



# Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 09:41 am BST

PDB ID : 5NDE  
Title : Crystal structure of metallo-beta-lactamase SPM-1 in space group P4222  
Authors : Hinchliffe, P.; Spencer, J.  
Deposited on : 2017-03-08  
Resolution : 1.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

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:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

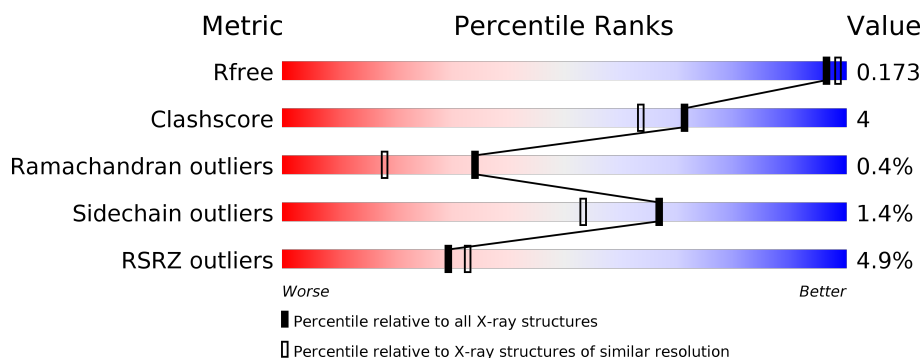
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.70 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))
$R_{free}$	130704	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	250	<div> <div>5%</div> <div> <div></div> <div>89%</div> <div>8%</div> <div></div> </div> </div>
1	B	250	<div> <div>5%</div> <div> <div></div> <div>94%</div> <div>5%</div> <div></div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	GOL	A	409	-	-	X	-

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4449 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 Beta-lactamase IMP-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	244	Total	C	N	O	S	0	1	0
			1933	1239	328	358	8			
1	B	247	Total	C	N	O	S	0	1	0
			1953	1251	331	362	9			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	29	GLY	-	expression tag	UNP Q8G9Q0
A	30	PRO	-	expression tag	UNP Q8G9Q0
B	29	GLY	-	expression tag	UNP Q8G9Q0
B	30	PRO	-	expression tag	UNP Q8G9Q0

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Zn	0	0
			2	2		
2	A	4	Total	Zn	0	0
			4	4		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 6 3 3	0	0
4	A	1	Total C O 6 3 3	0	0
4	A	1	Total C O 6 3 3	0	0
4	A	1	Total C O 6 3 3	0	0
4	B	1	Total C O 6 3 3	0	0

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total Cl 2 2	0	0

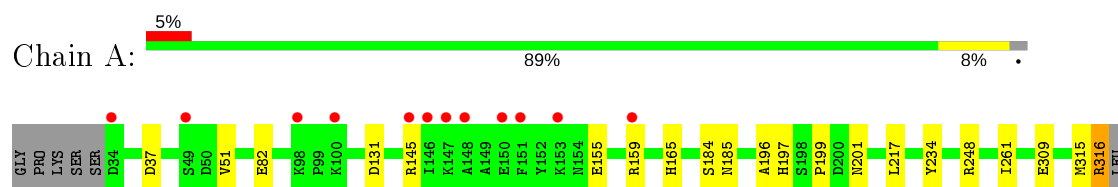
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	255	Total O 255 255	0	0
6	B	250	Total O 250 250	0	0

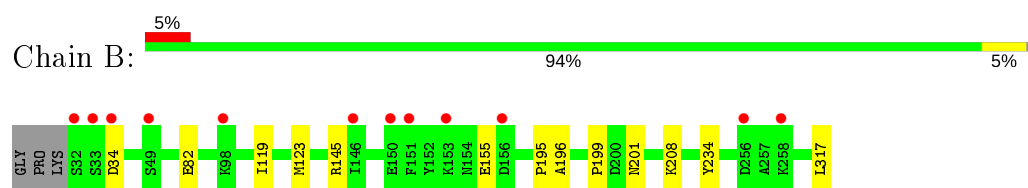
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-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. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

