



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:24 PM UTC

PDB ID : 2PFV / pdb_00002pfv
Title : S. cerevisiae Exo70 with additional residues to 2.1 Angstrom resolution
Authors : Moore, B.A.; Robinson, H.H.; Xu, Z.
Deposited on : 2007-04-05
Resolution : 2.10 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

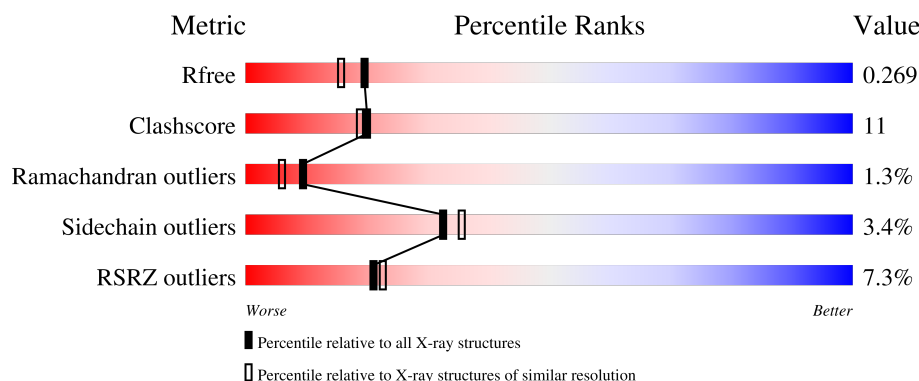
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	563	7% 72% 23% ...

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4415 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 Exocyst complex component EXO70.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	549	Total	C	N	O	S	0	0	0
			4304	2733	727	826	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	61	GLY	-	cloning artifact	UNP P19658
A	62	SER	-	cloning artifact	UNP P19658

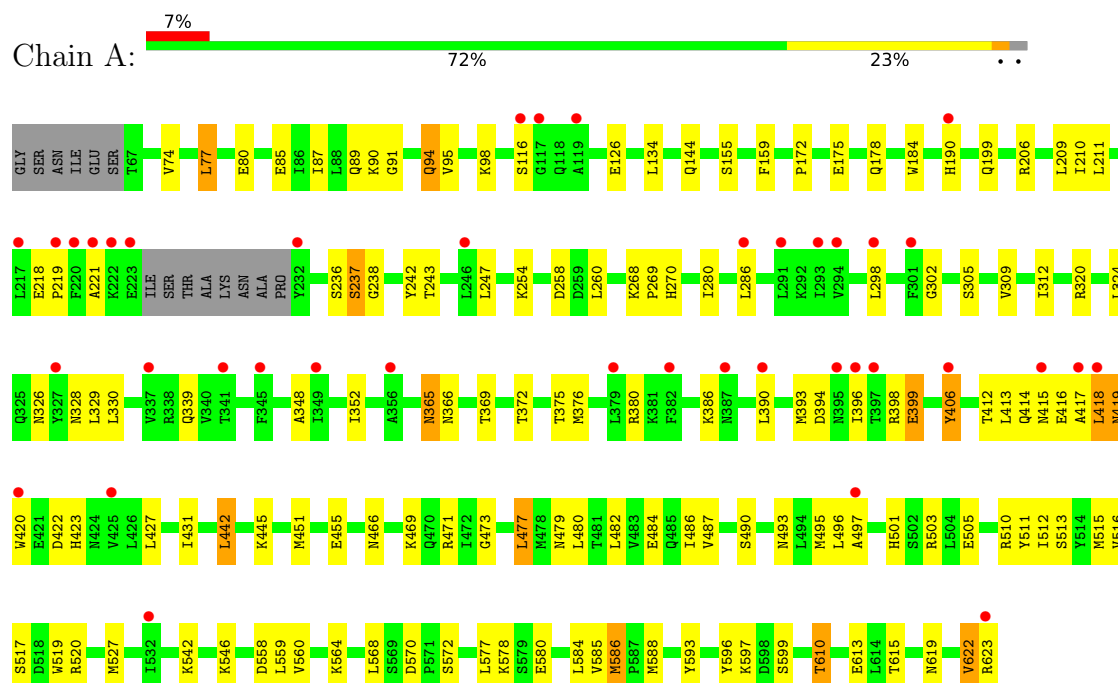
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	111	Total	O	0	0
			111	111		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Exocyst complex component EXO70



