



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 09:05 AM GMT

PDB ID : 2F9Z  
Title : Complex between the chemotaxis deamidase CheD and the chemotaxis phosphatase CheC from *Thermotoga maritima*  
Authors : Chao, X.; Park, S.Y; Bilwes, A.M.; Crane, B.R.  
Deposited on : 2005-12-06  
Resolution : 2.40 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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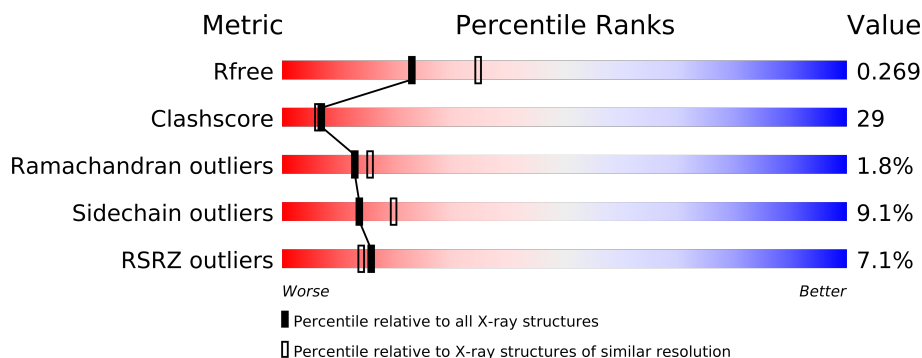
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.40 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))
$R_{free}$	66092	2207 (2.40-2.40)
Clashscore	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)
RSRZ outliers	66119	2210 (2.40-2.40)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	205	
1	B	205	
2	C	159	
2	D	159	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5552 atoms, of which 0 are hydrogen and 0 are deuterium.

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 chemotaxis protein CheC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	193	Total	C	N	O	S	0	0	0
			1452	939	222	281	10			
1	B	190	Total	C	N	O	S	0	0	0
			1433	930	224	268	11			

- Molecule 2 is a protein called PROTEIN (chemotaxis methylation protein).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	154	Total	C	N	O	S	0	0	0
			1140	719	201	210	10			
2	D	154	Total	C	N	O	S	0	0	0
			1135	715	202	209	9			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	ALA	-	CLONING ARTIFACT	UNP Q9X005
C	0	HIS	-	CLONING ARTIFACT	UNP Q9X005
D	-1	ALA	-	CLONING ARTIFACT	UNP Q9X005
D	0	HIS	-	CLONING ARTIFACT	UNP Q9X005

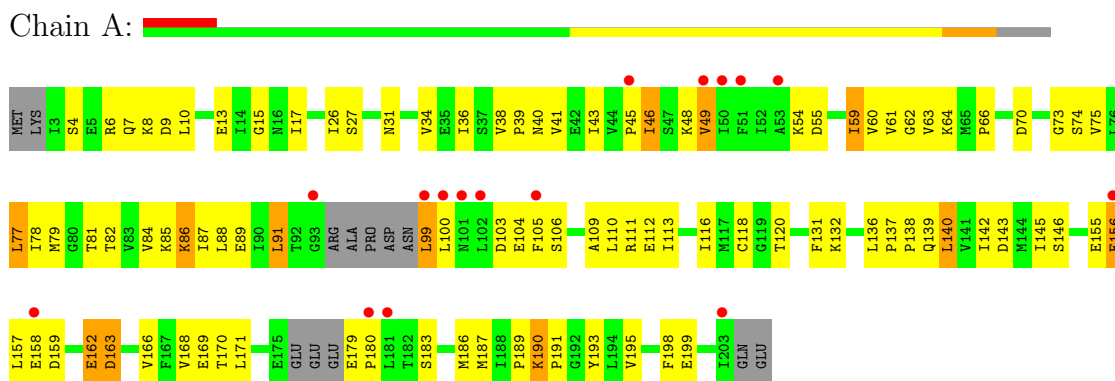
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	111	Total	O	0	0
			111	111		
3	B	100	Total	O	0	0
			100	100		
3	C	94	Total	O	0	0
			94	94		
3	D	87	Total	O	0	0
			87	87		

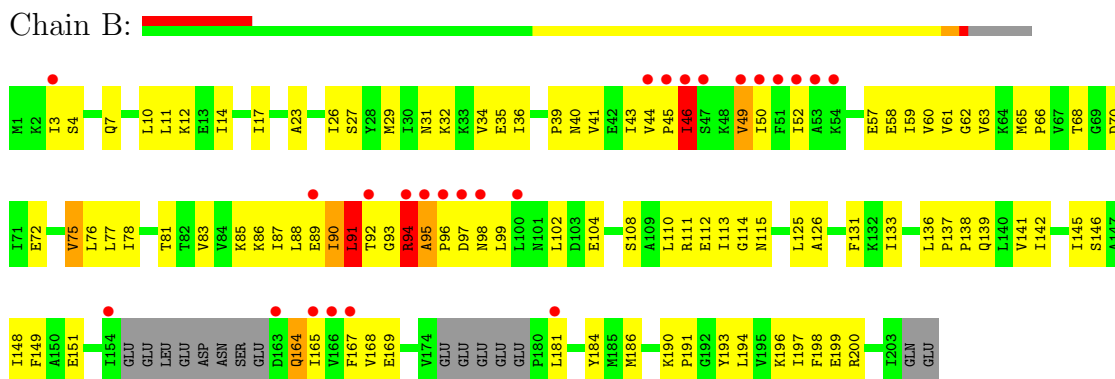
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

