

Full wwPDB X-ray Structure Validation Report (i)

Apr 18, 2024 – 10:38 AM EDT

PDB ID	:	7H2U
Title	:	Group deposition for crystallographic fragment screening of Coxsackievirus
		A16 (G-10) 2A protease – Crystal structure of Coxsackievirus A16 (G-10) 2A
		protease in complex with $Z355202286$ (A71EV2A-x0188)
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		X.; Thompson, W.; Tomlinson, C.W.E.; Wild, C.; Winokan, M.; Xavier, M
		A.E.; Fearon, D.; von Delft, F.
Deposited on	:	2024-04-04
Resolution	:	1.11 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* A user guide is available at https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

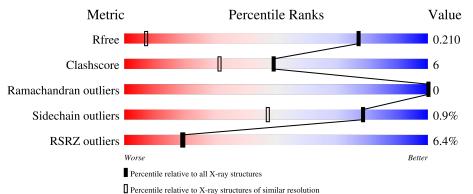
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\;DIFFRACTION$

The reported resolution of this entry is 1.11 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	$\begin{array}{c} \textbf{Whole archive} \\ \textbf{(\#Entries)} \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	1168 (1.14-1.10)
Clashscore	141614	1205 (1.14-1.10)
Ramachandran outliers	138981	1168 (1.14-1.10)
Sidechain outliers	138945	1165 (1.14-1.10)
RSRZ outliers	127900	1146 (1.14-1.10)

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 <=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	
			6%	
1	А	150	81%	10% •• 7%

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:

Validation Pipeline (wwPDB-VP) : 2.36.1



Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	SO4	А	208	-	-	Х	-



2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 1373 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.

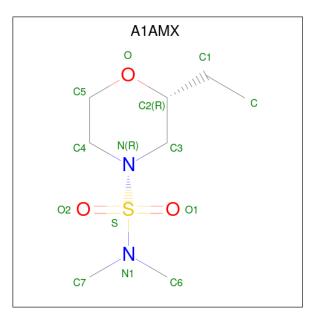
• Molecule 1 is a protein called Protease 2A.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	А	140	Total 1097	C 683	N 195	0 212	${ m S} 7$	0	2	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	1	GLN	-	expression tag	UNP Q65900
А	2	GLU	-	expression tag	UNP Q65900
A	3	GLN	-	expression tag	UNP Q65900
А	4	THR	-	expression tag	UNP Q65900
А	5	GLY	-	expression tag	UNP Q65900
А	6	GLY	-	expression tag	UNP Q65900

• Molecule 2 is (2R)-2-ethyl-N,N-dimethylmorpholine-4-sulfonamide (three-letter code: A1AMX) (formula: $C_8H_{18}N_2O_3S$).



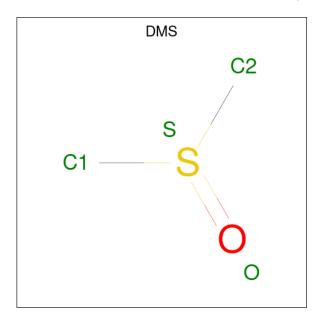


Mol	Chain	Residues		Ato	oms			ZeroOcc	AltConf
2	А	1	Total 14	C 8	N 2	0 3	S 1	0	0

• Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

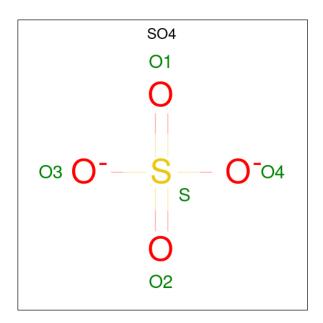
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	Total Zn 1 1	0	0

• Molecule 4 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	1	$\begin{array}{cccc} \text{Total} & \text{C} & \text{O} & \text{S} \\ 4 & 2 & 1 & 1 \end{array}$	0	0
4	А	1	$\begin{array}{cccc} \text{Total} & \text{C} & \text{O} & \text{S} \\ 4 & 2 & 1 & 1 \end{array}$	0	0
4	А	1	$\begin{array}{cccc} \text{Total} & \text{C} & \text{O} & \text{S} \\ 4 & 2 & 1 & 1 \end{array}$	0	0
4	А	1	$\begin{array}{cccc} \text{Total} & \text{C} & \text{O} & \text{S} \\ 4 & 2 & 1 & 1 \end{array}$	0	0
4	А	1	$\begin{array}{cccc} \text{Total} & \text{C} & \text{O} & \text{S} \\ 4 & 2 & 1 & 1 \end{array}$	0	0





