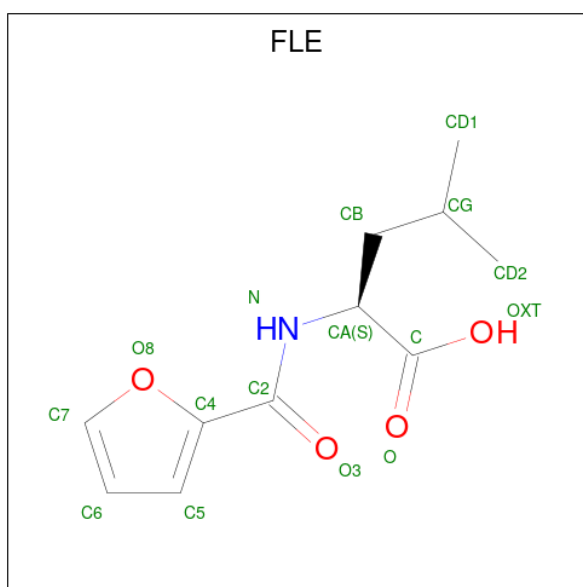
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	3	Total	Ca	0	0
			3	3		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

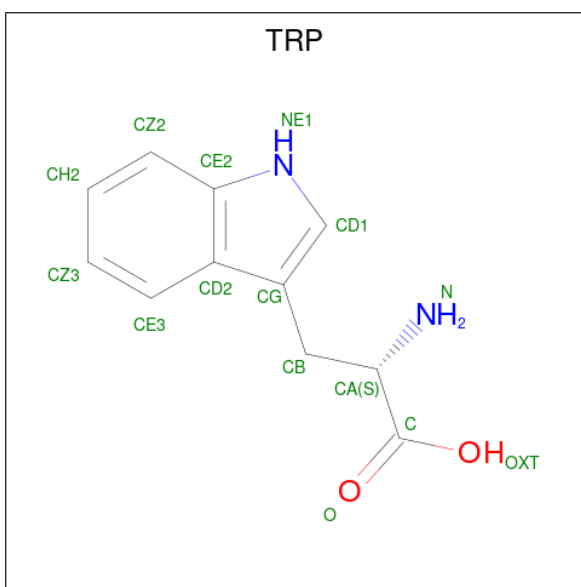
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		
4	B	1	Total	Zn	0	0
			1	1		

- Molecule 5 is FUROYL-LEUCINE (three-letter code: FLE) (formula: C₁₁H₁₅NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			15	11	1	3		
5	A	1	Total	C	N	O	0	0
			15	11	1	3		

- Molecule 6 is TRYPTOPHAN (three-letter code: TRP) (formula: C₁₁H₁₂N₂O₂).