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	45.54Å 60.07Å 222.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.84 – 2.10 40.84 – 2.10	Depositor EDS
% Data completeness (in resolution range)	88.3 (40.84-2.10) 90.0 (40.84-2.10)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 2.10Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.237 , 0.263 0.245 , 0.269	Depositor DCC
R_{free} test set	3128 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	30.5	Xtriage
Anisotropy	0.518	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 30.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4415	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.85% 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](#)

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	0.44	0/4376	0.88	12/5909 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	585	VAL	N-CA-C	9.88	119.73	110.74
1	A	610	THR	N-CA-C	-9.78	98.88	110.13
1	A	512	ILE	N-CA-C	-9.50	101.13	110.72
1	A	599	SER	N-CA-C	-8.32	103.12	113.18
1	A	219	PRO	N-CA-CB	6.38	109.95	103.25
1	A	260	LEU	N-CA-C	5.78	118.52	111.82
1	A	442	LEU	N-CA-C	-5.53	105.34	111.36
1	A	155	SER	N-CA-C	-5.31	105.61	111.71
1	A	455	GLU	N-CA-C	-5.28	103.00	109.65
1	A	596	TYR	N-CA-C	5.07	120.31	113.72
1	A	190	HIS	N-CA-C	-5.04	105.91	111.71
1	A	586	MET	N-CA-C	5.00	119.55	112.75

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	4304	0	4222	97	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	111	0	0	0	0
All	All	4415	0	4222	97	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 11.