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	А	1	Total 5	0 4	S 1	0	0

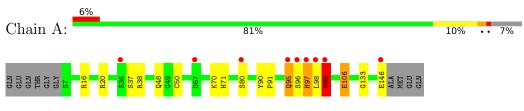
• Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	236	Total O 236 236	0	0



3 Residue-property plots (i)

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: Protease 2A



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	86.44Å 56.24Å 32.45Å	Depositor
a, b, c, α , β , γ	90.00° 95.67° 90.00°	Depositor
Resolution (Å)	47.07 - 1.11	Depositor
Resolution (A)	47.07 - 1.11	EDS
% Data completeness	97.0 (47.07-1.11)	Depositor
(in resolution range)	$97.0\ (47.07-1.11)$	EDS
R _{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.20 (at 1.11 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0267, REFMAC5	Depositor
D D.	0.188 , 0.194	Depositor
R, R_{free}	0.206 , 0.210	DCC
R_{free} test set	2917 reflections $(4.92%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	11.5	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.29, 37.2	EDS
L-test for twinning ²	$ L > = 0.50, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	1373	wwPDB-VP
Average B, all atoms $(Å^2)$	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 12.19% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: ZN, SO4, A1AMX, DMS

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		Bo	nd lengths	Bond angles	
Mol Ch	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	1.08	6/1123~(0.5%)	1.17	11/1527~(0.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers	
1	А	0	1	

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	А	95	GLN	C-O	12.75	1.47	1.23
1	А	99	MET	C-O	9.02	1.40	1.23
1	А	96	SER	CA-CB	8.56	1.65	1.52
1	А	96	SER	CB-OG	-7.13	1.32	1.42
1	А	106	GLU	CD-OE2	-5.85	1.19	1.25
1	А	98	LEU	C-O	5.48	1.33	1.23

All (6) bond length outliers are listed below:

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$\mathbf{Ideal}(^{o})$
1	А	99	MET	N-CA-CB	-8.94	94.51	110.60
1	А	98	LEU	CB-CG-CD2	-8.40	96.72	111.00
1	А	97	HIS	CB-CA-C	6.94	124.28	110.40
1	А	97	HIS	O-C-N	6.34	132.85	122.70
1	А	16	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	А	20	ARG	NE-CZ-NH2	-5.91	117.34	120.30
1	А	95	GLN	CB-CA-C	5.87	122.14	110.40

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Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	20	ARG	NE-CZ-NH1	5.66	123.13	120.30
1	А	97	HIS	CA-C-O	-5.51	108.53	120.10
1	А	90	TYR	CB-CG-CD2	-5.10	117.94	121.00
1	А	80	SER	CB-CA-C	5.07	119.73	110.10

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There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	99	MET	Mainchain

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	А	1097	0	1030	9	0
2	А	14	0	0	1	0
3	А	1	0	0	0	0
4	А	20	0	30	4	0
5	А	5	0	0	0	6
6	А	236	0	0	8	1
All	All	1373	0	1060	13	7

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 (13) 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:50:CYS:HB2	6:A:309:HOH:O	1.61	0.99
1:A:106:GLU:OE1	6:A:301:HOH:O	1.99	0.81
1:A:37:SER:O	1:A:95:GLN:HG3	1.90	0.71
4:A:203:DMS:H23	6:A:428:HOH:O	1.96	0.65
1:A:71:HIS:HE1	6:A:463:HOH:O	1.86	0.58
1:A:70:LYS:NZ	6:A:304:HOH:O	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:PRO:HD3	4:A:207:DMS:H21	1.90	0.53
4:A:203:DMS:C2	6:A:428:HOH:O	2.54	0.52
1:A:48:GLN:OE1	6:A:302:HOH:O	2.20	0.47
4:A:203:DMS:H22	6:A:416:HOH:O	2.14	0.47
1:A:38:ARG:CZ	1:A:97:HIS:CG	3.01	0.44
2:A:201:A1AMX:C3	2:A:201:A1AMX:C6	2.96	0.43
1:A:99:MET:O	1:A:133:GLY:HA2	2.18	0.43

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All (7) 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:208:SO4:O3	5:A:208:SO4:O3[2_556]	0.68	1.52
5:A:208:SO4:O4	5:A:208:SO4:O4[2_556]	1.19	1.01
5:A:208:SO4:S	5:A:208:SO4:O4[2_556]	1.63	0.57
5:A:208:SO4:S	5:A:208:SO4:O3[2_556]	1.82	0.38
5:A:208:SO4:O1	5:A:208:SO4:O4[2_556]	1.88	0.32
6:A:385:HOH:O	6:A:499:HOH:O[4_556]	2.02	0.18
5:A:208:SO4:S	5:A:208:SO4:S[2_556]	2.04	0.16

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	А	140/150~(93%)	137~(98%)	3~(2%)	0	100 100

There are no Ramachandran outliers to report.