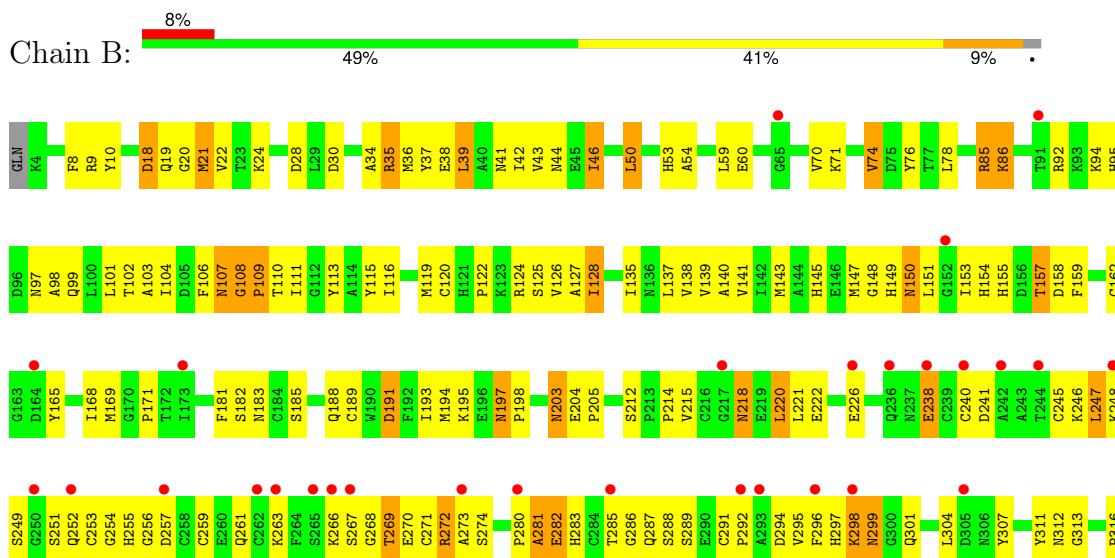


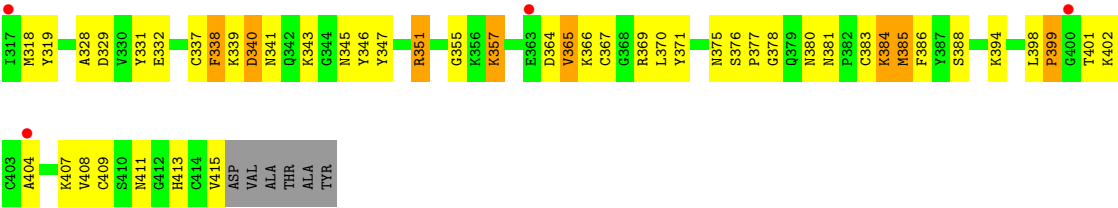
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			15	11	2	2		
6	A	1	Total	C	N	O	0	0
			15	11	2	2		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	113	Total	O	0	0
			113	113		
7	B	71	Total	O	0	0
			71	71		

- Molecule 1: Zinc metalloproteinase-disintegrin bothropasin





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	73.76Å 100.27Å 133.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.44 – 2.70 30.44 – 2.70	Depositor EDS
% Data completeness (in resolution range)	88.9 (30.44-2.70) 91.9 (30.44-2.70)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.68Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.210 , 0.294 0.232 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	46.7	Xtriage
Anisotropy	0.142	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 84.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	6675	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.10% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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 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	0.42	0/3294	0.58	0/4458
1	B	0.43	1/3265 (0.0%)	0.58	1/4418 (0.0%)
All	All	0.43	1/6559 (0.0%)	0.58	1/8876 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	355	GLY	C-N	-12.36	1.05	1.34

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	355	GLY	O-C-N	5.76	131.91	122.70

There are no chirality outliers.

There are no planarity outliers.

5.2 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 the 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 within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3219	0	2997	157	0
1	B	3190	0	2971	199	0
2	A	14	0	13	1	0
3	A	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	3	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	30	0	28	3	0
6	A	30	0	20	1	0
7	A	113	0	0	3	0
7	B	71	0	0	7	0
All	All	6675	0	6029	350	0

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 28.

The worst 5 of 350 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:357:LYS:HD3	1:B:357:LYS:N	1.52	1.13
1:A:353:GLU:OE1	1:A:353:GLU:HA	1.42	1.06
1:A:121:HIS:CD2	1:A:123:LYS:H	1.76	1.03
1:A:121:HIS:HD2	1:A:123:LYS:H	1.01	0.99
1:B:357:LYS:HD3	1:B:357:LYS:H	1.12	0.99

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 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 X-ray entries followed by that with respect to entries of similar resolution.

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	414/419 (99%)	351 (85%)	54 (13%)	9 (2%)	5	15
1	B	410/419 (98%)	320 (78%)	70 (17%)	20 (5%)	2	3
All	All	824/838 (98%)	671 (81%)	124 (15%)	29 (4%)	3	7

5 of 29 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	ASP
1	A	249	SER
1	B	107	ASN
1	B	157	THR
1	B	281	ALA

5.3.2 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 X-ray entries followed by that with respect to entries of similar resolution.

