



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 04:50 AM UTC

PDB ID : 8QMG / pdb_00008qmg
Title : Tobacco lectin Nictaba in apo state including I and Eu SAD datasets.
Authors : Bloch, Y.; Savvides, S.N.
Deposited on : 2023-09-22
Resolution : 2.30 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

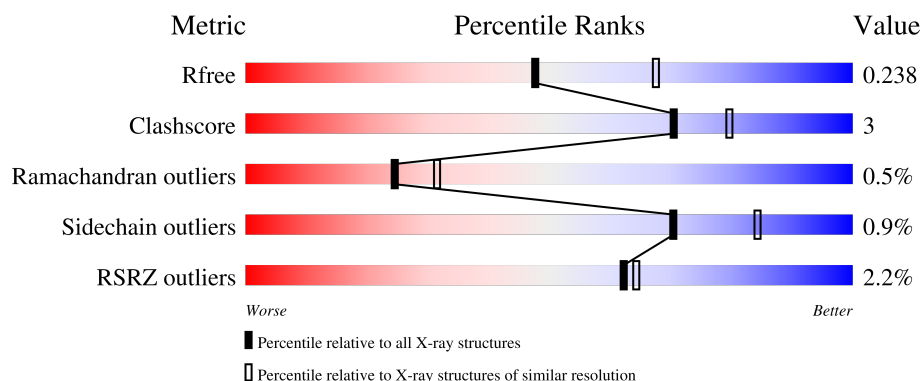
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	166	0% 93% 6% .
1	B	166	0% 92% 7% ..
1	C	166	3% 92% 7% .
1	D	166	2% 93% 6% ..
1	E	166	85% 14% .

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Mol	Chain	Length	Quality of chain
1	F	166	7% 90% 8% ..
1	G	166	3% 92% 7% ..
1	H	166	% 89% 10% .

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 22506 atoms, of which 10802 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 Nictaba.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	165	Total	C	H	N	O	S	0	2	0
			2708	869	1356	233	247	3			
1	B	165	Total	C	H	N	O	S	0	2	0
			2715	870	1360	236	247	2			
1	C	165	Total	C	H	N	O	S	0	0	0
			2690	864	1346	232	246	2			
1	D	165	Total	C	H	N	O	S	0	3	0
			2722	874	1361	234	250	3			
1	E	165	Total	C	H	N	O	S	0	0	0
			2690	864	1346	232	246	2			
1	F	165	Total	C	H	N	O	S	0	0	0
			2675	861	1335	231	246	2			
1	G	164	Total	C	H	N	O	S	0	0	0
			2682	862	1342	231	245	2			
1	H	165	Total	C	H	N	O	S	0	2	0
			2708	869	1356	233	247	3			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP Q94EW1
A	1	GLY	-	expression tag	UNP Q94EW1
B	0	MET	-	initiating methionine	UNP Q94EW1
B	1	GLY	-	expression tag	UNP Q94EW1
C	0	MET	-	initiating methionine	UNP Q94EW1
C	1	GLY	-	expression tag	UNP Q94EW1
D	0	MET	-	initiating methionine	UNP Q94EW1
D	1	GLY	-	expression tag	UNP Q94EW1
E	0	MET	-	initiating methionine	UNP Q94EW1
E	1	GLY	-	expression tag	UNP Q94EW1
F	0	MET	-	initiating methionine	UNP Q94EW1
F	1	GLY	-	expression tag	UNP Q94EW1
G	0	MET	-	initiating methionine	UNP Q94EW1

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Chain	Residue	Modelled	Actual	Comment	Reference
G	1	GLY	-	expression tag	UNP Q94EW1
H	0	MET	-	initiating methionine	UNP Q94EW1
H	1	GLY	-	expression tag	UNP Q94EW1

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Zn 1	0	0
2	B	4	Total 4	Zn 4	0	0
2	C	3	Total 3	Zn 3	0	0
2	D	4	Total 4	Zn 4	0	0
2	E	4	Total 4	Zn 4	0	0
2	F	5	Total 5	Zn 5	0	0
2	G	3	Total 3	Zn 3	0	0

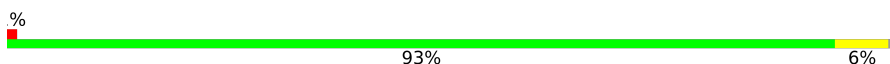
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	123	Total 123	O 123	0	0
3	B	113	Total 113	O 113	0	0
3	C	105	Total 105	O 105	0	0
3	D	111	Total 111	O 111	0	0
3	E	110	Total 110	O 110	0	0
3	F	107	Total 107	O 107	0	0
3	G	113	Total 113	O 113	0	0
3	H	110	Total 110	O 110	0	0

3 Residue-property plots [i](#)

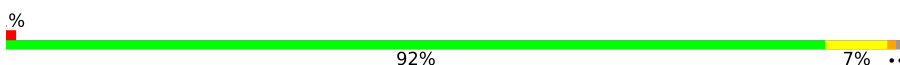
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Nictaba

