



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:53 AM UTC

PDB ID : 8D3F / pdb_00008d3f
Title : Crystal structure of human STAT1 in complex with the repeat region from Toxoplasma protein TgIST
Authors : Huang, Z.; Liu, H.; Nix, J.C.; Amarasinghe, G.K.; Sibley, L.D.
Deposited on : 2022-06-01
Resolution : 2.97 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

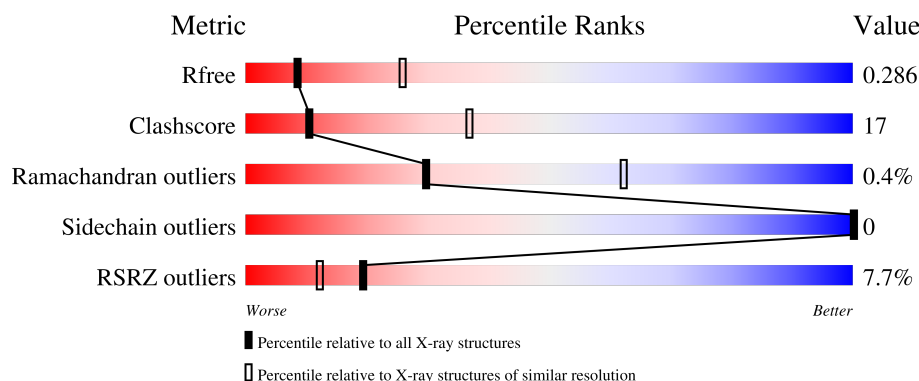
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.97 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	3580 (3.00-2.96)
Clashscore	190562	3904 (3.00-2.96)
Ramachandran outliers	187476	3761 (3.00-2.96)
Sidechain outliers	187428	3764 (3.00-2.96)
RSRZ outliers	180081	3579 (3.00-2.96)

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	642	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 4360 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 Signal transducer and activator of transcription 1-alpha/beta, Inhibitor of STAT1-dependent transcription TgIST.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	572	Total	C	N	O	S	0	0	0
			4360	2786	732	823	19			

There are 25 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	123	MET	-	initiating methionine	UNP P42224
A	124	TRP	-	expression tag	UNP P42224
A	125	SER	-	expression tag	UNP P42224
A	126	HIS	-	expression tag	UNP P42224
A	127	PRO	-	expression tag	UNP P42224
A	128	GLN	-	expression tag	UNP P42224
A	129	PHE	-	expression tag	UNP P42224
A	130	GLU	-	expression tag	UNP P42224
A	131	LYS	-	expression tag	UNP P42224
A	714	SER	-	linker	UNP P42224
A	715	GLY	-	linker	UNP P42224
A	716	GLY	-	linker	UNP P42224
A	717	GLY	-	linker	UNP P42224
A	718	GLY	-	linker	UNP P42224
A	719	SER	-	linker	UNP P42224
A	720	GLY	-	linker	UNP P42224
A	721	GLY	-	linker	UNP P42224
A	722	GLY	-	linker	UNP P42224
A	723	GLY	-	linker	UNP P42224
A	724	SER	-	linker	UNP P42224
A	725	GLY	-	linker	UNP P42224
A	726	GLY	-	linker	UNP P42224
A	727	GLY	-	linker	UNP P42224
A	728	GLY	-	linker	UNP P42224
A	729	GLY	-	linker	UNP P42224

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	59.96Å 163.92Å 226.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.64 – 2.97 46.64 – 2.97	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.64-2.97) 99.9 (46.64-2.97)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	0.18	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 2.96Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.248 , 0.288 0.248 , 0.286	Depositor DCC
R_{free} test set	1161 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	79.6	Xtriage
Anisotropy	0.881	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 85.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4360	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.46% 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	0.45	0/4450	0.67	0/6074

There are no bond length outliers.

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	4360	0	4062	142	0
All	All	4360	0	4062	142	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 17.

