



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 12, 2026 – 08:06 PM UTC

PDB ID : 2IJD / pdb\_00002ijd  
Title : Crystal Structure of the Poliovirus Precursor Protein 3CD  
Authors : Marcotte, L.L.; Gohara, D.W.; Filman, D.J.; Hogle, J.M.  
Deposited on : 2006-09-29  
Resolution : 3.40 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](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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

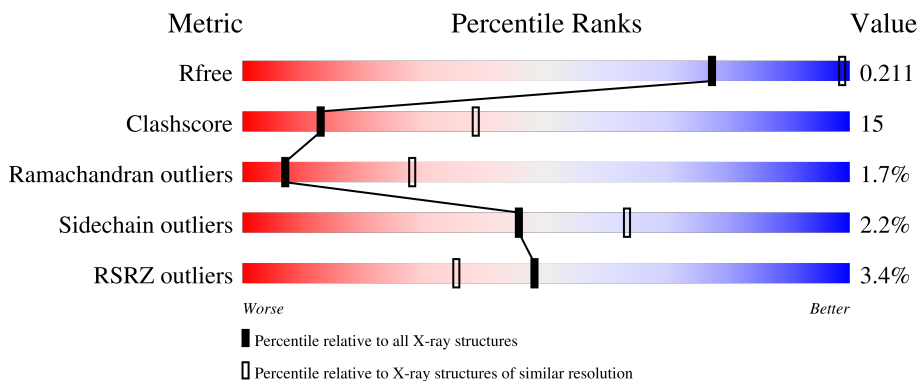
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


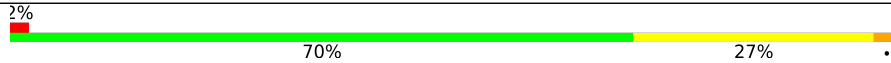
The reported resolution of this entry is 3.40 Å.

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))
$R_{free}$	180053	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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  $\geq 3$ , 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	1	644	
1	2	644	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 10170 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 Picornain 3C, RNA-directed RNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	1	644	5060	3226	850	956	28	0	0	0
1	2	644	5060	3226	850	956	28	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
1	55	ALA	GLU	engineered mutation	UNP P03300
1	58	ALA	ASP	engineered mutation	UNP P03300
1	63	ALA	GLU	engineered mutation	UNP P03300
1	147	ALA	CYS	engineered mutation	UNP P03300
1	629	ASP	LEU	engineered mutation	UNP P03300
1	638	ASP	ARG	engineered mutation	UNP P03300
2	55	ALA	GLU	engineered mutation	UNP P03300
2	58	ALA	ASP	engineered mutation	UNP P03300
2	63	ALA	GLU	engineered mutation	UNP P03300
2	147	ALA	CYS	engineered mutation	UNP P03300
2	629	ASP	LEU	engineered mutation	UNP P03300
2	638	ASP	ARG	engineered mutation	UNP P03300

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	1	1	Total	Zn	0	0
			1	1		
2	2	1	Total	Zn	0	0
			1	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	1	1	Total O S 5 4 1	0	0
3	1	1	Total O S 5 4 1	0	0
3	1	1	Total O S 5 4 1	0	0
3	1	1	Total O S 5 4 1	0	0
3	2	1	Total O S 5 4 1	0	0
3	2	1	Total O S 5 4 1	0	0
3	2	1	Total O S 5 4 1	0	0
3	2	1	Total O S 5 4 1	0	0

