



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:18 AM GMT

PDB ID : 3E4C  
Title : Procaspase-1 zymogen domain crystal structure  
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Deposited on : 2008-08-11  
Resolution : 2.05 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  
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

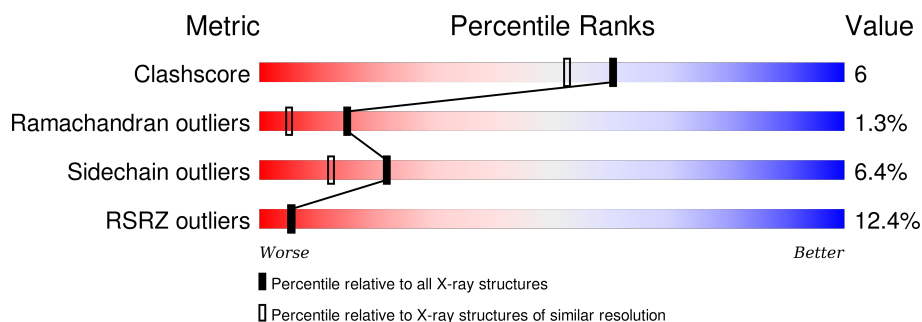
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.05 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	102246	1269 (2.04-2.04)
Ramachandran outliers	100387	1258 (2.04-2.04)
Sidechain outliers	100360	1258 (2.04-2.04)
RSRZ outliers	91569	1194 (2.04-2.04)

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. 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	302	
1	B	302	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4423 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 Caspase-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	Total	C	N	O	S	0	12	0
			2149	1351	371	410	17			
1	B	269	Total	C	N	O	S	0	12	0
			2137	1345	369	406	17			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	103	GLN	-	EXPRESSION TAG	UNP P29466
A	285	ALA	CYS	ENGINEERED	UNP P29466
B	103	GLN	-	EXPRESSION TAG	UNP P29466
B	285	ALA	CYS	ENGINEERED	UNP P29466

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		

- Molecule 3 is water.

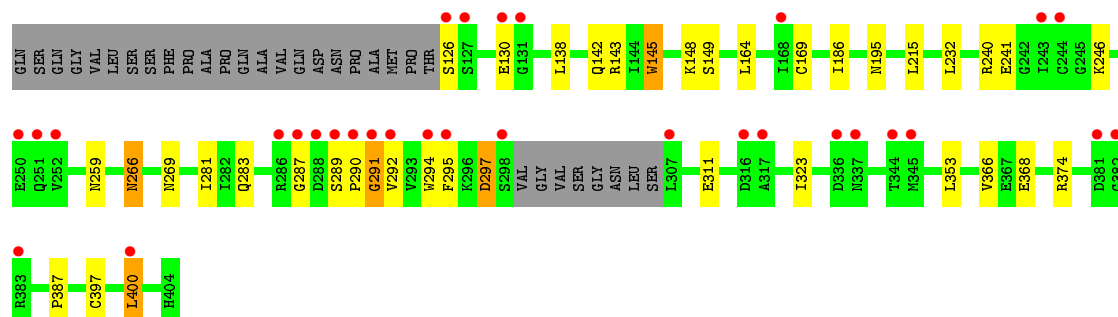
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	69	Total	O	0	0
			69	69		
3	B	67	Total	O	0	0
			67	67		

### 3 Residue-property plots


These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

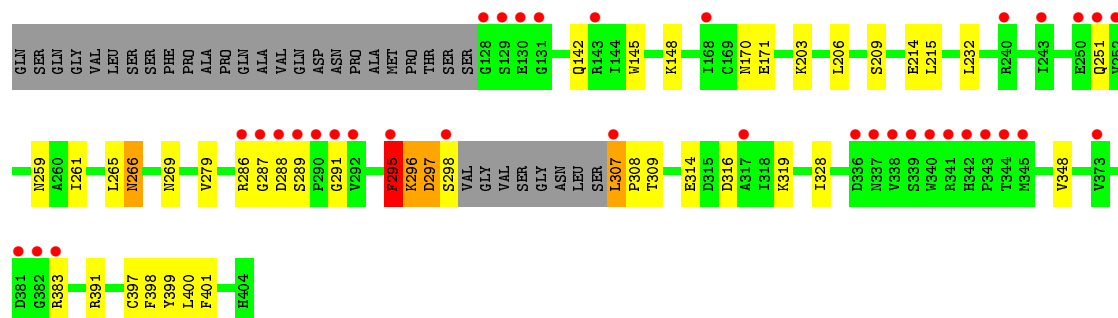
#### • Molecule 1: Caspase-1

Chain A: 



#### • Molecule 1: Caspase-1

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	56.10Å 58.00Å 60.12Å 105.95° 119.60° 92.77°	Depositor
Resolution (Å)	20.00 – 2.05 19.78 – 2.05	Depositor EDS
% Data completeness (in resolution range)	96.6 (20.00-2.05) 86.3 (19.78-2.05)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.78 (at 2.06Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.200 , 0.263 0.212 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	35.6	Xtriage
Anisotropy	0.133	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 59.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 37221 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4423	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.58% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: MG

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	0.56	0/2193	0.69	0/2957
1	B	0.61	0/2181	0.73	0/2941
All	All	0.59	0/4374	0.71	0/5898

There are no bond length outliers.