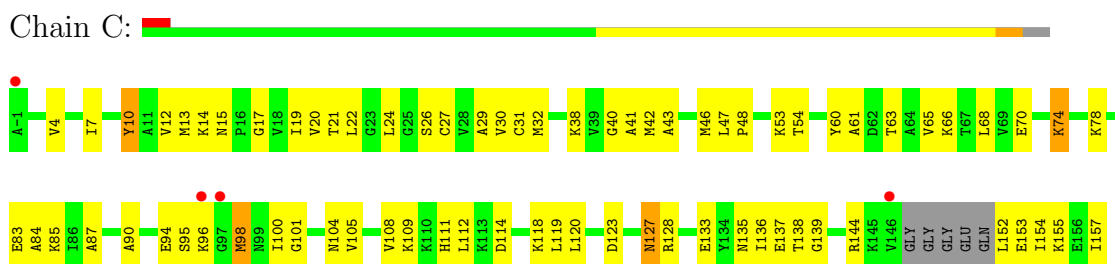
- Molecule 1: chemotaxis protein CheC



- Molecule 1: chemotaxis protein CheC

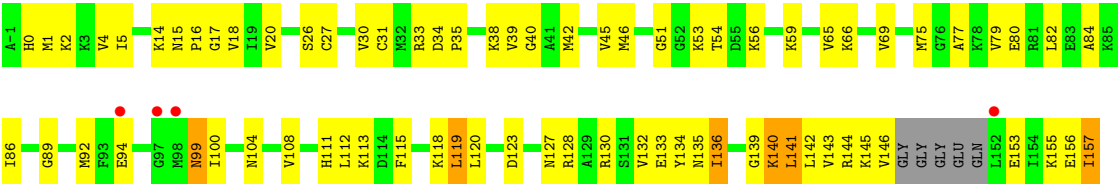


- Molecule 2: PROTEIN (chemotaxis methylation protein)



- Molecule 2: PROTEIN (chemotaxis methylation protein)





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.08Å 66.08Å 161.76Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.40 28.61 – 2.40	Depositor EDS
% Data completeness (in resolution range)	93.4 (30.00-2.40) 53.0 (28.61-2.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.63 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.211 , 0.275 0.216 , 0.269	Depositor DCC
$R_{free}$ test set	1630 reflections (9.98%)	DCC
Wilson B-factor (Å <sup>2</sup> )	41.9	Xtriage
Anisotropy	0.193	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 34.7	EDS
Estimated twinning fraction	0.186 for -h,-k,l 0.319 for h,-h-k,-l 0.217 for -k,-h,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.37$ , $\langle L^2 \rangle = 0.20$	Xtriage
Outliers	0 of 16378 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5552	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 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	0.36	0/1468	0.65	0/1987
1	B	0.42	0/1450	0.67	1/1961 (0.1%)
2	C	0.42	1/1149 (0.1%)	0.68	1/1531 (0.1%)
2	D	0.38	0/1145	0.69	1/1527 (0.1%)
All	All	0.39	1/5212 (0.0%)	0.67	3/7006 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	157	ILE	C-OXT	-5.57	1.12	1.23

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	114	ASP	CB-CA-C	-6.43	97.54	110.40
1	B	164	GLN	N-CA-C	-6.12	94.48	111.00
2	D	17	GLY	N-CA-C	5.54	126.95	113.10

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1452	0	1504	90	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1433	0	1518	97	0
2	C	1140	0	1220	65	0
2	D	1135	0	1202	59	0
3	A	111	0	0	10	0
3	B	100	0	0	18	0
3	C	94	0	0	6	0
3	D	87	0	0	6	0
All	All	5552	0	5444	305	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 29.

