



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 12:05 AM GMT

PDB ID : 2VHF
Title : STRUCTURE OF THE CYLD USP DOMAIN
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Barford, D.
Deposited on : 2007-11-21
Resolution : 2.80 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

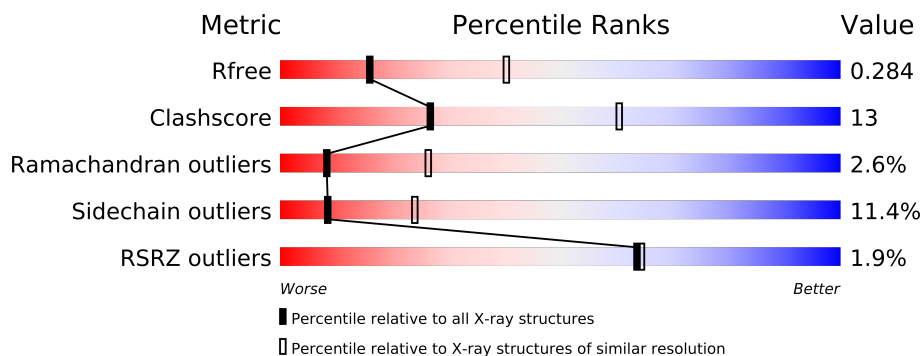
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.80 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}	66092	1799 (2.80-2.80)
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (2.80-2.80)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	374	
1	B	374	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5265 atoms, of which 0 are hydrogen and 0 are deuterium.

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 UBIQUITIN CARBOXYL-TERMINAL HYDROLASE CYLD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	0	0	0
			2653	1702	443	481	27			
1	B	334	Total	C	N	O	S	0	0	0
			2594	1665	427	475	27			

- 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.

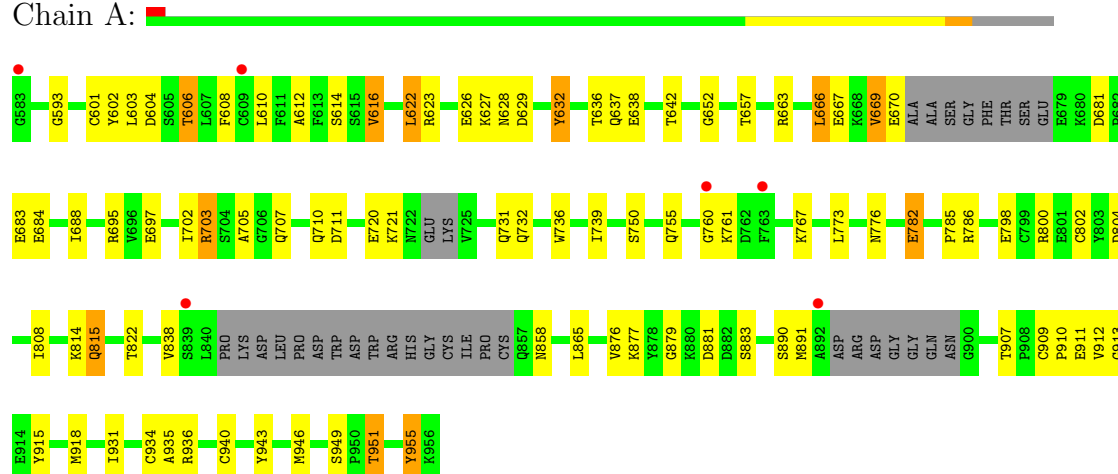
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	9	Total	O	0	0
			9	9		
3	B	5	Total	O	0	0
			5	5		

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.

