



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:01 AM UTC

PDB ID : 4YSC / pdb_00004ysc
Title : Completely oxidized structure of copper nitrite reductase from *Alcaligenes faecalis*
Authors : Fukuda, Y.; Tse, K.M.; Suzuki, M.; Diederichs, K.; Hirata, K.; Nakane, T.; Sugahara, M.; Nango, E.; Tono, K.; Joti, Y.; Kameshima, T.; Song, C.; Hatsui, T.; Yabashi, M.; Nureki, O.; Matsumura, H.; Inoue, T.; Iwata, S.; Mizohata, E.
Deposited on : 2015-03-17
Resolution : 2.03 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
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)	:	Engl & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

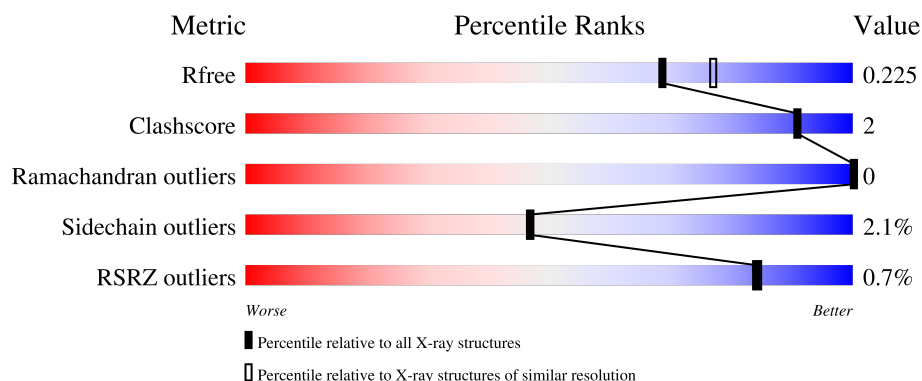
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.03 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	13299 (2.04-2.00)
Clashscore	190562	1022 (2.02-2.02)
Ramachandran outliers	187476	1014 (2.02-2.02)
Sidechain outliers	187428	1014 (2.02-2.02)
RSRZ outliers	180081	13314 (2.04-2.00)

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 $\leq 5\%$. 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	342	% 89% 8% ..
1	B	342	% 89% 9% .
1	C	342	92% 6% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7931 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 Copper-containing nitrite reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	336	Total	C	N	O	S	0	0	0
			2560	1639	430	480	11			
1	B	336	Total	C	N	O	S	0	0	0
			2560	1639	430	480	11			
1	C	335	Total	C	N	O	S	0	0	0
			2556	1637	429	479	11			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	3	MET	-	initiating methionine	UNP P38501
A	341	LEU	-	expression tag	UNP P38501
A	342	VAL	-	expression tag	UNP P38501
A	343	PRO	-	expression tag	UNP P38501
A	344	ARG	-	expression tag	UNP P38501
B	3	MET	-	initiating methionine	UNP P38501
B	341	LEU	-	expression tag	UNP P38501
B	342	VAL	-	expression tag	UNP P38501
B	343	PRO	-	expression tag	UNP P38501
B	344	ARG	-	expression tag	UNP P38501
C	3	MET	-	initiating methionine	UNP P38501
C	341	LEU	-	expression tag	UNP P38501
C	342	VAL	-	expression tag	UNP P38501
C	343	PRO	-	expression tag	UNP P38501
C	344	ARG	-	expression tag	UNP P38501

- Molecule 2 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Cu	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total 2	Cu 2	0	0
2	C	2	Total 2	Cu 2	0	0

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total 2	Cl 2	0	0
3	B	1	Total 1	Cl 1	0	0

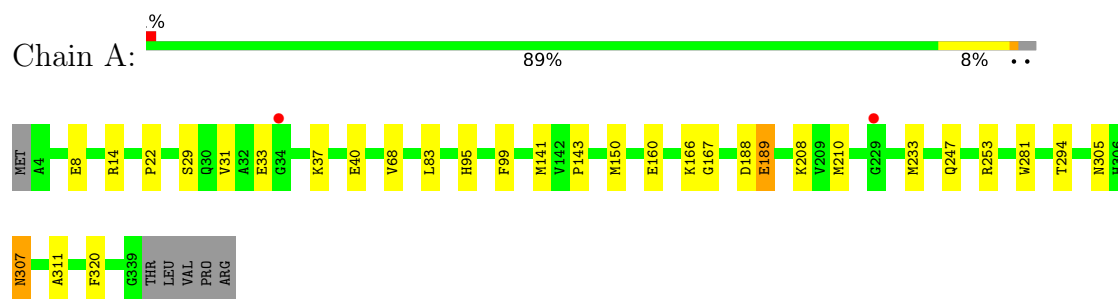
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	79	Total 79	O 79	0	0
4	B	86	Total 86	O 86	0	0
4	C	81	Total 81	O 81	0	0

