



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 08:41 PM UTC

PDB ID : 8A4J / pdb_00008a4j
Title : Human GDAP1, A247V mutant
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Deposited on : 2022-06-12
Resolution : 2.68 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

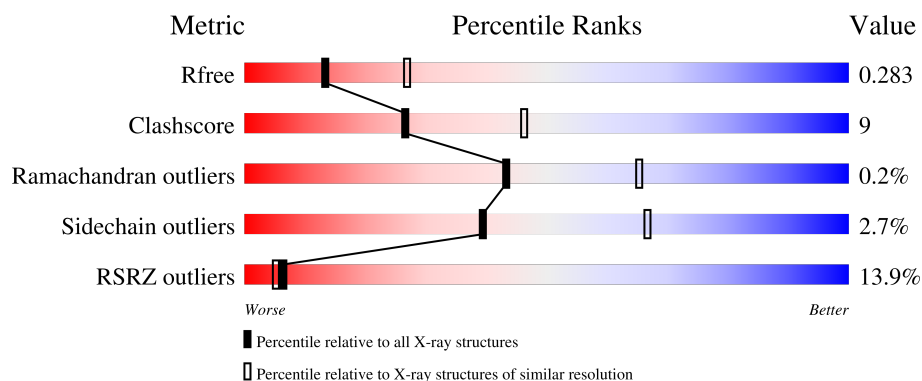
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.68 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	5070 (2.70-2.66)
Clashscore	190562	5409 (2.70-2.66)
Ramachandran outliers	187476	5324 (2.70-2.66)
Sidechain outliers	187428	5324 (2.70-2.66)
RSRZ outliers	180081	5070 (2.70-2.66)

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	280	4% 71% 16% • 11%
1	B	280	20% 64% 19% • 17%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8010 atoms, of which 4006 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 Ganglioside-induced differentiation-associated protein 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	249	Total	C	H	N	O	S	0	0	0
			4140	1319	2073	360	379	9			
1	B	232	Total	C	H	N	O	S	0	0	0
			3859	1232	1933	336	349	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	247	VAL	ALA	variant	UNP Q8TB36
B	247	VAL	ALA	variant	UNP Q8TB36

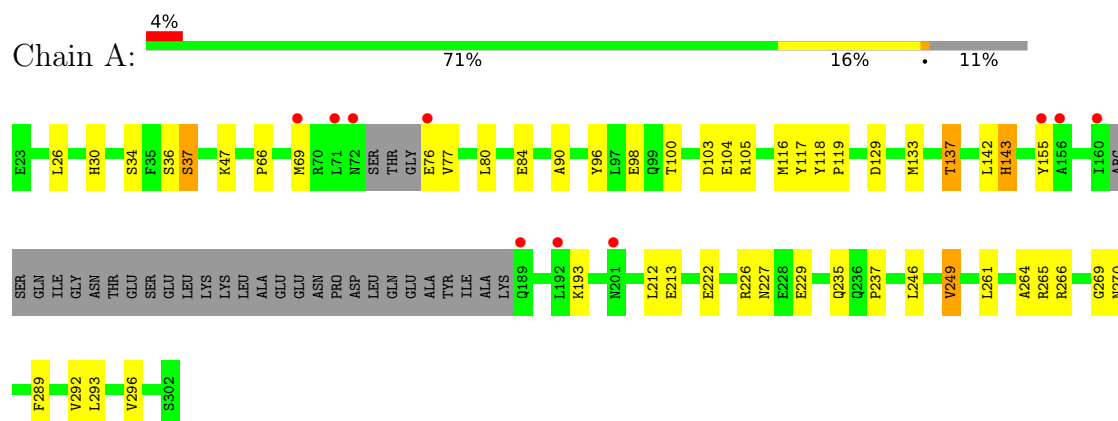
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	10	Total	O	0	0
			10	10		
2	B	1	Total	O	0	0
			1	1		

