



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 5ZW0 / pdb_00005zw0
Title : Apo-form PigA
Authors : Lee, C.-C.; Ko, T.-P.; Wang, A.H.J.
Deposited on : 2018-05-14
Resolution : 2.54 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

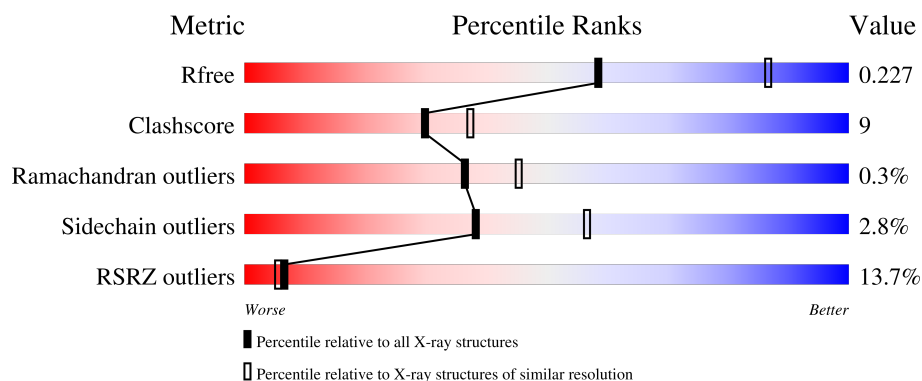
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.54 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	402	14% 74% 21% 5%
1	B	402	12% 78% 15% 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6043 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 L-prolyl-[peptidyl-carrier protein] dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	382	Total	C	N	O	S	0	0	0
			2927	1858	498	551	20			
1	B	375	Total	C	N	O	S	0	0	0
			2871	1822	488	541	20			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	387	GLY	-	expression tag	UNP Q5W271
A	388	SER	-	expression tag	UNP Q5W271
A	389	LEU	-	expression tag	UNP Q5W271
A	390	VAL	-	expression tag	UNP Q5W271
A	391	PRO	-	expression tag	UNP Q5W271
A	392	ARG	-	expression tag	UNP Q5W271
A	393	GLY	-	expression tag	UNP Q5W271
A	394	SER	-	expression tag	UNP Q5W271
A	395	HIS	-	expression tag	UNP Q5W271
A	396	HIS	-	expression tag	UNP Q5W271
A	397	HIS	-	expression tag	UNP Q5W271
A	398	HIS	-	expression tag	UNP Q5W271
A	399	HIS	-	expression tag	UNP Q5W271
A	400	HIS	-	expression tag	UNP Q5W271
A	401	HIS	-	expression tag	UNP Q5W271
A	402	HIS	-	expression tag	UNP Q5W271
B	387	GLY	-	expression tag	UNP Q5W271
B	388	SER	-	expression tag	UNP Q5W271
B	389	LEU	-	expression tag	UNP Q5W271
B	390	VAL	-	expression tag	UNP Q5W271
B	391	PRO	-	expression tag	UNP Q5W271
B	392	ARG	-	expression tag	UNP Q5W271
B	393	GLY	-	expression tag	UNP Q5W271
B	394	SER	-	expression tag	UNP Q5W271
B	395	HIS	-	expression tag	UNP Q5W271

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Chain	Residue	Modelled	Actual	Comment	Reference
B	396	HIS	-	expression tag	UNP Q5W271
B	397	HIS	-	expression tag	UNP Q5W271
B	398	HIS	-	expression tag	UNP Q5W271
B	399	HIS	-	expression tag	UNP Q5W271
B	400	HIS	-	expression tag	UNP Q5W271
B	401	HIS	-	expression tag	UNP Q5W271
B	402	HIS	-	expression tag	UNP Q5W271

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	143	Total 143	O 143	0	0
2	B	102	Total 102	O 102	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	79.22Å 163.02Å 146.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.90 – 2.54 19.90 – 2.54	Depositor EDS
% Data completeness (in resolution range)	90.4 (19.90-2.54) 90.3 (19.90-2.54)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.89 (at 2.53Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, R_{free}	0.179 , 0.227 0.180 , 0.227	Depositor DCC
R_{free} test set	1429 reflections (4.51%)	wwPDB-VP
Wilson B-factor (Å ²)	33.8	Xtriage
Anisotropy	0.240	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 74.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6043	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.23% 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

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	0.38	0/2981	0.58	0/4022
1	B	0.38	0/2921	0.57	0/3938
All	All	0.38	0/5902	0.58	0/7960

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

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	2927	0	2921	62	0
1	B	2871	0	2873	38	0
2	A	143	0	0	1	0
2	B	102	0	0	3	0
All	All	6043	0	5794	99	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 9.