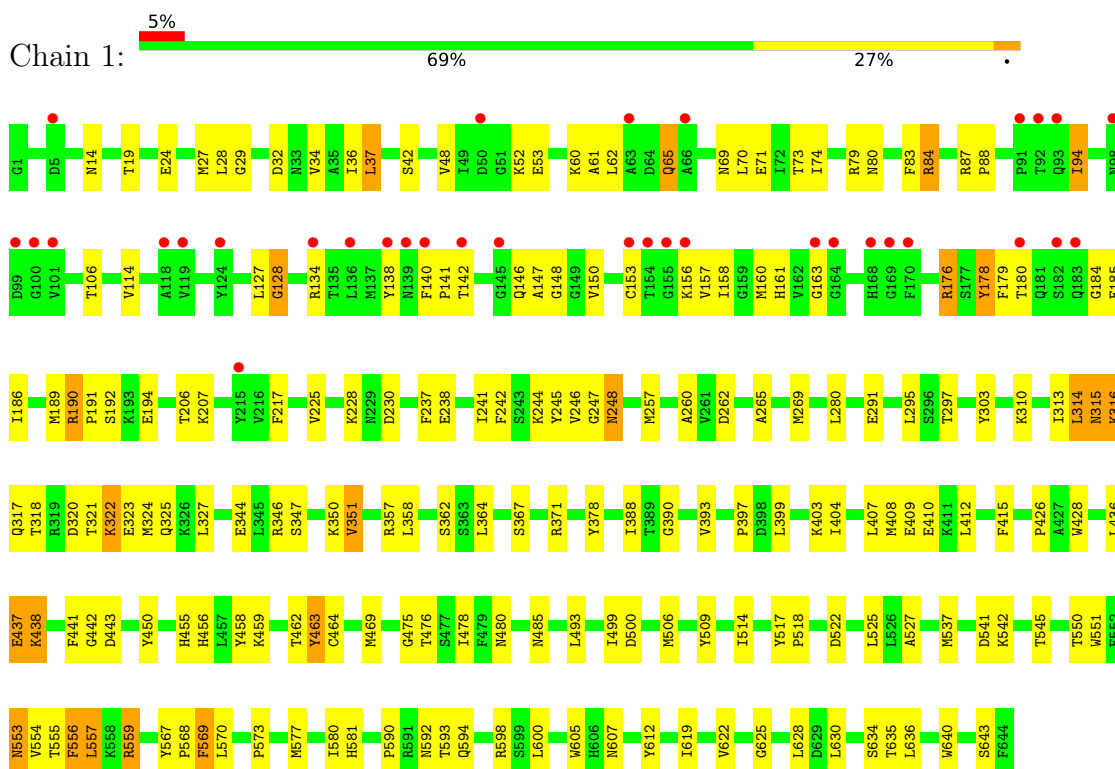
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	1	4	Total O 4 4	0	0
4	2	4	Total O 4 4	0	0

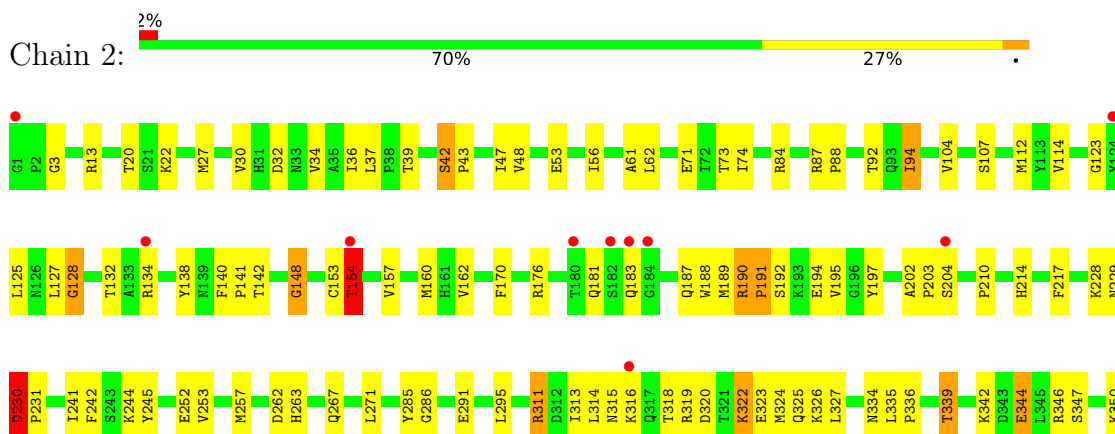
### 3 Residue-property plots

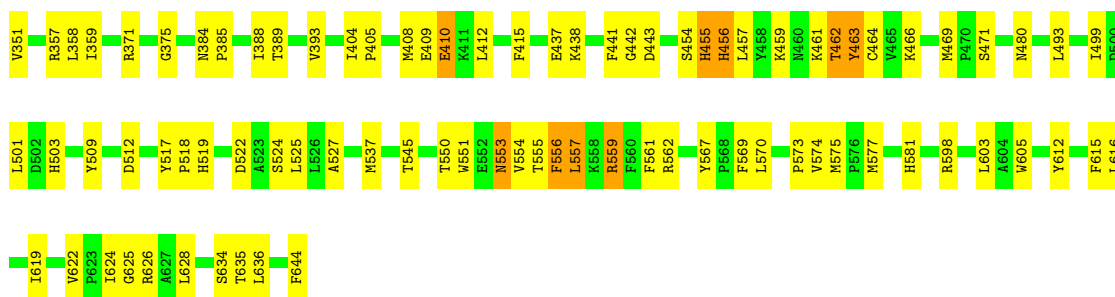
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Picornain 3C, RNA-directed RNA polymerase



- Molecule 1: Picornain 3C, RNA-directed RNA polymerase





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	208.54Å 230.44Å 151.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.90 – 3.40 19.90 – 3.40	Depositor EDS
% Data completeness (in resolution range)	98.5 (19.90-3.40) 97.5 (19.90-3.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.90 (at 3.44Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.199 , 0.231 0.177 , 0.211	Depositor DCC
$R_{free}$ test set	2510 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	94.1	Xtrriage
Anisotropy	0.051	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 61.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	10170	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.65% 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: SO4, ZN

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	1	0.85	0/5175	1.28	43/7006 (0.6%)
1	2	0.85	0/5175	1.26	45/7006 (0.6%)
All	All	0.85	0/10350	1.27	88/14012 (0.6%)

There are no bond length outliers.

