Full wwPDB X-ray Structure Validation Report (i)

Oct 27, 2026 – 04:09 PM JST

PDB ID	:	9XGD / pdb_00009xgd
Title	:	Crystal structure of dKeima570 at pH 4.0
Authors	:	Nam, K.H.
Deposited on	:	2025-10-30
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 (i) 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 (1)) were used in the production of this report:

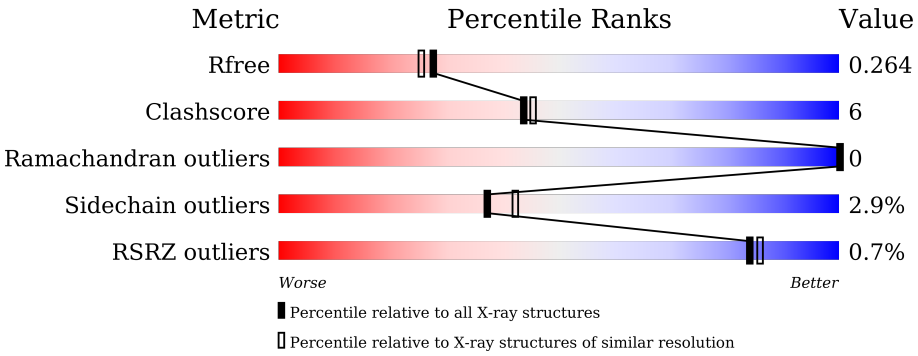
MolProbity	:	4-5-2 with Phenix2.2
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	2.2
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

1 Overall quality at a glance (i)

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 <=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	220	82% 16%
1	B	220	84% 14%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3607 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 Large stokes shift fluorescent protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	1	0
			1756	1122	289	331	14			
1	B	217	Total	C	N	O	S	0	1	0
			1756	1122	289	331	14			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	63	GYC	CYS	chromophore	UNP Q1JU63
A	63	GYC	TYR	chromophore	UNP Q1JU63
A	63	GYC	GLY	chromophore	UNP Q1JU63
A	192	VAL	ILE	conflict	UNP Q1JU63
B	63	GYC	CYS	chromophore	UNP Q1JU63
B	63	GYC	TYR	chromophore	UNP Q1JU63
B	63	GYC	GLY	chromophore	UNP Q1JU63
B	192	VAL	ILE	conflict	UNP Q1JU63

- Molecule 2 is water.

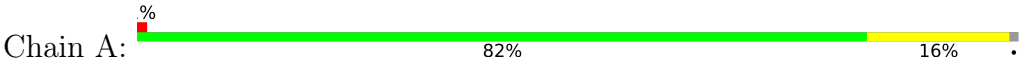
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	58	Total	O	0	0
			58	58		
2	B	37	Total	O	0	0
			37	37		

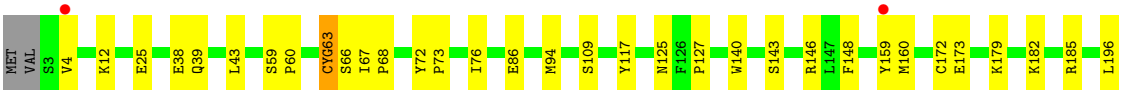
3 Residue-property plots (i)

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: Large stokes shift fluorescent protein

Chain A:



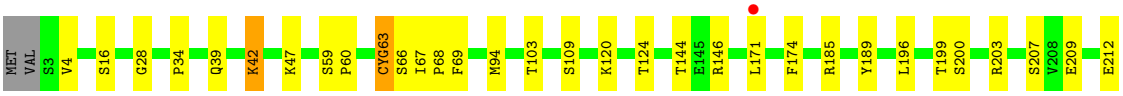




• Molecule 1: Large stokes shift fluorescent protein

Chain B:







4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	84.58Å 84.58Å 105.30Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	31.67 – 2.10 31.67 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.8 (31.67-2.10) 99.8 (31.67-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.33 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8.0430	Depositor
R, R_{free}	0.209 , 0.259 0.215 , 0.264	Depositor DCC
R_{free} test set	2002 reflections (7.71%)	wwPDB-VP
Wilson B-factor (Å^2)	39.0	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 29.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.041 for -h,-k,l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	3607	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.86% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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

Bond lengths and bond angles in the following residue types are not validated in this section: GYC

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	0.67	0/1760	1.19	3/2381 (0.1%)
1	B	0.65	0/1760	1.16	5/2381 (0.2%)
All	All	0.66	0/3520	1.17	8/4762 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	38	GLU	CB-CG-CD	6.74	124.06	112.60
1	B	103	THR	CA-CB-OG1	-6.30	100.15	109.60
1	B	124	THR	CA-CB-OG1	-6.18	100.33	109.60
1	B	144	THR	CA-CB-OG1	-6.02	100.57	109.60
1	B	199	THR	CA-CB-OG1	-5.89	100.77	109.60
1	A	72	TYR	CB-CA-C	5.43	115.99	109.31
1	A	12	LYS	CB-CA-C	5.10	119.05	109.71
1	B	185	ARG	CB-CA-C	5.08	117.12	110.06

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	185	ARG	Sidechain

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	1756	0	1680	19	0
1	B	1756	0	1680	25	0
2	A	58	0	0	3	0
2	B	37	0	0	0	0
All	All	3607	0	3360	42	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 6.

