



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:34 AM GMT

PDB ID : 2AUK
Title : Structure of E. coli RNA polymerase beta' G/G' insert
Authors : Chlenov, M.; Masuda, S.; Murakami, K.S.; Nikiforov, V.; Darst, S.A.; Mustaev, A.
Deposited on : 2005-08-28
Resolution : 2.30 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

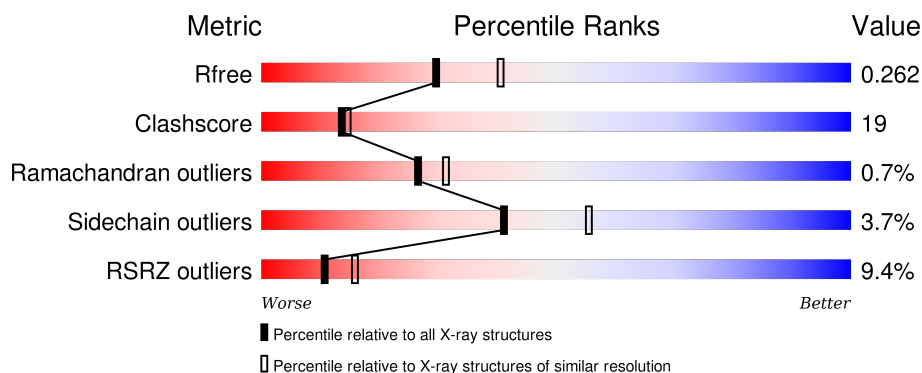
1 Overall quality at a glance

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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	190	<div> <div>6%</div> <div> <div></div> <div>65%</div> <div>31%</div> <div>• •</div> </div> </div>
1	B	190	<div> <div>4%</div> <div> <div></div> <div>69%</div> <div>24%</div> <div>• 5%</div> </div> </div>
1	C	190	<div> <div>4%</div> <div> <div></div> <div>75%</div> <div>19%</div> <div>• •</div> </div> </div>
1	D	190	<div> <div>15%</div> <div> <div></div> <div>64%</div> <div>29%</div> <div>• •</div> </div> </div>
1	E	190	<div> <div>16%</div> <div> <div></div> <div>62%</div> <div>35%</div> <div>• •</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7306 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 DNA-directed RNA polymerase beta' chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	189	Total	C	N	O	S	0	0	0
			1401	872	240	285	4			
1	B	181	Total	C	N	O	S	0	0	0
			1354	845	232	274	3			
1	C	183	Total	C	N	O	S	0	0	0
			1364	851	234	276	3			
1	D	183	Total	C	N	O	S	0	0	0
			1351	842	233	273	3			
1	E	188	Total	C	N	O	S	0	0	0
			1379	861	235	279	4			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	CLONING ARTIFACT	UNP P0A8T7
A	2	SER	-	CLONING ARTIFACT	UNP P0A8T7
A	3	HIS	-	CLONING ARTIFACT	UNP P0A8T7
A	4	MET	-	CLONING ARTIFACT	UNP P0A8T7
B	1	GLY	-	CLONING ARTIFACT	UNP P0A8T7
B	2	SER	-	CLONING ARTIFACT	UNP P0A8T7
B	3	HIS	-	CLONING ARTIFACT	UNP P0A8T7
B	4	MET	-	CLONING ARTIFACT	UNP P0A8T7
C	1	GLY	-	CLONING ARTIFACT	UNP P0A8T7
C	2	SER	-	CLONING ARTIFACT	UNP P0A8T7
C	3	HIS	-	CLONING ARTIFACT	UNP P0A8T7
C	4	MET	-	CLONING ARTIFACT	UNP P0A8T7
D	1	GLY	-	CLONING ARTIFACT	UNP P0A8T7
D	2	SER	-	CLONING ARTIFACT	UNP P0A8T7
D	3	HIS	-	CLONING ARTIFACT	UNP P0A8T7
D	4	MET	-	CLONING ARTIFACT	UNP P0A8T7
E	1	GLY	-	CLONING ARTIFACT	UNP P0A8T7
E	2	SER	-	CLONING ARTIFACT	UNP P0A8T7
E	3	HIS	-	CLONING ARTIFACT	UNP P0A8T7

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Chain	Residue	Modelled	Actual	Comment	Reference
E	4	MET	-	CLONING ARTIFACT	UNP P0A8T7