The worst 5 of 88 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	388	ILE	N-CA-C	-9.04	103.79	111.56
1	1	351	VAL	N-CA-C	-9.02	103.10	111.67
1	1	190	ARG	CA-C-N	8.82	130.87	119.84
1	1	190	ARG	C-N-CA	8.82	130.87	119.84
1	2	622	VAL	CA-C-N	8.41	127.98	119.24

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	1	5060	0	5028	153	0
1	2	5060	0	5028	154	0
2	1	1	0	0	0	0
2	2	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	1	20	0	0	1	0
3	2	20	0	0	2	0
4	1	4	0	0	0	0
4	2	4	0	0	0	0
All	All	10170	0	10056	306	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 15.

The worst 5 of 306 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)
1:1:228:LYS:H	1:1:228:LYS:HD2	1.10	1.16
1:2:39:THR:HG22	1:2:73:THR:HG23	1.27	1.12
1:2:228:LYS:H	1:2:228:LYS:HD2	0.96	1.07
1:1:37:LEU:HA	1:1:160:MET:HE1	1.28	1.06
1:2:409:GLU:HG3	1:2:410:GLU:H	1.20	1.02

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	1	642/644 (100%)	574 (89%)	56 (9%)	12 (2%)	6	26
1	2	642/644 (100%)	569 (89%)	63 (10%)	10 (2%)	7	29
All	All	1284/1288 (100%)	1143 (89%)	119 (9%)	22 (2%)	7	28

5 of 22 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1	128	GLY
1	1	178	TYR
1	1	316	LYS
1	2	134	ARG
1	2	322	LYS

### 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	1	546/546 (100%)	534 (98%)	12 (2%)	45 63
1	2	546/546 (100%)	534 (98%)	12 (2%)	45 63
All	All	1092/1092 (100%)	1068 (98%)	24 (2%)	45 63

5 of 24 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	2	138	TYR
1	2	311	ARG
1	2	187	GLN
1	2	334	ASN
1	1	397	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 27 such sidechains are listed below:

Mol	Chain	Res	Type
1	1	607	ASN
1	2	122	GLN
1	2	456	HIS
1	2	105	ASN
1	2	131	GLN

### 5.3.3 RNA [i](#)

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 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 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	$\# Z  > 2$
3	SO4	1	649	-	4,4,4	0.28	0	6,6,6	0.28	0
3	SO4	2	647	-	4,4,4	0.29	0	6,6,6	0.22	0
3	SO4	1	646	-	4,4,4	0.25	0	6,6,6	0.24	0
3	SO4	2	648	-	4,4,4	0.37	0	6,6,6	0.29	0
3	SO4	1	648	-	4,4,4	0.30	0	6,6,6	0.27	0
3	SO4	2	649	-	4,4,4	0.31	0	6,6,6	0.19	0
3	SO4	2	646	-	4,4,4	0.22	0	6,6,6	0.17	0
3	SO4	1	647	-	4,4,4	0.30	0	6,6,6	0.34	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	1	649	SO4	1	0
3	2	648	SO4	1	0
3	2	649	SO4	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 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, 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	1	644/644 (100%)	0.13	34 (5%) 32 24	11, 37, 61, 102	0
1	2	644/644 (100%)	-0.21	10 (1%) 70 56	13, 37, 61, 110	0
All	All	1288/1288 (100%)	-0.04	44 (3%) 48 35	11, 37, 61, 110	0

The worst 5 of 44 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	1	98	ASN	5.5
1	2	182	SER	4.6
1	2	180	THR	4.6
1	1	169	GLY	4.2
1	1	154	THR	4.1

### 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 oligosaccharides 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
3	SO4	1	648	5/5	0.53	0.25	109,112,114,116	0
3	SO4	2	648	5/5	0.76	0.18	77,78,88,89	0
3	SO4	1	647	5/5	0.77	0.18	84,85,88,89	0
3	SO4	1	646	5/5	0.79	0.15	104,104,106,106	0
3	SO4	2	647	5/5	0.80	0.18	101,104,106,108	0
3	SO4	2	646	5/5	0.83	0.15	118,120,122,124	0
3	SO4	1	649	5/5	0.90	0.18	87,88,92,92	0
3	SO4	2	649	5/5	0.92	0.17	88,95,96,97	0
2	ZN	2	645	1/1	0.99	0.09	35,35,35,35	0
2	ZN	1	645	1/1	1.00	0.13	25,25,25,25	0

## 6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.