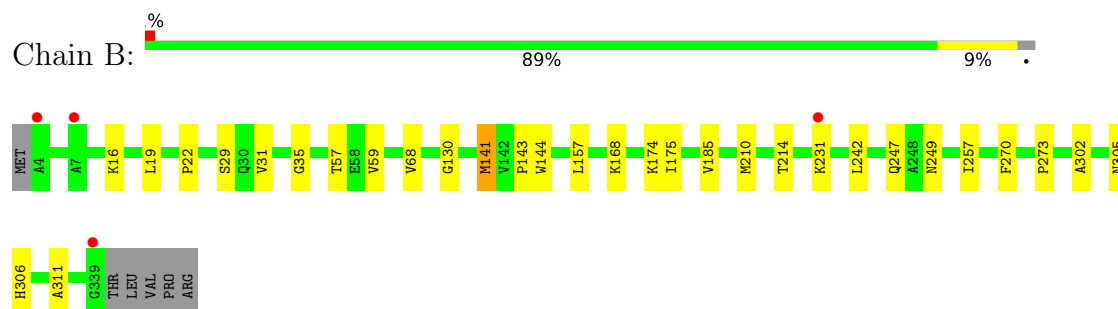
3 Residue-property plots [i](#)

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.

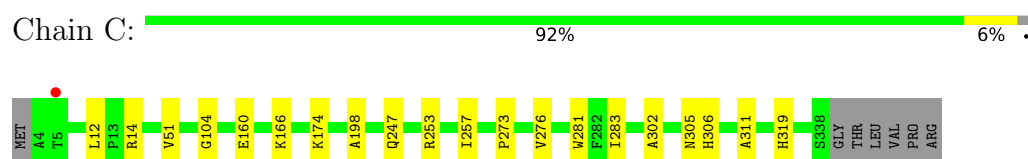
- Molecule 1: Copper-containing nitrite reductase



- Molecule 1: Copper-containing nitrite reductase



- Molecule 1: Copper-containing nitrite reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.10Å 103.80Å 147.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	73.90 – 2.03 73.90 – 2.03	Depositor EDS
% Data completeness (in resolution range)	100.0 (73.90-2.03) 100.0 (73.90-2.03)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 2.03Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.164 , 0.203 (Not available) , 0.225	Depositor DCC
R_{free} test set	2499 reflections (3.93%)	wwPDB-VP
Wilson B-factor (Å ²)	38.5	Xtriage
Anisotropy	0.022	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 35.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7931	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

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: CL, CU

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	1.27	3/2631 (0.1%)	1.12	2/3588 (0.1%)
1	B	1.29	5/2631 (0.2%)	1.11	6/3588 (0.2%)
1	C	1.31	2/2627 (0.1%)	1.13	10/3583 (0.3%)
All	All	1.29	10/7889 (0.1%)	1.12	18/10759 (0.2%)

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	B	0	1
1	C	0	1
All	All	0	2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	68	VAL	N-CA	-7.10	1.40	1.46
1	B	214	THR	C-O	-6.80	1.18	1.25
1	C	319	HIS	N-CA	6.46	1.53	1.45
1	A	68	VAL	N-CA	6.20	1.51	1.46
1	B	22	PRO	CA-C	6.13	1.55	1.51
1	B	273	PRO	CA-C	5.73	1.57	1.52
1	A	22	PRO	CA-C	5.51	1.54	1.51
1	B	57	THR	CA-C	5.49	1.59	1.52
1	A	294	THR	CA-CB	5.37	1.60	1.53
1	C	51	VAL	CA-CB	5.22	1.60	1.54

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	185	VAL	N-CA-C	6.28	113.27	107.56
1	A	307	ASN	N-CA-C	-6.21	99.04	108.67
1	B	35	GLY	CA-C-N	5.78	127.06	119.84
1	B	35	GLY	C-N-CA	5.78	127.06	119.84
1	C	283	ILE	CA-C-N	-5.52	114.27	119.85
1	C	283	ILE	C-N-CA	-5.52	114.27	119.85
1	C	273	PRO	CA-C-N	-5.45	114.63	120.03
1	C	273	PRO	C-N-CA	-5.45	114.63	120.03
1	B	157	LEU	CA-C-N	-5.44	114.17	119.78
1	B	157	LEU	C-N-CA	-5.44	114.17	119.78
1	B	168	LYS	N-CA-C	5.24	117.90	110.10
1	C	198	ALA	CA-C-N	-5.22	113.26	119.05
1	C	198	ALA	C-N-CA	-5.22	113.26	119.05
1	C	104	GLY	N-CA-C	5.22	116.91	111.95
1	A	167	GLY	N-CA-C	5.18	122.24	115.40
1	C	12	LEU	CA-C-N	-5.10	114.70	119.85
1	C	12	LEU	C-N-CA	-5.10	114.70	119.85
1	C	276	VAL	N-CA-C	5.08	116.67	108.85

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	306	HIS	Peptide
1	C	306	HIS	Peptide

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	2560	0	2485	14	0
1	B	2560	0	2485	10	0
1	C	2556	0	2482	3	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	79	0	0	1	0
4	B	86	0	0	0	0
4	C	81	0	0	0	0
All	All	7931	0	7452	26	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 2.