All (99) 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:129:GLU:HB2	1:A:132:ALA:HB3	1.55	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:129:GLU:HB2	1:B:132:ALA:HB3	1.57	0.85
1:A:56:SER:HA	1:A:63:GLU:HG2	1.61	0.83
1:B:372:LYS:NZ	2:B:501:HOH:O	2.16	0.77
1:A:366:GLY:HA2	1:A:370:ILE:HD12	1.70	0.73
1:B:139:MET:H	1:B:173:LYS:HE3	1.61	0.66
1:A:139:MET:H	1:A:173:LYS:NZ	1.96	0.64
1:B:203:LYS:HD3	1:B:211:TRP:HE1	1.65	0.62
1:B:340:PHE:HB2	1:B:344:ALA:HB2	1.83	0.60
1:B:49:GLY:HA2	1:B:52:HIS:CE1	2.37	0.59
1:A:340:PHE:HB2	1:A:344:ALA:HB2	1.85	0.59
1:A:175:ASN:O	1:A:177:ASP:N	2.36	0.59
1:B:203:LYS:HD3	1:B:211:TRP:NE1	2.17	0.59
1:B:366:GLY:HA2	1:B:370:ILE:HD12	1.85	0.58
1:A:125:ASN:O	1:A:159:ILE:HG23	2.04	0.58
1:A:203:LYS:HB2	1:A:211:TRP:HE1	1.69	0.57
1:A:346:ASP:HB3	1:A:349:LEU:HB2	1.88	0.55
1:A:206:LEU:HD12	1:A:360:PRO:HG2	1.88	0.55
1:B:132:ALA:HB1	1:B:139:MET:HG2	1.88	0.55
1:B:103:PHE:CG	1:B:236:ILE:HG12	2.40	0.55
1:A:127:ALA:O	1:A:136:ILE:HD13	2.07	0.55
1:B:128:THR:HA	1:B:133:GLY:HA2	1.88	0.54
1:B:170:ILE:HG13	1:B:188:ILE:HD12	1.88	0.54
1:B:206:LEU:HD12	1:B:360:PRO:HG2	1.89	0.54
1:A:3:PHE:HA	1:B:1:MET:HE3	1.89	0.54
1:A:353:ARG:HG2	1:A:357:ASN:HD21	1.73	0.53
1:A:139:MET:HE3	1:A:156:LYS:NZ	2.24	0.53
1:A:130:PRO:HD2	1:A:140:GLN:NE2	2.26	0.51
1:A:136:ILE:HD12	1:A:237:PHE:CE2	2.46	0.51
1:B:340:PHE:CD1	1:B:351:LEU:HD22	2.46	0.50
1:A:130:PRO:HD3	1:A:157:ILE:HG12	1.93	0.50
1:A:126:ALA:HA	1:A:159:ILE:HD12	1.92	0.50
1:A:249:SER:HB3	1:A:325:ILE:HD13	1.92	0.50
1:A:346:ASP:HB3	1:A:349:LEU:CB	2.41	0.50
1:A:42:TRP:CZ3	1:A:122:ILE:HG12	2.47	0.49
1:B:98:ILE:HG23	1:B:102:ARG:HD2	1.93	0.49
1:A:32:ILE:HG21	1:A:353:ARG:CZ	2.43	0.49
1:A:55:VAL:HG11	1:A:117:VAL:HG21	1.95	0.49
1:A:139:MET:H	1:A:173:LYS:HZ3	1.58	0.49
1:A:186:PHE:CE2	1:A:226:ARG:HG3	2.47	0.49
1:A:298:ARG:HD2	1:A:302:TYR:OH	2.13	0.49
1:B:132:ALA:HB2	1:B:139:MET:HA	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:ILE:HD12	1:B:237:PHE:CZ	2.48	0.49
1:A:269:ALA:O	1:A:271:GLN:HG3	2.13	0.48
1:A:340:PHE:CD1	1:A:351:LEU:HD22	2.49	0.48
1:B:264:LEU:HD22	1:B:268:LYS:HE3	1.96	0.48
1:B:298:ARG:HG3	1:B:302:TYR:CE2	2.49	0.48
1:B:150:TYR:OH	1:B:226:ARG:NH1	2.46	0.48
1:A:203:LYS:HE2	1:A:205:CYS:O	2.13	0.47
1:B:156:LYS:HB3	1:B:159:ILE:HD11	1.97	0.47