#### • Molecule 1: Beta-lactamase IMP-1



#### • Molecule 1: Beta-lactamase IMP-1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 42 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	130.74Å 130.74Å 100.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.26 – 1.70 36.26 – 1.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (36.26-1.70) 100.0 (36.26-1.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.13 (at 1.70Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.162 , 0.172 0.164 , 0.173	Depositor DCC
$R_{free}$ test set	4772 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.2	Xtriage
Anisotropy	0.351	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 48.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	4449	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.86% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: GOL, ZN, SO4, CL

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	0.38	0/1977	0.54	1/2670 (0.0%)
1	B	0.36	0/1997	0.53	0/2695
All	All	0.37	0/3974	0.54	1/5365 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	131	ASP	CB-CG-OD1	5.04	122.84	118.30

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	1933	0	1950	16	0
1	B	1953	0	1974	10	0
2	A	4	0	0	0	0
2	B	2	0	0	0	0
3	A	15	0	0	1	0
3	B	5	0	0	0	0
4	A	24	0	31	8	0
4	B	6	0	8	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	2	0	0	0	0
6	A	255	0	0	6	2
6	B	250	0	0	7	1
All	All	4449	0	3963	28	2

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 4.

All (28) 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:A:248:ARG:NH1	1:A:309[A]:GLU:OE2	2.21	0.74
1:A:315:MET:HA	6:B:519:HOH:O	1.88	0.73
1:A:37:ASP:OD1	6:A:501:HOH:O	2.06	0.72
4:A:409:GOL:H2	1:B:199:PRO:HB3	1.72	0.70
1:A:234:TYR:H	4:A:411:GOL:H32	1.55	0.70
1:B:234:TYR:H	4:B:404:GOL:H31	1.59	0.67
4:A:409:GOL:H11	6:B:519:HOH:O	1.97	0.64
1:A:315:MET:HB2	4:A:409:GOL:H12	1.84	0.58
1:A:316:ARG:HB2	6:A:537:HOH:O	2.04	0.56
1:A:315:MET:HG3	1:B:195:PRO:HG3	1.88	0.56
3:A:407:SO4:O1	6:A:502:HOH:O	2.13	0.55
1:A:217:LEU:HD11	1:A:261:ILE:HD11	1.91	0.52
1:A:315:MET:HE2	4:A:409:GOL:H12	1.91	0.51
1:A:155:GLU:HG3	1:A:159:ARG:HE	1.78	0.49
1:B:145:ARG:NH1	6:B:506:HOH:O	2.45	0.49
1:A:145:ARG:NH2	6:A:504:HOH:O	2.32	0.48
1:A:37:ASP:HB2	6:A:512:HOH:O	2.14	0.47
1:B:34:ASP:HA	6:B:620:HOH:O	2.13	0.47
4:A:410:GOL:O1	4:A:410:GOL:O3	2.31	0.46
1:A:184:SER:O	1:A:185:ASN:HB2	2.17	0.45
4:A:408:GOL:H12	6:A:647:HOH:O	2.15	0.45
1:A:199:PRO:HG3	1:B:317:LEU:HD23	1.98	0.45
1:A:197:HIS:CE1	4:A:411:GOL:H12	2.53	0.44
1:A:51:VAL:HG11	1:A:217:LEU:HD21	2.00	0.44
1:B:119:ILE:O	1:B:123[B]:MET:HG3	2.19	0.43
1:B:155:GLU:HG3	6:B:518:HOH:O	2.18	0.43
1:B:208:LYS:HE3	6:B:636:HOH:O	2.19	0.41
1:B:145:ARG:HD3	6:B:695:HOH:O	2.21	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:698:HOH:O	6:A:717:HOH:O[7_554]	2.07	0.13
6:A:586:HOH:O	6:B:646:HOH:O[6_555]	2.19	0.01

## 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	243/250 (97%)	238 (98%)	4 (2%)	1 (0%)	34	18
1	B	246/250 (98%)	240 (98%)	5 (2%)	1 (0%)	34	18
All	All	489/500 (98%)	478 (98%)	9 (2%)	2 (0%)	34	18

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	196	ALA
1	B	196	ALA

### 5.3.2 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 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	213/217 (98%)	209 (98%)	4 (2%)	57	41
1	B	216/217 (100%)	214 (99%)	2 (1%)	78	70
All	All	429/434 (99%)	423 (99%)	6 (1%)	67	53

