



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:37 PM GMT

PDB ID : 1PV9
Title : Prolidase from *Pyrococcus furiosus*
Authors : Maher, M.J.; Ghosh, M.; Grunden, A.M.; Menon, A.L.; Adams, M.W.; Freeman, H.C.; Guss, J.M.
Deposited on : 2003-06-27
Resolution : 2.00 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

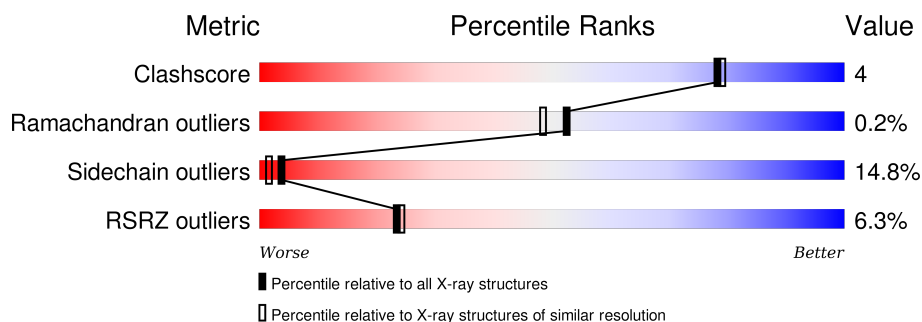
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	348	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 11% 81% 15% • • </div> </div>
1	B	348	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 11% 65% 24% • 9% </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5376 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 Xaa-Pro dipeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	337	Total	C	N	O	S	0	0	0
			2673	1706	439	519	9			
1	B	318	Total	C	N	O	S	0	0	0
			2523	1610	417	487	9			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Zn	0	0
			2	2		
2	A	2	Total	Zn	0	0
			2	2		

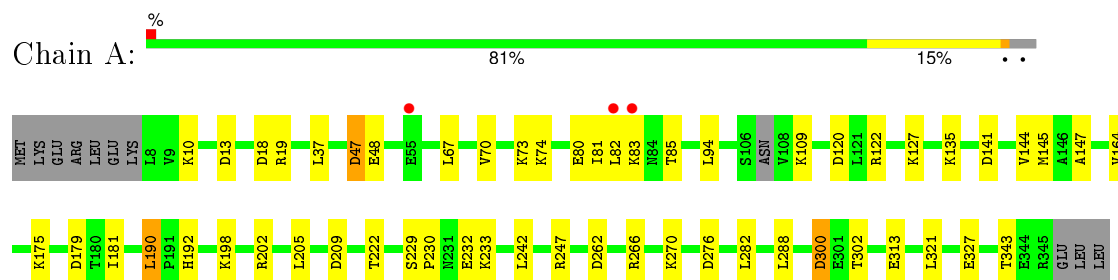
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	100	Total	O	0	0
			100	100		
3	B	76	Total	O	0	0
			76	76		

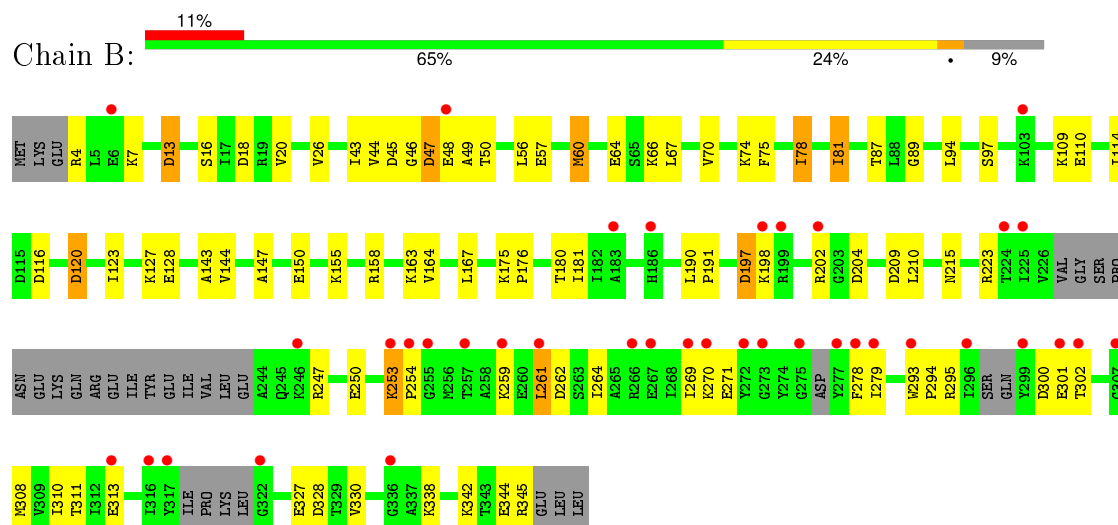
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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: Xaa-Pro dipeptidase