Chain A: 



- Molecule 1: Nictaba

Chain B: 

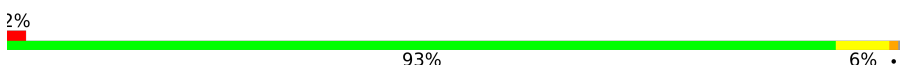


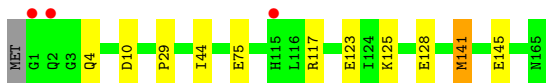
- Molecule 1: Nictaba

Chain C: 



- Molecule 1: Nictaba

Chain D: 

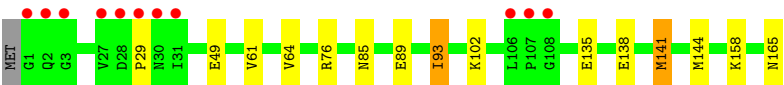
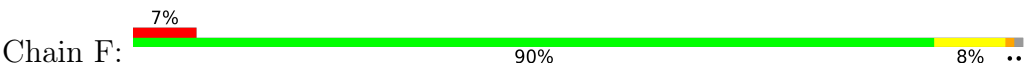


- Molecule 1: Nictaba

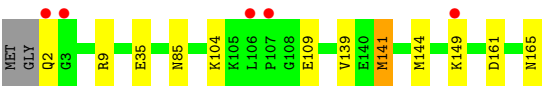
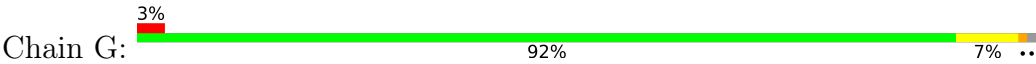
Chain E: 



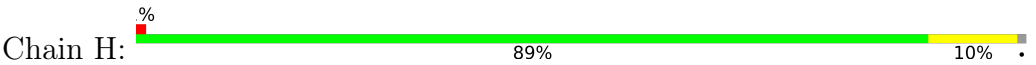
- Molecule 1: Nictaba



• Molecule 1: Nictaba



• Molecule 1: Nictaba



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.51Å 145.97Å 82.05Å 90.00° 95.95° 90.00°	Depositor
Resolution (Å)	71.23 – 2.30 71.23 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.9 (71.23-2.30) 94.2 (71.23-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.208 , 0.246 0.201 , 0.238	Depositor DCC
R_{free} test set	2274 reflections (3.50%)	wwPDB-VP
Wilson B-factor (Å ²)	33.9	Xtriage
Anisotropy	0.406	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 47.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	22506	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% 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](#)

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

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	0.23	0/1394	0.31	0/1883
1	B	0.22	0/1398	0.30	0/1889
1	C	0.21	0/1375	0.31	0/1859
1	D	0.22	0/1403	0.31	0/1895
1	E	0.22	0/1375	0.32	0/1859
1	F	0.23	0/1371	0.32	0/1855
1	G	0.21	0/1371	0.31	0/1854
1	H	0.20	0/1394	0.30	0/1883
All	All	0.22	0/11081	0.31	0/14977

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	1352	1356	1344	5	0
1	B	1355	1360	1352	8	0
1	C	1344	1346	1348	7	0
1	D	1361	1361	1349	7	0
1	E	1344	1346	1348	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	1340	1335	1337	12	0
1	G	1340	1342	1342	9	0
1	H	1352	1356	1344	11	0
2	A	1	0	0	0	0
2	B	4	0	0	0	0
2	C	3	0	0	0	0
2	D	4	0	0	0	0
2	E	4	0	0	0	0
2	F	5	0	0	0	0
2	G	3	0	0	0	0
3	A	123	0	0	1	3
3	B	113	0	0	4	1
3	C	105	0	0	0	0
3	D	111	0	0	3	0
3	E	110	0	0	5	1
3	F	107	0	0	5	2
3	G	113	0	0	5	1
3	H	110	0	0	1	0
All	All	11704	10802	10764	73	4

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 3.

The worst 5 of 73 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:49:GLU:OE2	3:B:301:HOH:O	1.88	0.92
1:F:138:GLU:OE2	3:F:301:HOH:O	1.98	0.82
1:H:52:ASN:ND2	3:H:201:HOH:O	2.15	0.79
1:B:75:GLU:OE2	3:B:302:HOH:O	2.02	0.76
1:D:75:GLU:OE2	3:D:301:HOH:O	2.04	0.76

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:354:HOH:O	3:G:309:HOH:O[1_655]	1.96	0.24
3:A:415:HOH:O	3:F:398:HOH:O[2_646]	2.12	0.08
3:A:305:HOH:O	3:E:305:HOH:O[2_645]	2.15	0.05
3:A:301:HOH:O	3:B:355:HOH:O[1_655]	2.16	0.04

