



# wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 13, 2024 – 10:16 am BST

PDB ID : 2V3Q  
Title : Serendipitous discovery and X-ray structure of a human phosphate binding apolipoprotein  
Authors : Morales, R.; Berna, A.; Carpentier, P.; Elias, M.; Contreras-Martel, C.; Renault, F.; Nicodeme, M.; Chesne-Seck, M.-L.; Bernier, F.; Dupuy, J.; Schaeffer, C.; Diemer, H.; Van Dorselaer, A.; Fontecilla, J.C.; Masson, P.; Rochu, D.; Chabriere, E.  
Deposited on : 2007-06-20  
Resolution : 1.89 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

<https://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	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
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)

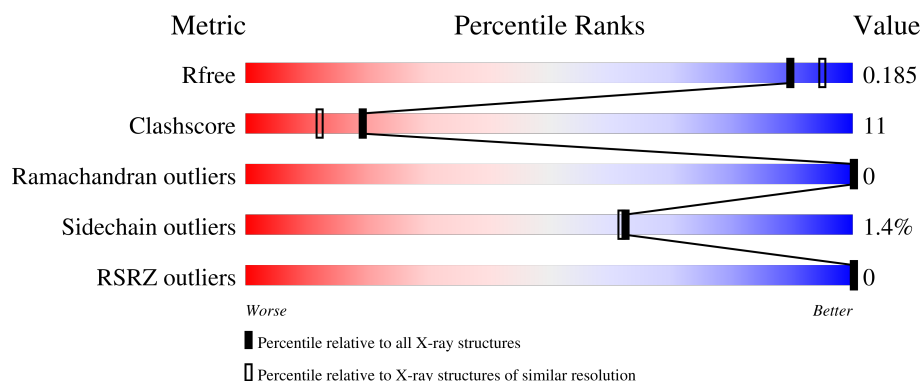
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*


The reported resolution of this entry is 1.89 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))
$R_{free}$	164625	7293 (1.90-1.90)
Clashscore	180529	8090 (1.90-1.90)
Ramachandran outliers	177936	8022 (1.90-1.90)
Sidechain outliers	177891	8022 (1.90-1.90)
RSRZ outliers	164620	7292 (1.90-1.90)

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	A	376	

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
2	GOL	A	1377	-	X	-	-
2	GOL	A	1378	-	-	X	-
2	GOL	A	1382	-	X	X	-
4	EDO	A	1390	-	-	X	-
4	EDO	A	1391	-	X	-	-

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3559 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 HUMAN PHOSPHATE BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	376	Total	C	N	O	S	0	15	0
			2788	1768	471	544	5			

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



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

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

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

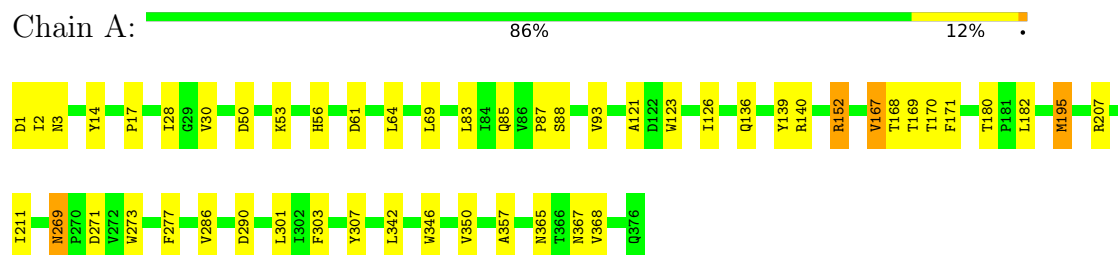
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	694	Total	O	0	0
			694	694		

### 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: HUMAN PHOSPHATE BINDING PROTEIN



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	96.25Å 86.84Å 89.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.95 – 1.89 44.95 – 1.89	Depositor EDS
% Data completeness (in resolution range)	99.6 (44.95-1.89) 99.6 (44.95-1.89)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.53 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.127 , 0.186 0.127 , 0.185	Depositor DCC
$R_{free}$ test set	1495 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	13.4	Xtriage
Anisotropy	0.008	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 63.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	3559	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.00% 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 ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PO4, EDO