There are no bond angle outliers.

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	2149	0	2116	18	0
1	B	2137	0	2103	31	0
2	B	1	0	0	0	0
3	A	69	0	0	1	0
3	B	67	0	0	2	0
All	All	4423	0	4219	48	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 6.

All (48) 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:148:LYS:NZ	1:B:397:CYS:SG	2.26	1.08
1:A:148:LYS:NZ	1:A:397:CYS:SG	2.27	1.07
1:B:142:GLN:HG3	3:B:405:HOH:O	1.72	0.88
1:B:259:ASN:HD22	1:B:391:ARG:HH22	1.20	0.87
1:A:290[A]:PRO:O	1:A:291[A]:GLY:O	1.92	0.85
1:B:266:ASN:C	1:B:266:ASN:HD22	1.86	0.77
1:B:295[B]:PHE:O	1:B:296[B]:LYS:CB	2.37	0.72
1:B:297[B]:ASP:O	1:B:298[B]:SER:HB3	1.89	0.72
1:A:266:ASN:HD22	1:A:266:ASN:C	1.93	0.71
1:B:297[B]:ASP:O	1:B:298[B]:SER:CB	2.39	0.71
1:B:287[B]:GLY:O	1:B:289[B]:SER:N	2.26	0.69
1:B:295[B]:PHE:O	1:B:296[B]:LYS:HB2	1.93	0.68
1:A:266:ASN:ND2	1:A:269:ASN:H	1.93	0.66
1:B:295[B]:PHE:O	1:B:296[B]:LYS:HG2	1.98	0.64
1:B:142:GLN:HE21	1:B:142:GLN:HA	1.63	0.64
1:B:295[B]:PHE:O	1:B:296[B]:LYS:CG	2.48	0.62
1:B:251:GLN:OE1	1:B:251:GLN:N	2.34	0.61
1:A:266:ASN:HD21	1:A:269:ASN:H	1.51	0.58
1:B:266:ASN:C	1:B:266:ASN:ND2	2.56	0.58
1:B:297[B]:ASP:OD1	1:B:297[B]:ASP:N	2.36	0.58
1:A:145:TRP:O	1:A:149:SER:HB3	2.07	0.54
1:B:286:ARG:HG2	1:B:291[B]:GLY:HA2	1.89	0.54
1:A:374:ARG:HD3	1:A:387:PRO:O	2.09	0.52
1:B:328:ILE:O	1:B:328:ILE:HG23	2.09	0.51
1:B:142:GLN:NE2	1:B:142:GLN:HA	2.27	0.50
1:A:368:GLU:OE1	1:B:148:LYS:NZ	2.40	0.49
1:B:142:GLN:CG	3:B:405:HOH:O	2.43	0.49
1:A:289[A]:SER:HB2	1:A:290[A]:PRO:HD2	1.93	0.49
1:A:366:VAL:HG12	1:A:400:LEU:CD2	2.44	0.48
1:A:126:SER:N	3:A:434:HOH:O	2.47	0.48
1:B:266:ASN:ND2	1:B:269:ASN:H	2.11	0.47
1:B:399:TYR:HB2	1:B:401:PHE:CE2	2.48	0.47
1:B:287[B]:GLY:C	1:B:289[B]:SER:N	2.69	0.46
1:B:279:VAL:CG2	1:B:398:PHE:HE2	2.28	0.46
1:B:314:GLU:O	1:B:319:LYS:HE3	2.16	0.46
1:B:148:LYS:CE	1:B:397:CYS:SG	3.04	0.45
1:B:170:ASN:ND2	1:B:206:LEU:H	2.16	0.44
1:B:307:LEU:N	1:B:308:PRO:CD	2.82	0.43
1:A:169:CYS:HB3	1:A:186:ILE:HD11	2.01	0.42
1:B:203:LYS:HB3	1:B:206:LEU:HD11	2.00	0.42
1:A:130:GLU:O	1:A:195:ASN:HB3	2.19	0.42
1:A:266:ASN:HB2	1:A:323:ILE:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:ARG:HB3	1:A:241:GLU:OE1	2.19	0.42
1:A:266:ASN:ND2	1:A:266:ASN:C	2.66	0.42
1:A:148:LYS:CE	1:A:397:CYS:SG	3.08	0.41
1:B:328:ILE:CG2	1:B:328:ILE:O	2.67	0.41
1:B:261:ILE:O	1:B:265:LEU:HD23	2.21	0.41
1:A:281:ILE:HD13	1:A:353:LEU:HD21	2.03	0.41