All (305) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:45:PRO:HB3	1:A:163:ASP:HB3	1.33	1.11
1:A:190:LYS:HB3	1:A:191:PRO:HD2	1.41	1.00
1:A:66:PRO:HG2	1:A:136:LEU:HB2	1.49	0.94
2:D:39:VAL:HG21	2:D:75:MET:HG2	1.51	0.91
2:C:7:ILE:H	2:C:7:ILE:HD12	1.36	0.90
2:C:13:MET:HG2	2:C:17:GLY:HA3	1.53	0.89
1:A:59:ILE:H	1:A:59:ILE:HD13	1.36	0.89
1:B:49:VAL:HG21	1:B:165:ILE:HD11	1.54	0.88
1:A:158:GLU:HG2	1:A:159:ASP:H	1.39	0.85
2:D:1:MET:SD	2:D:16:PRO:HB2	2.17	0.84
1:B:26:ILE:HD12	1:B:36:ILE:HD11	1.59	0.83
2:C:135:ASN:HD21	2:C:137:GLU:HB3	1.44	0.82
1:B:165:ILE:HD12	1:B:186:MET:SD	2.19	0.81
2:D:42:MET:HG2	3:D:1237:HOH:O	1.82	0.80
1:B:77:LEU:HD23	1:B:114:GLY:HA2	1.63	0.80
2:C:127:ASN:HD22	2:C:127:ASN:H	1.29	0.79
1:B:78:ILE:HB	1:B:184:TYR:HB2	1.65	0.78
1:B:63:VAL:HG11	1:B:114:GLY:HA3	1.68	0.75
1:B:41:VAL:HG12	1:B:168:VAL:HG22	1.69	0.75
1:A:13:GLU:O	1:A:17:ILE:HG12	1.87	0.75
2:D:99:ASN:H	2:D:99:ASN:ND2	1.82	0.75
1:B:190:LYS:HB3	1:B:191:PRO:HD2	1.69	0.74
1:A:171:LEU:HD11	1:A:180:PRO:HB2	1.67	0.74
1:B:12:LYS:HD3	1:B:40:ASN:HA	1.70	0.73
1:B:46:ILE:H	1:B:46:ILE:HD13	1.53	0.73
2:D:99:ASN:HD22	2:D:99:ASN:N	1.86	0.72
1:A:89:GLU:HB3	3:A:1187:HOH:O	1.91	0.70
2:C:43:ALA:HB1	3:C:1067:HOH:O	1.92	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:133:ILE:HG13	3:B:1357:HOH:O	1.92	0.70
2:D:86:ILE:HG23	2:D:119:LEU:HD21	1.72	0.70
1:B:34:VAL:HB	3:B:1104:HOH:O	1.91	0.70
2:D:14:LYS:C	2:D:16:PRO:HD2	2.12	0.70
1:A:190:LYS:HB3	1:A:191:PRO:CD	2.20	0.69
1:B:77:LEU:CD2	1:B:114:GLY:HA2	2.22	0.69
1:B:99:LEU:HB3	1:B:142:ILE:HD13	1.74	0.69
1:B:167:PHE:CD1	1:B:186:MET:HB3	2.27	0.69
1:A:60:VAL:HG21	1:A:78:ILE:HG23	1.75	0.69
1:A:45:PRO:CB	1:A:163:ASP:HB3	2.18	0.69
1:A:41:VAL:HG22	1:A:168:VAL:HG13	1.75	0.68
1:B:151:GLU:HG3	2:D:92:MET:SD	2.34	0.68
2:D:99:ASN:HD22	2:D:99:ASN:H	1.40	0.68
2:C:135:ASN:HD22	2:C:138:THR:H	1.40	0.68
1:A:136:LEU:HB3	1:A:137:PRO:HD2	1.76	0.68
1:B:44:VAL:HG23	3:B:1151:HOH:O	1.93	0.68
2:C:10:TYR:HD2	2:C:68:LEU:HD12	1.60	0.67
1:B:81:THR:HG22	1:B:85:LYS:HE2	1.75	0.67
1:A:74:SER:HB3	3:A:1365:HOH:O	1.94	0.67
2:C:127:ASN:N	2:C:127:ASN:HD22	1.92	0.67
1:B:190:LYS:HB2	1:B:193:TYR:CD2	2.31	0.66
1:A:64:LYS:HE2	3:A:1365:HOH:O	1.96	0.65
2:D:18:VAL:HG11	2:D:133:GLU:HB3	1.79	0.65
1:A:109:ALA:O	1:A:113:ILE:HG13	1.96	0.65
2:D:27:CYS:HA	3:D:1165:HOH:O	1.95	0.65
1:B:89:GLU:O	1:B:91:LEU:N	2.30	0.65
2:C:19:ILE:HG13	3:C:1057:HOH:O	1.97	0.65
1:A:88:LEU:HD21	1:A:106:SER:OG	1.97	0.64
2:C:24:LEU:HD11	2:C:29:ALA:HB2	1.78	0.64
1:B:148:ILE:HB	3:B:1253:HOH:O	1.98	0.64
2:D:15:ASN:N	2:D:16:PRO:HD2	2.13	0.64
1:A:39:PRO:HD3	3:A:1220:HOH:O	1.97	0.63