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	А	119/124~(96%)	118 (99%)	1 (1%)	81 51	

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	146	GLU

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such side chains are listed below:

Mol	Chain	Res	Type
1	А	71	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 8 ligands modelled in this entry, 1 is monoatomic - leaving 7 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	Turne	Chain	Res	Link	Bo	ond leng	ths	В	ond ang	les
10101	Type	Unam	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	DMS	А	204	-	3,3,3	0.28	0	3,3,3	0.44	0
2	A1AMX	А	201	-	14,14,14	5.41	9 (64%)	$15,\!20,\!20$	<mark>3.21</mark>	7 (46%)
4	DMS	А	205	-	3,3,3	0.26	0	3,3,3	0.32	0
4	DMS	А	203	-	3,3,3	0.70	0	3,3,3	0.72	0
5	SO4	А	208	-	4,4,4	0.37	0	$6,\!6,\!6$	0.12	0
4	DMS	А	207	-	3,3,3	0.55	0	3,3,3	0.68	0
4	DMS	А	206	-	3,3,3	0.07	0	3,3,3	0.04	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	A1AMX	А	201	-	-	4/12/24/24	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	$\mathrm{Ideal}(\mathrm{\AA})$
2	А	201	A1AMX	S-N1	-16.17	1.30	1.60
2	А	201	A1AMX	C1-C2	-5.67	1.23	1.51
2	А	201	A1AMX	S-N	5.01	1.69	1.60
2	А	201	A1AMX	C6-N1	-4.52	1.40	1.47
2	А	201	A1AMX	C7-N1	-4.18	1.40	1.47
2	А	201	A1AMX	O1-S	-3.88	1.34	1.42
2	А	201	A1AMX	O2-S	3.68	1.50	1.42
2	А	201	A1AMX	C-C1	3.55	1.72	1.51
2	А	201	A1AMX	C3-N	-3.01	1.45	1.47

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
2	А	201	A1AMX	02-S-01	-7.32	106.52	121.72
2	А	201	A1AMX	C5-C4-N	6.07	111.75	108.18
2	А	201	A1AMX	O2-S-N1	-4.36	102.87	108.89
2	А	201	A1AMX	C6-N1-S	-3.51	110.66	117.94
2	А	201	A1AMX	N-S-N1	2.99	116.39	105.97
2	А	201	A1AMX	01-S-N1	2.35	112.13	108.89
2	А	201	A1AMX	O-C2-C1	-2.29	100.94	107.02

There are no chirality outliers.



Mol	Chain	Res	Type	Atoms
2	А	201	A1AMX	C7-N1-S-O1
2	А	201	A1AMX	C3-N-S-O2
2	А	201	A1AMX	C-C1-C2-O
2	А	201	A1AMX	C-C1-C2-C3

All (4) torsion outliers are listed below:

There are no ring outliers.

4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	А	201	A1AMX	1	0
4	А	203	DMS	3	0
5	А	208	SO4	0	6
4	А	207	DMS	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^{th} 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 $>$	#RS	SRZ>	>2	$OWAB(Å^2)$	Q<0.9
1	А	140/150~(93%)	0.64	9~(6%)	19	19	8, 12, 19, 48	8 (5%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	98	LEU	23.6
1	А	99	MET	15.0
1	А	96	SER	14.8
1	А	97	HIS	14.0
1	А	95	GLN	13.3
1	А	80	SER	13.0
1	А	36[A]	SER	6.3
1	А	57[A]	ASP	5.2
1	А	146	GLU	2.2

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^{th} 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	\mathbf{RSR}	$\mathbf{B} ext{-factors}(\mathbf{A}^2)$	Q<0.9
2	A1AMX	A	201	14/14	0.75	0.40	20,20,20,20	14
4	DMS	А	205	4/4	0.84	0.16	45,55,57,58	0
5	SO4	А	208	5/5	0.86	0.49	$36,\!38,\!42,\!45$	0
4	DMS	А	203	4/4	0.90	0.14	20,22,24,33	0
4	DMS	А	206	4/4	0.95	0.15	26,26,27,28	0
4	DMS	А	204	4/4	0.97	0.12	21,21,22,23	0
4	DMS	А	207	4/4	0.98	0.17	18,21,23,27	0
3	ZN	А	202	1/1	1.00	0.03	$9,\!9,\!9,\!9$	0

6.5 Other polymers (i)

There are no such residues in this entry.