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	267/302 (88%)	246 (92%)	17 (6%)	4 (2%)	13	3
1	B	265/302 (88%)	249 (94%)	13 (5%)	3 (1%)	17	7
All	All	532/604 (88%)	495 (93%)	30 (6%)	7 (1%)	15	5

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	291[A]	GLY
1	A	297[A]	ASP
1	A	287[A]	GLY
1	B	288[B]	ASP
1	B	296[B]	LYS
1	A	294[A]	TRP
1	B	295[B]	PHE



### 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	242/267 (91%)	226 (93%)	16 (7%)	21	11
1	B	240/267 (90%)	225 (94%)	15 (6%)	22	12
All	All	482/534 (90%)	451 (94%)	31 (6%)	22	12

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	138	LEU
1	A	142	GLN
1	A	143	ARG
1	A	145	TRP
1	A	164	LEU
1	A	215	LEU
1	A	232	LEU
1	A	246	LYS
1	A	259	ASN
1	A	266	ASN
1	A	283	GLN
1	A	292[A]	VAL
1	A	295[A]	PHE
1	A	297[A]	ASP
1	A	311	GLU
1	A	400	LEU
1	B	145	TRP
1	B	171	GLU
1	B	209	SER
1	B	214	GLU
1	B	215	LEU
1	B	232	LEU
1	B	266	ASN
1	B	295[B]	PHE
1	B	297[B]	ASP
1	B	307	LEU
1	B	309	THR

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Mol	Chain	Res	Type
1	B	316	ASP
1	B	348	VAL
1	B	383	ARG
1	B	400	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	170	ASN
1	A	195	ASN
1	A	263	ASN
1	A	266	ASN
1	A	269	ASN
1	A	385	GLN
1	B	142	GLN
1	B	147	GLN
1	B	170	ASN
1	B	259	ASN
1	B	263	ASN
1	B	266	ASN
1	B	269	ASN
1	B	283	GLN

### 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 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

## 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 ⓘ

### 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	271/302 (89%)	0.51	31 (11%) 7 7	22, 36, 69, 96	12 (4%)
1	B	269/302 (89%)	0.74	36 (13%) 4 4	8, 34, 77, 123	12 (4%)
All	All	540/604 (89%)	0.63	67 (12%) 5 5	8, 35, 75, 123	24 (4%)

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	340	TRP	13.6
1	B	338	VAL	12.0
1	B	343	PRO	11.1
1	B	341	ARG	8.7
1	B	344	THR	7.5
1	B	342	HIS	7.0
1	A	344	THR	6.9
1	B	130	GLU	6.7
1	B	298[B]	SER	6.5
1	B	337	ASN	6.4
1	B	290[B]	PRO	6.1
1	B	345	MET	5.9
1	A	288[A]	ASP	5.7
1	B	129	SER	5.7
1	B	288[B]	ASP	5.5
1	A	298[A]	SER	5.4
1	A	290[A]	PRO	5.3
1	B	128	GLY	5.1
1	A	382	GLY	4.8
1	B	252	VAL	4.5
1	A	317	ALA	4.3
1	B	307	LEU	4.3
1	B	289[B]	SER	4.2
1	A	294[A]	TRP	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	286	ARG	4.1
1	A	250	GLU	4.0
1	B	339	SER	3.8
1	A	287[A]	GLY	3.7
1	B	382	GLY	3.7
1	B	381	ASP	3.7
1	B	291[B]	GLY	3.7
1	A	337	ASN	3.6
1	A	286	ARG	3.5
1	B	131	GLY	3.5
1	A	292[A]	VAL	3.5
1	A	126	SER	3.4
1	B	250	GLU	3.4
1	A	291[A]	GLY	3.4
1	A	251	GLN	3.3
1	B	243	ILE	3.3
1	A	243	ILE	3.2
1	A	295[A]	PHE	3.2
1	B	251	GLN	3.1
1	A	383	ARG	3.1
1	B	292[B]	VAL	3.1
1	A	127	SER	3.0
1	B	336	ASP	3.0
1	B	295[B]	PHE	2.9
1	A	252	VAL	2.8
1	A	381	ASP	2.8
1	A	400	LEU	2.7
1	A	316	ASP	2.7
1	B	317	ALA	2.6
1	B	287[B]	GLY	2.5
1	A	289[A]	SER	2.5
1	A	131	GLY	2.5
1	B	168	ILE	2.5
1	A	244	CYS	2.4
1	A	336	ASP	2.4
1	A	307	LEU	2.4
1	B	383	ARG	2.3
1	A	345	MET	2.3
1	A	168	ILE	2.2
1	B	240	ARG	2.1
1	B	143	ARG	2.1
1	A	130	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	373	VAL	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	MG	B	1	1/1	0.98	0.02	-2.58	34,34,34,34	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.