All (142) 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:564:LEU:HD22	1:A:569:LEU:HD11	1.45	0.98
1:A:564:LEU:HD12	1:A:610:GLY:O	1.70	0.91
1:A:564:LEU:HD22	1:A:569:LEU:CD1	2.08	0.83
1:A:465:PRO:HB3	1:A:559:GLU:HG3	1.61	0.82
1:A:313:ILE:HG22	1:A:349:VAL:HG11	1.58	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:LEU:HB3	1:A:471:ILE:HD12	1.64	0.77
1:A:586:ARG:O	1:A:590:LEU:HD12	1.85	0.77
1:A:380:PHE:HB3	1:A:409:LEU:HD21	1.66	0.76
1:A:378:ARG:CZ	1:A:426:VAL:HG23	2.22	0.70
1:A:561:ILE:O	1:A:565:ILE:HG13	1.92	0.69
1:A:430:LEU:HD23	1:A:454:PRO:HB2	1.74	0.68
1:A:598:THR:HA	1:A:666:TYR:O	1.94	0.68
1:A:746:ALA:HA	1:A:749:VAL:HG12	1.75	0.68
1:A:164:GLU:HG2	1:A:283:LEU:HD11	1.75	0.68
1:A:511:LYS:H	1:A:574:ASN:HD21	1.43	0.67
1:A:747:LEU:HD13	1:A:751:ARG:NH2	2.09	0.67
1:A:511:LYS:H	1:A:574:ASN:ND2	1.93	0.66
1:A:495:TRP:O	1:A:499:SER:HB3	1.95	0.66
1:A:492:CYS:HB2	1:A:536:LEU:HD22	1.77	0.65
1:A:162:SER:HA	1:A:165:ASP:HB2	1.79	0.64
1:A:598:THR:HG22	1:A:666:TYR:HB2	1.80	0.64
1:A:239:TRP:CZ2	1:A:256:LEU:HG	2.33	0.63
1:A:743:VAL:HG12	1:A:744:ARG:H	1.64	0.63
1:A:580:GLY:HA2	1:A:602:ARG:HA	1.81	0.62
1:A:584:LYS:O	1:A:588:ARG:HG3	2.00	0.61
1:A:498:LEU:HD12	1:A:537:ILE:HD13	1.82	0.61
1:A:586:ARG:HD2	1:A:590:LEU:HD11	1.82	0.60
1:A:600:LEU:HA	1:A:667:LEU:HD11	1.83	0.59
1:A:141:GLU:OE2	1:A:142:LEU:N	2.35	0.59
1:A:649:ARG:HB3	1:A:681:TYR:CE1	2.38	0.58
1:A:248:ILE:HB	1:A:472:LEU:CD2	2.34	0.58
1:A:223:VAL:O	1:A:227:THR:HG23	2.04	0.58
1:A:461:VAL:HG13	1:A:464:LEU:HD12	1.85	0.57
1:A:291:HIS:HA	1:A:296:LYS:HE3	1.86	0.57
1:A:634:TYR:CE2	1:A:647:ILE:HG21	2.40	0.56
1:A:667:LEU:HD22	1:A:677:PHE:HZ	1.69	0.56
1:A:599:PHE:O	1:A:667:LEU:HD12	2.05	0.56
1:A:214:VAL:HG11	1:A:297:ASN:OD1	2.06	0.56
1:A:378:ARG:NH1	1:A:426:VAL:HG23	2.21	0.55
1:A:141:GLU:OE1	1:A:145:LYS:HD2	2.05	0.55
1:A:366:LYS:HG2	1:A:367:ASP:OD1	2.07	0.55
1:A:301:LEU:O	1:A:305:THR:HG23	2.07	0.55
1:A:635:THR:O	1:A:639:LEU:HD12	2.06	0.55
1:A:170:TYR:HA	1:A:202:MET:HE1	1.88	0.54
1:A:322:GLN:CD	1:A:453:LEU:HD12	2.33	0.54
1:A:620:SER:O	1:A:620:SER:OG	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:515:ASN:OD1	1:A:518:GLN:HG3	2.08	0.54
1:A:667:LEU:HD22	1:A:677:PHE:CZ	2.42	0.53
1:A:651:TYR:O	1:A:652:LYS:NZ	2.42	0.53
1:A:510:THR:HB	1:A:574:ASN:HD21	1.72	0.53
1:A:153:VAL:HG13	1:A:220:LEU:HD11	1.90	0.53
1:A:587:GLU:HG3	1:A:600:LEU:HD11	1.91	0.52
1:A:276:GLN:HA	1:A:279:LYS:HD3	1.91	0.52
1:A:456:VAL:HG23	1:A:471:ILE:HD11	1.92	0.52
1:A:554:PHE:O	1:A:558:ILE:HD12	2.09	0.52
1:A:248:ILE:HD13	1:A:475:ASN:HB2	1.92	0.52
1:A:166:LEU:HD11	1:A:205:MET:HE3	1.92	0.51
1:A:256:LEU:HD22	1:A:348:LEU:HD11	1.92	0.51
1:A:239:TRP:CH2	1:A:256:LEU:HG	2.45	0.51
1:A:378:ARG:HD3	1:A:429:GLU:CD	2.35	0.51
1:A:360:VAL:HG12	1:A:437:THR:HB	1.92	0.51
1:A:145:LYS:HE3	1:A:227:THR:HG22	1.91	0.50
1:A:565:ILE:HA	1:A:569:LEU:HB2	1.93	0.50
1:A:391:ASN:OD1	1:A:394:GLU:HB2	2.12	0.50
1:A:488:LEU:C	1:A:490:PRO:HD3	2.36	0.50
1:A:595:GLN:O	1:A:598:THR:OG1	2.24	0.50
1:A:330:GLN:C	1:A:331:ARG:HG2	2.36	0.50
1:A:365:ASP:HB3	1:A:368:VAL:HG11	1.93	0.50
1:A:311:GLN:HA	1:A:314:GLN:HB2	1.92	0.50
1:A:141:GLU:CD	1:A:142:LEU:N	2.70	0.50
1:A:671:ILE:HD11	1:A:675:HIS:ND1	2.27	0.49