1:A:45:PRO:HA	1:A:163:ASP:O	1.98	0.63
2:D:31:CYS:SG	2:D:42:MET:HG3	2.40	0.62
1:B:3:ILE:HA	3:B:1116:HOH:O	1.99	0.62
1:B:102:LEU:HD13	3:B:1236:HOH:O	2.00	0.62
2:C:15:ASN:HB2	2:C:38:LYS:HB3	1.80	0.62
1:B:89:GLU:HG3	1:B:94:ARG:HA	1.82	0.61
2:D:141:LEU:HD12	2:D:157:ILE:HD11	1.82	0.61
1:B:141:VAL:HG13	1:B:148:ILE:HD13	1.82	0.61
1:A:106:SER:O	1:A:110:LEU:HB2	2.01	0.61
1:A:45:PRO:HG2	1:A:48:LYS:CB	2.31	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:34:VAL:HG22	1:B:35:GLU:H	1.66	0.61
1:B:193:TYR:HA	1:B:196:LYS:HB2	1.83	0.61
1:B:86:LYS:HG3	1:B:90:ILE:HD13	1.82	0.60
1:B:60:VAL:O	1:B:142:ILE:HA	2.01	0.60
1:A:190:LYS:CB	1:A:191:PRO:HD2	2.24	0.60
1:B:34:VAL:HG22	1:B:35:GLU:N	2.17	0.60
1:B:95:ALA:N	1:B:96:PRO:CD	2.65	0.60
1:A:59:ILE:N	1:A:59:ILE:HD13	2.11	0.60
1:B:29:MET:HG2	3:B:1304:HOH:O	2.02	0.60
1:A:82:THR:O	1:A:86:LYS:HG2	2.02	0.59
2:D:84:ALA:H	2:D:120:LEU:HD13	1.67	0.59
2:D:65:VAL:O	2:D:69:VAL:HG13	2.03	0.59
1:B:39:PRO:HG2	1:B:169:GLU:HG3	1.85	0.59
1:A:189:PRO:HB2	1:A:193:TYR:HB2	1.84	0.58
2:C:22:LEU:HD23	3:C:1356:HOH:O	2.03	0.58
2:C:7:ILE:H	2:C:7:ILE:CD1	2.12	0.58
1:A:99:LEU:HG	1:A:142:ILE:HG21	1.85	0.58
1:B:164:GLN:O	1:B:194:LEU:HD22	2.03	0.58
1:A:156:GLU:HG3	1:A:157:LEU:H	1.67	0.58
1:A:41:VAL:HG13	1:A:168:VAL:HG22	1.85	0.58
1:B:44:VAL:HB	1:B:45:PRO:HD2	1.83	0.58
2:D:77:ALA:HB1	2:D:82:LEU:HD11	1.85	0.58
2:D:133:GLU:HB2	2:D:142:LEU:HB3	1.86	0.58
2:C:32:MET:HA	2:C:83:GLU:O	2.03	0.58
1:B:11:LEU:HB2	1:B:41:VAL:HG21	1.85	0.58
2:C:128:ARG:HH12	2:C:152:LEU:HD13	1.69	0.58
2:C:19:ILE:HG21	2:C:42:MET:HB3	1.85	0.57
1:B:190:LYS:HB2	1:B:193:TYR:HD2	1.68	0.57
1:B:131:PHE:CZ	1:B:200:ARG:HG3	2.39	0.57
2:C:109:LYS:HE3	2:C:119:LEU:HD12	1.86	0.57
1:B:27:SER:HB2	1:B:34:VAL:HG12	1.86	0.57
1:A:63:VAL:CG1	1:A:138:PRO:HB2	2.34	0.57
1:A:64:LYS:HD3	1:A:139:GLN:NE2	2.19	0.57
2:D:30:VAL:HG22	2:D:108:VAL:HG21	1.86	0.57
1:A:156:GLU:HG3	1:A:157:LEU:N	2.20	0.57
1:A:59:ILE:H	1:A:59:ILE:CD1	2.07	0.56
1:B:89:GLU:C	1:B:91:LEU:H	2.08	0.56
2:D:145:LYS:HG2	3:D:1031:HOH:O	2.05	0.56
1:B:26:ILE:HG22	3:B:1104:HOH:O	2.03	0.56
1:A:9:ASP:O	1:A:13:GLU:HG3	2.05	0.56
1:B:95:ALA:N	1:B:96:PRO:HD2	2.20	0.56
1:B:50:ILE:HA	3:B:1303:HOH:O	2.05	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:43:ILE:HG12	1:A:198:PHE:HZ	1.70	0.56
2:D:145:LYS:O	2:D:153:GLU:HG3	2.06	0.56
1:B:96:PRO:HA	3:B:1236:HOH:O	2.06	0.55
2:D:123:ASP:OD2	2:D:143:VAL:HG11	2.05	0.55
1:B:39:PRO:HD2	1:B:169:GLU:O	2.06	0.55
1:A:60:VAL:HG23	1:A:79:MET:O	2.06	0.55
1:B:60:VAL:HG21	3:B:1253:HOH:O	2.07	0.55
2:D:135:ASN:O	2:D:139:GLY:N	2.40	0.55
1:A:171:LEU:N	3:A:1220:HOH:O	2.40	0.55
1:B:91:LEU:HD23	1:B:92:THR:HG23	1.89	0.54