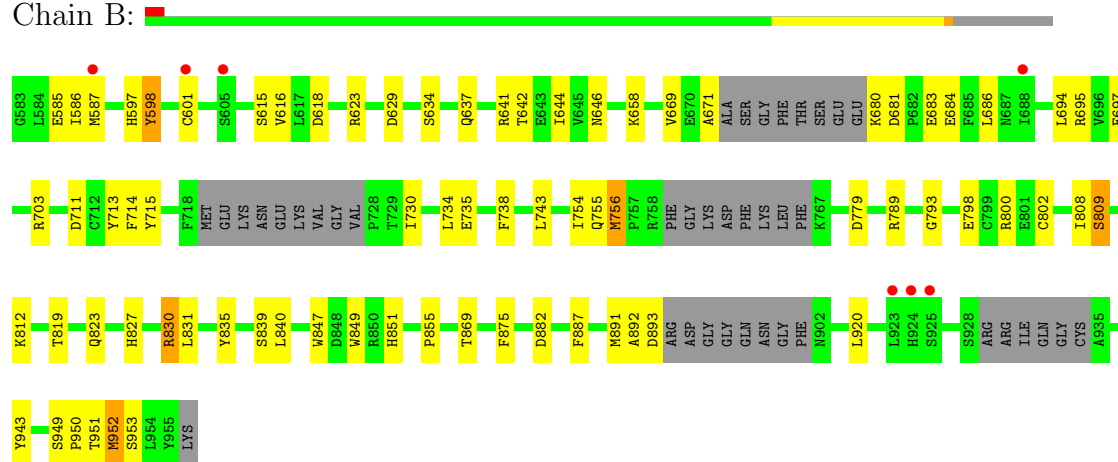
• Molecule 1: UBIQUITIN CARBOXYL-TERMINAL HYDROLASE CYLD

Chain A:



• Molecule 1: UBIQUITIN CARBOXYL-TERMINAL HYDROLASE CYLD

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	60.49Å 89.08Å 171.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.79 – 2.80 29.79 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.7 (29.79-2.80) 99.1 (29.79-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.233 , 0.281 0.239 , 0.284	Depositor DCC
R_{free} test set	764 reflections (3.38%)	DCC
Wilson B-factor (Å ²)	59.5	Xtriage
Anisotropy	0.672	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 63.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.20$	Xtriage
Outliers	0 of 23373 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	5265	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.93% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.46	1/2709 (0.0%)	0.56	0/3665
1	B	0.41	1/2651 (0.0%)	0.55	0/3597
All	All	0.44	2/5360 (0.0%)	0.56	0/7262

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	798	GLU	CD-OE2	7.02	1.33	1.25
1	B	798	GLU	CD-OE2	6.81	1.33	1.25

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2653	0	0	39	0
1	B	2594	0	0	29	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	9	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	5	0	0	0	0
All	All	5265	0	0	68	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 13.

