



wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 18, 2025 – 04:10 PM EDT

PDB ID : 9ONM / pdb_00009onm
Title : near-apo rubrerythrin
Authors : Budziszewski, G.R.; Snell, M.E.; Monteiro, D.C.F.; Lynch, M.L.; Bowman, S.E.J.
Deposited on : 2025-05-15
Resolution : 1.57 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

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 Phenix2.0rc1
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.006 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.44

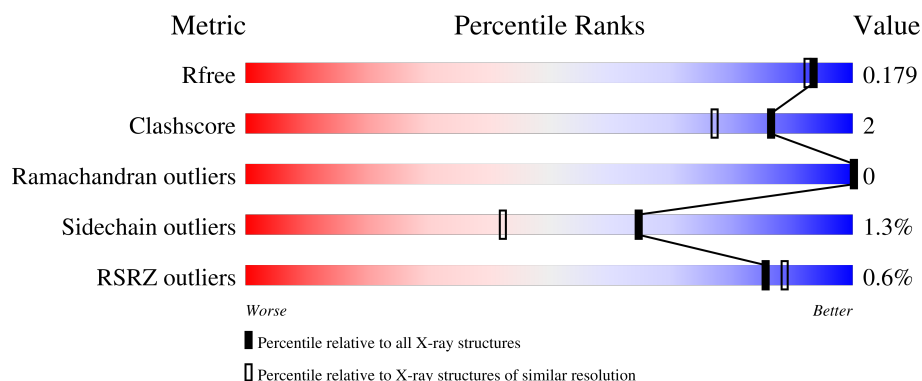
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.57 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	7165 (1.60-1.56)
Clashscore	180529	1026 (1.58-1.58)
Ramachandran outliers	177936	1005 (1.58-1.58)
Sidechain outliers	177891	1004 (1.58-1.58)
RSRZ outliers	164620	7163 (1.60-1.56)

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	140	<div> <div>%</div> <div> <div></div> <div>94%</div> <div>6%</div> <div>.</div> </div> </div>
1	B	140	<div> <div>%</div> <div> <div></div> <div>94%</div> <div>6%</div> <div>.</div> </div> </div>
1	C	140	<div> <div></div> <div> <div></div> <div>95%</div> <div></div> <div>.</div> <div>..</div> </div> </div>
1	D	140	<div> <div>%</div> <div> <div></div> <div>96%</div> <div></div> <div>.</div> <div>..</div> </div> </div>
1	E	140	<div> <div>%</div> <div> <div></div> <div>94%</div> <div>5%</div> <div>.</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	140	<div><div><div>%</div><div><div></div></div><div>91%</div><div>8%<div></div></div></div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14226 atoms, of which 6486 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 Rubrerythrin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	139	Total	C	H	N	O	S	0	8	0
			2148	684	1039	197	226	2			
1	B	139	Total	C	H	N	O	S	0	10	0
			2171	692	1052	199	226	2			
1	C	139	Total	C	H	N	O	S	0	13	0
			2209	704	1076	199	228	2			
1	D	139	Total	C	H	N	O	S	0	17	0
			2233	709	1085	204	233	2			
1	E	139	Total	C	H	N	O	S	0	11	0
			2183	695	1062	196	228	2			
1	F	139	Total	C	H	N	O	S	0	11	0
			2183	695	1062	196	228	2			

- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Fe	0	0
			2	2		
2	B	1	Total	Fe	0	0
			1	1		
2	C	3	Total	Fe	0	0
			3	3		
2	D	2	Total	Fe	0	0
			2	2		
2	E	2	Total	Fe	0	0
			2	2		
2	F	2	Total	Fe	0	0
			2	2		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	H	O	0	0
			17	4	10	3		
3	B	1	Total	C	H	O	0	0
			17	4	10	3		
3	C	1	Total	C	H	O	0	0
			17	4	10	3		
3	C	1	Total	C	H	O	0	0
			17	4	10	3		
3	D	1	Total	C	H	O	0	0
			17	4	10	3		
3	D	1	Total	C	H	O	0	0
			17	4	10	3		
3	E	1	Total	C	H	O	0	0
			17	4	10	3		
3	F	1	Total	C	H	O	0	0
			17	4	10	3		
3	F	1	Total	C	H	O	0	0
			17	4	10	3		
3	F	1	Total	C	H	O	0	0
			17	4	10	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	145	Total	O	0	10
			155	155		

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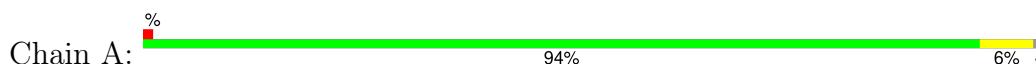
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	156	Total 162	O 162	0	6
4	C	136	Total 142	O 142	0	6
4	D	134	Total 138	O 138	0	4
4	E	143	Total 152	O 152	0	9
4	F	145	Total 151	O 151	0	6

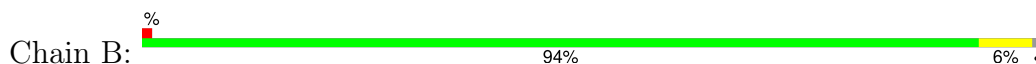
3 Residue-property plots [i](#)

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.