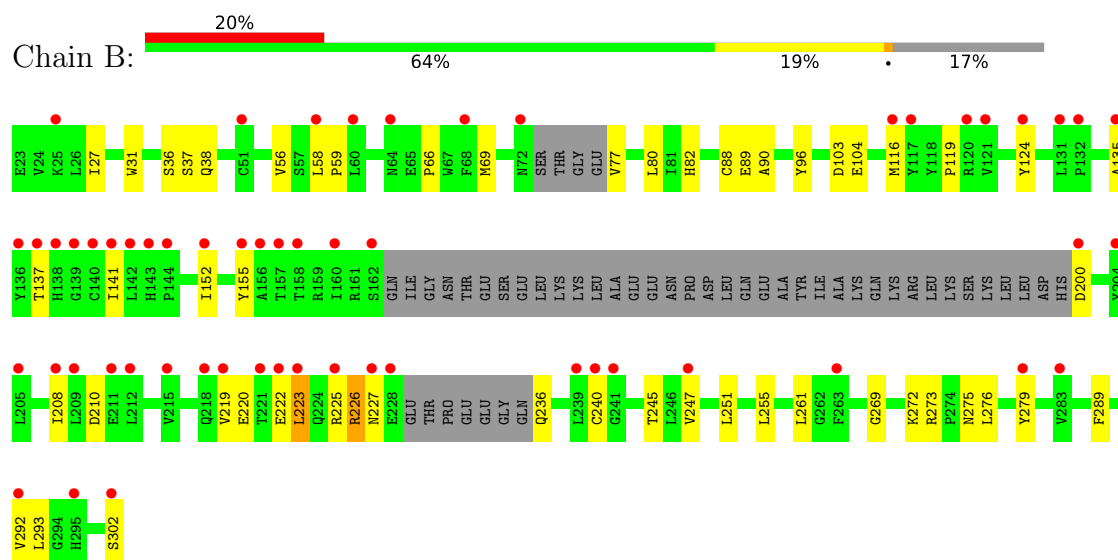
3 Residue-property plots [i](#)

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: Ganglioside-induced differentiation-associated protein 1



- Molecule 1: Ganglioside-induced differentiation-associated protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	73.29Å 115.69Å 115.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.28 – 2.68 42.28 – 2.68	Depositor EDS
% Data completeness (in resolution range)	99.6 (42.28-2.68) 99.5 (42.28-2.68)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.20_4459	Depositor
R, R_{free}	0.252 , 0.284 0.252 , 0.283	Depositor DCC
R_{free} test set	2000 reflections (7.06%)	wwPDB-VP
Wilson B-factor (Å ²)	85.0	Xtriage
Anisotropy	0.528	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 78.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8010	wwPDB-VP
Average B, all atoms (Å ²)	127.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.64% 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	0.80	1/2112 (0.0%)	0.94	0/2857
1	B	0.58	1/1968 (0.1%)	0.72	0/2663
All	All	0.71	2/4080 (0.0%)	0.84	0/5520

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	269	GLY	C-O	-9.71	1.18	1.24
1	A	143	HIS	C-O	5.23	1.26	1.23