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	82	GLU
1	A	165	HIS
1	A	201	ASN
1	A	316	ARG
1	B	82	GLU
1	B	201	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 17 ligands modelled in this entry, 8 are monoatomic - leaving 9 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	$\# Z  > 2$
4	GOL	A	408	-	5,5,5	0.35	0	5,5,5	0.70	0
4	GOL	A	410	-	5,5,5	0.36	0	5,5,5	0.25	0
3	SO4	A	405	-	4,4,4	0.14	0	6,6,6	0.12	0
3	SO4	A	406	-	4,4,4	0.14	0	6,6,6	0.08	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	GOL	A	409	-	5,5,5	0.25	0	5,5,5	0.69	0
3	SO4	A	407	-	4,4,4	0.14	0	6,6,6	0.06	0
3	SO4	B	403	-	4,4,4	0.14	0	6,6,6	0.09	0
4	GOL	A	411	2	5,5,5	0.30	0	5,5,5	0.24	0
4	GOL	B	404	-	5,5,5	0.32	0	5,5,5	0.54	0

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
4	GOL	A	408	-	-	2/4/4/4	-
4	GOL	A	410	-	-	2/4/4/4	-
4	GOL	A	409	-	-	1/4/4/4	-
4	GOL	B	404	-	-	0/4/4/4	-
4	GOL	A	411	2	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	410	GOL	O1-C1-C2-C3
4	A	408	GOL	O1-C1-C2-C3
4	A	410	GOL	O1-C1-C2-O2
4	A	409	GOL	O1-C1-C2-O2
4	A	408	GOL	O1-C1-C2-O2

There are no ring outliers.

6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	408	GOL	1	0
4	A	410	GOL	1	0
4	A	409	GOL	4	0
3	A	407	SO4	1	0
4	A	411	GOL	2	0
4	B	404	GOL	1	0

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	244/250 (97%)	-0.02	12 (4%)	29 33	17, 26, 65, 98	0
1	B	247/250 (98%)	-0.01	12 (4%)	29 33	16, 26, 56, 92	1 (0%)
All	All	491/500 (98%)	-0.01	24 (4%)	29 33	16, 26, 61, 98	1 (0%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	148	ALA	5.9
1	B	32	SER	5.9
1	A	146	ILE	5.6
1	B	153	LYS	4.4
1	B	146	ILE	4.0
1	A	151	PHE	3.9
1	A	150	GLU	3.7
1	B	33	SER	3.3
1	A	147	LYS	3.1
1	A	145	ARG	2.9
1	B	34	ASP	2.8
1	A	100	LYS	2.8
1	B	49	SER	2.5
1	B	150	GLU	2.5
1	A	159	ARG	2.4
1	A	153	LYS	2.4
1	A	34	ASP	2.3
1	B	151	PHE	2.2
1	B	258	LYS	2.2
1	A	98	LYS	2.1
1	B	98	LYS	2.1
1	B	156	ASP	2.1
1	B	256	ASP	2.1
1	A	49	SER	2.0

## 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 carbohydrates 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
4	GOL	A	411	6/6	0.71	0.23	28,49,61,67	0
4	GOL	A	409	6/6	0.78	0.26	25,28,32,36	6
4	GOL	A	410	6/6	0.80	0.17	75,79,80,83	0
3	SO4	A	406	5/5	0.81	0.19	75,78,93,95	0
5	CL	A	412	1/1	0.82	0.10	74,74,74,74	0
3	SO4	A	407	5/5	0.84	0.15	101,102,108,109	0
4	GOL	A	408	6/6	0.84	0.17	29,45,54,56	0
4	GOL	B	404	6/6	0.88	0.20	30,66,72,76	0
5	CL	A	413	1/1	0.90	0.09	68,68,68,68	0
3	SO4	B	403	5/5	0.96	0.10	75,79,82,85	0
3	SO4	A	405	5/5	0.96	0.12	53,61,67,70	0
2	ZN	A	404	1/1	0.98	0.06	25,25,25,25	0
2	ZN	A	403	1/1	0.99	0.08	23,23,23,23	1
2	ZN	B	402	1/1	0.99	0.06	27,27,27,27	1
2	ZN	A	402	1/1	1.00	0.09	21,21,21,21	0
2	ZN	A	401	1/1	1.00	0.04	20,20,20,20	0
2	ZN	B	401	1/1	1.00	0.08	21,21,21,21	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.