All (42) 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:B:63[B]:GYC:O2	1:B:63[B]:GYC:HD1	1.61	0.99
1:A:66:SER:OG	1:A:117:TYR:OH	1.87	0.81
1:A:63[A]:GYC:O2	1:A:63[A]:GYC:HD1	1.82	0.78
1:B:63[B]:GYC:N2	1:B:212:GLU:OE2	2.26	0.69
1:A:63[B]:GYC:HE1	1:A:143:SER:OG	1.93	0.67
1:A:207:SER:OG	2:A:301:HOH:O	2.15	0.65
1:A:39:GLN:OE1	1:A:66:SER:HB3	2.03	0.58
1:B:67:ILE:N	1:B:68:PRO:CD	2.67	0.58
1:B:200:SER:HB3	1:B:209:GLU:HB2	1.85	0.57
1:A:159:TYR:CE1	1:A:173:GLU:HG2	2.43	0.54
1:B:59:SER:N	1:B:60:PRO:CD	2.71	0.53
1:B:39:GLN:OE1	1:B:66:SER:HB3	2.09	0.53
1:A:109:SER:HB2	2:A:309:HOH:O	2.09	0.52
1:A:86:GLU:OE1	1:A:182:LYS:NZ	2.40	0.52
1:B:63[A]:GYC:N2	1:B:212:GLU:OE2	2.42	0.52
1:B:63[A]:GYC:HE2	1:B:196:LEU:HB2	1.92	0.51
1:B:146:ARG:HD3	1:B:189:TYR:CE1	2.44	0.51
1:A:125:ASN:O	1:A:127:PRO:HD3	2.13	0.49
1:A:59:SER:N	1:A:60:PRO:CD	2.77	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:LEU:HD11	1:B:220:LEU:HD11	1.95	0.47
1:B:66:SER:C	1:B:68:PRO:HD2	2.39	0.47
1:B:63[B]:GYC:SG1	1:B:212:GLU:HB2	2.54	0.47
1:B:94:MET:HE3	1:B:174:PHE:CE1	2.49	0.47
1:A:203:ARG:HB3	1:A:203:ARG:CZ	2.45	0.46
1:B:63[B]:GYC:HD1	1:B:63[B]:GYC:C2	2.41	0.46
1:B:63[A]:GYC:SG1	1:B:212:GLU:HB2	2.57	0.45
1:A:148:PHE:CE1	1:B:171:LEU:HD21	2.52	0.44
1:B:63[A]:GYC:N2	1:B:63[A]:GYC:HD2	2.33	0.44
1:A:63[B]:GYC:HE2	1:A:196:LEU:HB3	2.00	0.43
1:A:140:TRP:CE3	1:A:160:MET:HE3	2.52	0.43
1:A:73:PRO:HG2	1:A:76:ILE:HD12	2.00	0.43
1:B:63[A]:GYC:O3	1:B:66:SER:C	2.60	0.43
1:A:67:ILE:N	1:A:68:PRO:CD	2.82	0.42
1:B:67:ILE:N	1:B:68:PRO:HD3	2.33	0.42
1:B:171:LEU:N	1:B:171:LEU:HD23	2.34	0.42
1:B:16:SER:O	1:B:120:LYS:HA	2.19	0.42
1:B:34:PRO:HA	1:B:69:PHE:HA	2.02	0.42
1:A:94:MET:HB3	1:A:172:CYS:SG	2.60	0.41
1:B:66:SER:C	1:B:68:PRO:CD	2.93	0.41
1:A:25:GLU:OE1	2:A:302:HOH:O	2.21	0.41
1:B:28:GLY:HA2	1:B:42:LYS:O	2.21	0.41
1:B:63[B]:GYC:O3	1:B:66:SER:C	2.63	0.41

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	212/220 (96%)	210 (99%)	2 (1%)	0	100	100
1	B	212/220 (96%)	206 (97%)	6 (3%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	424/440 (96%)	416 (98%)	8 (2%)	0	100 100

There are no Ramachandran outliers to report.