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	165/166 (99%)	161 (98%)	3 (2%)	1 (1%)	21	27
1	B	165/166 (99%)	158 (96%)	6 (4%)	1 (1%)	21	27
1	C	163/166 (98%)	156 (96%)	6 (4%)	1 (1%)	21	27
1	D	166/166 (100%)	159 (96%)	6 (4%)	1 (1%)	21	27
1	E	163/166 (98%)	157 (96%)	6 (4%)	0	100	100
1	F	163/166 (98%)	157 (96%)	5 (3%)	1 (1%)	21	27
1	G	162/166 (98%)	157 (97%)	5 (3%)	0	100	100
1	H	165/166 (99%)	159 (96%)	5 (3%)	1 (1%)	21	27
All	All	1312/1328 (99%)	1264 (96%)	42 (3%)	6 (0%)	24	31

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	29	PRO
1	D	29	PRO
1	H	29	PRO
1	A	29	PRO
1	B	29	PRO

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	144/143 (101%)	144 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	144/143 (101%)	142 (99%)	2 (1%)	59	76
1	C	142/143 (99%)	142 (100%)	0	100	100
1	D	145/143 (101%)	142 (98%)	3 (2%)	47	66
1	E	142/143 (99%)	141 (99%)	1 (1%)	76	87
1	F	141/143 (99%)	139 (99%)	2 (1%)	59	76
1	G	142/143 (99%)	139 (98%)	3 (2%)	47	66
1	H	144/143 (101%)	144 (100%)	0	100	100
All	All	1144/1144 (100%)	1133 (99%)	11 (1%)	70	82

5 of 11 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	141	MET
1	G	2	GLN
1	G	144	MET
1	G	141	MET
1	D	141[B]	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	165	ASN
1	H	4	GLN
1	H	85	ASN
1	H	20	GLN
1	F	52	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 24 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	165/166 (99%)	-0.08	2 (1%) 76 77	16, 36, 72, 101	1 (0%)
1	B	165/166 (99%)	-0.16	2 (1%) 76 77	17, 36, 63, 104	1 (0%)
1	C	165/166 (99%)	0.10	5 (3%) 52 54	30, 41, 71, 103	0
1	D	165/166 (99%)	0.06	3 (1%) 67 69	17, 39, 71, 116	2 (1%)
1	E	165/166 (99%)	-0.04	0 100 100	29, 39, 64, 91	0
1	F	165/166 (99%)	0.01	11 (6%) 24 26	27, 34, 74, 110	0
1	G	164/166 (98%)	-0.02	5 (3%) 52 54	28, 36, 69, 118	0
1	H	165/166 (99%)	0.04	1 (0%) 85 86	17, 41, 70, 97	1 (0%)
All	All	1319/1328 (99%)	-0.01	29 (2%) 62 64	16, 38, 71, 118	5 (0%)

The worst 5 of 29 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	2	GLN	4.5
1	A	1	GLY	4.1
1	F	1	GLY	3.9
1	F	28	ASP	3.5
1	F	29	PRO	3.4

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 ⓘ

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, 95th 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
2	ZN	D	203	1/1	0.82	0.16	141,141,141,141	0
2	ZN	D	204	1/1	0.85	0.15	102,102,102,102	0
2	ZN	B	204	1/1	0.93	0.14	102,102,102,102	0
2	ZN	E	201	1/1	0.95	0.07	79,79,79,79	0
2	ZN	D	201	1/1	0.96	0.12	98,98,98,98	0
2	ZN	B	203	1/1	0.96	0.05	59,59,59,59	0
2	ZN	E	202	1/1	0.97	0.05	82,82,82,82	0
2	ZN	E	203	1/1	0.97	0.07	66,66,66,66	0
2	ZN	F	205	1/1	0.97	0.08	67,67,67,67	0
2	ZN	G	203	1/1	0.97	0.05	70,70,70,70	0
2	ZN	A	201	1/1	0.98	0.06	51,51,51,51	0
2	ZN	G	202	1/1	0.98	0.04	58,58,58,58	0
2	ZN	C	203	1/1	0.98	0.04	52,52,52,52	0
2	ZN	C	202	1/1	0.99	0.02	38,38,38,38	0
2	ZN	E	204	1/1	0.99	0.08	78,78,78,78	0
2	ZN	F	201	1/1	0.99	0.02	27,27,27,27	0
2	ZN	F	203	1/1	0.99	0.03	47,47,47,47	0
2	ZN	F	204	1/1	0.99	0.04	53,53,53,53	0
2	ZN	B	201	1/1	0.99	0.06	38,38,38,38	0
2	ZN	B	202	1/1	0.99	0.02	41,41,41,41	0
2	ZN	D	202	1/1	0.99	0.08	52,52,52,52	0
2	ZN	G	201	1/1	1.00	0.01	28,28,28,28	0
2	ZN	F	202	1/1	1.00	0.03	37,37,37,37	0
2	ZN	C	201	1/1	1.00	0.04	32,32,32,32	0

6.5 Other polymers ⓘ

There are no such residues in this entry.