



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 08:13 PM UTC

PDB ID : 1Q67 / pdb_00001q67
Title : Crystal structure of Dcp1p
Authors : She, M.; Decker, C.J.; Liu, Y.; Chen, N.; Parker, R.; Song, H.
Deposited on : 2003-08-12
Resolution : 2.30 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)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

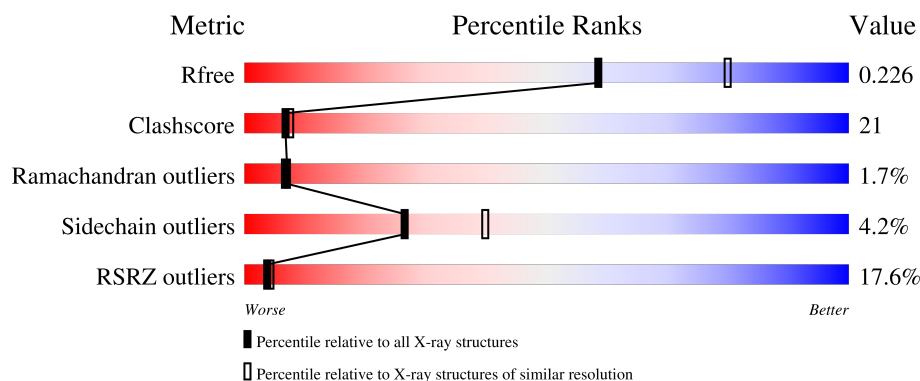
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	231	12% 34% 26% • 37%
1	B	231	11% 44% 24% • 30%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2726 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 Decapping protein involved in mRNA degradation-Dcp1p.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	146	Total	C	N	O	S	0	0	0
			1222	794	203	222	3			
1	B	161	Total	C	N	O	S	0	0	0
			1348	876	225	244	3			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	67	Total	O	0	0
			67	67		
2	B	89	Total	O	0	0
			89	89		

4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	123.55Å 123.55Å 77.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.30 20.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.8 (20.00-2.30) 97.6 (20.00-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.97 (at 2.20Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.227 , 0.268 0.228 , 0.226	Depositor DCC
R_{free} test set	1727 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	40.5	Xtriage
Anisotropy	0.301	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 51.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.033 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2726	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 32.41 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 9.4345e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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](#)

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	0.39	0/1250	0.94	10/1692 (0.6%)
1	B	0.38	0/1380	0.90	5/1867 (0.3%)
All	All	0.38	0/2630	0.92	15/3559 (0.4%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	41	THR	CA-C-N	7.19	126.45	118.97
1	B	41	THR	C-N-CA	7.19	126.45	118.97
1	B	191	VAL	N-CA-C	-6.26	98.86	107.99
1	A	41	THR	CA-C-N	5.69	125.37	119.05
1	A	41	THR	C-N-CA	5.69	125.37	119.05
1	A	227	LYS	N-CA-C	5.56	122.65	110.80
1	A	199	GLU	N-CA-C	-5.51	100.71	109.96
1	A	190	LEU	N-CA-C	5.40	118.62	109.76
1	A	191	VAL	N-CA-C	-5.35	99.38	107.73
1	A	31	ASP	CA-C-N	5.25	124.92	119.56
1	A	31	ASP	C-N-CA	5.25	124.92	119.56
1	B	46	LEU	N-CA-C	5.20	117.71	109.24
1	B	23	ASN	N-CA-C	-5.14	100.70	108.67
1	A	225	GLU	CA-C-N	5.04	126.14	119.84
1	A	225	GLU	C-N-CA	5.04	126.14	119.84

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	1222	0	1225	69	0
1	B	1348	0	1349	46	0
2	A	67	0	0	6	0
2	B	89	0	0	8	0
All	All	2726	0	2574	110	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 21.