1:A:746:ALA:HA	1:A:749:VAL:CG1	2.42	0.49
1:A:331:ARG:HD2	1:A:468:TRP:CD2	2.48	0.49
1:A:284:GLU:HB2	1:A:294:ILE:HG21	1.94	0.49
1:A:646:ASP:OD2	1:A:682:SER:OG	2.13	0.49
1:A:256:LEU:HD13	1:A:346:ARG:HD3	1.94	0.49
1:A:166:LEU:HA	1:A:169:GLU:HG2	1.95	0.48
1:A:170:TYR:HA	1:A:202:MET:CE	2.43	0.48
1:A:494:ARG:HA	1:A:535:GLY:O	2.14	0.48
1:A:167:GLN:NE2	1:A:287:TYR:HB3	2.28	0.47
1:A:232:ILE:HD11	1:A:312:LEU:HA	1.95	0.47
1:A:364:PHE:HD2	1:A:433:LEU:HD13	1.78	0.47
1:A:335:LEU:O	1:A:457:VAL:HA	2.14	0.47
1:A:380:PHE:HB3	1:A:409:LEU:CD2	2.40	0.47
1:A:141:GLU:OE2	1:A:142:LEU:HB2	2.14	0.47
1:A:221:LEU:HD21	1:A:305:THR:HG22	1.97	0.47
1:A:264:THR:OG1	1:A:349:VAL:O	2.28	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:495:TRP:HB2	1:A:527:LEU:HD11	1.97	0.47
1:A:320:GLU:OE2	1:A:346:ARG:NH1	2.47	0.47
1:A:579:MET:HE3	1:A:579:MET:HB2	1.66	0.46
1:A:347:LEU:HD12	1:A:347:LEU:HA	1.80	0.46
1:A:557:TRP:CE3	1:A:558:ILE:HG13	2.51	0.46
1:A:378:ARG:HD3	1:A:429:GLU:OE1	2.16	0.46
1:A:667:LEU:O	1:A:671:ILE:HG22	2.16	0.46
1:A:156:ILE:HD13	1:A:217:ILE:HG13	1.97	0.45
1:A:498:LEU:HA	1:A:501:VAL:HG22	1.98	0.45
1:A:652:LYS:NZ	1:A:662:ASN:HB2	2.32	0.45
1:A:194:GLN:C	1:A:196:GLN:H	2.25	0.45
1:A:210:ARG:HB3	1:A:293:PRO:HG2	2.00	0.45
1:A:345:LEU:HD11	1:A:435:PHE:CG	2.52	0.45
1:A:512:ARG:NH1	1:A:574:ASN:O	2.50	0.45
1:A:620:SER:O	1:A:622:ASN:N	2.51	0.44
1:A:649:ARG:HD2	1:A:681:TYR:CD1	2.52	0.44
1:A:498:LEU:HA	1:A:498:LEU:HD23	1.74	0.44
1:A:220:LEU:HA	1:A:223:VAL:HG12	1.99	0.44
1:A:335:LEU:HD22	1:A:341:PHE:CG	2.52	0.44
1:A:554:PHE:O	1:A:557:TRP:HB3	2.17	0.44
1:A:276:GLN:HA	1:A:279:LYS:CD	2.48	0.44
1:A:476:MET:SD	1:A:501:VAL:HG21	2.58	0.43
1:A:653:VAL:HG23	1:A:663:PRO:HD3	2.01	0.43
1:A:160:ILE:O	1:A:164:GLU:HG3	2.18	0.43
1:A:634:TYR:HE2	1:A:647:ILE:HG21	1.82	0.43
1:A:639:LEU:CD2	1:A:644:PHE:HB2	2.49	0.43
1:A:229:ASN:O	1:A:233:ASN:HB2	2.18	0.43
1:A:514:LEU:HA	1:A:514:LEU:HD23	1.69	0.43
1:A:639:LEU:HD21	1:A:644:PHE:HB2	2.01	0.43
1:A:202:MET:HB3	1:A:202:MET:HE3	1.79	0.43
1:A:270:LEU:HD23	1:A:270:LEU:HA	1.69	0.43
1:A:458:ILE:HB	1:A:463:GLN:HB2	2.01	0.43
1:A:655:ALA:HB2	1:A:712:VAL:HG22	2.00	0.42
1:A:146:VAL:HB	1:A:265:ILE:HD11	2.01	0.42
1:A:335:LEU:HD22	1:A:341:PHE:CB	2.49	0.42
1:A:430:LEU:HD11	1:A:486:PHE:CE2	2.54	0.42
1:A:483:ASN:O	1:A:484:LEU:HD23	2.20	0.42
1:A:316:SER:OG	1:A:348:LEU:N	2.51	0.42
1:A:231:LEU:HD12	1:A:231:LEU:HA	1.79	0.42
1:A:667:LEU:HD12	1:A:667:LEU:HA	1.62	0.42
1:A:474:TYR:OH	1:A:482:ARG:N	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:ARG:HH11	1:A:180:ARG:HG3	1.84	0.41
1:A:430:LEU:N	1:A:430:LEU:HD12	2.34	0.41
1:A:359:LYS:HB2	1:A:389:VAL:HG22	2.02	0.41
1:A:661:GLU:CD	1:A:747:LEU:HD21	2.44	0.41
1:A:604:SER:HB2	1:A:633:PRO:HG2	2.02	0.41
1:A:353:GLU:H	1:A:353:GLU:HG3	1.53	0.41
1:A:370:GLU:HA	1:A:373:THR:OG1	2.21	0.41
1:A:232:ILE:HD13	1:A:315:SER:HB2	2.02	0.41
1:A:366:LYS:NZ	1:A:449:GLU:OE1	2.41	0.41
1:A:390:MET:HG2	1:A:401:ALA:C	2.46	0.41
1:A:643:THR:OG1	1:A:645:PRO:HD2	2.20	0.41
1:A:439:LEU:HD11	1:A:441:GLN:HG3	2.02	0.40
1:A:244:GLN:HB2	1:A:475:ASN:HB3	2.03	0.40
1:A:747:LEU:O	1:A:751:ARG:HG3	2.21	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	562/642 (88%)	517 (92%)	43 (8%)	2 (0%)	30 62