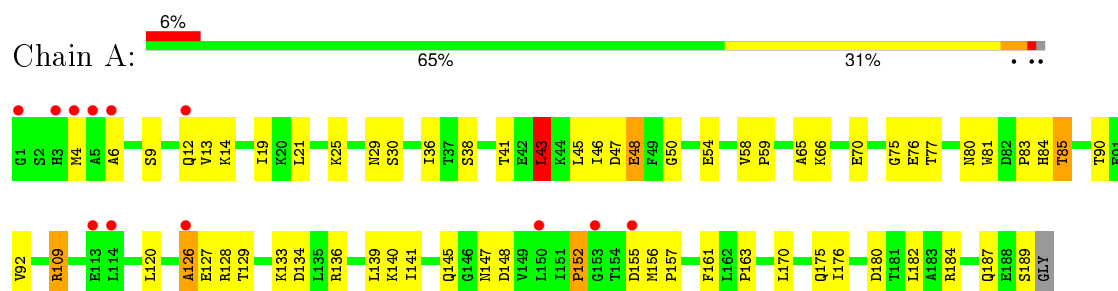
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	97	Total O 97 97	0	0
2	B	111	Total O 111 111	0	0
2	C	91	Total O 91 91	0	0
2	D	87	Total O 87 87	0	0
2	E	71	Total O 71 71	0	0

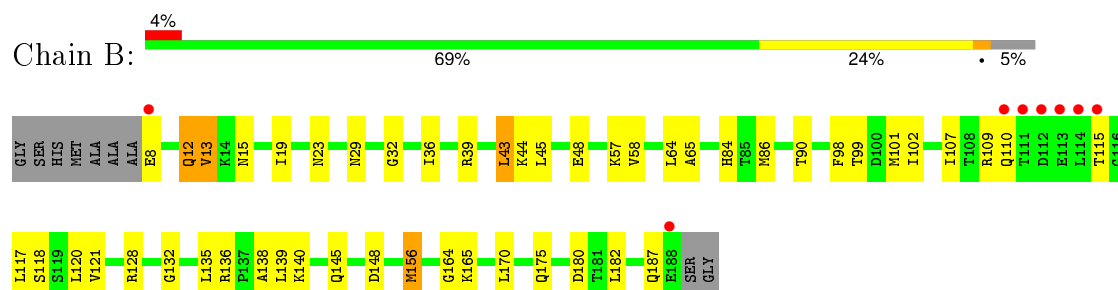
3 Residue-property plots

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.

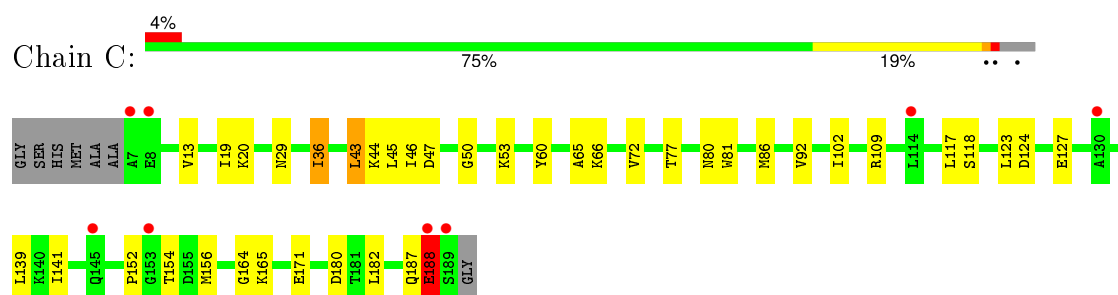
- Molecule 1: DNA-directed RNA polymerase beta' chain



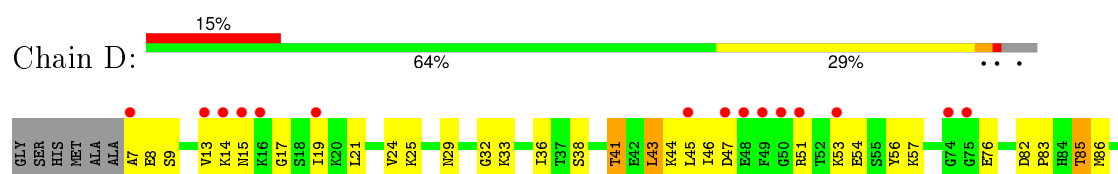
- Molecule 1: DNA-directed RNA polymerase beta' chain

