



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 01:27 am BST

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

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.11
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)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

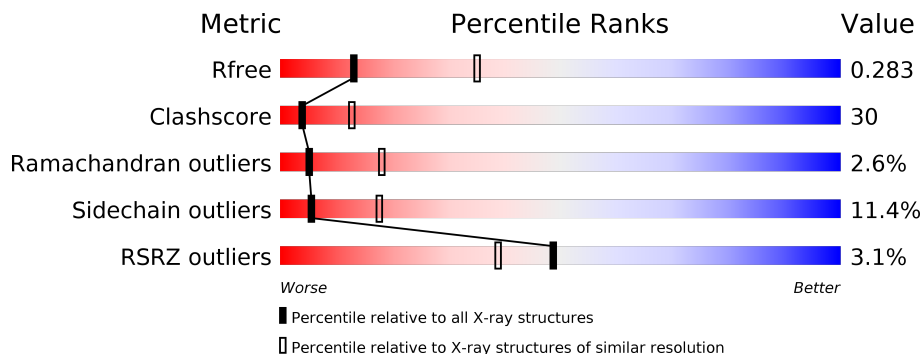
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (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 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	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 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 UBIQUITIN CARBOXYL-TERMINAL HYDROLASE CYLD.

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

- 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		

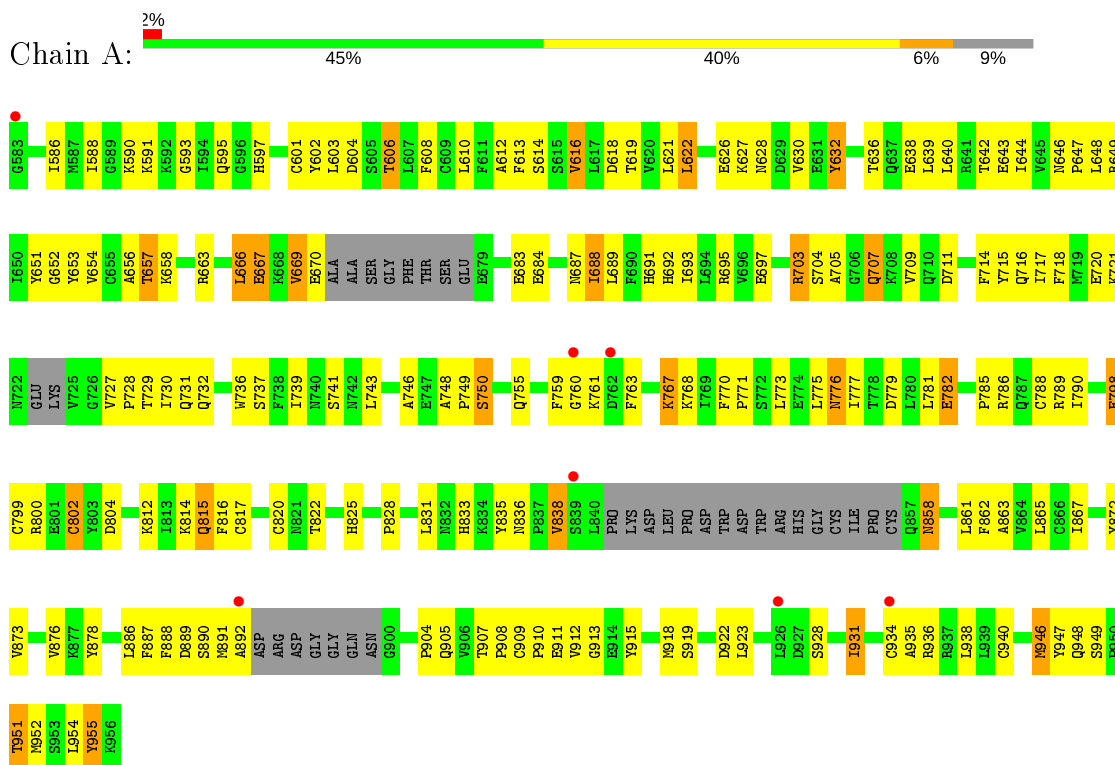
- Molecule 3 is water.