1:A:84:VAL:HG13	1:A:110:LEU:HD11	1.89	0.54
1:B:145:ILE:HA	3:B:1253:HOH:O	2.07	0.54
2:D:4:VAL:HA	2:D:20:VAL:O	2.07	0.54
1:A:70:ASP:HB3	1:A:131:PHE:CD2	2.43	0.54
1:A:170:THR:OG1	1:A:183:SER:HB2	2.08	0.54
1:A:55:ASP:O	1:A:145:ILE:HG21	2.07	0.53
2:C:61:ALA:O	2:C:65:VAL:HG12	2.07	0.53
1:A:46:ILE:H	1:A:46:ILE:HD13	1.74	0.53
1:A:70:ASP:HB3	1:A:131:PHE:HD2	1.73	0.53
1:A:159:ASP:HA	2:C:46:MET:CE	2.39	0.53
1:A:112:GLU:O	1:A:116:ILE:HG13	2.09	0.53
2:D:111:HIS:O	2:D:115:PHE:HD1	1.91	0.53
2:D:1:MET:SD	2:D:16:PRO:CB	2.93	0.52
1:A:158:GLU:HG2	1:A:159:ASP:N	2.16	0.52
1:A:171:LEU:HB2	3:A:1220:HOH:O	2.09	0.52
1:A:4:SER:O	1:A:8:LYS:HG3	2.09	0.52
1:B:89:GLU:C	1:B:91:LEU:N	2.63	0.52
1:A:111:ARG:HD3	1:A:140:LEU:HB2	1.91	0.52
1:A:91:LEU:HD22	1:A:105:PHE:CE1	2.45	0.52
2:D:51:GLY:H	2:D:53:LYS:HE2	1.74	0.52
1:A:63:VAL:HG12	1:A:138:PRO:HB2	1.92	0.52
2:C:40:GLY:HA3	2:C:136:ILE:HD12	1.90	0.51
1:A:59:ILE:HB	1:A:99:LEU:HD11	1.91	0.51
1:B:88:LEU:HD13	1:B:102:LEU:HD21	1.91	0.51
2:C:47:LEU:HA	2:C:104:ASN:HD21	1.75	0.51
2:C:47:LEU:HB3	2:C:48:PRO:HD2	1.92	0.51
1:A:7:GLN:HB3	3:A:1207:HOH:O	2.10	0.51
2:C:65:VAL:HG11	2:C:111:HIS:HB2	1.91	0.51
2:C:144:ARG:HH21	2:C:154:ILE:HD13	1.75	0.51
1:B:60:VAL:HB	3:B:1169:HOH:O	2.09	0.51
1:B:115:ASN:HB2	1:B:138:PRO:HG3	1.93	0.51
2:C:4:VAL:HA	2:C:20:VAL:O	2.10	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:135:ASN:ND2	2:C:137:GLU:HB3	2.21	0.51
2:C:14:LYS:HG2	2:C:15:ASN:H	1.75	0.51
1:A:45:PRO:HA	1:A:163:ASP:C	2.31	0.50
1:A:45:PRO:O	1:A:49:VAL:HB	2.11	0.50
2:C:87:ALA:HA	2:C:123:ASP:O	2.11	0.50
1:B:148:ILE:HA	2:D:92:MET:HE2	1.92	0.50
2:D:30:VAL:O	2:D:42:MET:HA	2.11	0.50
1:A:39:PRO:CD	3:A:1220:HOH:O	2.59	0.50
1:A:73:GLY:HA2	1:A:190:LYS:HG3	1.94	0.49
2:C:21:THR:O	2:C:21:THR:HG23	2.11	0.49
1:A:6:ARG:O	1:A:10:LEU:HD23	2.11	0.49
2:D:35:PRO:O	2:D:38:LYS:HG3	2.11	0.49
1:A:81:THR:OG1	1:A:99:LEU:HD13	2.12	0.49
2:D:104:ASN:O	2:D:108:VAL:HG12	2.12	0.49
1:A:70:ASP:CG	1:A:132:LYS:H	2.16	0.49
2:C:101:GLY:O	2:C:105:VAL:HG23	2.13	0.49
2:C:13:MET:HB3	3:C:1057:HOH:O	2.12	0.49
2:C:20:VAL:HG12	2:C:21:THR:N	2.28	0.49
1:A:39:PRO:HD2	1:A:169:GLU:O	2.13	0.49
2:C:60:TYR:HB2	2:C:63:THR:HG22	1.94	0.49
1:B:149:PHE:CZ	1:B:186:MET:HE3	2.48	0.48
2:D:4:VAL:HG12	2:D:20:VAL:HB	1.95	0.48
2:C:144:ARG:HE	2:C:154:ILE:HG12	1.77	0.48
2:D:94:GLU:HG2	3:D:1294:HOH:O	2.13	0.48
1:A:46:ILE:HA	1:A:49:VAL:CG1	2.43	0.48
2:D:113:LYS:HD2	3:D:1037:HOH:O	2.12	0.48
1:A:45:PRO:O	1:A:49:VAL:N	2.46	0.48
2:C:65:VAL:HG11	2:C:111:HIS:CB	2.44	0.48
2:C:155:LYS:O	2:C:155:LYS:HG3	2.13	0.48
2:D:140:LYS:HG3	2:D:156:GLU:HB3	1.95	0.48
1:B:108:SER:HA	1:B:111:ARG:NH1	2.29	0.48