• Molecule 1: Xaa-Pro dipeptidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.48Å 97.23Å 69.88Å 90.00° 97.09° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 48.61 – 2.00	Depositor EDS
% Data completeness (in resolution range)	96.8 (20.00-2.00) 95.1 (48.61-2.00)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.69 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.245 , 0.277 0.247 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	36.1	Xtriage
Anisotropy	0.429	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 51.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 48110 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5376	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.31% 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, 0.375 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: ZN

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	0.29	0/2715	0.63	9/3655 (0.2%)
1	B	0.29	0/2559	0.63	10/3437 (0.3%)
All	All	0.29	0/5274	0.63	19/7092 (0.3%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	262	ASP	CB-CG-OD2	5.50	123.25	118.30
1	B	47	ASP	CB-CG-OD2	5.49	123.24	118.30
1	B	262	ASP	CB-CG-OD2	5.34	123.10	118.30
1	A	141	ASP	CB-CG-OD2	5.33	123.10	118.30
1	B	120	ASP	CB-CG-OD2	5.27	123.04	118.30
1	A	13	ASP	CB-CG-OD2	5.25	123.02	118.30
1	A	179	ASP	CB-CG-OD2	5.24	123.02	118.30
1	B	197	ASP	CB-CG-OD2	5.24	123.02	118.30
1	A	18	ASP	CB-CG-OD2	5.22	123.00	118.30
1	A	300	ASP	CB-CG-OD2	5.19	122.97	118.30
1	A	120	ASP	CB-CG-OD2	5.19	122.97	118.30
1	A	47	ASP	CB-CG-OD2	5.18	122.96	118.30
1	A	276	ASP	CB-CG-OD2	5.13	122.92	118.30
1	B	45	ASP	CB-CG-OD2	5.11	122.90	118.30
1	B	204	ASP	CB-CG-OD2	5.07	122.86	118.30
1	B	13	ASP	CB-CG-OD2	5.04	122.84	118.30
1	B	18	ASP	CB-CG-OD2	5.03	122.83	118.30
1	B	328	ASP	CB-CG-OD2	5.03	122.83	118.30
1	B	300	ASP	CB-CG-OD2	5.02	122.82	118.30

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	2673	0	2715	14	0
1	B	2523	0	2563	23	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	100	0	0	1	0
3	B	76	0	0	0	0
All	All	5376	0	5278	35	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 4.

All (35) 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:75:PHE:O	1:B:78:ILE:HG22	1.75	0.84
1:B:181:ILE:HB	1:B:209:ASP:HB3	1.64	0.79
1:A:300:ASP:OD1	1:A:302:THR:HG23	1.91	0.70
1:B:313:GLU:HB3	1:B:327:GLU:HB2	1.76	0.66
1:A:313:GLU:HB3	1:A:327:GLU:HB2	1.78	0.65
1:A:181:ILE:HB	1:A:209:ASP:HB3	1.78	0.64
1:B:143:ALA:HB1	1:B:164:VAL:HG13	1.84	0.59
1:A:122:ARG:HD3	1:A:288:LEU:HD12	1.87	0.56
1:B:43:ILE:HG21	1:B:81:ILE:HD11	1.88	0.55
1:B:164:VAL:HG11	1:B:210:LEU:HD22	1.93	0.51
1:A:190:LEU:HD13	1:B:60:MET:HG3	1.93	0.49
1:B:261:LEU:HA	1:B:264:ILE:HD12	1.94	0.48
1:A:313:GLU:CB	1:A:327:GLU:HB2	2.44	0.48
1:B:147:ALA:HB2	1:B:164:VAL:HG21	1.96	0.47
1:A:266:ARG:HG2	1:A:270:LYS:HD2	1.96	0.47
1:A:222:THR:HB	1:A:327:GLU:HB3	1.97	0.47
1:B:313:GLU:CB	1:B:327:GLU:HB2	2.45	0.47
1:A:122:ARG:HD2	3:A:417:HOH:O	2.16	0.45
1:A:147:ALA:HB2	1:A:164:VAL:HG11	1.99	0.45
1:B:20:VAL:HG22	1:B:89:GLY:HA3	1.98	0.44
1:B:175:LYS:HB2	1:B:176:PRO:CD	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:190:LEU:HA	1:B:191:PRO:HD3	1.84	0.44
1:B:310:ILE:HG13	1:B:330:VAL:HG23	1.99	0.43
1:A:229:SER:HA	1:A:230:PRO:HD3	1.88	0.43
1:A:313:GLU:HB3	1:A:327:GLU:CB	2.45	0.43
1:A:19:ARG:HD2	1:A:82:LEU:HD23	2.00	0.43
1:B:44:VAL:HG22	1:B:49:ALA:HB2	2.00	0.42
1:A:192:HIS:HD2	1:B:57:GLU:OE2	2.03	0.41
1:B:87:THR:HG23	1:B:110:GLU:HG3	2.02	0.41
1:B:123:ILE:O	1:B:308:MET:HG2	2.21	0.41
1:B:180:THR:HG23	1:B:210:LEU:HB3	2.02	0.41
1:B:175:LYS:HB2	1:B:176:PRO:HD2	2.03	0.41
1:B:253:LYS:HA	1:B:254:PRO:HD2	1.83	0.40
1:B:293:TRP:HA	1:B:294:PRO:HA	1.85	0.40
1:B:223:ARG:NH2	1:B:327:GLU:O	2.53	0.40