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

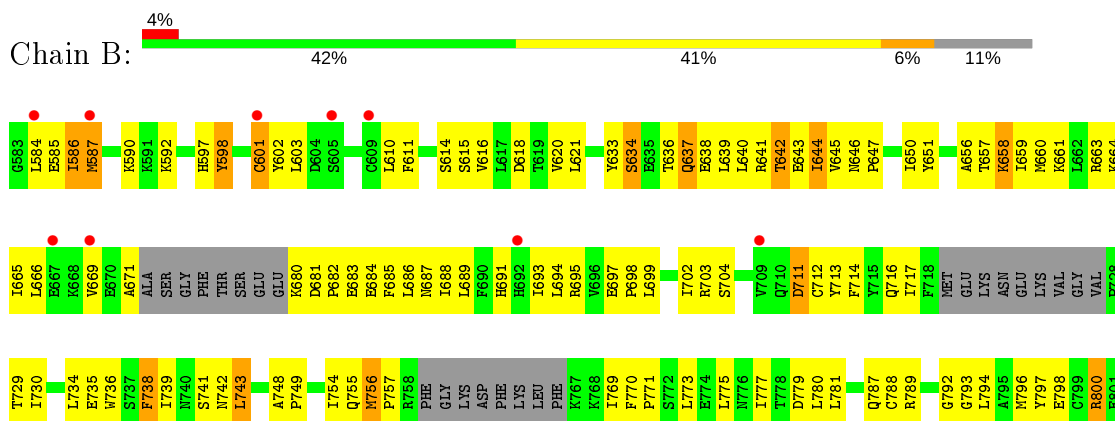
3 Residue-property plots [i](#)

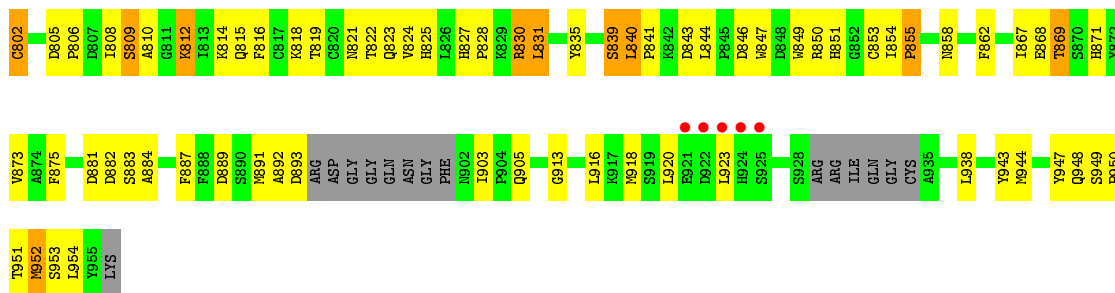
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: UBIQUITIN CARBOXYL-TERMINAL HYDROLASE CYLD



• Molecule 1: UBIQUITIN CARBOXYL-TERMINAL HYDROLASE CYLD





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)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.233 , 0.281 0.239 , 0.283	Depositor DCC
R_{free} test set	764 reflections (3.27%)	wwPDB-VP
Wilson B-factor (Å ²)	59.5	Xtrriage
Anisotropy	0.672	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 62.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.20$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5265	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtrriage'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.

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:
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 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	2653	0	2542	141	0
1	B	2594	0	2480	172	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	9	0	0	0	0
3	B	5	0	0	0	0
All	All	5265	0	5022	310	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 30.