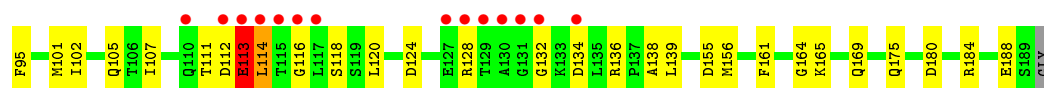


- Molecule 1: DNA-directed RNA polymerase beta' chain

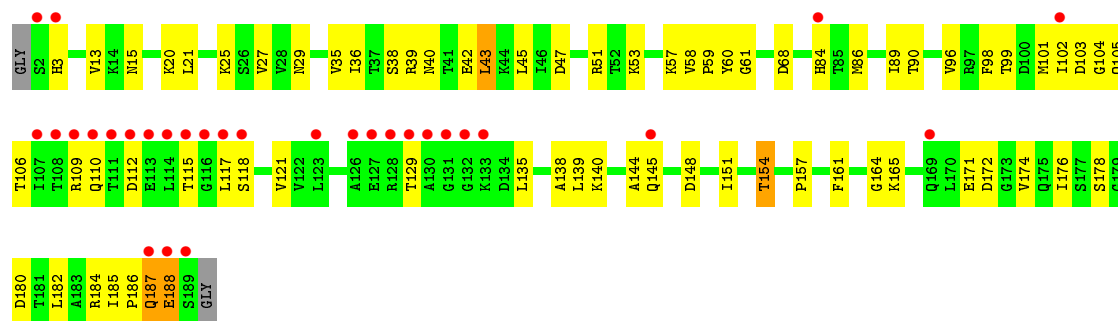


- Molecule 1: DNA-directed RNA polymerase beta' chain





- Molecule 1: DNA-directed RNA polymerase beta' chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.03Å 95.13Å 86.07Å 90.00° 92.06° 90.00°	Depositor
Resolution (Å)	35.00 – 2.30 34.10 – 2.19	Depositor EDS
% Data completeness (in resolution range)	(Not available) (35.00-2.30) 90.3 (34.10-2.19)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.76 (at 2.20Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.225 , 0.264 0.220 , 0.262	Depositor DCC
R_{free} test set	3069 reflections (8.04%)	DCC
Wilson B-factor (Å ²)	29.4	Xtriage
Anisotropy	0.129	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 51.3	EDS
Estimated twinning fraction	0.027 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 43459 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7306	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.23% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.37	0/1418	0.68	1/1920 (0.1%)
1	B	0.39	0/1371	0.69	0/1857
1	C	0.40	0/1381	0.70	1/1871 (0.1%)
1	D	0.41	1/1368 (0.1%)	0.67	0/1855
1	E	0.36	0/1396	0.65	1/1893 (0.1%)
All	All	0.39	1/6934 (0.0%)	0.68	3/9396 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	113	GLU	CA-CB	6.00	1.67	1.53

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	188	GLU	N-CA-C	7.30	130.71	111.00
1	A	43	LEU	CA-CB-CG	5.46	127.86	115.30
1	E	154	THR	N-CA-C	-5.19	96.99	111.00

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	1401	0	1415	58	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1354	0	1372	55	0
1	C	1364	0	1379	35	0
1	D	1351	0	1350	64	0
1	E	1379	0	1384	62	0
2	A	97	0	0	7	0
2	B	111	0	0	7	0
2	C	91	0	0	3	0
2	D	87	0	0	9	0
2	E	71	0	0	7	0
All	All	7306	0	6900	255	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 19.