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	2067	2073	2072	39	0
1	B	1926	1933	1931	37	0
2	A	10	0	0	1	0
2	B	1	0	0	0	0
All	All	4004	4006	4003	74	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 (74) 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:ASP:OD2	2:A:401:HOH:O	1.93	0.87
1:A:69:MET:HE1	1:A:155:TYR:HE1	1.43	0.83
1:B:226:ARG:HH21	1:B:227:ASN:HB2	1.48	0.77
1:A:26:LEU:CD2	1:A:80:LEU:HD11	2.16	0.75
1:A:103:ASP:OD1	1:A:104:GLU:N	2.22	0.72
1:A:69:MET:HE1	1:A:155:TYR:CE1	2.25	0.69
1:A:222:GLU:OE2	1:A:226:ARG:NH2	2.27	0.67
1:B:245:THR:OG1	1:B:247:VAL:HG12	1.96	0.66
1:A:26:LEU:HD22	1:A:80:LEU:HD11	1.79	0.64
1:B:251:LEU:HD23	1:B:279:TYR:CE1	2.32	0.63
1:A:66:PRO:HD2	1:A:69:MET:HG3	1.81	0.62
1:B:251:LEU:HD23	1:B:279:TYR:CZ	2.35	0.62
1:B:36:SER:HB2	1:B:90:ALA:HB2	1.83	0.60
1:A:34:SER:OG	1:A:37:SER:HB3	2.03	0.59
1:B:77:VAL:HG13	1:B:77:VAL:O	2.01	0.59
1:B:141:ILE:HD11	1:B:152:ILE:HD13	1.85	0.58
1:A:30:HIS:HB2	1:A:37:SER:HB2	1.86	0.58
1:B:219:VAL:HG12	1:B:223:LEU:HD12	1.86	0.58
1:B:226:ARG:NH1	1:B:236:GLN:NE2	2.52	0.57
1:B:226:ARG:NH1	1:B:236:GLN:HE22	2.03	0.55
1:B:226:ARG:NH2	1:B:227:ASN:HB2	2.21	0.54
1:A:116:MET:HE3	1:A:117:TYR:CE1	2.42	0.54
1:A:36:SER:HB2	1:A:90:ALA:HB2	1.90	0.54
1:B:66:PRO:HD2	1:B:69:MET:HG3	1.89	0.53
1:A:66:PRO:HD2	1:A:69:MET:CG	2.39	0.53
1:A:96:TYR:O	1:A:100:THR:HG23	2.09	0.52
1:A:292:VAL:HG12	1:A:293:LEU:HD23	1.91	0.52
1:B:31:TRP:CH2	1:B:77:VAL:N	2.78	0.52
1:A:293:LEU:O	1:A:296:VAL:HG22	2.10	0.50
1:B:59:PRO:HB3	1:B:302:SER:HB3	1.93	0.50
1:A:133:MET:O	1:A:137:THR:HG23	2.12	0.50
1:B:289:PHE:C	1:B:289:PHE:CD1	2.90	0.49
1:A:264:ALA:O	1:A:265:ARG:C	2.55	0.49
1:A:26:LEU:HD23	1:A:80:LEU:HD11	1.94	0.49
1:B:240:CYS:SG	1:B:247:VAL:HG11	2.54	0.48
1:A:103:ASP:OD1	1:A:105:ARG:N	2.39	0.48
1:A:84:GLU:OE2	1:B:58:LEU:N	2.45	0.48
1:B:219:VAL:HG12	1:B:223:LEU:CD1	2.44	0.48
1:B:272:LYS:HG2	1:B:273:ARG:HG3	1.96	0.47
1:B:124:TYR:CE2	1:B:222:GLU:HG2	2.50	0.47
1:B:220:GLU:OE1	1:B:275:ASN:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:255:LEU:HD11	1:B:276:LEU:HD11	1.97	0.46
1:B:135:ALA:HB1	1:B:208:ILE:CD1	2.46	0.46
1:A:103:ASP:OD1	1:A:103:ASP:C	2.59	0.46
1:A:155:TYR:C	1:A:155:TYR:CD1	2.95	0.45
1:A:96:TYR:CE1	1:A:100:THR:HG21	2.51	0.45
1:B:103:ASP:OD1	1:B:104:GLU:N	2.50	0.45
1:B:137:THR:HG22	1:B:261:LEU:HD11	1.98	0.45
1:B:38:GLN:HG2	1:B:293:LEU:CD2	2.47	0.45
1:B:82:HIS:ND1	1:B:96:TYR:OH	2.36	0.44
1:A:292:VAL:HG12	1:A:293:LEU:CD2	2.48	0.44
1:A:246:LEU:HA	1:A:249:VAL:HG13	2.00	0.44
1:A:69:MET:CE	1:A:155:TYR:HE1	2.23	0.44
1:A:118:TYR:HB3	1:A:119:PRO:HD3	2.00	0.43
1:B:292:VAL:HG12	1:B:293:LEU:HG	2.00	0.43
1:B:116:MET:O	1:B:119:PRO:HD2	2.19	0.43
1:B:38:GLN:HG2	1:B:293:LEU:HD22	1.99	0.43
1:A:235:GLN:O	1:A:235:GLN:HG3	2.18	0.43
1:B:220:GLU:OE1	1:B:220:GLU:HA	2.19	0.43
1:A:226:ARG:HB3	1:A:237:PRO:HG2	2.00	0.42
1:B:69:MET:CE	1:B:155:TYR:CE2	3.02	0.42
1:A:269:GLY:O	1:A:270:ASN:C	2.62	0.42
1:A:289:PHE:CE2	1:A:293:LEU:HD12	2.55	0.42
1:B:31:TRP:HH2	1:B:77:VAL:N	2.16	0.42
1:A:212:LEU:O	1:A:213:GLU:C	2.63	0.42
1:B:88:CYS:O	1:B:89:GLU:HB2	2.20	0.42
1:A:84:GLU:HG2	1:B:56:VAL:O	2.20	0.41
1:A:226:ARG:NH1	1:A:229:GLU:OE1	2.54	0.41
1:A:137:THR:HG22	1:A:261:LEU:HD11	2.02	0.41
1:B:69:MET:HE3	1:B:155:TYR:CE2	2.56	0.41
1:A:142:LEU:O	1:A:143:HIS:C	2.63	0.40
1:A:213:GLU:HG3	1:A:266:ARG:HH12	1.87	0.40
1:A:47:LYS:NZ	1:A:98:GLU:OE1	2.42	0.40
1:B:27:ILE:O	1:B:80:LEU:HD12	2.22	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	243/280 (87%)	235 (97%)	8 (3%)	0	100	100
1	B	224/280 (80%)	215 (96%)	8 (4%)	1 (0%)	30	51
All	All	467/560 (83%)	450 (96%)	16 (3%)	1 (0%)	43	65