All (97) 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:477:LEU:HD23	1:A:515:MET:HE2	1.45	0.99
1:A:326:ASN:HB2	1:A:329:LEU:HD23	1.45	0.95
1:A:366:ASN:HD21	1:A:471:ARG:HE	1.11	0.94
1:A:477:LEU:CD2	1:A:515:MET:HE2	2.13	0.77
1:A:366:ASN:ND2	1:A:471:ARG:HE	1.84	0.76
1:A:375:THR:HG22	1:A:376:MET:HE2	1.69	0.74
1:A:87:ILE:HA	1:A:90:LYS:HE3	1.68	0.74
1:A:380:ARG:HH11	1:A:380:ARG:HG3	1.53	0.73
1:A:366:ASN:HD21	1:A:471:ARG:NE	1.87	0.70
1:A:515:MET:HE1	1:A:568:LEU:HD11	1.73	0.68
1:A:268:LYS:HB2	1:A:269:PRO:HD3	1.76	0.67
1:A:90:LYS:HZ2	1:A:94:GLN:HE22	1.41	0.67
1:A:394:ASP:C	1:A:396:ILE:H	2.03	0.67
1:A:221:ALA:CB	1:A:286:LEU:HD21	2.24	0.66
1:A:74:VAL:HG11	1:A:126:GLU:HG2	1.76	0.66
1:A:298:LEU:HD12	1:A:302:GLY:HA2	1.78	0.64
1:A:431:ILE:HD11	1:A:495:MET:HE2	1.79	0.62
1:A:326:ASN:HB2	1:A:329:LEU:CD2	2.25	0.61
1:A:90:LYS:NZ	1:A:94:GLN:HE22	1.97	0.61
1:A:320:ARG:HG2	1:A:320:ARG:HH11	1.66	0.61
1:A:427:LEU:HD21	1:A:495:MET:HE3	1.83	0.60
1:A:206:ARG:O	1:A:210:ILE:HG12	2.01	0.59
1:A:221:ALA:HB3	1:A:286:LEU:HD21	1.84	0.58
1:A:380:ARG:HG3	1:A:380:ARG:NH1	2.15	0.58
1:A:90:LYS:HZ2	1:A:94:GLN:NE2	2.01	0.58
1:A:610:THR:OG1	1:A:613:GLU:HG3	2.03	0.58
1:A:243:THR:O	1:A:247:LEU:HG	2.04	0.58
1:A:414:GLN:O	1:A:416:GLU:N	2.36	0.57
1:A:211:LEU:C	1:A:211:LEU:HD23	2.30	0.56
1:A:375:THR:HG22	1:A:376:MET:CE	2.35	0.56
1:A:328:ASN:OD1	1:A:329:LEU:HD22	2.05	0.55
1:A:580:GLU:O	1:A:584:LEU:HG	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:GLU:O	1:A:89:GLN:HG3	2.07	0.55
1:A:560:VAL:HG12	1:A:564:LYS:HE3	1.88	0.55
1:A:221:ALA:HB1	1:A:286:LEU:HD21	1.89	0.54
1:A:473:GLY:O	1:A:477:LEU:HD22	2.07	0.54
1:A:199:GLN:HE22	1:A:270:HIS:CD2	2.26	0.54
1:A:419:ASN:HD22	1:A:422:ASP:HB2	1.74	0.52
1:A:390:LEU:HG	1:A:423:HIS:HB3	1.91	0.52
1:A:369:THR:H	1:A:479:ASN:HD21	1.57	0.51
1:A:469:LYS:HE2	1:A:517:SER:OG	2.11	0.51
1:A:622:VAL:O	1:A:623:ARG:C	2.53	0.51
1:A:236:SER:O	1:A:237:SER:HB2	2.11	0.51
1:A:380:ARG:NH1	1:A:490:SER:HB2	2.27	0.50
1:A:91:GLY:O	1:A:95:VAL:HG23	2.12	0.50
1:A:348:ALA:O	1:A:352:ILE:HG13	2.11	0.49
1:A:445:LYS:HD2	1:A:445:LYS:C	2.37	0.49
1:A:372:THR:HG22	1:A:482:LEU:HD23	1.94	0.49
1:A:175:GLU:CG	1:A:178:GLN:HG3	2.43	0.49
1:A:496:LEU:O	1:A:497:ALA:C	2.57	0.48
1:A:513:SER:HA	1:A:516:VAL:HG12	1.96	0.48
1:A:542:LYS:O	1:A:546:LYS:HG3	2.13	0.47
1:A:418:LEU:HD12	1:A:418:LEU:O	2.13	0.47
1:A:365:ASN:HD22	1:A:365:ASN:N	2.12	0.47
1:A:593:TYR:CE2	1:A:597:LYS:HD3	2.50	0.47
1:A:527:MET:HB2	1:A:588:MET:HE1	1.95	0.47
1:A:619:ASN:O	1:A:623:ARG:HG3	2.14	0.47
1:A:94:GLN:NE2	1:A:94:GLN:C	2.72	0.47
1:A:386:LYS:O	1:A:390:LEU:HB2	2.15	0.47
1:A:578:LYS:HG3	1:A:622:VAL:CB	2.45	0.47
1:A:305:SER:O	1:A:309:VAL:HG23	2.15	0.47
1:A:175:GLU:HG2	1:A:178:GLN:OE1	2.15	0.46
1:A:519:TRP:CD2	1:A:559:LEU:HD13	2.51	0.46
1:A:527:MET:HB2	1:A:588:MET:CE	2.46	0.46
1:A:98:LYS:HG2	1:A:184:TRP:CE2	2.50	0.46
1:A:159:PHE:CZ	1:A:172:PRO:HD2	2.50	0.46
1:A:398:ARG:NH1	1:A:413:LEU:HD23	2.32	0.45
1:A:320:ARG:HG2	1:A:320:ARG:NH1	2.31	0.45
1:A:412:THR:OG1	1:A:503:ARG:HB3	2.17	0.45
1:A:312:ILE:HD11	1:A:330:LEU:HD22	1.98	0.45
1:A:501:HIS:O	1:A:505:GLU:HG2	2.16	0.44
1:A:570:ASP:OD1	1:A:572:SER:HB2	2.18	0.44
1:A:517:SER:HA	1:A:520:ARG:NH1	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:477:LEU:HD22	1:A:511:TYR:HB3	2.00	0.43
1:A:399:GLU:N	1:A:399:GLU:OE1	2.52	0.43
1:A:238:GLY:O	1:A:242:TYR:HB2	2.18	0.43
1:A:369:THR:H	1:A:479:ASN:ND2	2.15	0.43
1:A:510:ARG:HD2	1:A:511:TYR:CE1	2.54	0.42
1:A:77:LEU:HD12	1:A:77:LEU:HA	1.91	0.42
1:A:254:LYS:HE3	1:A:258:ASP:OD2	2.19	0.42
1:A:418:LEU:O	1:A:420:TRP:N	2.51	0.42
1:A:199:GLN:HE22	1:A:270:HIS:HD2	1.66	0.42
1:A:280:ILE:HG13	1:A:324:LEU:HD21	2.02	0.42
1:A:312:ILE:CD1	1:A:330:LEU:HD22	2.50	0.42
1:A:380:ARG:NH1	1:A:490:SER:CB	2.83	0.42
1:A:298:LEU:CD1	1:A:302:GLY:HA2	2.47	0.41
1:A:451:MET:HE1	1:A:466:ASN:HA	2.02	0.41
1:A:480:LEU:HD23	1:A:480:LEU:HA	1.92	0.41
1:A:376:MET:HB3	1:A:486:ILE:HG21	2.03	0.41
1:A:486:ILE:O	1:A:490:SER:HB3	2.20	0.41
1:A:406:TYR:CD2	1:A:406:TYR:C	2.99	0.41
1:A:175:GLU:HG3	1:A:178:GLN:HG3	2.01	0.41
1:A:393:MET:HA	1:A:393:MET:HE2	2.02	0.41
1:A:516:VAL:O	1:A:516:VAL:HG22	2.20	0.41
1:A:487:VAL:O	1:A:493:ASN:HB2	2.20	0.40
1:A:586:MET:HE2	1:A:615:THR:HG1	1.86	0.40
1:A:394:ASP:C	1:A:396:ILE:N	2.70	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	545/563 (97%)	525 (96%)	13 (2%)	7 (1%)	9 6