All (26) 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:189:GLU:H	1:A:189:GLU:CD	1.95	0.72
1:A:141:MET:HG2	4:A:555:HOH:O	1.93	0.69
1:B:257:ILE:HD12	1:B:302:ALA:HB3	1.87	0.56
1:A:188:ASP:HB2	1:A:189:GLU:OE1	2.06	0.55
1:B:143:PRO:HB2	1:B:210:MET:CE	2.39	0.52
1:B:143:PRO:HB2	1:B:210:MET:HE1	1.91	0.51
1:A:29:SER:OG	1:A:31:VAL:O	2.28	0.51
1:C:257:ILE:HD12	1:C:302:ALA:HB3	1.95	0.49
1:C:305:ASN:O	1:C:311:ALA:HB2	2.15	0.46
1:A:95:HIS:HB3	1:A:150:MET:HE1	1.97	0.46
1:B:29:SER:OG	1:B:31:VAL:O	2.32	0.46
1:B:141:MET:HE3	1:B:144:TRP:CE3	2.51	0.45
1:A:305:ASN:O	1:A:311:ALA:HB2	2.17	0.45
1:A:14:ARG:NH2	1:A:40:GLU:OE2	2.50	0.45
1:A:307:ASN:HA	1:B:249:ASN:O	2.18	0.44
1:A:33:GLU:O	1:A:37:LYS:NZ	2.52	0.43
1:A:253:ARG:HA	1:A:281:TRP:O	2.19	0.43
1:B:175:ILE:HD12	1:B:242:LEU:HD23	2.01	0.43
1:B:130:GLY:HA2	1:B:270:PHE:CD2	2.55	0.42
1:A:83:LEU:HD22	1:A:99:PHE:HZ	1.84	0.41
1:A:189:GLU:CD	1:A:189:GLU:N	2.72	0.41
1:A:143:PRO:HB2	1:A:210:MET:HE1	2.03	0.41
1:A:233:MET:O	1:A:320:PHE:HA	2.21	0.40
1:B:19:LEU:HD23	1:B:19:LEU:HA	1.95	0.40
1:B:305:ASN:O	1:B:311:ALA:HB2	2.22	0.40
1:C:253:ARG:HA	1:C:281:TRP:O	2.21	0.40

There are no symmetry-related clashes.

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	334/342 (98%)	328 (98%)	6 (2%)	0	100	100
1	B	334/342 (98%)	328 (98%)	6 (2%)	0	100	100
1	C	333/342 (97%)	330 (99%)	3 (1%)	0	100	100
All	All	1001/1026 (98%)	986 (98%)	15 (2%)	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	264/270 (98%)	258 (98%)	6 (2%)	44	43
1	B	264/270 (98%)	258 (98%)	6 (2%)	44	43
1	C	264/270 (98%)	259 (98%)	5 (2%)	50	50
All	All	792/810 (98%)	775 (98%)	17 (2%)	47	47

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	GLU
1	A	160	GLU
1	A	166	LYS
1	A	189	GLU
1	A	208	LYS

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Mol	Chain	Res	Type
1	A	247	GLN
1	B	16	LYS
1	B	59	VAL
1	B	141	MET
1	B	174	LYS
1	B	231	LYS
1	B	247	GLN
1	C	14	ARG
1	C	160	GLU
1	C	166	LYS
1	C	174	LYS
1	C	247	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	HIS
1	B	115	ASN
1	B	190	ASN

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 9 ligands modelled in this entry, 9 are 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

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 [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, 95th 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 ²)	Q<0.9
1	A	336/342 (98%)	-0.15	2 (0%) 85 86	26, 36, 59, 86	0
1	B	336/342 (98%)	-0.16	4 (1%) 76 76	24, 34, 62, 83	0
1	C	335/342 (97%)	-0.34	1 (0%) 90 90	23, 31, 53, 73	0
All	All	1007/1026 (98%)	-0.22	7 (0%) 84 84	23, 34, 59, 86	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	4	ALA	4.2
1	B	339	GLY	3.2
1	A	229	GLY	2.7
1	B	7	ALA	2.5
1	C	5	THR	2.2
1	B	231	LYS	2.2
1	A	34	GLY	2.1

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, 95th 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 ²)	Q<0.9
3	CL	A	404	1/1	0.97	0.08	40,40,40,40	0
3	CL	A	403	1/1	0.98	0.08	38,38,38,38	0
3	CL	B	403	1/1	0.99	0.03	35,35,35,35	0
2	CU	B	402	1/1	1.00	0.02	34,34,34,34	0
2	CU	C	401	1/1	1.00	0.01	29,29,29,29	0
2	CU	C	402	1/1	1.00	0.01	29,29,29,29	0
2	CU	A	401	1/1	1.00	0.01	35,35,35,35	0
2	CU	A	402	1/1	1.00	0.01	33,33,33,33	0
2	CU	B	401	1/1	1.00	0.02	36,36,36,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.