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	225	ARG

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	233/259 (90%)	226 (97%)	7 (3%)	36	63
1	B	217/259 (84%)	212 (98%)	5 (2%)	44	71
All	All	450/518 (87%)	438 (97%)	12 (3%)	39	67

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

Mol	Chain	Res	Type
1	A	37	SER
1	A	76	GLU
1	A	77	VAL
1	A	137	THR
1	A	193	LYS

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Mol	Chain	Res	Type
1	A	227	ASN
1	A	249	VAL
1	B	37	SER
1	B	200	ASP
1	B	210	ASP
1	B	223	LEU
1	B	226	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	290	ASN
1	B	33	HIS
1	B	38	GLN
1	B	143	HIS
1	B	201	ASN
1	B	236	GLN
1	B	267	ASN

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 ⓘ

There are no oligosaccharides 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	249/280 (88%)	0.19	10 (4%) 42 37	63, 89, 183, 227	0
1	B	232/280 (82%)	1.23	57 (24%) 2 1	77, 140, 264, 351	0
All	All	481/560 (85%)	0.69	67 (13%) 6 5	63, 102, 245, 351	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	142	LEU	6.5
1	B	160	ILE	6.5
1	B	200	ASP	6.4
1	A	72	ASN	6.3
1	B	141	ILE	5.6
1	B	139	GLY	5.6
1	B	240	CYS	5.4
1	B	228	GLU	4.4
1	B	239	LEU	4.3
1	B	158	THR	4.3
1	A	71	LEU	4.2
1	B	120	ARG	4.2
1	B	138	HIS	3.8
1	B	155	TYR	3.8
1	B	204	TYR	3.8
1	B	143	HIS	3.5
1	A	189	GLN	3.5
1	B	137	THR	3.5
1	B	156	ALA	3.4
1	A	76	GLU	3.4
1	A	160	ILE	3.4
1	B	140	CYS	3.4
1	B	223	LEU	3.4
1	B	295	HIS	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	135	ALA	3.3
1	A	192	LEU	3.3
1	B	144	PRO	3.3
1	B	136	TYR	3.2
1	B	131	LEU	3.2
1	B	205	LEU	3.2
1	B	162	SER	3.1
1	B	60	LEU	3.1
1	B	208	ILE	3.1
1	B	68	PHE	3.1
1	B	117	TYR	3.0
1	B	263	PHE	2.9
1	B	218	GLN	2.9
1	B	72	ASN	2.9
1	A	201	ASN	2.9
1	B	157	THR	2.8
1	B	211	GLU	2.8
1	B	247	VAL	2.8
1	B	209	LEU	2.7
1	B	51	CYS	2.7
1	B	215	VAL	2.7
1	B	221	THR	2.7
1	B	132	PRO	2.7
1	B	302	SER	2.7
1	B	279	TYR	2.6
1	B	241	GLY	2.6
1	B	121	VAL	2.6
1	B	212	LEU	2.5
1	B	292	VAL	2.4
1	B	25	LYS	2.3
1	B	225	ARG	2.3
1	A	156	ALA	2.3
1	B	58	LEU	2.3
1	B	227	ASN	2.2
1	B	219	VAL	2.2
1	B	64	ASN	2.2
1	B	124	TYR	2.2
1	B	283	VAL	2.1
1	B	116	MET	2.1
1	A	69	MET	2.1
1	B	152	ILE	2.1
1	B	222	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	155	TYR	2.1

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.