1:B:298:ARG:HD2	1:B:302:TYR:OH	2.14	0.47
1:B:272:GLN:HB2	1:B:277:ILE:HG22	1.97	0.46
1:A:290:MET:HE2	1:A:290:MET:HB3	1.74	0.46
1:B:346:ASP:HB3	1:B:349:LEU:HB2	1.98	0.46
1:B:203:LYS:HE2	1:B:205:CYS:O	2.15	0.46
1:A:128:THR:HA	1:A:133:GLY:HA2	1.98	0.46
1:A:175:ASN:C	1:A:177:ASP:H	2.24	0.46
1:A:362:ARG:H	1:A:362:ARG:HG2	1.59	0.45
1:A:249:SER:HB3	1:A:325:ILE:CD1	2.46	0.45
1:A:345:MET:HG2	1:A:352:VAL:HG21	1.99	0.45
1:A:139:MET:HE3	1:A:156:LYS:HZ3	1.80	0.45
1:B:17:TYR:OH	2:B:502:HOH:O	2.18	0.45
1:A:95:ALA:HB1	1:A:125:ASN:CG	2.41	0.45
1:B:340:PHE:HD1	1:B:351:LEU:HD22	1.82	0.45
1:B:245:LYS:NZ	2:B:504:HOH:O	2.50	0.44
1:A:331:GLN:HG3	2:A:607:HOH:O	2.18	0.44
1:A:355:LEU:O	1:A:359:ILE:HG13	2.18	0.43
1:A:40:GLU:OE2	1:A:44:LYS:NZ	2.50	0.43
1:A:178:HIS:HB3	1:A:179:GLY:H	1.71	0.43
1:B:290:MET:HE2	1:B:290:MET:HB3	1.65	0.43
1:B:362:ARG:H	1:B:362:ARG:HG2	1.61	0.42
1:A:171:TYR:CE1	1:A:237:PHE:HB2	2.53	0.42
1:A:327:GLU:HG2	1:A:362:ARG:HH22	1.84	0.42
1:B:116:LEU:HD23	1:B:121:LEU:HB2	2.02	0.42
1:A:203:LYS:HD3	1:A:211:TRP:NE1	2.35	0.42
1:A:129:GLU:HG2	1:A:155:LYS:O	2.20	0.42
1:A:129:GLU:CG	1:A:156:LYS:HD3	2.50	0.42
1:A:354:HIS:O	1:A:358:MET:HG3	2.19	0.42
1:A:22:ASP:HA	1:A:26:GLN:HE21	1.85	0.42
1:A:57:HIS:ND1	1:A:63:GLU:OE2	2.49	0.41
1:B:229:MET:HE3	1:B:229:MET:HB2	1.73	0.41
1:A:159:ILE:O	1:A:211:TRP:HA	2.20	0.41
1:B:315:GLU:O	1:B:318:ILE:HG22	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:GLY:HA2	1:A:52:HIS:CE1	2.55	0.41
1:A:353:ARG:O	1:A:357:ASN:ND2	2.54	0.41
1:B:222:PRO:HD2	1:B:225:GLN:HG3	2.02	0.41
1:A:168:PHE:HB3	1:A:170:ILE:HD11	2.02	0.41
1:A:143:ALA:HA	1:A:151:ILE:O	2.21	0.41
1:B:353:ARG:O	1:B:357:ASN:ND2	2.53	0.41
1:A:136:ILE:HD12	1:A:237:PHE:CZ	2.56	0.41
1:A:266:TYR:CE2	1:A:277:ILE:HD11	2.55	0.41
1:B:74:GLU:OE1	1:B:302:TYR:OH	2.27	0.41
1:B:171:TYR:CE1	1:B:237:PHE:HB2	2.56	0.41
1:A:256:LEU:HD12	1:A:256:LEU:HA	1.78	0.41
1:A:129:GLU:HG3	1:A:156:LYS:HD3	2.01	0.40
1:A:139:MET:HE1	1:A:171:TYR:O	2.20	0.40
1:A:159:ILE:CG2	1:A:162:ALA:HB2	2.52	0.40
1:A:340:PHE:CE1	1:A:351:LEU:HD22	2.56	0.40
1:A:272:GLN:HB2	1:A:277:ILE:HG22	2.04	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	380/402 (94%)	356 (94%)	22 (6%)	2 (0%)	24	34
1	B	371/402 (92%)	351 (95%)	20 (5%)	0	100	100
All	All	751/804 (93%)	707 (94%)	42 (6%)	2 (0%)	36	45