All (310) 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:616:VAL:CA	1:B:952:MET:CE	1.93	1.45
1:B:616:VAL:HA	1:B:952:MET:CE	0.99	1.45
1:B:730:ILE:HD11	1:B:756:MET:CE	1.68	1.24
1:B:827:HIS:HB3	1:B:830:ARG:HH12	1.06	1.21
1:B:616:VAL:HA	1:B:952:MET:HE3	1.20	1.18
1:A:779:ASP:HA	1:A:786:ARG:HD3	1.20	1.15
1:B:730:ILE:HD11	1:B:756:MET:HE1	1.24	1.10
1:B:616:VAL:HA	1:B:952:MET:HE1	1.34	1.08
1:B:830:ARG:HG2	1:B:830:ARG:HH11	0.94	1.06
1:B:616:VAL:CA	1:B:952:MET:HE2	1.71	1.04
1:A:632:TYR:N	1:A:632:TYR:HD2	1.60	1.00
1:A:789:ARG:HH21	1:A:812:LYS:HB3	1.26	0.97
1:A:915:TYR:O	1:A:918:MET:HG3	1.66	0.94
1:B:730:ILE:CD1	1:B:756:MET:HE1	1.98	0.94
1:B:615:SER:O	1:B:952:MET:HE1	1.69	0.92
1:A:632:TYR:N	1:A:632:TYR:CD2	2.30	0.92
1:B:730:ILE:HD11	1:B:756:MET:HE3	1.50	0.92
1:A:616:VAL:HG13	1:A:952:MET:HG2	1.53	0.91
1:B:827:HIS:HB3	1:B:830:ARG:NH1	1.86	0.91
1:A:632:TYR:H	1:A:632:TYR:HD2	1.11	0.90
1:A:729:THR:HB	1:A:732:GLN:HG3	1.52	0.90
1:B:830:ARG:HG2	1:B:830:ARG:NH1	1.75	0.88
1:B:616:VAL:CA	1:B:952:MET:HE3	1.80	0.87
1:B:616:VAL:HA	1:B:952:MET:HE2	0.86	0.85
1:A:614:SER:OG	1:A:616:VAL:HG22	1.78	0.84
1:B:616:VAL:N	1:B:952:MET:HE3	1.93	0.83
1:A:936:ARG:HH11	1:A:936:ARG:HG2	1.41	0.83
1:A:789:ARG:NH2	1:A:812:LYS:HB3	1.93	0.82
1:A:595:GLN:HA	1:A:904:PRO:HG3	1.61	0.81
1:B:827:HIS:CB	1:B:830:ARG:HH12	1.92	0.80
1:B:793:GLY:CA	1:B:851:HIS:HD2	1.95	0.79
1:B:646:ASN:O	1:B:650:ILE:HG13	1.81	0.78
1:B:730:ILE:HG12	1:B:771:PRO:HA	1.65	0.78
1:A:931:ILE:HD11	1:A:936:ARG:HA	1.65	0.78
1:B:769:ILE:O	1:B:771:PRO:HD3	1.84	0.77
1:B:697:GLU:H	1:B:697:GLU:CD	1.86	0.76
1:B:730:ILE:CD1	1:B:756:MET:CE	2.56	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:830:ARG:CG	1:B:830:ARG:HH11	1.85	0.76
1:B:669:VAL:HG21	1:B:693:ILE:HD11	1.67	0.76
1:B:699:LEU:HB2	1:B:713:TYR:HA	1.68	0.75
1:B:633:TYR:O	1:B:637:GLN:HB2	1.86	0.75
1:B:680:LYS:NZ	1:B:681:ASP:HB2	2.02	0.74
1:B:639:LEU:HD12	1:B:643:GLU:HB2	1.69	0.74
1:B:843:ASP:O	1:B:846:ASP:HB2	1.87	0.74
1:A:918:MET:HB2	1:A:923:LEU:HD21	1.69	0.73
1:B:615:SER:O	1:B:952:MET:CE	2.36	0.73
1:A:786:ARG:HH11	1:A:815:GLN:HE21	1.36	0.73
1:A:863:ALA:HB3	1:A:946:MET:HG3	1.71	0.73
1:B:621:LEU:HD22	1:B:641:ARG:HA	1.69	0.73
1:A:878:TYR:HD2	1:A:886:LEU:HG	1.54	0.72
1:B:640:LEU:O	1:B:645:VAL:HG23	1.89	0.71
1:A:878:TYR:CD2	1:A:886:LEU:HG	2.25	0.71
1:B:616:VAL:N	1:B:952:MET:CE	2.53	0.71
1:B:788:CYS:SG	1:B:816:PHE:HA	2.30	0.71
1:B:881:ASP:HB3	1:B:884:ALA:HB2	1.71	0.71
1:A:684:GLU:O	1:A:688:ILE:HG22	1.90	0.70
1:A:779:ASP:CA	1:A:786:ARG:HD3	2.12	0.69
1:B:773:LEU:HD11	1:B:916:LEU:HD13	1.72	0.69
1:A:919:SER:O	1:A:922:ASP:HB2	1.91	0.69
1:B:602:TYR:HB2	1:B:682:PRO:HD3	1.73	0.69
1:B:680:LYS:HG2	1:B:681:ASP:N	2.07	0.68
1:B:730:ILE:CG1	1:B:771:PRO:HA	2.23	0.68