All (255) 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:85:THR:HG21	1:A:184:ARG:HD3	1.38	1.03
1:B:107:ILE:HD12	1:B:120:LEU:HD22	1.50	0.93
1:D:21:LEU:HB3	1:D:24:VAL:HG11	1.54	0.88
1:D:45:LEU:HB2	1:D:54:GLU:HB2	1.56	0.87
1:E:3:HIS:HA	2:E:192:HOH:O	1.78	0.82
1:B:13:VAL:HG21	1:B:45:LEU:HD13	1.64	0.80
1:D:114:LEU:HD23	1:D:114:LEU:N	1.99	0.78
1:C:47:ASP:OD2	1:C:53:LYS:HE2	1.82	0.78
1:E:187:GLN:HG2	1:E:188:GLU:H	1.49	0.77
1:D:9:SER:HB3	1:D:83:PRO:HG3	1.65	0.77
1:E:98:PHE:HA	1:E:101:MET:HE3	1.67	0.76
1:C:154:THR:HG22	1:C:156:MET:HB2	1.66	0.75
1:D:47:ASP:OD2	1:D:53:LYS:HE3	1.87	0.74
1:D:114:LEU:HD23	1:D:114:LEU:H	1.53	0.74
1:B:57:LYS:HD3	1:C:171:GLU:HG2	1.67	0.74
1:E:13:VAL:HG21	1:E:45:LEU:HD13	1.69	0.74
1:A:77:THR:OG1	1:D:155:ASP:HB3	1.88	0.73
1:D:85:THR:HG21	1:D:184:ARG:HD3	1.70	0.73
1:B:101:MET:HE2	1:B:138:ALA:O	1.87	0.73
1:B:29:ASN:HB2	1:B:180:ASP:OD1	1.88	0.73
1:B:156:MET:HE3	1:C:77:THR:HG22	1.69	0.73
1:E:21:LEU:HD21	1:E:43:LEU:HD22	1.71	0.72
1:B:140:LYS:HZ1	1:B:148:ASP:CG	1.92	0.72
1:A:84:HIS:O	1:A:187:GLN:HG3	1.90	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:13:VAL:HG22	1:B:45:LEU:HD22	1.72	0.71
1:B:136:ARG:HG2	1:B:136:ARG:HH11	1.55	0.71
1:E:25:LYS:HB2	1:E:38:SER:HB3	1.72	0.70
1:D:111:THR:HG23	1:D:116:GLY:O	1.92	0.70
1:D:111:THR:HG22	1:D:113:GLU:H	1.57	0.69
1:E:117:LEU:HD23	2:E:238:HOH:O	1.91	0.69
1:A:136:ARG:CZ	2:A:229:HOH:O	2.40	0.69
1:D:43:LEU:HD12	1:D:44:LYS:N	2.07	0.69
1:E:99:THR:H	1:E:101:MET:HE2	1.59	0.68
1:D:9:SER:CB	1:D:83:PRO:HG3	2.24	0.68
1:E:29:ASN:HB2	1:E:180:ASP:OD1	1.94	0.67
1:D:21:LEU:HB3	1:D:24:VAL:CG1	2.24	0.66
1:B:156:MET:HE1	1:C:65:ALA:HB2	1.78	0.66
1:A:85:THR:HG22	2:A:208:HOH:O	1.95	0.66
1:E:151:ILE:O	1:E:154:THR:O	2.14	0.66
1:C:86:MET:HB2	1:C:187:GLN:NE2	2.09	0.65
1:B:156:MET:CE	1:C:77:THR:HG22	2.26	0.65
1:C:29:ASN:HB2	1:C:180:ASP:OD1	1.96	0.65
1:A:21:LEU:HD21	1:A:43:LEU:HD22	1.79	0.65
1:C:164:GLY:O	1:C:165:LYS:HB2	1.97	0.64
1:C:46:ILE:CG2	1:C:50:GLY:HA2	2.27	0.64
1:A:65:ALA:HB2	1:D:156:MET:CE	2.27	0.64
1:A:65:ALA:HB2	1:D:156:MET:HE1	1.80	0.64
1:D:169:GLN:HA	2:D:221:HOH:O	1.97	0.64
1:A:155:ASP:O	1:A:157:PRO:HD3	1.98	0.64
1:E:47:ASP:OD2	1:E:51:ARG:HB2	1.98	0.64
1:D:14:LYS:HG3	2:D:248:HOH:O	1.98	0.63
1:E:13:VAL:HG22	1:E:45:LEU:HD22	1.81	0.63
1:B:98:PHE:HA	1:B:101:MET:CE	2.29	0.63
1:B:57:LYS:HD3	1:C:171:GLU:CG	2.29	0.63
1:E:15:ASN:HB2	1:E:45:LEU:HD21	1.81	0.62
1:E:27:VAL:HG21	1:E:178:SER:HB2	1.81	0.62
1:B:15:ASN:HB2	1:B:45:LEU:HD21	1.81	0.62
1:D:111:THR:HA	1:D:116:GLY:O	2.00	0.62
1:A:136:ARG:NH1	2:A:229:HOH:O	2.32	0.62
1:C:66:LYS:HD2	1:C:72:VAL:HG12	1.82	0.62
1:E:112:ASP:HB2	1:E:115:THR:HB	1.81	0.62
1:A:13:VAL:HG21	1:A:45:LEU:HD13	1.81	0.62
1:C:13:VAL:HG13	1:C:45:LEU:HD22	1.81	0.61