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	333/348 (96%)	320 (96%)	13 (4%)	0	100	100
1	B	308/348 (88%)	298 (97%)	9 (3%)	1 (0%)	46	41
All	All	641/696 (92%)	618 (96%)	22 (3%)	1 (0%)	52	48

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	46	GLY

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	289/300 (96%)	259 (90%)	30 (10%)	9	4
1	B	271/300 (90%)	218 (80%)	53 (20%)	1	0
All	All	560/600 (93%)	477 (85%)	83 (15%)	4	1

All (83) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	LYS
1	A	37	LEU
1	A	47	ASP
1	A	48	GLU
1	A	67	LEU
1	A	70	VAL
1	A	73	LYS
1	A	74	LYS
1	A	80	GLU
1	A	81	ILE
1	A	83	LYS
1	A	85	THR
1	A	94	LEU
1	A	109	LYS
1	A	127	LYS
1	A	135	LYS
1	A	144	VAL
1	A	145	MET
1	A	175	LYS
1	A	190	LEU
1	A	198	LYS
1	A	202	ARG
1	A	205	LEU
1	A	232	GLU
1	A	233	LYS
1	A	242	LEU
1	A	247	ARG

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Mol	Chain	Res	Type
1	A	282	LEU
1	A	321	LEU
1	A	343	THR
1	B	4	ARG
1	B	7	LYS
1	B	13	ASP
1	B	16	SER
1	B	26	VAL
1	B	47	ASP
1	B	48	GLU
1	B	50	THR
1	B	56	LEU
1	B	60	MET
1	B	64	GLU
1	B	66	LYS
1	B	67	LEU
1	B	70	VAL
1	B	74	LYS
1	B	78	ILE
1	B	81	ILE
1	B	94	LEU
1	B	97	SER
1	B	109	LYS
1	B	114	ILE
1	B	116	ASP
1	B	120	ASP
1	B	127	LYS
1	B	128	GLU
1	B	144	VAL
1	B	150	GLU
1	B	155	LYS
1	B	158	ARG
1	B	163	LYS
1	B	167	LEU
1	B	197	ASP
1	B	198	LYS
1	B	202	ARG
1	B	215	ASN
1	B	247	ARG
1	B	250	GLU
1	B	253	LYS
1	B	259	LYS

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Mol	Chain	Res	Type
1	B	261	LEU
1	B	269	ILE
1	B	270	LYS
1	B	271	GLU
1	B	278	PHE
1	B	279	ILE
1	B	295	ARG
1	B	301	GLU
1	B	302	THR
1	B	311	THR
1	B	338	LYS
1	B	342	LYS
1	B	344	GLU
1	B	345	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	192	HIS
1	B	215	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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 ⓘ

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, 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	337/348 (96%)	0.19	3 (0%) 85 86	15, 21, 25, 29	33 (9%)
1	B	318/348 (91%)	0.80	38 (11%) 6 6	13, 21, 25, 29	38 (11%)
All	All	655/696 (94%)	0.49	41 (6%) 23 24	13, 21, 25, 29	71 (10%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	255	GLY	6.1
1	B	293	TRP	6.1
1	B	278	PHE	5.9
1	B	269	ILE	5.1
1	B	202	ARG	4.7
1	B	266	ARG	4.5
1	B	259	LYS	4.2
1	A	83	LYS	4.1
1	B	273	GLY	4.0
1	B	322	GLY	3.8
1	B	257	THR	3.3
1	B	267	GLU	3.2
1	B	253	LYS	3.2
1	B	279	ILE	3.1
1	B	316	ILE	2.9
1	B	302	THR	2.9
1	B	275	GLY	2.8
1	B	336	GLY	2.8
1	B	48	GLU	2.7
1	B	186	HIS	2.7
1	B	299	TYR	2.6
1	B	307	GLY	2.6
1	B	313	GLU	2.5
1	B	272	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	301	GLU	2.4
1	B	254	PRO	2.3
1	B	317	TYR	2.3
1	A	55	GLU	2.3
1	B	198	LYS	2.2
1	B	261	LEU	2.2
1	B	224	THR	2.2
1	B	183	ALA	2.1
1	B	296	ILE	2.1
1	B	270	LYS	2.1
1	B	225	ILE	2.1
1	A	82	LEU	2.1
1	B	6	GLU	2.1
1	B	277	TYR	2.1
1	B	199	ARG	2.0
1	B	103	LYS	2.0
1	B	246	LYS	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, 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	LLDF	B-factors(Å ²)	Q<0.9
2	ZN	B	401	1/1	0.86	0.10	-0.94	55,55,55,55	0
2	ZN	B	402	1/1	0.95	0.07	-1.35	53,53,53,53	0
2	ZN	A	401	1/1	0.99	0.08	-2.68	37,37,37,37	0
2	ZN	A	402	1/1	0.99	0.07	-3.06	38,38,38,38	0

6.5 Other polymers ⓘ

There are no such residues in this entry.