All (110) 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:184:VAL:HG11	1:A:218:ILE:HG21	1.50	0.92
1:B:76:THR:HG22	1:B:78:LEU:H	1.32	0.92
1:A:72:VAL:HG23	1:A:136:ASP:HB3	1.53	0.88
1:A:227:LYS:HG2	1:A:228:ASP:H	1.42	0.83
1:B:58:LYS:HZ1	1:B:62:GLN:HE21	1.23	0.82
1:A:69:LEU:HD11	1:A:227:LYS:HG3	1.62	0.80
1:B:184:VAL:HG11	1:B:218:ILE:HG21	1.63	0.80
1:A:219:LYS:O	1:A:223:GLU:HG2	1.83	0.79
1:A:184:VAL:HG11	1:A:218:ILE:CG2	2.14	0.78
1:A:195:ASN:ND2	1:A:199:GLU:H	1.83	0.77
1:A:195:ASN:HD21	1:A:199:GLU:H	1.33	0.75
1:A:72:VAL:CG2	1:A:136:ASP:HB3	2.20	0.72
1:A:72:VAL:HG21	1:A:136:ASP:O	1.94	0.67
1:B:184:VAL:HG11	1:B:218:ILE:CG2	2.25	0.66
1:A:227:LYS:CG	1:A:228:ASP:H	2.08	0.66
1:A:146:ASN:ND2	1:A:151:ASP:HB2	2.10	0.66
1:B:223:GLU:HG3	2:B:301:HOH:O	1.95	0.65
1:B:172:GLU:HB2	2:B:266:HOH:O	1.98	0.64
1:A:227:LYS:HG2	1:A:228:ASP:N	2.12	0.63
1:A:206:HIS:CD2	1:B:206:HIS:NE2	2.67	0.63
1:A:62:GLN:NE2	1:B:58:LYS:HG2	2.14	0.63
1:A:145:LEU:HD23	1:A:152:ASN:OD1	1.99	0.62
1:A:184:VAL:HG12	1:A:193:ILE:HG12	1.82	0.61
1:A:177:ASN:ND2	1:A:197:LYS:NZ	2.48	0.60
1:A:72:VAL:HG22	1:A:72:VAL:O	2.02	0.59
1:B:230:PHE:O	1:B:231:ALA:CB	2.51	0.59
1:A:41:THR:HG22	1:A:65:LEU:HB3	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:204:TRP:HH2	1:B:206:HIS:HD1	1.51	0.58
1:A:148:ILE:HD11	1:B:56:TRP:HB3	1.85	0.58
1:A:177:ASN:HD21	1:A:197:LYS:HZ3	1.52	0.57
1:B:184:VAL:HG12	1:B:193:ILE:HG12	1.86	0.57
1:B:50:ASP:OD2	1:B:53:LYS:HB2	2.04	0.57
1:B:146:ASN:ND2	1:B:151:ASP:HB2	2.19	0.57
1:A:158:VAL:HG21	1:A:196:LEU:HD11	1.87	0.56
1:B:166:ARG:NE	1:B:179:LEU:HD12	2.21	0.56
1:B:58:LYS:NZ	1:B:62:GLN:HE21	1.99	0.55
1:A:227:LYS:HB2	2:A:265:HOH:O	2.06	0.54
1:B:225:GLU:HG3	1:B:226:PRO:HD2	1.89	0.54
1:A:148:ILE:HG23	2:B:235:HOH:O	2.07	0.54
1:A:195:ASN:HB2	2:A:261:HOH:O	2.07	0.54
1:B:227:LYS:HD3	1:B:231:ALA:HB3	1.89	0.53
1:A:62:GLN:HE22	1:B:58:LYS:HG2	1.74	0.53
1:A:166:ARG:NE	1:A:179:LEU:HD12	2.23	0.53
1:B:170:ASN:ND2	1:B:173:GLU:HB2	2.23	0.53
1:A:177:ASN:HD21	1:A:197:LYS:NZ	2.07	0.52
1:A:23:ASN:ND2	1:A:27:ILE:HG13	2.24	0.52
1:B:171:ALA:O	1:B:175:THR:HG23	2.10	0.52
1:B:58:LYS:HZ1	1:B:62:GLN:NE2	2.01	0.52
1:B:140:TYR:HB2	1:B:157:ILE:HB	1.92	0.52
1:A:177:ASN:ND2	1:A:197:LYS:HZ3	2.07	0.51
1:B:40:HIS:HD2	2:B:299:HOH:O	1.92	0.51
1:B:149:ASN:HB2	1:B:150:PRO:HD2	1.92	0.51
1:B:219:LYS:HD2	2:B:311:HOH:O	2.11	0.51
1:A:140:TYR:HB2	1:A:157:ILE:HB	1.93	0.51
1:A:148:ILE:HD11	1:B:56:TRP:CB	2.41	0.51
1:A:22:LEU:O	1:A:23:ASN:HB2	2.10	0.51