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	361/363 (99%)	323 (90%)	38 (10%)	5	14
1	B	358/363 (99%)	326 (91%)	32 (9%)	8	20
All	All	719/726 (99%)	649 (90%)	70 (10%)	6	17

5 of 70 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	253	CYS
1	B	272	ARG
1	B	357	LYS
1	A	260	GLU
1	A	252	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 42 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	99	GLN
1	B	255	HIS
1	B	150	ASN
1	B	197	ASN
1	B	299	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 8 are monoatomic - leaving 5 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or 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 length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	TRP	A	507	4,5	14,16,16	0.92	1 (7%)	13,22,22	1.18	1 (7%)
6	TRP	A	505	4,5	14,16,16	0.98	1 (7%)	13,22,22	1.28	1 (7%)
5	FLE	A	504	6	11,15,16	1.71	3 (27%)	12,19,21	3.73	6 (50%)
5	FLE	A	506	6	11,15,16	1.96	4 (36%)	12,19,21	2.12	4 (33%)
2	NAG	A	501	1	14,14,15	1.69	2 (14%)	17,19,21	2.09	5 (29%)

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
6	TRP	A	507	4,5	-	1/7/8/8	0/2/2/2
6	TRP	A	505	4,5	-	1/7/8/8	0/2/2/2
5	FLE	A	504	6	-	2/10/14/16	0/1/1/1
5	FLE	A	506	6	-	3/10/14/16	0/1/1/1
2	NAG	A	501	1	-	4/6/23/26	0/1/1/1

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	NAG	C1-C2	4.48	1.58	1.52
5	A	504	FLE	CB-CA	3.50	1.58	1.53
5	A	506	FLE	CB-CA	3.30	1.57	1.53
5	A	506	FLE	C2-N	3.30	1.41	1.34
5	A	506	FLE	C4-C2	3.27	1.54	1.49

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	504	FLE	C4-C2-N	8.01	130.58	115.19
5	A	504	FLE	CB-CA-N	6.60	120.65	110.69
2	A	501	NAG	C1-O5-C5	5.46	119.50	112.19
5	A	506	FLE	C4-C2-N	5.40	125.56	115.19
5	A	504	FLE	CB-CA-C	4.50	117.91	110.99

There are no chirality outliers.

5 of 11 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	504	FLE	N-CA-CB-CG
5	A	504	FLE	C-CA-CB-CG
5	A	506	FLE	N-CA-CB-CG
5	A	506	FLE	C-CA-CB-CG
5	A	506	FLE	O-C-CA-CB

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	505	TRP	1	0
5	A	504	FLE	1	0
5	A	506	FLE	2	0
2	A	501	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	355:GLY	C	356:LYS	N	1.05

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	416/419 (99%)	0.02	11 (2%) 57 56	22, 38, 78, 180	0
1	B	412/419 (98%)	0.36	33 (8%) 20 18	24, 49, 129, 235	0
All	All	828/838 (98%)	0.19	44 (5%) 33 31	22, 43, 105, 235	0

The worst 5 of 44 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	240	CYS	6.4
1	A	361	ALA	6.1
1	B	273	ALA	5.8
1	B	262	CYS	4.5
1	B	363	GLU	4.3

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NAG	A	501	14/15	0.82	0.12	42,45,52,61	0
5	FLE	A	506	15/16	0.88	0.10	42,42,42,42	0
6	TRP	A	507	15/15	0.88	0.13	53,53,53,53	0
5	FLE	A	504	15/16	0.92	0.10	24,24,24,24	0
3	CA	B	424	1/1	0.97	0.04	72,72,72,72	0
6	TRP	A	505	15/15	0.97	0.06	16,16,16,16	0
3	CA	A	1	1/1	0.97	0.07	19,19,19,19	0
3	CA	B	423	1/1	0.98	0.04	34,34,34,34	0
3	CA	A	502	1/1	0.98	0.02	34,34,34,34	0
3	CA	A	2	1/1	0.99	0.05	26,26,26,26	0
3	CA	B	422	1/1	0.99	0.03	24,24,24,24	0
4	ZN	A	503	1/1	0.99	0.11	33,33,33,33	0
4	ZN	B	1	1/1	0.99	0.06	48,48,48,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.