1:A:786:ARG:NH1	1:A:815:GLN:HE21	1.92	0.68
1:B:918:MET:HB2	1:B:923:LEU:HD13	1.75	0.68
1:A:588:ILE:HG23	1:A:908:PRO:HD3	1.74	0.67
1:B:634:SER:O	1:B:638:GLU:HB2	1.94	0.67
1:A:786:ARG:HH11	1:A:815:GLN:NE2	1.92	0.67
1:B:716:GLN:HB2	1:B:755:GLN:NE2	2.10	0.66
1:B:680:LYS:HZ3	1:B:681:ASP:HB2	1.61	0.66
1:B:687:ASN:O	1:B:691:HIS:HB3	1.95	0.66
1:B:951:THR:C	1:B:953:SER:H	1.99	0.66
1:A:704:SER:O	1:A:707:GLN:HB2	1.96	0.65
1:A:867:ILE:HD13	1:A:872:TYR:CD2	2.31	0.65
1:B:616:VAL:CA	1:B:952:MET:HE1	2.04	0.65
1:A:909:CYS:N	1:A:910:PRO:HD3	2.12	0.65
1:B:793:GLY:HA2	1:B:851:HIS:HD2	1.60	0.65
1:B:949:SER:C	1:B:951:THR:H	2.00	0.65
1:B:646:ASN:HB2	1:B:647:PRO:HD3	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:886:LEU:HD23	1:A:908:PRO:HA	1.78	0.64
1:A:715:TYR:OH	1:A:737:SER:HB3	1.97	0.64
1:A:928:SER:O	1:A:936:ARG:HG3	1.97	0.64
1:B:716:GLN:HA	1:B:755:GLN:HB2	1.80	0.64
1:A:591:LYS:HG2	1:A:653:TYR:CD1	2.33	0.63
1:B:867:ILE:HA	1:B:871:HIS:O	1.98	0.63
1:A:638:GLU:O	1:A:642:THR:HB	1.97	0.63
1:B:808:ILE:O	1:B:814:LYS:HE2	1.99	0.63
1:B:657:THR:HG22	1:B:658:LYS:HD2	1.81	0.63
1:B:738:PHE:CD1	1:B:780:LEU:HD22	2.33	0.63
1:A:759:PHE:CB	1:A:767:LYS:HE2	2.28	0.63
1:A:788:CYS:HB2	1:A:817:CYS:H	1.64	0.63
1:A:704:SER:HB3	1:A:743:LEU:HD23	1.80	0.63
1:B:615:SER:C	1:B:952:MET:CE	2.68	0.62
1:B:615:SER:C	1:B:952:MET:HE3	2.19	0.62
1:A:688:ILE:O	1:A:692:HIS:HB2	2.00	0.62
1:A:862:PHE:CE2	1:A:946:MET:HB2	2.34	0.62
1:A:936:ARG:HG2	1:A:936:ARG:NH1	2.12	0.61
1:B:610:LEU:O	1:B:614:SER:HB3	1.99	0.61
1:A:812:LYS:O	1:A:814:LYS:HE3	2.01	0.61
1:B:658:LYS:HD2	1:B:658:LYS:N	2.15	0.61
1:A:911:GLU:HA	1:A:911:GLU:OE2	2.01	0.61
1:A:590:LYS:HG2	1:A:905:GLN:HG3	1.83	0.60
1:A:657:THR:HG22	1:A:658:LYS:HD2	1.82	0.60
1:A:750:SER:O	1:A:948:GLN:HA	2.02	0.60
1:B:639:LEU:HD12	1:B:643:GLU:CB	2.31	0.60
1:B:640:LEU:HA	1:B:644:ILE:HD12	1.84	0.59
1:B:656:ALA:HA	1:B:659:ILE:HD13	1.83	0.59
1:B:660:MET:O	1:B:664:LYS:HG2	2.01	0.59
1:A:703:ARG:HH11	1:A:782:GLU:CD	2.05	0.59
1:B:611:PHE:HB3	1:B:645:VAL:HG13	1.84	0.59
1:A:640:LEU:HA	1:A:644:ILE:HD12	1.83	0.59
1:A:889:ASP:HB3	1:A:892:ALA:HB2	1.84	0.59
1:B:841:PRO:HD2	1:B:844:LEU:HD12	1.85	0.59
1:B:793:GLY:CA	1:B:851:HIS:CD2	2.84	0.58
1:A:618:ASP:O	1:A:622:LEU:HD13	2.03	0.58
1:B:893:ASP:HB3	1:B:903:ILE:HB	1.86	0.58
1:A:746:ALA:HA	1:A:782:GLU:HG3	1.85	0.58
1:B:666:LEU:HD21	1:B:685:PHE:CE1	2.38	0.58
1:A:799:CYS:HB2	1:A:835:TYR:HB3	1.86	0.57
1:A:616:VAL:CG1	1:A:952:MET:HG2	2.32	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:642:THR:HG22	1:A:643:GLU:HG2	1.87	0.57
1:A:776:ASN:HA	1:A:858:ASN:ND2	2.20	0.57
1:A:715:TYR:HE2	1:A:717:ILE:HG22	1.69	0.56
1:B:891:MET:O	1:B:893:ASP:N	2.38	0.56
1:B:775:LEU:HG	1:B:947:TYR:CE2	2.40	0.56