1:A:65:LEU:HD21	1:A:214:ILE:CD1	2.42	0.50
1:A:195:ASN:HD22	1:A:195:ASN:C	2.19	0.50
1:B:163:VAL:O	1:B:167:LYS:HG3	2.12	0.50
1:A:187:LYS:O	1:A:188:ASP:HB2	2.11	0.49
1:A:148:ILE:HG13	1:A:148:ILE:O	2.12	0.49
1:A:58:LYS:NZ	1:A:62:GLN:HG3	2.28	0.49
1:B:229:SER:O	1:B:230:PHE:C	2.57	0.48
1:B:80:PRO:HD3	2:B:283:HOH:O	2.12	0.48
1:A:40:HIS:HA	1:A:65:LEU:O	2.14	0.47
1:A:182:MET:HG2	1:A:195:ASN:HA	1.95	0.47
1:A:139:ASN:HB3	2:A:265:HOH:O	2.14	0.47
1:A:179:LEU:HD22	1:A:196:LEU:HB3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:33:LYS:O	1:B:34:ILE:C	2.58	0.47
1:B:177:ASN:HD21	1:B:197:LYS:HE2	1.80	0.47
1:A:58:LYS:HZ1	1:A:62:GLN:HE21	1.62	0.46
1:A:211:ARG:NH2	2:A:263:HOH:O	2.47	0.46
1:A:227:LYS:N	2:A:265:HOH:O	2.47	0.46
1:B:142:LEU:C	1:B:142:LEU:HD23	2.41	0.46
1:A:153:PHE:HE2	1:A:155:MET:HE3	1.81	0.46
1:B:138:TYR:CD2	1:B:156:GLY:HA3	2.50	0.46
1:A:72:VAL:O	1:A:73:SER:C	2.59	0.45
1:B:62:GLN:O	1:B:146:ASN:HA	2.16	0.45
1:B:172:GLU:O	1:B:176:LEU:HG	2.16	0.45
1:B:184:VAL:CG1	1:B:218:ILE:HG21	2.40	0.45
1:B:194:LYS:NZ	2:B:265:HOH:O	2.49	0.45
1:B:188:ASP:OD2	1:B:189:GLU:HG3	2.17	0.45
1:A:140:TYR:CZ	1:A:221:LEU:HD22	2.52	0.44
1:A:62:GLN:O	1:A:146:ASN:HA	2.17	0.44
1:B:62:GLN:N	1:B:62:GLN:CD	2.75	0.44
1:A:153:PHE:CE2	1:A:155:MET:HE3	2.52	0.44
1:B:40:HIS:HA	1:B:65:LEU:O	2.18	0.44
1:A:42:PRO:HB3	1:A:147:ARG:HH21	1.83	0.43
1:A:206:HIS:O	1:A:207:THR:C	2.61	0.43
1:A:62:GLN:N	1:A:62:GLN:CD	2.77	0.43
1:B:19:ARG:HG3	1:B:19:ARG:NH1	2.32	0.43
1:A:159:PRO:HG2	1:A:226:PRO:HG3	2.00	0.43
1:A:195:ASN:ND2	1:A:195:ASN:C	2.77	0.43
1:A:50:ASP:HB2	1:A:57:ASN:ND2	2.35	0.42
1:B:61:TYR:C	1:B:62:GLN:CD	2.87	0.42
1:A:177:ASN:ND2	1:A:197:LYS:HZ1	2.15	0.42
1:A:142:LEU:C	1:A:142:LEU:HD23	2.45	0.42
1:B:212:GLN:HG3	2:B:303:HOH:O	2.19	0.42
1:A:24:PHE:HB3	2:A:270:HOH:O	2.18	0.41
1:A:58:LYS:HE3	1:A:62:GLN:HE21	1.86	0.41
1:B:187:LYS:HE3	1:B:187:LYS:HB2	1.86	0.41
1:A:41:THR:HA	1:A:42:PRO:HD3	1.85	0.41
1:A:138:TYR:CD2	1:A:156:GLY:HA3	2.56	0.41
1:A:227:LYS:CG	1:A:228:ASP:N	2.74	0.41
1:A:165:LYS:HE3	1:A:165:LYS:HB2	1.90	0.41
1:B:79:LEU:HA	1:B:80:PRO:HD3	1.89	0.41
1:A:33:LYS:HD2	1:A:70:ARG:NH2	2.35	0.40
1:A:184:VAL:CG1	1:A:218:ILE:HG21	2.36	0.40
1:A:140:TYR:CE2	1:A:221:LEU:HD22	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:33:LYS:HD2	1:A:70:ARG:HH22	1.86	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	142/231 (62%)	132 (93%)	6 (4%)	4 (3%)	4	2
1	B	157/231 (68%)	150 (96%)	6 (4%)	1 (1%)	21	27
All	All	299/462 (65%)	282 (94%)	12 (4%)	5 (2%)	7	7