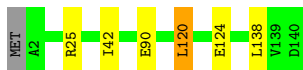
- Molecule 1: Rubrerythrin



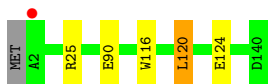
- Molecule 1: Rubrerythrin



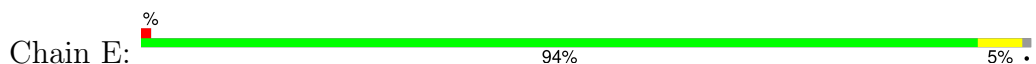
- Molecule 1: Rubrerythrin



- Molecule 1: Rubrerythrin



- Molecule 1: Rubrerythrin



- Molecule 1: Rubrerythrin



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	203.92Å 203.92Å 68.78Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	33.76 – 1.57 33.76 – 1.57	Depositor EDS
% Data completeness (in resolution range)	78.2 (33.76-1.57) 78.5 (33.76-1.57)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 1.57Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, R_{free}	0.148 , 0.179 0.148 , 0.179	Depositor DCC
R_{free} test set	7465 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	18.7	Xtriage
Anisotropy	0.021	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.009 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	14226	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% 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: FE, PEG

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	0.41	0/1171	0.54	0/1579
1	B	0.42	0/1175	0.53	0/1583
1	C	0.43	0/1194	0.52	0/1608
1	D	0.39	0/1229	0.50	0/1655
1	E	0.40	0/1179	0.52	0/1588
1	F	0.40	0/1183	0.50	0/1593
All	All	0.41	0/7131	0.52	0/9606

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 [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	1109	1039	1001	5	0
1	B	1119	1052	1015	5	0
1	C	1133	1076	1032	4	0
1	D	1148	1085	1028	3	0
1	E	1121	1062	1020	4	0
1	F	1121	1062	1019	8	0
2	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	3	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
3	B	14	20	20	0	0
3	C	14	20	20	1	0
3	D	14	20	20	0	0
3	E	7	10	10	0	0
3	F	28	40	40	3	0
4	A	155	0	0	0	0
4	B	162	0	0	1	0
4	C	142	0	0	0	0
4	D	138	0	0	0	0
4	E	152	0	0	0	0
4	F	151	0	0	0	0
All	All	7740	6486	6225	27	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.

The worst 5 of 27 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:F:97:MET:HG2	1:F:101:MET:HE3	1.84	0.60
1:F:97:MET:CG	1:F:101:MET:HE3	2.36	0.55
1:F:65:GLU:HG3	1:F:76:PHE:CE2	2.43	0.53
1:F:17:PHE:CE1	1:F:61[A]:LEU:HG	2.43	0.53
1:F:90:GLU:OE1	1:F:124:GLU:OE1	2.27	0.52

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	145/140 (104%)	141 (97%)	4 (3%)	0	100	100
1	B	147/140 (105%)	144 (98%)	3 (2%)	0	100	100
1	C	150/140 (107%)	145 (97%)	5 (3%)	0	100	100
1	D	154/140 (110%)	149 (97%)	5 (3%)	0	100	100
1	E	148/140 (106%)	143 (97%)	5 (3%)	0	100	100
1	F	148/140 (106%)	143 (97%)	5 (3%)	0	100	100
All	All	892/840 (106%)	865 (97%)	27 (3%)	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	113/106 (107%)	112 (99%)	1 (1%)	75	59
1	B	113/106 (107%)	112 (99%)	1 (1%)	75	59
1	C	114/106 (108%)	111 (97%)	3 (3%)	41	12
1	D	119/106 (112%)	116 (98%)	3 (2%)	42	13
1	E	114/106 (108%)	113 (99%)	1 (1%)	75	59
1	F	115/106 (108%)	114 (99%)	1 (1%)	75	59
All	All	688/636 (108%)	678 (98%)	10 (2%)	65	36