All (68) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:616:VAL:CA	1:B:952:MET:CE	1.93	1.46
1:A:632:TYR:N	1:A:632:TYR:CD2	2.30	0.95
1:B:730:ILE:CD1	1:B:756:MET:CE	2.56	0.84
1:B:615:SER:O	1:B:952:MET:CE	2.36	0.73
1:B:616:VAL:N	1:B:952:MET:CE	2.53	0.71
1:B:830:ARG:CG	1:B:830:ARG:NH1	2.55	0.70
1:B:615:SER:C	1:B:952:MET:CE	2.68	0.62
1:B:793:GLY:CA	1:B:851:HIS:CD2	2.84	0.61
1:B:891:MET:O	1:B:893:ASP:N	2.38	0.56
1:A:915:TYR:O	1:A:918:MET:CG	2.54	0.56
1:A:666:LEU:O	1:A:670:GLU:N	2.40	0.55
1:A:626:GLU:N	1:A:629:ASP:OD2	2.40	0.54
1:B:808:ILE:O	1:B:809:SER:C	2.46	0.54
1:B:713:TYR:CD1	1:B:713:TYR:N	2.75	0.54
1:A:593:GLY:N	1:A:652:GLY:O	2.44	0.51
1:A:683:GLU:N	1:A:683:GLU:OE1	2.44	0.50
1:B:734:LEU:O	1:B:735:GLU:C	2.49	0.50
1:A:683:GLU:OE2	1:A:755:GLN:NE2	2.44	0.50
1:B:949:SER:O	1:B:951:THR:N	2.44	0.50
1:B:949:SER:C	1:B:951:THR:N	2.65	0.50
1:A:955:TYR:N	1:A:955:TYR:CD2	2.79	0.50
1:B:847:TRP:O	1:B:849:TRP:CD1	2.65	0.49
1:A:632:TYR:O	1:A:636:THR:OG1	2.30	0.49
1:B:697:GLU:CD	1:B:697:GLU:N	2.65	0.49
1:A:603:LEU:O	1:A:604:ASP:C	2.52	0.48
1:A:936:ARG:O	1:A:940:CYS:SG	2.71	0.48
1:B:951:THR:C	1:B:953:SER:N	2.65	0.48
1:A:936:ARG:CG	1:A:936:ARG:NH1	2.75	0.48
1:B:623:ARG:NH1	1:B:629:ASP:OD2	2.46	0.48
1:B:827:HIS:CB	1:B:830:ARG:NH1	2.77	0.47
1:B:715:TYR:O	1:B:755:GLN:N	2.47	0.47
1:B:835:TYR:O	1:B:835:TYR:CD1	2.67	0.47
1:A:702:ILE:O	1:A:710:GLN:N	2.47	0.47
1:A:951:THR:O	1:A:951:THR:OG1	2.32	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:681:ASP:OD1	1:A:684:GLU:N	2.48	0.47
1:A:720:GLU:O	1:A:721:LYS:C	2.54	0.47
1:B:597:HIS:O	1:B:598:TYR:C	2.52	0.46
1:A:760:GLY:O	1:A:761:LYS:C	2.54	0.46
1:A:909:CYS:SG	1:A:909:CYS:O	2.73	0.46
1:A:786:ARG:NH1	1:A:815:GLN:NE2	2.64	0.45
1:A:911:GLU:O	1:A:913:GLY:N	2.50	0.45
1:A:703:ARG:NH1	1:A:782:GLU:OE2	2.50	0.45
1:A:877:LYS:NZ	1:A:881:ASP:O	2.49	0.45
1:A:881:ASP:OD1	1:A:883:SER:N	2.50	0.43
1:A:683:GLU:OE2	1:A:943:TYR:OH	2.36	0.43
1:A:602:TYR:O	1:A:606:THR:CG2	2.66	0.43
1:A:626:GLU:O	1:A:628:ASN:N	2.51	0.43
1:A:935:ALA:O	1:A:936:ARG:C	2.57	0.43
1:B:875:PHE:CZ	1:B:887:PHE:CD1	3.07	0.43
1:A:669:VAL:O	1:A:670:GLU:C	2.57	0.42
1:B:680:LYS:CG	1:B:681:ASP:N	2.82	0.42
1:B:683:GLU:O	1:B:684:GLU:C	2.57	0.42
1:B:779:ASP:OD2	1:B:789:ARG:NH2	2.52	0.42
1:A:608:PHE:O	1:A:612:ALA:N	2.52	0.42
1:A:614:SER:OG	1:A:616:VAL:CG2	2.68	0.42
1:B:683:GLU:OE1	1:B:943:TYR:OH	2.38	0.42
1:A:736:TRP:N	1:A:736:TRP:CD1	2.87	0.42
1:A:622:LEU:N	1:A:622:LEU:CD1	2.83	0.42
1:A:638:GLU:O	1:A:642:THR:N	2.52	0.42
1:A:911:GLU:C	1:A:913:GLY:N	2.74	0.41
1:A:877:LYS:NZ	1:A:879:GLY:O	2.53	0.41
1:A:623:ARG:O	1:A:637:GLN:NE2	2.53	0.41
1:A:731:GLN:O	1:A:732:GLN:C	2.59	0.41
1:A:909:CYS:N	1:A:910:PRO:CD	2.83	0.41
1:B:694:LEU:O	1:B:695:ARG:C	2.59	0.41
1:A:808:ILE:O	1:A:814:LYS:NZ	2.54	0.41
1:B:669:VAL:O	1:B:671:ALA:N	2.54	0.41
1:B:641:ARG:O	1:B:646:ASN:CG	2.59	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	331/374 (88%)	271 (82%)	53 (16%)	7 (2%)	11	33
1	B	322/374 (86%)	268 (83%)	44 (14%)	10 (3%)	7	21
All	All	653/748 (87%)	539 (82%)	97 (15%)	17 (3%)	8	26