1:B:787:GLN:O	1:B:815:GLN:HG3	2.06	0.56
1:A:911:GLU:C	1:A:913:GLY:H	2.09	0.56
1:A:586:ILE:HG22	1:A:651:TYR:O	2.06	0.55
1:B:736:TRP:O	1:B:739:ILE:HG22	2.07	0.55
1:B:850:ARG:HB3	1:B:853:CYS:HB3	1.87	0.55
1:A:716:GLN:HA	1:A:755:GLN:HB2	1.89	0.55
1:A:760:GLY:HA3	1:A:763:PHE:O	2.07	0.55
1:A:867:ILE:HD13	1:A:872:TYR:CE2	2.41	0.55
1:A:666:LEU:O	1:A:669:VAL:HG13	2.05	0.55
1:A:588:ILE:CG2	1:A:908:PRO:HD3	2.37	0.54
1:A:608:PHE:O	1:A:612:ALA:HB2	2.07	0.54
1:B:642:THR:HA	1:B:646:ASN:ND2	2.22	0.54
1:A:689:LEU:HA	1:A:693:ILE:HD12	1.89	0.54
1:B:610:LEU:HB2	1:B:611:PHE:CD1	2.42	0.54
1:B:586:ILE:HG23	1:B:651:TYR:O	2.08	0.54
1:A:748:ALA:HB2	1:A:781:LEU:HD11	1.90	0.54
1:B:713:TYR:CD1	1:B:713:TYR:N	2.75	0.54
1:B:806:PRO:HA	1:B:810:ALA:HB2	1.90	0.54
1:B:808:ILE:O	1:B:809:SER:C	2.46	0.54
1:B:793:GLY:HA2	1:B:851:HIS:CD2	2.43	0.54
1:B:666:LEU:HD21	1:B:685:PHE:CD1	2.43	0.53
1:A:931:ILE:HD11	1:A:936:ARG:CA	2.36	0.53
1:A:727:VAL:HG23	1:A:768:LYS:O	2.09	0.53
1:A:788:CYS:HB2	1:A:817:CYS:N	2.24	0.53
1:A:718:PHE:HZ	1:A:741:SER:HB2	1.74	0.52
1:B:854:ILE:HD12	1:B:854:ILE:N	2.24	0.52
1:A:802:CYS:HB3	1:A:814:LYS:HG3	1.91	0.52
1:A:687:ASN:O	1:A:691:HIS:HB3	2.10	0.52
1:B:602:TYR:HB2	1:B:682:PRO:CD	2.40	0.52
1:A:915:TYR:C	1:A:918:MET:HG3	2.26	0.52
1:A:776:ASN:HD22	1:A:777:ILE:N	2.07	0.52
1:A:649:ARG:NE	1:A:954:LEU:HD11	2.25	0.51
1:A:828:PRO:HA	1:A:831:LEU:HG	1.91	0.51
1:B:616:VAL:CB	1:B:952:MET:HE2	2.39	0.51
1:B:640:LEU:HD22	1:B:689:LEU:HD21	1.91	0.51
1:A:798:GLU:HG2	1:A:838:VAL:HG12	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:584:LEU:C	1:B:586:ILE:H	2.14	0.51
1:B:867:ILE:HG13	1:B:871:HIS:O	2.11	0.51
1:B:738:PHE:CE1	1:B:780:LEU:HB3	2.46	0.50
1:A:591:LYS:HG2	1:A:653:TYR:CE1	2.46	0.50
1:A:729:THR:HG22	1:A:731:GLN:H	1.76	0.50
1:A:798:GLU:HG3	1:A:836:ASN:HD21	1.76	0.50
1:A:667:GLU:HA	1:A:670:GLU:HB3	1.93	0.50
1:B:734:LEU:O	1:B:735:GLU:C	2.49	0.50
1:A:730:ILE:CD1	1:A:771:PRO:HA	2.42	0.50
1:A:873:VAL:HG21	1:A:887:PHE:CE1	2.47	0.50
1:A:683:GLU:OE2	1:A:755:GLN:NE2	2.44	0.50
1:B:873:VAL:HG12	1:B:889:ASP:OD1	2.11	0.50
1:B:729:THR:HA	1:B:770:PHE:O	2.11	0.50
1:A:955:TYR:N	1:A:955:TYR:CD2	2.79	0.50
1:B:704:SER:HB3	1:B:743:LEU:HD12	1.94	0.50
1:B:789:ARG:HG2	1:B:815:GLN:HG2	1.94	0.50
1:A:613:PHE:HB3	1:A:955:TYR:HE2	1.77	0.49
1:A:789:ARG:HH21	1:A:812:LYS:CB	2.12	0.49
1:B:777:ILE:HD12	1:B:781:LEU:HD11	1.93	0.49
1:A:886:LEU:CD2	1:A:908:PRO:HA	2.41	0.49
1:B:847:TRP:O	1:B:849:TRP:CD1	2.65	0.49
1:B:695:ARG:HB3	1:B:695:ARG:NH1	2.28	0.49
1:B:839:SER:O	1:B:840:LEU:HD13	2.13	0.49
1:A:825:HIS:CD2	1:A:833:HIS:HB2	2.48	0.49
1:B:639:LEU:HD12	1:B:643:GLU:CD	2.33	0.49
1:A:728:PRO:HD2	1:A:768:LYS:O	2.12	0.49
1:A:861:LEU:HB2	1:A:947:TYR:CE2	2.47	0.49