1:B:91:LEU:HB3	1:B:92:THR:H	1.34	0.48
1:B:65:MET:HB3	1:B:75:VAL:HG23	1.94	0.48
1:B:66:PRO:HG2	1:B:136:LEU:HB2	1.95	0.48
1:A:46:ILE:HA	1:A:49:VAL:HG12	1.94	0.48
1:B:72:GLU:HB2	1:B:193:TYR:CE2	2.49	0.48
2:D:157:ILE:HD13	2:D:157:ILE:H	1.79	0.48
1:B:193:TYR:O	1:B:197:ILE:HG13	2.14	0.47
1:B:45:PRO:HA	1:B:164:GLN:HG2	1.96	0.47
1:A:166:VAL:O	1:A:186:MET:HA	2.14	0.47
2:C:12:VAL:HG22	2:C:41:ALA:HB2	1.96	0.47
2:C:14:LYS:HG2	2:C:15:ASN:N	2.29	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:46:MET:SD	2:D:59:LYS:HE3	2.54	0.47
2:C:90:ALA:HB2	3:C:1330:HOH:O	2.14	0.47
1:B:86:LYS:HG3	1:B:90:ILE:CD1	2.44	0.47
1:A:27:SER:HB2	1:A:34:VAL:HG12	1.97	0.47
1:A:60:VAL:HG21	1:A:78:ILE:CG2	2.42	0.47
1:A:60:VAL:HG22	1:A:61:VAL:N	2.30	0.47
2:C:27:CYS:SG	2:C:46:MET:HA	2.55	0.46
1:B:26:ILE:HD12	1:B:36:ILE:CD1	2.39	0.46
2:D:132:VAL:HA	2:D:142:LEU:O	2.14	0.46
1:A:99:LEU:HD23	1:A:100:LEU:H	1.80	0.46
1:B:167:PHE:HD1	1:B:186:MET:HB3	1.77	0.46
1:A:99:LEU:HD23	1:A:100:LEU:N	2.31	0.46
1:B:86:LYS:CG	1:B:90:ILE:HD13	2.46	0.46
2:C:83:GLU:HB3	2:C:120:LEU:HD22	1.98	0.46
2:C:20:VAL:HG13	2:C:133:GLU:HG2	1.98	0.46
1:B:17:ILE:HD11	3:B:1287:HOH:O	2.15	0.46
2:C:108:VAL:O	2:C:112:LEU:HG	2.16	0.46
2:D:0:HIS:HB3	2:D:2:LYS:NZ	2.31	0.46
1:B:49:VAL:O	1:B:52:ILE:HG22	2.15	0.46
2:D:5:ILE:HD11	2:D:42:MET:HE2	1.98	0.46
1:A:45:PRO:HB3	1:A:163:ASP:CB	2.24	0.45
2:D:14:LYS:O	2:D:136:ILE:HD11	2.17	0.45
1:B:60:VAL:HG21	1:B:145:ILE:HA	1.99	0.45
2:C:53:LYS:HG3	2:C:54:THR:N	2.31	0.45
1:A:60:VAL:O	1:A:142:ILE:HA	2.15	0.45
1:A:179:GLU:CB	1:A:180:PRO:HD2	2.46	0.45
2:C:7:ILE:N	2:C:7:ILE:HD12	2.19	0.45
2:D:128:ARG:HH11	2:D:145:LYS:HE3	1.81	0.45
2:D:33:ARG:HA	2:D:40:GLY:HA2	1.99	0.45
1:B:94:ARG:O	1:B:95:ALA:HB3	2.17	0.45
1:A:162:GLU:HA	1:A:162:GLU:OE1	2.16	0.45
1:A:159:ASP:HA	2:C:46:MET:HE1	1.98	0.45
2:C:48:PRO:O	2:C:61:ALA:HB3	2.17	0.45
2:C:41:ALA:CB	2:C:68:LEU:HG	2.47	0.45
1:A:116:ILE:O	1:A:120:THR:HG23	2.17	0.44
2:C:118:LYS:HA	2:C:118:LYS:HD2	1.81	0.44
2:D:84:ALA:O	2:D:119:LEU:HD23	2.17	0.44
1:A:75:VAL:CG1	1:A:118:CYS:HB3	2.47	0.44
1:B:190:LYS:HB3	1:B:191:PRO:CD	2.44	0.44
2:C:65:VAL:HG13	2:C:66:LYS:N	2.33	0.44
2:C:66:LYS:HD2	3:C:1143:HOH:O	2.17	0.44
1:B:14:ILE:HD13	1:B:125:LEU:HD12	1.99	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:30:VAL:HB	2:C:43:ALA:HB3	1.99	0.44
2:D:120:LEU:HD12	2:D:120:LEU:N	2.32	0.44
1:B:167:PHE:CE1	1:B:186:MET:HB3	2.52	0.44
1:B:136:LEU:HB3	1:B:137:PRO:HD2	1.99	0.44
2:C:70:GLU:O	2:C:74:LYS:HG3	2.18	0.44
1:B:63:VAL:O	1:B:76:LEU:HD12	2.18	0.44
1:B:58:GLU:HG2	1:B:59:ILE:N	2.33	0.44
1:A:85:LYS:O	1:A:87:ILE:N	2.51	0.44
1:B:4:SER:HB3	1:B:7:GLN:HG3	1.99	0.44