1:A:19:ILE:HD12	1:A:43:LEU:HD11	1.82	0.61
1:A:129:THR:O	1:A:133:LYS:HG2	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:169:GLN:HB2	2:D:206:HOH:O	2.00	0.61
1:A:152:PRO:HB2	2:A:217:HOH:O	2.01	0.61
1:D:17:GLY:HA3	1:D:46:ILE:O	2.01	0.60
1:B:39:ARG:HD3	2:B:264:HOH:O	2.00	0.60
1:D:19:ILE:HG23	1:D:43:LEU:HD11	1.82	0.60
1:D:36:ILE:HG21	1:D:41:THR:HG21	1.85	0.59
1:A:170:LEU:HD13	1:A:176:ILE:HG22	1.85	0.59
1:D:112:ASP:O	1:D:114:LEU:N	2.30	0.59
1:B:43:LEU:HD12	1:B:44:LYS:N	2.18	0.58
1:B:136:ARG:HG2	1:B:136:ARG:NH1	2.16	0.58
1:E:98:PHE:HA	1:E:101:MET:CE	2.33	0.58
1:A:19:ILE:HG23	1:A:43:LEU:HD13	1.85	0.58
1:B:140:LYS:NZ	1:B:148:ASP:CG	2.56	0.58
1:B:132:GLY:HA2	1:B:135:LEU:HD12	1.86	0.57
1:E:21:LEU:CD2	1:E:43:LEU:HD22	2.33	0.57
1:C:19:ILE:HG23	1:C:43:LEU:CD1	2.34	0.56
1:B:101:MET:HE1	1:B:139:LEU:HD23	1.88	0.56
1:D:128:ARG:HB3	1:D:132:GLY:HA3	1.87	0.56
1:E:115:THR:HA	2:E:232:HOH:O	2.05	0.55
1:E:47:ASP:OD2	1:E:53:LYS:HE2	2.07	0.55
1:A:36:ILE:HD12	1:A:58:VAL:HG11	1.89	0.55
1:D:41:THR:HG23	2:D:191:HOH:O	2.05	0.55
1:C:92:VAL:HG23	1:C:141:ILE:HG21	1.88	0.55
1:E:35:VAL:HG11	1:E:89:ILE:HD13	1.89	0.55
1:B:43:LEU:C	1:B:43:LEU:HD12	2.27	0.55
1:C:154:THR:CG2	1:C:156:MET:HB2	2.35	0.55
1:E:140:LYS:HE3	2:E:225:HOH:O	2.06	0.55
1:D:13:VAL:HG23	1:D:76:GLU:O	2.08	0.54
1:B:156:MET:CE	1:C:65:ALA:HB2	2.37	0.54
1:C:43:LEU:HD12	1:C:44:LYS:N	2.22	0.54
1:B:84:HIS:O	1:B:187:GLN:HG3	2.08	0.54
1:B:13:VAL:CG2	1:B:45:LEU:HD13	2.34	0.53
1:D:15:ASN:ND2	1:D:53:LYS:HD3	2.23	0.53
1:A:29:ASN:HB2	1:A:180:ASP:OD1	2.09	0.53
1:E:86:MET:HB3	2:E:204:HOH:O	2.08	0.53
1:E:144:ALA:HB3	1:E:145:GLN:NE2	2.23	0.53
1:A:85:THR:HG21	1:A:184:ARG:HB3	1.91	0.53
1:C:156:MET:HG2	2:C:257:HOH:O	2.08	0.53
1:D:107:ILE:HD12	1:D:120:LEU:HD22	1.90	0.53
1:D:118:SER:CB	1:E:42:GLU:OE2	2.57	0.52
1:B:115:THR:HG22	1:B:117:LEU:HG	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:15:ASN:HB2	1:B:45:LEU:CD2	2.38	0.52
1:A:47:ASP:HB2	1:A:48:GLU:OE2	2.10	0.52
1:E:106:THR:HG23	1:E:135:LEU:HD12	1.92	0.52
1:E:101:MET:HE1	1:E:138:ALA:O	2.10	0.52
1:D:102:ILE:N	1:D:102:ILE:HD12	2.25	0.52
1:A:85:THR:CG2	1:A:184:ARG:HB3	2.39	0.51
1:A:141:ILE:HD12	1:A:176:ILE:HD11	1.92	0.51
1:A:136:ARG:NH1	2:A:283:HOH:O	2.43	0.51
1:A:13:VAL:CG2	1:A:45:LEU:HD22	2.41	0.51
1:A:156:MET:SD	1:B:65:ALA:HB2	2.50	0.51
1:C:19:ILE:HG23	1:C:43:LEU:HD11	1.93	0.51
1:D:32:GLY:HA2	1:E:161:PHE:O	2.11	0.51
1:C:102:ILE:HD12	1:C:102:ILE:N	2.25	0.51
1:B:110:GLN:NE2	1:B:121:VAL:HG21	2.26	0.51
1:D:47:ASP:OD2	1:D:51:ARG:HB2	2.10	0.51
1:A:139:LEU:HD12	1:A:182:LEU:HB3	1.93	0.51
1:A:84:HIS:HA	1:A:187:GLN:CD	2.30	0.50
1:D:25:LYS:HB2	1:D:38:SER:HB3	1.93	0.50
1:D:101:MET:HE1	1:D:139:LEU:HD23	1.93	0.50
1:E:84:HIS:CD2	1:E:187:GLN:HE22	2.29	0.50