1:A:602:TYR:CE2	1:A:603:LEU:HD13	2.48	0.48
1:A:597:HIS:CD2	1:A:656:ALA:HB1	2.48	0.48
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:621:LEU:HB3	1:B:641:ARG:HG3	1.95	0.48
1:A:714:PHE:HE1	1:A:755:GLN:OE1	1.96	0.48
1:B:610:LEU:HB2	1:B:611:PHE:HD1	1.77	0.48
1:B:739:ILE:HG12	1:B:739:ILE:O	2.13	0.48
1:B:681:ASP:HB3	1:B:683:GLU:OE2	2.13	0.48
1:B:830:ARG:CG	1:B:830:ARG:NH1	2.55	0.48
1:A:622:LEU:HD11	1:B:849:TRP:CE3	2.49	0.48
1:B:841:PRO:HD2	1:B:844:LEU:CD1	2.43	0.48
1:A:717:ILE:HG12	1:A:755:GLN:O	2.13	0.48
1:A:936:ARG:CG	1:A:936:ARG:NH1	2.75	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:787:GLN:HG2	1:B:792:GLY:HA2	1.95	0.48
1:B:949:SER:C	1:B:951:THR:N	2.65	0.48
1:B:793:GLY:N	1:B:851:HIS:HD2	2.11	0.48
1:B:802:CYS:HB3	1:B:814:LYS:HG3	1.96	0.48
1:A:771:PRO:HG2	1:A:938:LEU:HD21	1.95	0.48
1:A:648:LEU:HB2	1:A:654:VAL:CG2	2.44	0.48
1:B:841:PRO:HB2	1:B:844:LEU:HG	1.94	0.47
1:A:788:CYS:SG	1:A:816:PHE:HA	2.54	0.47
1:A:619:THR:CG2	1:B:855:PRO:HG3	2.44	0.47
1:B:951:THR:C	1:B:953:SER:N	2.65	0.47
1:B:835:TYR:CD1	1:B:835:TYR:O	2.67	0.47
1:A:639:LEU:HD12	1:A:643:GLU:HB2	1.95	0.47
1:B:620:VAL:HG13	1:B:694:LEU:HD13	1.96	0.47
1:B:827:HIS:CG	1:B:828:PRO:HD2	2.50	0.47
1:A:951:THR:OG1	1:A:951:THR:O	2.32	0.47
1:A:825:HIS:HD2	1:A:833:HIS:HB2	1.80	0.47
1:B:949:SER:O	1:B:951:THR:N	2.44	0.47
1:A:889:ASP:O	1:A:892:ALA:HB2	2.14	0.47
1:B:704:SER:HA	1:B:742:ASN:O	2.14	0.47
1:B:699:LEU:HD12	1:B:713:TYR:O	2.15	0.47
1:A:593:GLY:HA3	1:A:888:PHE:CE1	2.50	0.47
1:A:720:GLU:O	1:A:721:LYS:C	2.54	0.47
1:B:875:PHE:CG	1:B:938:LEU:HD13	2.50	0.47
1:A:619:THR:HG22	1:B:855:PRO:HG3	1.97	0.47
1:A:630:VAL:HB	1:A:632:TYR:CE2	2.50	0.46
1:B:698:PRO:HB3	1:B:711:ASP:HB3	1.98	0.46
1:B:775:LEU:O	1:B:858:ASN:HA	2.16	0.46
1:B:597:HIS:O	1:B:598:TYR:C	2.52	0.46
1:A:730:ILE:HG13	1:A:770:PHE:O	2.15	0.46
1:A:760:GLY:O	1:A:761:LYS:C	2.54	0.46
1:A:646:ASN:HB2	1:A:647:PRO:HD3	1.98	0.46
1:B:797:TYR:HE1	1:B:818:LYS:HD3	1.80	0.46
1:A:595:GLN:HB3	1:A:656:ALA:HB2	1.98	0.46
1:A:909:CYS:O	1:A:909:CYS:SG	2.73	0.46
1:B:789:ARG:NH1	1:B:812:LYS:HD3	2.32	0.45
1:B:659:ILE:O	1:B:663:ARG:HB2	2.16	0.45
1:B:587:MET:O	1:B:592:LYS:HB3	2.17	0.45
1:B:816:PHE:HB2	1:B:821:ASN:HB2	1.99	0.45
1:B:702:ILE:HD12	1:B:712:CYS:SG	2.57	0.45
1:B:794:LEU:HD23	1:B:796:MET:CE	2.47	0.45
1:B:875:PHE:CD1	1:B:938:LEU:HD13	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:717:ILE:HG13	1:B:757:PRO:HD3	1.98	0.45
1:B:590:LYS:HG3	1:B:905:GLN:NE2	2.32	0.44
1:A:667:GLU:HA	1:A:670:GLU:CB	2.46	0.44
1:A:703:ARG:NH1	1:A:782:GLU:OE2	2.50	0.44
1:B:586:ILE:O	1:B:592:LYS:HD2	2.17	0.44
1:B:825:HIS:O	1:B:831:LEU:HB3	2.17	0.44
1:B:748:ALA:HA	1:B:749:PRO:HD2	1.83	0.43
1:B:683:GLU:HG3	1:B:944:MET:CE	2.48	0.43
1:A:602:TYR:O	1:A:606:THR:CG2	2.66	0.43