1:B:148:ILE:HA	2:D:92:MET:CE	2.48	0.43
1:A:60:VAL:CG2	1:A:78:ILE:HG23	2.46	0.43
1:B:76:LEU:HB3	1:B:186:MET:HG3	1.99	0.43
2:C:98:MET:O	2:C:98:MET:HG3	2.18	0.43
2:C:30:VAL:HG22	2:C:108:VAL:HG11	2.00	0.43
1:A:63:VAL:HA	1:A:139:GLN:O	2.17	0.43
2:D:69:VAL:HG21	2:D:115:PHE:CE2	2.54	0.43
1:A:145:ILE:HG23	1:A:146:SER:N	2.33	0.43
1:B:23:ALA:HB1	1:B:34:VAL:HG13	2.01	0.43
1:B:91:LEU:C	1:B:93:GLY:H	2.22	0.43
2:D:144:ARG:HA	2:D:153:GLU:O	2.18	0.43
1:A:15:GLY:C	1:A:38:VAL:HG21	2.40	0.43
2:D:18:VAL:HA	2:D:134:TYR:O	2.18	0.43
2:C:135:ASN:O	2:C:139:GLY:N	2.51	0.42
1:B:49:VAL:HG12	3:B:1103:HOH:O	2.19	0.42
1:B:62:GLY:N	1:B:141:VAL:O	2.52	0.42
1:B:59:ILE:HD12	2:D:146:VAL:HG12	2.01	0.42
2:C:127:ASN:N	2:C:127:ASN:ND2	2.64	0.42
1:B:94:ARG:C	1:B:96:PRO:CD	2.87	0.42
1:B:98:ASN:O	1:B:102:LEU:HB2	2.20	0.42
1:B:83:VAL:HG22	1:B:181:LEU:HD13	2.02	0.42
1:B:7:GLN:NE2	1:B:198:PHE:CG	2.87	0.42
1:A:145:ILE:HD13	3:A:1363:HOH:O	2.19	0.42
1:B:87:ILE:HG21	1:B:113:ILE:CD1	2.49	0.42
2:C:26:SER:HB2	2:C:101:GLY:CA	2.50	0.42
2:D:26:SER:O	2:D:89:GLY:HA2	2.20	0.42
1:B:191:PRO:HA	3:B:1336:HOH:O	2.18	0.42
1:B:61:VAL:HG22	1:B:99:LEU:HD13	2.02	0.42
2:D:66:LYS:HA	2:D:69:VAL:HG22	2.02	0.42
2:D:53:LYS:HG2	2:D:53:LYS:H	1.62	0.42
2:D:100:ILE:HG13	2:D:100:ILE:H	1.54	0.42
1:A:195:VAL:O	1:A:199:GLU:HG2	2.19	0.42
2:C:26:SER:HB2	2:C:101:GLY:HA2	2.02	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:62:GLY:HA2	1:A:77:LEU:O	2.19	0.41
1:A:189:PRO:HB2	1:A:193:TYR:CB	2.50	0.41
1:B:193:TYR:HA	1:B:196:LYS:CB	2.50	0.41
1:A:171:LEU:CB	3:A:1220:HOH:O	2.66	0.41
2:C:4:VAL:HG23	2:C:21:THR:HA	2.03	0.41
1:A:84:VAL:O	1:A:88:LEU:HB2	2.21	0.41
2:D:79:VAL:HA	2:D:82:LEU:HD13	2.02	0.41
2:C:48:PRO:HD3	2:C:104:ASN:ND2	2.35	0.41
1:A:137:PRO:HA	1:A:138:PRO:HD2	1.89	0.41
2:D:143:VAL:HB	2:D:155:LYS:HG3	2.02	0.41
2:C:53:LYS:HG3	2:C:54:THR:H	1.86	0.41
1:B:32:LYS:HG2	3:B:1018:HOH:O	2.20	0.41
1:B:145:ILE:HG23	1:B:146:SER:N	2.36	0.41
2:C:100:ILE:HG12	2:C:100:ILE:H	1.71	0.41
2:C:94:GLU:OE2	2:C:94:GLU:HA	2.21	0.41
2:D:128:ARG:HE	2:D:145:LYS:CE	2.33	0.40
1:A:26:ILE:HD12	1:A:36:ILE:HD11	2.03	0.40
1:B:68:THR:HG22	3:B:1285:HOH:O	2.21	0.40
1:B:104:GLU:HG2	1:B:104:GLU:H	1.55	0.40
1:A:179:GLU:HB3	1:A:180:PRO:HD2	2.03	0.40
2:D:34:ASP:OD1	2:D:77:ALA:HA	2.21	0.40
1:B:65:MET:SD	1:B:75:VAL:CG2	3.10	0.40
2:D:56:LYS:HG2	3:D:1228:HOH:O	2.21	0.40
1:B:46:ILE:HG21	1:B:165:ILE:HD13	2.03	0.40
2:C:31:CYS:O	2:C:84:ALA:HA	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	187/205 (91%)	169 (90%)	14 (8%)	4 (2%)	11	12
1	B	184/205 (90%)	159 (86%)	19 (10%)	6 (3%)	6	5
2	C	150/159 (94%)	131 (87%)	17 (11%)	2 (1%)	18	24

