



Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 02:34 AM GMT

PDB ID : 1QZ7
Title : Beta-catenin binding domain of Axin in complex with beta-catenin
Authors : Xing, Y.; Clements, W.K.; Kimelman, D.; Xu, W.
Deposited on : 2003-09-15
Resolution : 2.20 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>

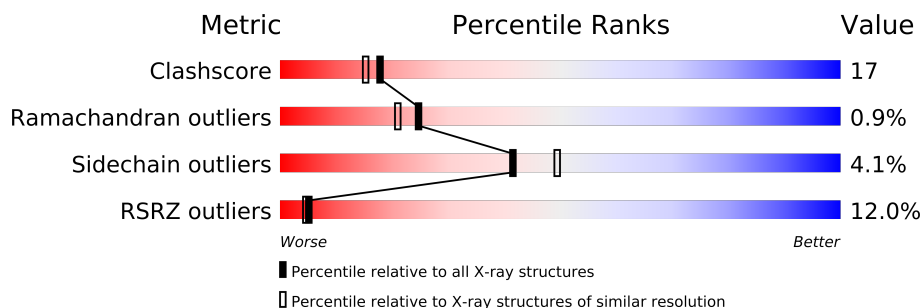
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.20 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	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (2.20-2.20)

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	533	
2	B	70	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4148 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 Beta-catenin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	524	Total	C	N	O	S	0	0	0
			3979	2496	721	735	27			

- Molecule 2 is a protein called Axin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	17	Total	C	N	O	S	0	0	0
			138	83	24	30	1			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	31	Total	O	0	0
			31	31		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	85.14Å 75.04Å 101.40Å 90.00° 97.47° 90.00°	Depositor
Resolution (Å)	42.21 – 2.20 42.21 – 2.09	Depositor EDS
% Data completeness (in resolution range)	81.1 (42.21-2.20) 93.5 (42.21-2.09)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.10Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.221 , 0.256 0.231 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	35.7	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 37253 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4148	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.75% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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	0.31	0/4035	0.53	0/5476
2	B	0.28	0/139	0.87	1/187 (0.5%)
All	All	0.31	0/4174	0.55	1/5663 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	475	GLU	CB-CA-C	9.99	130.38	110.40

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	3979	0	4094	129	0
2	B	138	0	124	21	0
3	A	31	0	0	0	0
All	All	4148	0	4218	143	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 17.