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	621	GLN
1	A	582	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	444/577 (77%)	444 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	158	HIS
1	A	233	ASN
1	A	260	GLN
1	A	272	GLN
1	A	357	ASN
1	A	548	ASN
1	A	568	HIS
1	A	574	ASN
1	A	670	ASN

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	572/642 (89%)	0.62	44 (7%) 19 12	63, 89, 144, 187	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	488	LEU	5.3
1	A	483	ASN	5.2
1	A	487	PHE	5.1
1	A	368	VAL	3.9
1	A	543	CYS	3.7
1	A	622	ASN	3.6
1	A	763	THR	3.6
1	A	746	ALA	3.6
1	A	485	SER	3.6
1	A	415	ALA	3.5
1	A	420	ASN	3.3
1	A	702	ILE	3.2
1	A	628	PHE	3.2
1	A	481	PRO	3.2
1	A	683	ARG	3.0
1	A	764	ALA	2.9
1	A	179	ASN	2.9
1	A	136	LEU	2.8
1	A	482	ARG	2.7
1	A	596	PRO	2.7
1	A	349	VAL	2.7
1	A	742	GLN	2.6
1	A	530	ASN	2.6
1	A	484	LEU	2.5
1	A	414	ASN	2.5
1	A	177	LEU	2.5
1	A	419	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	237	VAL	2.4
1	A	712	VAL	2.4
1	A	500	GLU	2.4
1	A	489	THR	2.4
1	A	668	TYR	2.3
1	A	531	ALA	2.3
1	A	616	TRP	2.3
1	A	195	GLU	2.3
1	A	551	ASN	2.3
1	A	374	VAL	2.2
1	A	624	GLY	2.2
1	A	197	LEU	2.2
1	A	433	LEU	2.2
1	A	327	THR	2.2
1	A	255	CYS	2.0
1	A	579	MET	2.0
1	A	542	PHE	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.