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	176	PRO
1	A	136	ILE

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	308/325 (95%)	299 (97%)	9 (3%)	37	55
1	B	302/325 (93%)	294 (97%)	8 (3%)	40	59
All	All	610/650 (94%)	593 (97%)	17 (3%)	38	57

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	37	LEU
1	A	58	GLN
1	A	136	ILE
1	A	142	THR
1	A	178	HIS
1	A	230	GLU
1	A	318	ILE
1	A	354	HIS
1	B	1	MET
1	B	10	SER
1	B	37	LEU
1	B	62	SER
1	B	63	GLU
1	B	136	ILE
1	B	354	HIS
1	B	373	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	34	GLN
1	A	217	ASN
1	A	225	GLN
1	A	296	GLN

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Mol	Chain	Res	Type
1	A	310	GLN
1	A	354	HIS
1	A	368	ASN
1	B	57	HIS
1	B	238	HIS

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	382/402 (95%)	0.26	55 (14%) 6 5	10, 44, 200, 287	0
1	B	375/402 (93%)	0.20	49 (13%) 7 6	13, 45, 217, 263	0
All	All	757/804 (94%)	0.23	104 (13%) 6 5	10, 44, 212, 287	0

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	278	GLY	7.3
1	A	278	GLY	7.1
1	A	361	SER	6.2
1	A	277	ILE	6.1
1	B	136	ILE	5.6
1	A	349	LEU	5.6
1	B	137	TYR	5.4
1	B	277	ILE	5.3
1	A	351	LEU	5.3
1	B	128	THR	5.1
1	B	349	LEU	5.1
1	B	343	ALA	5.0
1	B	351	LEU	4.8
1	B	127	ALA	4.7
1	B	345	MET	4.6
1	B	358	MET	4.6
1	B	139	MET	4.4
1	A	200	VAL	4.4
1	B	365	SER	4.4
1	B	341	GLY	4.3
1	A	206	LEU	4.3
1	A	358	MET	4.2
1	A	276	ALA	4.2
1	B	276	ALA	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	359	ILE	4.1
1	A	352	VAL	4.1
1	A	205	CYS	4.0
1	A	343	ALA	4.0
1	B	361	SER	4.0
1	B	206	LEU	3.9
1	B	342	GLY	3.9
1	B	364	PHE	3.9
1	A	143	ALA	3.8
1	A	355	LEU	3.8
1	A	141	ALA	3.7
1	A	364	PHE	3.7
1	B	133	GLY	3.7
1	B	134	SER	3.7
1	A	357	ASN	3.6
1	A	362	ARG	3.6
1	A	204	ASP	3.4
1	B	352	VAL	3.4
1	B	355	LEU	3.4
1	A	341	GLY	3.4
1	A	270	ARG	3.4
1	A	347	GLN	3.4
1	A	356	LEU	3.4
1	B	126	ALA	3.3
1	B	344	ALA	3.3
1	B	359	ILE	3.3
1	B	200	VAL	3.2
1	B	353	ARG	3.2
1	A	340	PHE	3.2
1	B	201	ILE	3.2
1	B	130	PRO	3.1
1	A	201	ILE	3.1
1	A	180	PHE	3.1
1	A	354	HIS	3.0
1	B	280	PHE	3.0
1	B	273	PHE	3.0
1	B	362	ARG	3.0
1	B	202	PRO	3.0
1	B	340	PHE	3.0
1	B	363	ILE	2.9
1	B	135	ASP	2.9
1	A	182	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	354	HIS	2.8
1	A	134	SER	2.7
1	A	353	ARG	2.7
1	B	205	CYS	2.7
1	A	273	PHE	2.7
1	A	148	GLY	2.7
1	A	360	PRO	2.6
1	A	348	GLU	2.6
1	A	280	PHE	2.6
1	B	203	LYS	2.5
1	A	342	GLY	2.5
1	A	158	PHE	2.5
1	A	271	GLN	2.5
1	A	132	ALA	2.5
1	B	231	GLY	2.4
1	A	346	ASP	2.4
1	A	269	ALA	2.4
1	A	183	VAL	2.4
1	A	345	MET	2.4
1	B	357	ASN	2.4
1	A	202	PRO	2.3
1	B	339	THR	2.3
1	A	133	GLY	2.3
1	A	140	GLN	2.3
1	B	204	ASP	2.3
1	A	365	SER	2.3
1	A	131	ASP	2.3
1	B	158	PHE	2.3
1	A	272	GLN	2.2
1	B	140	GLN	2.2
1	A	181	LEU	2.2
1	A	366	GLY	2.2
1	B	141	ALA	2.1
1	B	271	GLN	2.1
1	A	142	THR	2.1
1	B	279	GLN	2.1
1	A	154	GLY	2.1
1	A	130	PRO	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 [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.