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	116	SER
1	A	415	ASN
1	A	417	ALA
1	A	622	VAL
1	A	406	TYR
1	A	419	ASN
1	A	237	SER

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	466/514 (91%)	450 (97%)	16 (3%)	32	35

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	77	LEU
1	A	80	GLU
1	A	94	GLN
1	A	134	LEU
1	A	144	GLN
1	A	209	LEU
1	A	218	GLU
1	A	339	GLN
1	A	365	ASN
1	A	399	GLU
1	A	418	LEU
1	A	442	LEU
1	A	477	LEU
1	A	484	GLU
1	A	558	ASP
1	A	577	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	ASN
1	A	94	GLN
1	A	102	GLN
1	A	120	ASN
1	A	136	GLN
1	A	164	ASN
1	A	190	HIS
1	A	199	GLN
1	A	240	ASN
1	A	252	ASN
1	A	263	GLN
1	A	274	GLN
1	A	300	ASN
1	A	313	ASN
1	A	339	GLN
1	A	342	GLN
1	A	357	ASN
1	A	365	ASN
1	A	366	ASN
1	A	400	ASN
1	A	419	ASN
1	A	441	ASN
1	A	454	GLN
1	A	460	ASN
1	A	462	ASN
1	A	479	ASN
1	A	485	GLN
1	A	501	HIS
1	A	602	ASN
1	A	619	ASN
1	A	620	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 [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	549/563 (97%)	0.71	40 (7%)	21 22	23, 42, 72, 82	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	232	TYR	5.1
1	A	220	PHE	5.0
1	A	291	LEU	3.6
1	A	117	GLY	3.4
1	A	390	LEU	3.4
1	A	420	TRP	3.4
1	A	293	ILE	3.3
1	A	532	ILE	3.2
1	A	294	VAL	3.2
1	A	623	ARG	3.2
1	A	345	PHE	2.9
1	A	406	TYR	2.9
1	A	221	ALA	2.9
1	A	301	PHE	2.8
1	A	417	ALA	2.7
1	A	395	ASN	2.6
1	A	379	LEU	2.6
1	A	190	HIS	2.6
1	A	387	ASN	2.6
1	A	415	ASN	2.5
1	A	382	PHE	2.5
1	A	119	ALA	2.5
1	A	116	SER	2.4
1	A	418	LEU	2.4
1	A	246	LEU	2.4
1	A	286	LEU	2.4
1	A	337	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	223	GLU	2.3
1	A	222	LYS	2.3
1	A	219	PRO	2.2
1	A	327	TYR	2.2
1	A	397	THR	2.2
1	A	298	LEU	2.2
1	A	217	LEU	2.1
1	A	349	ILE	2.1
1	A	425	VAL	2.1
1	A	356	ALA	2.1
1	A	396	ILE	2.1
1	A	497	ALA	2.1
1	A	341	THR	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.