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	23	ASN
1	A	227	LYS
1	B	230	PHE
1	A	207	THR
1	A	60	GLU

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	137/208 (66%)	129 (94%)	8 (6%)	18	26
1	B	150/208 (72%)	146 (97%)	4 (3%)	39	58
All	All	287/416 (69%)	275 (96%)	12 (4%)	26	40

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	145	LEU
1	A	148	ILE
1	A	154	SER
1	A	170	ASN
1	A	172	GLU
1	A	176	LEU
1	A	195	ASN
1	A	212	GLN
1	B	23	ASN
1	B	54	ASP
1	B	167	LYS
1	B	223	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	ASN
1	A	62	GLN
1	A	177	ASN
1	A	195	ASN
1	A	206	HIS
1	B	23	ASN
1	B	40	HIS
1	B	43	HIS
1	B	62	GLN
1	B	170	ASN
1	B	177	ASN
1	B	212	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	146/231 (63%)	0.97	28 (19%) 3 3	31, 46, 69, 79	0
1	B	161/231 (69%)	0.82	26 (16%) 4 5	31, 44, 71, 78	0
All	All	307/462 (66%)	0.89	54 (17%) 4 4	31, 45, 71, 79	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	228	ASP	6.3
1	B	80	PRO	6.2
1	A	22	LEU	5.1
1	B	230	PHE	4.7
1	B	176	LEU	4.7
1	A	206	HIS	4.5
1	B	195	ASN	4.2
1	B	79	LEU	4.1
1	A	227	LYS	4.1
1	A	23	ASN	3.8
1	B	17	PHE	3.7
1	B	76	THR	3.6
1	B	175	THR	3.4
1	B	170	ASN	3.3
1	A	184	VAL	3.2
1	B	174	ASP	3.2
1	A	189	GLU	3.1
1	A	175	THR	3.1
1	B	172	GLU	2.9
1	A	186	VAL	2.9
1	A	53	LYS	2.9
1	A	174	ASP	2.7
1	B	228	ASP	2.7
1	B	206	HIS	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	208	VAL	2.7
1	A	176	LEU	2.7
1	B	73	SER	2.6
1	A	55	GLU	2.6
1	B	75	ASN	2.6
1	B	78	LEU	2.6
1	A	226	PRO	2.6
1	A	25	ASN	2.5
1	A	209	SER	2.5
1	A	213	ASN	2.5
1	B	198	HIS	2.5
1	B	208	VAL	2.4
1	A	198	HIS	2.4
1	A	135	LYS	2.4
1	A	207	THR	2.4
1	B	177	ASN	2.3
1	B	74	GLN	2.3
1	B	19	ARG	2.3
1	A	178	PRO	2.2
1	B	173	GLU	2.2
1	B	188	ASP	2.2
1	A	173	GLU	2.2
1	B	171	ALA	2.2
1	B	71	ASP	2.2
1	B	189	GLU	2.1
1	A	71	ASP	2.0
1	A	152	ASN	2.0
1	A	51	PHE	2.0
1	A	188	ASP	2.0
1	A	148	ILE	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.