All (143) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:159:LEU:HD23	1:A:171:ALA:HB2	1.48	0.96
1:A:271:MET:O	1:A:275:LEU:HG	1.84	0.78
1:A:550:ARG:HH22	1:A:561:VAL:CG2	1.99	0.76
1:A:272:ALA:HA	1:A:275:LEU:HD12	1.68	0.75
1:A:364:GLN:H	1:A:364:GLN:HE21	1.35	0.74
1:A:364:GLN:H	1:A:364:GLN:NE2	1.85	0.74
1:A:355:PRO:O	1:A:359:GLU:HG2	1.87	0.74
1:A:270:LYS:HB3	1:A:274:ARG:HH12	1.55	0.70
2:B:468:ASN:HD22	2:B:469:PRO:CD	2.05	0.70
1:A:154:PRO:O	1:A:158:LYS:HG3	1.91	0.70
1:A:257:THR:HG23	2:B:476:HIS:ND1	2.07	0.69
1:A:153:ILE:HB	1:A:154:PRO:HD3	1.75	0.69
1:A:508:LYS:HE3	1:A:568:GLU:OE1	1.92	0.68
1:A:201:THR:O	1:A:205:THR:HG22	1.93	0.68
1:A:188:ILE:HG23	1:A:194:MET:HB3	1.77	0.67
2:B:468:ASN:HD22	2:B:469:PRO:HD2	1.60	0.66
1:A:371:THR:O	1:A:371:THR:HG22	1.95	0.66
1:A:200:ARG:HA	1:A:203:GLN:HE21	1.61	0.66
1:A:292:LYS:O	1:A:296:ILE:HG12	1.95	0.66
1:A:278:GLY:O	1:A:282:MET:HG3	1.97	0.65
1:A:189:MET:SD	1:A:226:GLU:HB2	2.38	0.63
1:A:483:ASN:ND2	1:A:486:ARG:NH2	2.48	0.62
1:A:270:LYS:HB3	1:A:274:ARG:NH1	2.14	0.61
1:A:189:MET:HB2	1:A:221:LEU:HD22	1.80	0.61
1:A:533:ILE:HB	1:A:534:PRO:HD3	1.83	0.60
1:A:190:ARG:O	1:A:192:PRO:HD3	2.02	0.60
2:B:468:ASN:HD22	2:B:469:PRO:N	2.00	0.59
1:A:173:VAL:HG12	1:A:177:GLN:HE21	1.68	0.58
1:A:458:GLU:OE2	1:A:503:HIS:HD2	1.86	0.58
1:A:557:GLN:O	1:A:558:GLN:HG2	2.03	0.58
1:A:550:ARG:HD3	1:A:558:GLN:OE1	2.04	0.58
1:A:159:LEU:HD11	1:A:167:VAL:HG12	1.85	0.57
1:A:296:ILE:HD12	2:B:476:HIS:HD2	1.69	0.57
1:A:181:LYS:HB2	1:A:184:SER:OG	2.03	0.57
1:A:399:GLU:H	1:A:399:GLU:CD	2.08	0.56
1:A:518:ALA:CB	1:A:580:LEU:HD13	2.36	0.56
1:A:200:ARG:HA	1:A:203:GLN:NE2	2.22	0.55
1:A:159:LEU:CD2	1:A:171:ALA:HB2	2.29	0.54
1:A:371:THR:CG2	1:A:371:THR:O	2.55	0.54
1:A:525:ALA:HB3	1:A:526:PRO:HD3	1.90	0.53
1:A:347:LEU:O	1:A:350:CYS:HB3	2.09	0.53
1:A:174:MET:HA	1:A:177:GLN:NE2	2.23	0.53
1:A:188:ILE:HG23	1:A:194:MET:CB	2.38	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:559:GLN:HG2	1:A:560:PHE:N	2.23	0.53
1:A:171:ALA:O	1:A:175:VAL:HG23	2.09	0.53
1:A:547:THR:HA	1:A:558:GLN:NE2	2.23	0.52
2:B:472:ILE:C	2:B:472:ILE:HD13	2.29	0.52
1:A:164:ASP:OD2	1:A:167:VAL:HG23	2.09	0.52
1:A:558:GLN:OE1	1:A:558:GLN:HA	2.10	0.52
1:A:354:LYS:HZ3	1:A:354:LYS:HB2	1.74	0.52
1:A:205:THR:O	1:A:205:THR:HG23	2.10	0.52
1:A:550:ARG:HH22	1:A:561:VAL:HG21	1.73	0.51
1:A:543:ALA:O	1:A:547:THR:HG23	2.10	0.51
1:A:542:ARG:HG2	1:A:542:ARG:HH11	1.75	0.51
1:A:535:ARG:HH11	1:A:538:GLN:NE2	2.09	0.51
1:A:200:ARG:O	1:A:204:ASN:ND2	2.44	0.51
1:A:149:ALA:O	1:A:153:ILE:HG12	2.11	0.50
1:A:535:ARG:NH1	1:A:538:GLN:NE2	2.59	0.50
1:A:171:ALA:HA	1:A:174:MET:HE3	1.94	0.50
1:A:186:HIS:HA	1:A:189:MET:HE2	1.93	0.50
1:A:230:ALA:HA	1:A:233:LYS:CE	2.41	0.50
2:B:468:ASN:C	2:B:470:GLU:H	2.14	0.49
1:A:227:GLY:O	1:A:231:ILE:HG13	2.12	0.49
1:A:658:VAL:HG12	1:A:662:MET:CE	2.42	0.49
1:A:257:THR:HG23	2:B:476:HIS:CE1	2.46	0.49
1:A:483:ASN:HD21	1:A:486:ARG:NH2	2.11	0.49
2:B:479:ARG:CB	2:B:479:ARG:HH11	2.26	0.48
1:A:159:LEU:HD21	1:A:167:VAL:HG12	1.94	0.48
1:A:178:LEU:HB3	1:A:184:SER:HB2	1.95	0.48
1:A:160:LEU:HD12	1:A:201:THR:OG1	2.14	0.48
1:A:518:ALA:HB2	1:A:580:LEU:HD13	1.95	0.48
1:A:431:ASN:O	1:A:435:LYS:HG3	2.13	0.48
1:A:448:VAL:O	1:A:452:LEU:HD22	2.15	0.47
1:A:552:SER:O	1:A:553:MET:HB2	2.15	0.47