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	188/190 (99%)	183 (97%)	5 (3%)	39 45
1	B	188/190 (99%)	182 (97%)	6 (3%)	34 38
All	All	376/380 (99%)	365 (97%)	11 (3%)	37 42

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	VAL
1	A	43	LEU
1	A	146	ARG
1	A	179	LYS
1	A	211	CYS
1	B	4	VAL
1	B	42	LYS
1	B	47	LYS
1	B	109	SER
1	B	203	ARG
1	B	207	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	157	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)

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
1	GYC	A	63[A]	1	22,22,23	1.08	2 (9%)	26,30,32	2.12	6 (23%)
1	GYC	A	63[B]	1	22,22,23	1.12	3 (13%)	26,30,32	1.33	4 (15%)
1	GYC	B	63[A]	1	22,22,23	0.92	1 (4%)	26,30,32	1.49	2 (7%)
1	GYC	B	63[B]	1	22,22,23	0.90	1 (4%)	26,30,32	1.69	4 (15%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	GYC	A	63[A]	1	-	4/9/29/30	0/2/2/2
1	GYC	A	63[B]	1	-	2/9/29/30	0/2/2/2
1	GYC	B	63[A]	1	-	1/9/29/30	0/2/2/2
1	GYC	B	63[B]	1	-	4/9/29/30	0/2/2/2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	63[A]	GYC	C1-N3	3.30	1.42	1.37
1	B	63[B]	GYC	C1-N3	3.06	1.42	1.37
1	A	63[B]	GYC	C1-N3	2.64	1.41	1.37
1	A	63[A]	GYC	C1-N2	2.59	1.36	1.32
1	A	63[B]	GYC	CA1-N1	2.45	1.61	1.48
1	B	63[A]	GYC	O3-C3	2.15	1.32	1.19

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	63[B]	GYC	C1-N2	2.05	1.35	1.32

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	63[A]	GYC	C2-N3-C1	5.08	110.54	107.97
1	A	63[A]	GYC	N3-C1-N2	-4.88	108.07	111.45
1	A	63[A]	GYC	CA2-N2-C1	4.61	109.17	105.77
1	B	63[A]	GYC	CA3-N3-C1	-4.47	121.80	127.16
1	B	63[A]	GYC	C2-N3-C1	4.38	110.18	107.97
1	B	63[B]	GYC	CA2-N2-C1	4.36	108.98	105.77
1	B	63[B]	GYC	C2-N3-C1	4.31	110.14	107.97
1	A	63[A]	GYC	CA3-N3-C1	-3.50	122.96	127.16
1	A	63[B]	GYC	O3-C3-CA3	-3.44	115.99	126.39
1	B	63[B]	GYC	N3-C1-N2	-3.18	109.25	111.45
1	A	63[A]	GYC	C2-CA2-N2	-2.72	107.03	108.93
1	A	63[B]	GYC	C2-N3-C1	2.42	109.19	107.97
1	A	63[A]	GYC	CG2-CB2-CA2	2.31	132.77	129.94
1	A	63[B]	GYC	CD1-CE1-CZ	-2.19	117.48	119.88
1	A	63[B]	GYC	CA2-C2-N3	2.13	104.38	103.37
1	B	63[B]	GYC	CA1-C1-N3	2.02	127.48	124.85

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	63[A]	GYC	C1-CA1-CB1-SG1
1	B	63[B]	GYC	N1-CA1-CB1-SG1
1	B	63[B]	GYC	C1-CA1-CB1-SG1
1	A	63[A]	GYC	C2-CA2-CB2-CG2
1	B	63[B]	GYC	C2-CA2-CB2-CG2
1	A	63[A]	GYC	N2-CA2-CB2-CG2
1	B	63[B]	GYC	N2-CA2-CB2-CG2
1	A	63[A]	GYC	C3-CA3-N3-C1
1	A	63[B]	GYC	C3-CA3-N3-C1
1	A	63[A]	GYC	C1-CA1-CB1-SG1
1	A	63[B]	GYC	C1-CA1-CB1-SG1

There are no ring outliers.

4 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	63[A]	GYC	1	0
1	A	63[B]	GYC	2	0
1	B	63[A]	GYC	5	0
1	B	63[B]	GYC	5	0

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, 95^{th} 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	216/220 (98%)	-0.06	2 (0%)	81	83	31, 40, 61, 82	0
1	B	216/220 (98%)	0.10	1 (0%)	87	88	31, 45, 63, 76	0
All	All	432/440 (98%)	0.02	3 (0%)	84	86	31, 43, 61, 82	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	4	VAL	2.5
1	B	171	LEU	2.4
1	A	159	TYR	2.2

6.2 Non-standard residues in protein, DNA, RNA chains

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
1	GYC	B	63[A]	21/22	0.71	0.25	40,89,111,113	21
1	GYC	B	63[B]	21/22	0.71	0.25	35,57,68,70	21
1	GYC	A	63[A]	21/22	0.76	0.17	30,62,73,75	21
1	GYC	A	63[B]	21/22	0.76	0.17	44,81,99,102	21

6.3 Carbohydrates

There are no oligosaccharides 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.