1:B:802:CYS:O	1:B:814:LYS:HG3	2.18	0.43
1:A:626:GLU:O	1:A:628:ASN:N	2.51	0.43
1:A:781:LEU:HD23	1:A:782:GLU:N	2.34	0.43
1:A:935:ALA:O	1:A:936:ARG:C	2.57	0.43
1:B:818:LYS:O	1:B:822:THR:HG23	2.17	0.43
1:A:798:GLU:HG3	1:A:836:ASN:ND2	2.33	0.43
1:B:680:LYS:HZ1	1:B:681:ASP:HB2	1.80	0.43
1:B:883:SER:HB3	1:B:913:GLY:HA3	2.01	0.43
1:A:790:ILE:HB	1:A:820:CYS:SG	2.58	0.43
1:B:875:PHE:CZ	1:B:887:PHE:CD1	3.07	0.43
1:B:601:CYS:C	1:B:603:LEU:H	2.22	0.42
1:A:586:ILE:HG22	1:A:586:ILE:O	2.20	0.42
1:A:669:VAL:O	1:A:670:GLU:C	2.57	0.42
1:A:748:ALA:HA	1:A:749:PRO:HD3	1.91	0.42
1:B:741:SER:HB3	1:B:743:LEU:HD13	2.00	0.42
1:B:805:ASP:HA	1:B:806:PRO:HD3	1.70	0.42
1:B:868:GLU:C	1:B:869:THR:HG1	2.22	0.42
1:B:614:SER:C	1:B:954:LEU:HD12	2.39	0.42
1:B:683:GLU:O	1:B:684:GLU:C	2.57	0.42
1:A:597:HIS:HD2	1:A:656:ALA:HB1	1.83	0.42
1:B:779:ASP:OD2	1:B:789:ARG:NH2	2.52	0.42
1:B:597:HIS:HE1	1:B:659:ILE:HG22	1.84	0.42
1:B:844:LEU:C	1:B:846:ASP:H	2.22	0.42
1:A:736:TRP:N	1:A:736:TRP:CD1	2.87	0.42
1:B:824:VAL:HG12	1:B:825:HIS:ND1	2.35	0.42
1:B:952:MET:HB3	1:B:952:MET:HE3	1.70	0.42
1:A:622:LEU:CD1	1:A:622:LEU:N	2.83	0.42
1:B:615:SER:HB3	1:B:954:LEU:HG	2.02	0.42
1:A:703:ARG:NH2	1:A:709:VAL:HG22	2.34	0.41
1:B:683:GLU:HG3	1:B:944:MET:HE1	2.02	0.41
1:A:775:LEU:HA	1:A:775:LEU:HD12	1.85	0.41
1:A:911:GLU:C	1:A:913:GLY:N	2.74	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:825:HIS:C	1:B:827:HIS:H	2.22	0.41
1:B:636:THR:OG1	1:B:665:ILE:HD13	2.21	0.41
1:B:685:PHE:O	1:B:688:ILE:HG22	2.20	0.41
1:B:862:PHE:CE2	1:B:948:GLN:HB3	2.56	0.41
1:A:669:VAL:CG2	1:A:688:ILE:HG13	2.50	0.41
1:A:918:MET:HE2	1:A:923:LEU:CD2	2.51	0.41
1:A:731:GLN:O	1:A:732:GLN:C	2.59	0.41
1:B:658:LYS:HE2	1:B:661:LYS:NZ	2.35	0.41
1:B:694:LEU:O	1:B:695:ARG:C	2.59	0.41
1:B:738:PHE:HD1	1:B:780:LEU:HD22	1.81	0.41
1:B:777:ILE:CD1	1:B:781:LEU:HD11	2.50	0.41
1:B:669:VAL:O	1:B:671:ALA:N	2.54	0.41
1:A:788:CYS:CB	1:A:817:CYS:H	2.31	0.41
1:A:909:CYS:N	1:A:910:PRO:CD	2.83	0.41
1:A:776:ASN:CG	1:A:858:ASN:HD21	2.24	0.40
1:B:756:MET:HB2	1:B:943:TYR:O	2.21	0.40
1:B:621:LEU:HD21	1:B:645:VAL:HG21	2.02	0.40
1:B:680:LYS:HG2	1:B:681:ASP:HB2	2.02	0.40
1:B:697:GLU:CD	1:B:697:GLU:N	2.65	0.40
1:A:618:ASP:HA	1:A:621:LEU:HD12	2.02	0.40
1:A:632:TYR:O	1:A:636:THR:OG1	2.30	0.40
1:A:715:TYR:CE2	1:A:717:ILE:HG22	2.54	0.40
1:B:800:ARG:C	1:B:800:ARG:HD3	2.42	0.40
1:A:595:GLN:HA	1:A:904:PRO:CG	2.41	0.40
1:A:593:GLY:N	1:A:652:GLY:O	2.44	0.40
1:B:641:ARG:O	1:B:646:ASN:CG	2.59	0.40
1:B:680:LYS:NZ	1:B:684:GLU:HB2	2.37	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	331/374 (88%)	271 (82%)	53 (16%)	7 (2%)	7	23
1	B	322/374 (86%)	268 (83%)	44 (14%)	10 (3%)	4	14
All	All	653/748 (87%)	539 (82%)	97 (15%)	17 (3%)	5	18