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	150/159 (94%)	137 (91%)	13 (9%)	0	100	100
All	All	671/728 (92%)	596 (89%)	63 (9%)	12 (2%)	13	15

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	90	ILE
1	B	91	LEU
1	B	94	ARG
1	B	95	ALA
1	A	163	ASP
1	A	54	LYS
1	A	86	LYS
1	B	126	ALA
1	A	190	LYS
1	B	46	ILE
2	C	74	LYS
2	C	153	GLU

### 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	162/179 (90%)	146 (90%)	16 (10%)	11	16
1	B	161/179 (90%)	146 (91%)	15 (9%)	13	19
2	C	119/122 (98%)	112 (94%)	7 (6%)	28	42
2	D	117/122 (96%)	104 (89%)	13 (11%)	9	12
All	All	559/602 (93%)	508 (91%)	51 (9%)	14	20

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	40	ASN
1	A	46	ILE

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Mol	Chain	Res	Type
1	A	49	VAL
1	A	59	ILE
1	A	77	LEU
1	A	91	LEU
1	A	99	LEU
1	A	103	ASP
1	A	104	GLU
1	A	140	LEU
1	A	143	ASP
1	A	155	GLU
1	A	156	GLU
1	A	162	GLU
1	A	187	MET
1	B	10	LEU
1	B	31	ASN
1	B	43	ILE
1	B	46	ILE
1	B	49	VAL
1	B	57	GLU
1	B	70	ASP
1	B	75	VAL
1	B	91	LEU
1	B	94	ARG
1	B	97	ASP
1	B	110	LEU
1	B	112	GLU
1	B	139	GLN
1	B	199	GLU
2	C	10	TYR
2	C	78	LYS
2	C	85	LYS
2	C	95	SER
2	C	96	LYS
2	C	98	MET
2	C	127	ASN
2	D	45	VAL
2	D	54	THR
2	D	80	GLU
2	D	99	ASN
2	D	112	LEU
2	D	118	LYS
2	D	119	LEU

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Mol	Chain	Res	Type
2	D	127	ASN
2	D	130	ARG
2	D	136	ILE
2	D	140	LYS
2	D	141	LEU
2	D	157	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
2	C	15	ASN
2	C	104	ASN
2	C	127	ASN
2	C	135	ASN
2	D	99	ASN
2	D	104	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	193/205 (94%)	0.36	16 (8%) 11 10	15, 50, 94, 115	0
1	B	190/205 (92%)	0.51	25 (13%) 4 4	20, 48, 113, 120	0
2	C	154/159 (96%)	-0.13	4 (2%) 53 51	8, 30, 63, 85	0
2	D	154/159 (96%)	-0.02	4 (2%) 53 51	14, 32, 78, 108	0
All	All	691/728 (94%)	0.21	49 (7%) 16 14	8, 40, 99, 120	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	46	ILE	7.1
1	B	50	ILE	6.5
2	D	97	GLY	6.2
1	B	96	PRO	5.3
2	C	97	GLY	5.0
1	A	156	GLU	4.9
1	B	53	ALA	4.3
1	B	49	VAL	4.2
1	A	99	LEU	4.2
1	A	101	ASN	4.1
2	C	96	LYS	4.0
2	C	-1	ALA	3.8
1	A	105	PHE	3.7
1	B	45	PRO	3.7
1	B	95	ALA	3.5
1	B	98	ASN	3.5
1	A	100	LEU	3.5
1	A	50	ILE	3.4
1	B	97	ASP	3.4
1	B	44	VAL	3.4
2	D	98	MET	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	51	PHE	3.4
1	B	94	ARG	3.4
1	B	47	SER	3.2
1	A	93	GLY	3.2
1	B	154	ILE	3.2
1	B	52	ILE	3.1
2	C	146	VAL	3.0
1	A	180	PRO	3.0
1	B	92	THR	2.9
1	B	163	ASP	2.9
1	B	181	LEU	2.9
1	B	54	LYS	2.8
1	B	100	LEU	2.6
1	A	49	VAL	2.6
1	B	167	PHE	2.6
1	A	203	ILE	2.5
1	B	89	GLU	2.5
1	A	53	ALA	2.4
1	A	181	LEU	2.4
1	B	165	ILE	2.4
1	A	102	LEU	2.3
1	A	45	PRO	2.3
2	D	152	LEU	2.2
1	A	158	GLU	2.2
2	D	94	GLU	2.2
1	B	51	PHE	2.2
1	B	3	ILE	2.1
1	B	166	VAL	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 ⓘ

There are no carbohydrates 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.