1:A:253:PHE:HB3	2:B:472:ILE:HD12	1.97	0.47
1:A:159:LEU:HD11	1:A:167:VAL:CG1	2.45	0.47
1:A:354:LYS:HE3	1:A:387:ASN:O	2.15	0.47
1:A:547:THR:HA	1:A:558:GLN:HE22	1.80	0.47
1:A:354:LYS:HB2	1:A:354:LYS:NZ	2.30	0.47
1:A:253:PHE:HD2	2:B:469:PRO:O	1.98	0.47
1:A:145:ASP:O	1:A:149:ALA:HB3	2.15	0.47
1:A:490:GLY:O	1:A:493:VAL:HG12	2.15	0.47
1:A:149:ALA:HA	1:A:152:ALA:HB3	1.97	0.46
2:B:468:ASN:HD22	2:B:468:ASN:C	2.18	0.46
1:A:428:THR:O	1:A:435:LYS:HE2	2.15	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:172:ALA:CB	1:A:210:THR:HG23	2.46	0.46
1:A:308:ASN:OD1	1:A:310:GLU:HB2	2.15	0.46
1:A:177:GLN:C	1:A:179:SER:H	2.18	0.46
2:B:475:GLU:O	2:B:478:GLN:HB2	2.15	0.46
1:A:179:SER:OG	1:A:217:THR:HG23	2.16	0.46
1:A:561:VAL:O	1:A:562:GLU:HB2	2.16	0.46
1:A:226:GLU:H	1:A:226:GLU:CD	2.19	0.46
1:A:346:VAL:O	1:A:349:VAL:HG22	2.16	0.46
1:A:491:LEU:HB2	1:A:492:PRO:HD3	1.98	0.45
1:A:152:ALA:O	1:A:155:GLU:HB3	2.16	0.45
1:A:173:VAL:HG12	1:A:177:GLN:NE2	2.30	0.45
1:A:164:ASP:O	1:A:168:VAL:HG23	2.16	0.45
1:A:538:GLN:NE2	1:A:542:ARG:HH21	2.15	0.45
1:A:447:LEU:O	1:A:451:VAL:HG23	2.16	0.45
1:A:414:ILE:H	1:A:414:ILE:HG13	1.60	0.45
1:A:198:ILE:HD13	1:A:217:THR:HG21	1.99	0.45
1:A:518:ALA:HB1	1:A:580:LEU:HD13	1.99	0.45
2:B:468:ASN:O	2:B:470:GLU:N	2.50	0.44
1:A:173:VAL:O	1:A:177:GLN:HG3	2.16	0.44
1:A:640:LEU:CD1	1:A:662:MET:HE1	2.48	0.44
1:A:308:ASN:O	1:A:312:LYS:HG3	2.16	0.44
1:A:523:ASN:O	1:A:527:LEU:HB2	2.17	0.44
1:A:195:VAL:O	1:A:199:VAL:HG23	2.17	0.44
1:A:354:LYS:HD3	1:A:391:ALA:HB2	2.00	0.44
1:A:560:PHE:N	1:A:560:PHE:CD1	2.86	0.44
1:A:458:GLU:OE2	1:A:503:HIS:CD2	2.70	0.44
1:A:567:GLU:H	1:A:567:GLU:CD	2.20	0.44
1:A:196:SER:O	1:A:200:ARG:HG3	2.18	0.43
1:A:350:CYS:HB3	1:A:353:ASN:HB2	1.99	0.43
2:B:479:ARG:HB3	2:B:479:ARG:HH11	1.82	0.43
1:A:162:ASP:HB3	1:A:168:VAL:HG22	2.01	0.43
2:B:479:ARG:HB2	2:B:479:ARG:NH1	2.33	0.43
1:A:225:ARG:HG3	1:A:225:ARG:HH11	1.84	0.43
1:A:545:GLN:O	1:A:549:ARG:HG3	2.19	0.43
1:A:257:THR:HG23	2:B:476:HIS:CG	2.53	0.43
1:A:658:VAL:HG12	1:A:662:MET:HE1	1.99	0.43
1:A:658:VAL:O	1:A:662:MET:HG3	2.19	0.43
1:A:237:ILE:HB	1:A:238:PRO:HD3	2.01	0.43
1:A:270:LYS:O	1:A:274:ARG:HG3	2.18	0.43
1:A:659:LEU:HD23	1:A:662:MET:HE3	2.00	0.42
1:A:549:ARG:HH11	1:A:549:ARG:HG2	1.83	0.42
1:A:659:LEU:HA	1:A:662:MET:HE2	2.01	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:479:ARG:CB	2:B:479:ARG:NH1	2.82	0.42
2:B:472:ILE:HG23	2:B:473:LEU:N	2.35	0.42
1:A:397:GLY:HA2	1:A:399:GLU:OE1	2.19	0.42
1:A:296:ILE:HD12	2:B:476:HIS:CD2	2.52	0.42
1:A:230:ALA:HA	1:A:233:LYS:HE3	2.00	0.41
1:A:153:ILE:HG21	1:A:191:SER:OG	2.21	0.41
1:A:160:LEU:HD22	1:A:171:ALA:CB	2.51	0.41
1:A:483:ASN:HD21	1:A:486:ARG:HH22	1.68	0.41
1:A:228:LEU:HD22	1:A:262:LEU:HD23	2.03	0.41
1:A:156:LEU:HD12	1:A:178:LEU:HD12	2.02	0.41
1:A:354:LYS:HB3	1:A:355:PRO:CD	2.51	0.40
1:A:188:ILE:HG12	1:A:194:MET:HE2	2.02	0.40
1:A:302:GLN:NE2	1:A:342:ARG:HH21	2.19	0.40
1:A:493:VAL:HG13	1:A:494:VAL:N	2.36	0.40
2:B:468:ASN:ND2	2:B:469:PRO:HD2	2.33	0.40
1:A:229:LEU:HG	1:A:233:LYS:HE2	2.02	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	522/533 (98%)	488 (94%)	30 (6%)	4 (1%)	27	24
2	B	15/70 (21%)	13 (87%)	1 (7%)	1 (7%)	2	0
All	All	537/603 (89%)	501 (93%)	31 (6%)	5 (1%)	25	21