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 [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	A	281/340 (83%)	243 (86%)	38 (14%)	4	11
1	B	279/340 (82%)	253 (91%)	26 (9%)	9	26
All	All	560/680 (82%)	496 (89%)	64 (11%)	5	18

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
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

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Mol	Chain	Res	Type
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. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	597	HIS
1	A	707	GLN
1	A	710	GLN
1	A	776	ASN
1	A	787	GLN
1	A	815	GLN
1	A	821	ASN
1	A	836	ASN
1	A	858	ASN
1	A	905	GLN
1	B	597	HIS
1	B	637	GLN
1	B	716	GLN
1	B	740	ASN
1	B	755	GLN
1	B	821	ASN
1	B	836	ASN

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Mol	Chain	Res	Type
1	B	851	HIS
1	B	905	GLN

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 [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

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%)	0.01	7 (2%) 63 54	26, 54, 97, 133	5 (1%)
1	B	334/374 (89%)	0.17	14 (4%) 36 26	32, 62, 105, 135	9 (2%)
All	All	675/748 (90%)	0.09	21 (3%) 49 39	26, 60, 101, 135	14 (2%)

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	892	ALA	4.2
1	B	922	ASP	3.9
1	B	601	CYS	3.3
1	B	605	SER	3.2
1	B	921	GLU	3.1
1	B	587	MET	2.9
1	B	709	VAL	2.8
1	B	692	HIS	2.6
1	B	584	LEU	2.6
1	A	760	GLY	2.5
1	B	924	HIS	2.5
1	A	839	SER	2.4
1	B	925	SER	2.4
1	A	583	GLY	2.3
1	B	609	CYS	2.2
1	A	762	ASP	2.2
1	B	667	GLU	2.2
1	A	926	LEU	2.1
1	B	923	LEU	2.1
1	A	934	CYS	2.0
1	B	669	VAL	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. 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(\AA^2)	Q<0.9
2	ZN	A	1957	1/1	0.98	0.13	63,63,63,63	0
2	ZN	B	1956	1/1	0.99	0.14	47,47,47,47	0
2	ZN	A	1958	1/1	0.99	0.12	43,43,43,43	0
2	ZN	B	1957	1/1	0.99	0.11	67,67,67,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.