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	920	LEU
1	A	627	LYS
1	A	705	ALA
1	A	912	VAL
1	B	598	TYR
1	B	892	ALA
1	B	809	SER
1	A	601	CYS
1	A	782	GLU
1	B	855	PRO
1	B	869	THR
1	B	952	MET
1	A	785	PRO
1	B	585	GLU
1	B	950	PRO
1	B	644	ILE
1	A	739	ILE

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	281/340 (83%)	243 (86%)	38 (14%)	6	16
1	B	279/340 (82%)	253 (91%)	26 (9%)	13	35
All	All	560/680 (82%)	496 (89%)	64 (11%)	8	24

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	606	THR
1	A	610	LEU
1	A	616	VAL
1	A	622	LEU
1	A	632	TYR
1	A	657	THR
1	A	663	ARG
1	A	666	LEU
1	A	667	GLU
1	A	669	VAL
1	A	688	ILE
1	A	695	ARG
1	A	697	GLU
1	A	703	ARG
1	A	707	GLN
1	A	711	ASP
1	A	750	SER
1	A	767	LYS
1	A	773	LEU
1	A	776	ASN
1	A	800	ARG
1	A	802	CYS
1	A	804	ASP
1	A	815	GLN
1	A	822	THR
1	A	838	VAL
1	A	858	ASN
1	A	865	LEU
1	A	876	VAL
1	A	890	SER
1	A	891	MET
1	A	907	THR
1	A	931	ILE
1	A	934	CYS
1	A	946	MET

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Mol	Chain	Res	Type
1	A	949	SER
1	A	951	THR
1	A	955	TYR
1	B	586	ILE
1	B	587	MET
1	B	601	CYS
1	B	618	ASP
1	B	634	SER
1	B	637	GLN
1	B	642	THR
1	B	658	LYS
1	B	686	LEU
1	B	703	ARG
1	B	711	ASP
1	B	714	PHE
1	B	738	PHE
1	B	743	LEU
1	B	754	ILE
1	B	756	MET
1	B	800	ARG
1	B	802	CYS
1	B	812	LYS
1	B	819	THR
1	B	823	GLN
1	B	830	ARG
1	B	831	LEU
1	B	839	SER
1	B	840	LEU
1	B	882	ASP

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

5.3.3 RNA ⓘ

There are no RNA chains 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 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.

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	341/374 (91%)	-0.22	6 (1%) 65 66	26, 54, 97, 133	5 (1%)
1	B	334/374 (89%)	-0.08	7 (2%) 60 61	32, 62, 105, 135	9 (2%)
All	All	675/748 (90%)	-0.15	13 (1%) 64 64	26, 60, 101, 135	14 (2%)

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	892	ALA	4.2
1	B	587	MET	3.6
1	B	601	CYS	3.0
1	B	924	HIS	2.7
1	A	583	GLY	2.5
1	B	605	SER	2.4
1	B	925	SER	2.3
1	A	609	CYS	2.2
1	B	923	LEU	2.2
1	B	688	ILE	2.1
1	A	839	SER	2.1
1	A	760	GLY	2.1
1	A	763	PHE	2.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	ZN	B	1956	1/1	0.17	0.39	47,47,47,47	0
2	ZN	A	1958	1/1	0.13	-0.76	43,43,43,43	0
2	ZN	A	1957	1/1	0.13	-0.85	63,63,63,63	0
2	ZN	B	1957	1/1	0.11	-1.18	67,67,67,67	0

6.5 Other polymers

There are no such residues in this entry.