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	553	MET
1	A	208	VAL
1	A	555	GLY
2	B	469	PRO

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Mol	Chain	Res	Type
1	A	166	VAL

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	428/442 (97%)	414 (97%)	14 (3%)	50	60
2	B	16/62 (26%)	12 (75%)	4 (25%)	1	0
All	All	444/504 (88%)	426 (96%)	18 (4%)	41	49

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	163	GLU
1	A	252	LEU
1	A	254	TYR
1	A	302	GLN
1	A	364	GLN
1	A	437	MET
1	A	452	LEU
1	A	462	GLU
1	A	480	MET
1	A	483	ASN
1	A	527	LEU
1	A	567	GLU
1	A	580	LEU
1	A	665	ASP
2	B	468	ASN
2	B	472	ILE
2	B	474	ASP
2	B	479	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	177	GLN

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Mol	Chain	Res	Type
1	A	186	HIS
1	A	193	GLN
1	A	203	GLN
1	A	220	ASN
1	A	302	GLN
1	A	322	GLN
1	A	326	ASN
1	A	364	GLN
1	A	415	ASN
1	A	483	ASN
1	A	499	HIS
1	A	503	HIS
1	A	538	GLN
1	A	601	GLN
1	A	611	GLN
2	B	468	ASN

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 ⓘ

There are no ligands in this entry.

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	524/533 (98%)	0.72	62 (11%) 5 5	22, 47, 99, 108	0
2	B	17/70 (24%)	1.17	3 (17%) 2 2	71, 81, 94, 94	0
All	All	541/603 (89%)	0.74	65 (12%) 5 4	22, 48, 99, 108	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	160	LEU	12.7
1	A	142	TYR	11.9
1	A	551	THR	8.9
1	A	144	ASP	8.3
1	A	146	ALA	8.3
1	A	143	GLN	8.3
1	A	148	LEU	7.9
1	A	178	LEU	7.7
1	A	149	ALA	7.6
1	A	552	SER	6.9
1	A	183	ALA	6.6
1	A	159	LEU	6.5
1	A	556	THR	6.3
2	B	481	MET	5.9
1	A	145	ASP	5.8
1	A	150	THR	5.7
1	A	174	MET	5.4
1	A	181	LYS	5.3
1	A	557	GLN	5.2
1	A	165	GLN	5.2
1	A	167	VAL	5.1
1	A	147	GLU	5.0
1	A	186	HIS	4.8
1	A	177	GLN	4.8

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Mol	Chain	Res	Type	RSRZ
1	A	156	LEU	4.6
1	A	153	ILE	4.5
1	A	184	SER	4.4
1	A	352	SER	4.4
1	A	185	ARG	4.2
1	A	166	VAL	4.2
1	A	168	VAL	4.2
1	A	189	MET	4.1
1	A	555	GLY	4.0
1	A	558	GLN	4.0
1	A	550	ARG	4.0
1	A	665	ASP	3.9
1	A	175	VAL	3.8
1	A	151	ARG	3.8
1	A	188	ILE	3.6
1	A	164	ASP	3.5
1	A	172	ALA	3.4
1	A	206	ASN	3.4
1	A	559	GLN	3.2
1	A	161	ASN	3.0
1	A	228	LEU	2.9
1	A	182	GLU	2.9
1	A	549	ARG	2.8
1	A	190	ARG	2.8
1	A	221	LEU	2.6
1	A	187	ALA	2.6
1	A	489	TYR	2.6
1	A	169	ASN	2.6
1	A	205	THR	2.6
2	B	480	VAL	2.6
1	A	208	VAL	2.4
2	B	475	GLU	2.4
1	A	152	ALA	2.3
1	A	560	PHE	2.3
1	A	193	GLN	2.2
1	A	171	ALA	2.2
1	A	213	CYS	2.2
1	A	232	PHE	2.1
1	A	350	CYS	2.1
1	A	197	ALA	2.1
1	A	542	ARG	2.0

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 ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.