1:A:13:VAL:HG22	1:A:45:LEU:HD22	1.94	0.50
1:D:24:VAL:HG12	1:D:41:THR:HB	1.94	0.50
1:D:118:SER:CB	1:E:57:LYS:NZ	2.75	0.50
1:D:101:MET:CE	1:D:139:LEU:HD23	2.42	0.50
1:B:102:ILE:HD12	1:B:102:ILE:N	2.26	0.50
1:B:109:ARG:HD3	1:B:120:LEU:CD2	2.42	0.50
1:B:36:ILE:HD11	1:B:64:LEU:HD21	1.93	0.50
1:B:101:MET:CE	1:B:139:LEU:HD23	2.41	0.50
1:A:85:THR:HG21	1:A:184:ARG:CD	2.27	0.49
1:A:48:GLU:CD	1:A:48:GLU:H	2.14	0.49
1:D:19:ILE:HG23	1:D:43:LEU:CD1	2.43	0.49
1:C:46:ILE:HG22	1:C:50:GLY:HA2	1.94	0.49
1:E:164:GLY:O	1:E:165:LYS:HB2	2.13	0.49
1:E:39:ARG:O	1:E:40:ASN:HB2	2.11	0.49
1:A:14:LYS:HB2	1:A:54:GLU:OE2	2.13	0.49
1:A:109:ARG:NH1	1:A:120:LEU:HD21	2.28	0.49
1:A:9:SER:HB3	1:A:83:PRO:HB3	1.94	0.48
1:E:140:LYS:HE2	1:E:148:ASP:OD1	2.14	0.48
1:C:43:LEU:HD12	1:C:43:LEU:C	2.34	0.48
1:D:41:THR:CG2	2:D:191:HOH:O	2.60	0.48
1:B:109:ARG:HD3	1:B:120:LEU:HD23	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:36:ILE:CG2	1:D:41:THR:HG21	2.44	0.48
1:B:164:GLY:O	1:B:165:LYS:HB2	2.13	0.48
1:E:101:MET:CE	1:E:139:LEU:HD23	2.44	0.48
1:B:140:LYS:NZ	1:B:148:ASP:OD1	2.47	0.48
1:C:109:ARG:NH1	1:C:118:SER:OG	2.46	0.48
1:B:109:ARG:CZ	1:B:120:LEU:HD21	2.43	0.47
1:A:4:MET:CE	1:A:6:ALA:HB3	2.44	0.47
1:B:98:PHE:HA	1:B:101:MET:HE1	1.96	0.47
1:B:8:GLU:N	2:B:235:HOH:O	2.46	0.47
1:E:157:PRO:HB2	2:E:229:HOH:O	2.14	0.47
1:D:136:ARG:HD3	2:D:249:HOH:O	2.15	0.47
1:B:90:THR:HB	1:B:182:LEU:HD11	1.97	0.47
1:A:134:ASP:C	2:A:229:HOH:O	2.52	0.47
1:D:7:ALA:C	1:D:9:SER:H	2.19	0.47
1:D:82:ASP:OD2	1:D:85:THR:HB	2.15	0.46
1:E:99:THR:N	1:E:101:MET:HE2	2.28	0.46
1:E:104:GLY:O	1:E:129:THR:HG23	2.16	0.46
1:E:36:ILE:HD11	1:E:60:TYR:HA	1.97	0.46
1:A:25:LYS:HB2	1:A:38:SER:HB3	1.96	0.46
1:D:43:LEU:C	1:D:43:LEU:HD12	2.35	0.46
1:B:36:ILE:HD12	1:B:58:VAL:HG11	1.98	0.46
1:B:170:LEU:C	1:B:170:LEU:HD12	2.36	0.46
1:D:29:ASN:HB2	1:D:180:ASP:OD1	2.16	0.46
1:A:19:ILE:HG23	1:A:43:LEU:CD1	2.45	0.46
1:E:61:GLY:HA2	1:E:89:ILE:HD12	1.97	0.46
1:B:19:ILE:HG23	1:B:43:LEU:CD1	2.45	0.46
1:D:76:GLU:OE1	1:E:157:PRO:HG2	2.16	0.46
1:D:102:ILE:HB	1:D:105:GLN:HB3	1.98	0.46
1:C:139:LEU:HD12	1:C:182:LEU:HB3	1.97	0.46
1:A:38:SER:OG	1:A:41:THR:OG1	2.34	0.46
1:E:61:GLY:HA2	1:E:89:ILE:CD1	2.47	0.45
1:A:145:GLN:HB2	1:A:147:ASN:ND2	2.31	0.45
1:E:13:VAL:CG2	1:E:45:LEU:HD22	2.46	0.45
1:B:98:PHE:HA	1:B:101:MET:HE3	1.99	0.45
1:A:163:PRO:HD3	2:B:236:HOH:O	2.16	0.45
1:E:98:PHE:HB2	1:E:172:ASP:OD2	2.16	0.45
1:D:95:PHE:HZ	1:E:187:GLN:HG3	1.81	0.45
1:A:90:THR:HG21	1:A:176:ILE:HD12	1.98	0.45
1:C:19:ILE:HD12	1:C:43:LEU:HD11	1.99	0.45
1:B:12:GLN:NE2	2:B:224:HOH:O	2.42	0.45
1:B:19:ILE:HD12	1:B:43:LEU:HD11	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:101:MET:C	1:E:102:ILE:HD12	2.36	0.45
1:A:36:ILE:CD1	1:A:58:VAL:HG11	2.47	0.45