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	1.37	8/2901 (0.3%)	1.00	6/3964 (0.2%)

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	307	TYR	CD1-CE1	5.67	1.47	1.39
1	A	350	VAL	CB-CG2	5.55	1.64	1.52
1	A	303	PHE	CE1-CZ	5.48	1.47	1.37
1	A	167	VAL	CB-CG2	5.32	1.64	1.52
1	A	14	TYR	CD2-CE2	5.25	1.47	1.39

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	50	ASP	CB-CG-OD2	-5.52	113.34	118.30
1	A	140	ARG	NE-CZ-NH2	-5.48	117.56	120.30
1	A	195	MET	CG-SD-CE	5.33	108.73	100.20
1	A	182[A]	LEU	CB-CA-C	-5.23	100.27	110.20
1	A	182[B]	LEU	CB-CA-C	-5.23	100.27	110.20

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts ⓘ

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	2788	0	2757	50	0
2	A	24	0	30	14	0
3	A	5	0	0	0	0
4	A	48	0	72	19	0
5	A	694	0	0	26	3
All	All	3559	0	2859	66	3

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 11.

The worst 5 of 66 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)
4:A:1390:EDO:H12	5:A:2684:HOH:O	1.33	1.25
2:A:1377:GOL:H12	5:A:2587:HOH:O	1.58	1.02
1:A:3:ASN:H	1:A:56:HIS:HD2	1.15	0.92
1:A:69:LEU:HB3	4:A:1387:EDO:H11	1.53	0.90
1:A:53:LYS:HG2	2:A:1382:GOL:H12	1.56	0.86

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:2341:HOH:O	5:A:2341:HOH:O[4_566]	1.51	0.69
5:A:2371:HOH:O	5:A:2683:HOH:O[6_555]	2.11	0.09
5:A:2294:HOH:O	5:A:2684:HOH:O[6_555]	2.18	0.02

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	389/376 (104%)	374 (96%)	15 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	293/278 (105%)	289 (99%)	4 (1%)	62 62

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	PRO
1	A	61	ASP
1	A	83	LEU
1	A	269	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	HIS
1	A	85	GLN
1	A	269	ASN
1	A	337	ASN
1	A	365	ASN

### 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 [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

17 ligands are modelled in this entry.

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	$\# Z  > 2$
2	GOL	A	1377	-	5,5,5	0.68	0	5,5,5	1.96	2 (40%)
2	GOL	A	1382	-	5,5,5	1.40	1 (20%)	5,5,5	1.94	3 (60%)
4	EDO	A	1395	-	3,3,3	0.49	0	2,2,2	0.22	0
4	EDO	A	1387	-	3,3,3	0.71	0	2,2,2	0.54	0
2	GOL	A	1378	-	5,5,5	0.41	0	5,5,5	0.58	0
4	EDO	A	1394	-	3,3,3	0.36	0	2,2,2	1.57	0
2	GOL	A	1393	-	5,5,5	0.41	0	5,5,5	1.27	1 (20%)
4	EDO	A	1388	-	3,3,3	0.60	0	2,2,2	0.41	0
4	EDO	A	1392	-	3,3,3	0.27	0	2,2,2	3.15	2 (100%)
4	EDO	A	1390	-	3,3,3	0.43	0	2,2,2	0.72	0
4	EDO	A	1384	-	3,3,3	0.59	0	2,2,2	0.63	0
3	PO4	A	1379	-	4,4,4	1.19	0	6,6,6	0.85	0
4	EDO	A	1389	-	3,3,3	0.42	0	2,2,2	0.04	0
4	EDO	A	1391	-	3,3,3	0.59	0	2,2,2	2.80	2 (100%)
4	EDO	A	1383	-	3,3,3	0.77	0	2,2,2	0.43	0
4	EDO	A	1386	-	3,3,3	0.61	0	2,2,2	0.42	0
4	EDO	A	1385	-	3,3,3	0.26	0	2,2,2	0.96	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
2	GOL	A	1377	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	1382	-	-	3/4/4/4	-
4	EDO	A	1395	-	-	1/1/1/1	-
4	EDO	A	1387	-	-	1/1/1/1	-
2	GOL	A	1378	-	-	2/4/4/4	-
4	EDO	A	1394	-	-	0/1/1/1	-
2	GOL	A	1393	-	-	4/4/4/4	-
4	EDO	A	1388	-	-	1/1/1/1	-
4	EDO	A	1392	-	-	0/1/1/1	-
4	EDO	A	1390	-	-	1/1/1/1	-
4	EDO	A	1384	-	-	1/1/1/1	-
4	EDO	A	1389	-	-	1/1/1/1	-
4	EDO	A	1391	-	-	1/1/1/1	-
4	EDO	A	1383	-	-	0/1/1/1	-
4	EDO	A	1386	-	-	1/1/1/1	-
4	EDO	A	1385	-	-	1/1/1/1	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1382	GOL	O2-C2	-2.25	1.36	1.43