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	120[B]	LEU
1	E	25	ARG
1	F	25	ARG
1	C	120[A]	LEU
1	C	120[B]	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 23 ligands modelled in this entry, 12 are monoatomic - leaving 11 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	PEG	D	202	-	6,6,6	0.27	0	5,5,5	0.43	0
3	PEG	D	201	-	6,6,6	0.25	0	5,5,5	0.36	0
3	PEG	B	201	-	6,6,6	0.26	0	5,5,5	0.32	0
3	PEG	C	201	-	6,6,6	0.30	0	5,5,5	0.31	0
3	PEG	F	201	-	6,6,6	0.33	0	5,5,5	0.38	0
3	PEG	E	201	-	6,6,6	0.25	0	5,5,5	0.35	0
3	PEG	F	202	-	6,6,6	0.37	0	5,5,5	0.98	0
3	PEG	F	204	-	6,6,6	0.24	0	5,5,5	0.25	0
3	PEG	F	203	-	6,6,6	0.25	0	5,5,5	0.35	0
3	PEG	B	202	-	6,6,6	0.25	0	5,5,5	0.49	0
3	PEG	C	202	-	6,6,6	0.24	0	5,5,5	0.37	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
3	PEG	D	202	-	-	2/4/4/4	-
3	PEG	D	201	-	-	0/4/4/4	-
3	PEG	B	201	-	-	3/4/4/4	-
3	PEG	C	201	-	-	2/4/4/4	-
3	PEG	F	201	-	-	1/4/4/4	-
3	PEG	E	201	-	-	2/4/4/4	-
3	PEG	F	202	-	-	4/4/4/4	-
3	PEG	F	204	-	-	3/4/4/4	-
3	PEG	F	203	-	-	0/4/4/4	-
3	PEG	B	202	-	-	1/4/4/4	-
3	PEG	C	202	-	-	1/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 19 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	F	202	PEG	C1-C2-O2-C3
3	F	204	PEG	O1-C1-C2-O2
3	F	204	PEG	O2-C3-C4-O4
3	E	201	PEG	O1-C1-C2-O2
3	B	201	PEG	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	202	PEG	2	0
3	F	203	PEG	1	0
3	C	202	PEG	1	0

5.7 Other polymers [i](#)

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	139/140 (99%)	-0.58	1 (0%) 84 88	8, 20, 42, 76	4 (2%)
1	B	139/140 (99%)	-0.59	1 (0%) 84 88	8, 19, 40, 70	5 (3%)
1	C	139/140 (99%)	-0.58	0 100 100	8, 20, 37, 62	7 (5%)
1	D	139/140 (99%)	-0.47	1 (0%) 84 88	8, 22, 43, 57	9 (6%)
1	E	139/140 (99%)	-0.52	1 (0%) 84 88	8, 21, 40, 74	6 (4%)
1	F	139/140 (99%)	-0.43	1 (0%) 84 88	8, 22, 43, 51	6 (4%)
All	All	834/840 (99%)	-0.53	5 (0%) 85 89	8, 21, 41, 76	37 (4%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	2	ALA	4.4
1	B	2	ALA	4.3
1	D	2	ALA	3.5
1	A	2	ALA	3.4
1	F	8[A]	LYS	3.3

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	PEG	F	202	7/7	0.76	0.20	33,40,47,57	17
3	PEG	C	201	7/7	0.80	0.16	46,55,70,70	0
3	PEG	D	202	7/7	0.82	0.12	43,55,69,71	0
3	PEG	B	201	7/7	0.83	0.17	35,45,54,63	17
3	PEG	C	202	7/7	0.83	0.15	43,55,77,77	0
3	PEG	B	202	7/7	0.84	0.15	44,53,62,64	0
3	PEG	F	203	7/7	0.84	0.12	35,49,62,74	0
3	PEG	F	204	7/7	0.84	0.14	49,61,73,79	0
3	PEG	E	201	7/7	0.91	0.12	27,35,53,53	17
2	FE	D	204	1/1	0.92	0.08	18,18,18,18	1
3	PEG	D	201	7/7	0.93	0.10	28,35,61,74	0
2	FE	B	203	1/1	0.93	0.10	20,20,20,20	1
2	FE	A	202	1/1	0.93	0.10	17,17,17,17	1
3	PEG	F	201	7/7	0.94	0.09	26,35,48,51	0
2	FE	E	202	1/1	0.94	0.05	19,19,19,19	1
2	FE	F	206	1/1	0.94	0.06	18,18,18,18	1
2	FE	C	205	1/1	0.94	0.06	19,19,19,19	1
2	FE	F	205	1/1	0.95	0.06	23,23,23,23	1
2	FE	E	203	1/1	0.96	0.06	17,17,17,17	1
2	FE	C	203	1/1	0.97	0.04	14,14,14,14	1
2	FE	C	204	1/1	0.97	0.06	21,21,21,21	1
2	FE	A	201	1/1	0.97	0.04	19,19,19,19	1
2	FE	D	203	1/1	0.97	0.05	19,19,19,19	1

6.5 Other polymers [i](#)

There are no such residues in this entry.