1:A:175:GLN:NE2	2:A:213:HOH:O	2.48	0.45
1:D:82:ASP:CG	1:D:85:THR:HB	2.36	0.44
1:A:83:PRO:O	1:A:187:GLN:NE2	2.50	0.44
1:E:27:VAL:HG21	1:E:178:SER:CB	2.44	0.44
1:D:54:GLU:HB3	1:D:56:TYR:CE2	2.51	0.44
1:E:176:ILE:O	1:E:176:ILE:HD12	2.16	0.44
1:A:13:VAL:HG12	1:A:76:GLU:O	2.17	0.44
1:B:23:ASN:HA	1:C:117:LEU:HD23	1.99	0.44
1:D:8:GLU:HA	2:D:264:HOH:O	2.16	0.44
1:B:99:THR:H	1:B:101:MET:HE3	1.83	0.44
1:E:144:ALA:HB3	1:E:145:GLN:HE22	1.83	0.44
1:A:161:PHE:O	1:B:32:GLY:HA2	2.18	0.44
1:E:139:LEU:HD12	1:E:182:LEU:HB3	2.00	0.44
1:A:12:GLN:HG3	1:A:75:GLY:O	2.18	0.43
1:D:54:GLU:HA	1:D:54:GLU:OE1	2.18	0.43
1:C:152:PRO:HB3	2:C:281:HOH:O	2.19	0.43
1:E:112:ASP:HB2	1:E:115:THR:CB	2.49	0.43
1:D:114:LEU:CD2	1:D:114:LEU:H	2.25	0.43
1:A:170:LEU:HD13	1:A:176:ILE:CG2	2.47	0.43
1:B:170:LEU:HD12	1:B:170:LEU:O	2.18	0.42
1:A:140:LYS:HZ1	1:A:148:ASP:CG	2.23	0.42
1:E:103:ASP:C	1:E:105:GLN:H	2.22	0.42
1:D:112:ASP:O	1:D:114:LEU:HD23	2.19	0.42
1:D:47:ASP:HB3	1:D:53:LYS:HD2	1.99	0.42
1:E:110:GLN:NE2	1:E:121:VAL:HB	2.34	0.42
1:D:38:SER:OG	1:D:41:THR:HG22	2.20	0.42
1:A:189:SER:O	1:B:145:GLN:NE2	2.52	0.42
1:E:58:VAL:HA	1:E:59:PRO:HD3	1.76	0.42
1:E:90:THR:HB	1:E:182:LEU:HD11	2.01	0.42
1:C:80:ASN:ND2	1:C:81:TRP:H	2.17	0.42
1:D:95:PHE:CE2	1:D:175:GLN:HB2	2.54	0.42
1:A:92:VAL:HG23	1:A:141:ILE:HG21	2.01	0.42
1:A:66:LYS:HB3	1:A:70:GLU:CD	2.40	0.42
1:E:96:VAL:HG23	1:E:176:ILE:HG12	2.01	0.42
1:E:109:ARG:HD2	1:E:118:SER:OG	2.19	0.42
1:A:126:ALA:O	1:A:128:ARG:N	2.52	0.42
1:C:20:LYS:HA	1:C:20:LYS:HD2	1.75	0.41
1:E:13:VAL:CG2	1:E:45:LEU:HD13	2.46	0.41
1:A:4:MET:HE3	1:A:6:ALA:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:VAL:HA	1:A:59:PRO:HD3	1.87	0.41
1:E:110:GLN:O	1:E:118:SER:HA	2.21	0.41
1:D:138:ALA:CB	1:D:161:PHE:HA	2.50	0.41
1:B:175:GLN:NE2	2:B:216:HOH:O	2.47	0.41
1:D:95:PHE:CZ	1:E:187:GLN:HG3	2.55	0.41
1:C:19:ILE:HG23	1:C:43:LEU:HD13	2.01	0.41
1:A:43:LEU:C	1:A:43:LEU:HD12	2.41	0.41
1:D:101:MET:C	1:D:102:ILE:HD12	2.41	0.41
1:D:164:GLY:O	1:D:165:LYS:HB2	2.21	0.41
1:E:171:GLU:O	1:E:174:VAL:HG12	2.21	0.41
1:E:20:LYS:HD2	1:E:68:ASP:OD2	2.21	0.41
1:E:185:ILE:HA	1:E:186:PRO:HD2	1.89	0.41
1:C:127:GLU:CD	2:C:217:HOH:O	2.58	0.41
1:D:114:LEU:HD21	2:D:238:HOH:O	2.21	0.41
1:D:85:THR:HG23	1:D:86:MET:N	2.34	0.41
1:C:86:MET:HB2	1:C:187:GLN:HE21	1.82	0.41
1:E:20:LYS:NZ	2:E:241:HOH:O	2.54	0.41
1:A:46:ILE:CG2	1:A:50:GLY:HA2	2.50	0.41
1:B:128:ARG:HD2	2:B:277:HOH:O	2.21	0.41
1:D:33:LYS:HA	2:D:220:HOH:O	2.20	0.40
1:B:15:ASN:HA	2:B:244:HOH:O	2.21	0.40
1:A:59:PRO:HG2	1:A:81:TRP:CE2	2.57	0.40
1:C:36:ILE:HD11	1:C:60:TYR:HA	2.03	0.40
1:B:110:GLN:O	1:B:118:SER:HA	2.22	0.40
1:C:187:GLN:HB2	1:C:188:GLU:HG2	2.02	0.40
1:B:19:ILE:HG23	1:B:43:LEU:HD11	2.02	0.40