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1392	EDO	O1-C1-C2	-3.19	88.98	111.91
4	A	1391	EDO	O2-C2-C1	-3.12	89.47	111.91
4	A	1392	EDO	O2-C2-C1	-3.11	89.56	111.91
2	A	1377	GOL	O1-C1-C2	-2.84	96.57	110.20
4	A	1391	EDO	O1-C1-C2	-2.44	94.36	111.91

There are no chirality outliers.

5 of 22 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1377	GOL	C1-C2-C3-O3
2	A	1393	GOL	C1-C2-C3-O3
2	A	1377	GOL	O1-C1-C2-C3
2	A	1378	GOL	C1-C2-C3-O3
2	A	1382	GOL	O1-C1-C2-C3

There are no ring outliers.

13 monomers are involved in 33 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1377	GOL	2	0
2	A	1382	GOL	7	0
4	A	1395	EDO	1	0
4	A	1387	EDO	1	0
2	A	1378	GOL	5	0
4	A	1394	EDO	2	0
4	A	1388	EDO	1	0
4	A	1390	EDO	5	0
4	A	1384	EDO	2	0
4	A	1389	EDO	3	0
4	A	1383	EDO	2	0
4	A	1386	EDO	1	0
4	A	1385	EDO	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	A	376/376 (100%)	-1.01	0 100 100	4, 10, 18, 30	15 (3%)

There are no RSRZ outliers to report.

### 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, 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(Å <sup>2</sup> )	Q<0.9
4	EDO	A	1391	4/4	0.81	0.16	24,26,29,39	0
2	GOL	A	1378	6/6	0.83	0.14	36,38,42,45	0
4	EDO	A	1389	4/4	0.85	0.14	40,40,43,44	0
2	GOL	A	1393	6/6	0.88	0.12	26,35,38,43	0
4	EDO	A	1395	4/4	0.89	0.15	24,28,28,34	4
2	GOL	A	1382	6/6	0.90	0.15	23,24,29,36	0
4	EDO	A	1386	4/4	0.90	0.12	34,40,41,48	0
4	EDO	A	1387	4/4	0.90	0.15	28,32,32,35	0
4	EDO	A	1384	4/4	0.91	0.14	26,31,40,45	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	EDO	A	1392	4/4	0.92	0.09	17,21,27,35	0
4	EDO	A	1390	4/4	0.93	0.11	27,30,31,32	0
4	EDO	A	1385	4/4	0.93	0.14	27,33,37,43	0
4	EDO	A	1388	4/4	0.93	0.08	20,28,30,41	0
4	EDO	A	1394	4/4	0.93	0.11	16,30,32,34	0
4	EDO	A	1383	4/4	0.93	0.09	17,20,26,29	0
2	GOL	A	1377	6/6	0.94	0.09	16,26,30,38	0
3	PO4	A	1379	5/5	1.00	0.03	4,5,5,7	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.