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	187/190 (98%)	179 (96%)	5 (3%)	3 (2%)	12	11
1	B	179/190 (94%)	173 (97%)	6 (3%)	0	100	100
1	C	181/190 (95%)	172 (95%)	8 (4%)	1 (1%)	30	36
1	D	181/190 (95%)	175 (97%)	5 (3%)	1 (1%)	30	36
1	E	186/190 (98%)	174 (94%)	11 (6%)	1 (0%)	34	41
All	All	914/950 (96%)	873 (96%)	35 (4%)	6 (1%)	26	31

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	188	GLU
1	A	127	GLU
1	D	113	GLU
1	E	187	GLN
1	A	152	PRO
1	A	126	ALA

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	154/156 (99%)	148 (96%)	6 (4%)	39	53
1	B	151/156 (97%)	145 (96%)	6 (4%)	38	52
1	C	151/156 (97%)	146 (97%)	5 (3%)	45	61
1	D	147/156 (94%)	139 (95%)	8 (5%)	27	36
1	E	150/156 (96%)	147 (98%)	3 (2%)	63	79
All	All	753/780 (96%)	725 (96%)	28 (4%)	41	55

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	SER
1	A	43	LEU

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Mol	Chain	Res	Type
1	A	48	GLU
1	A	80	ASN
1	A	85	THR
1	A	109	ARG
1	B	12	GLN
1	B	13	VAL
1	B	43	LEU
1	B	48	GLU
1	B	86	MET
1	B	156	MET
1	C	36	ILE
1	C	43	LEU
1	C	123	LEU
1	C	124	ASP
1	C	188	GLU
1	D	41	THR
1	D	43	LEU
1	D	57	LYS
1	D	85	THR
1	D	114	LEU
1	D	124	ASP
1	D	134	ASP
1	D	188	GLU
1	E	43	LEU
1	E	184	ARG
1	E	188	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	ASN
1	A	147	ASN
1	A	159	GLN
1	A	175	GLN
1	B	12	GLN
1	B	145	GLN
1	B	175	GLN
1	C	80	ASN
1	C	110	GLN
1	C	175	GLN
1	C	187	GLN
1	D	15	ASN

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Mol	Chain	Res	Type
1	E	15	ASN
1	E	80	ASN
1	E	84	HIS
1	E	110	GLN
1	E	145	GLN
1	E	159	GLN
1	E	169	GLN
1	E	175	GLN
1	E	187	GLN

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 carbohydrates 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	189/190 (99%)	0.27	12 (6%) 23 31	14, 31, 58, 73	0
1	B	181/190 (95%)	0.13	8 (4%) 38 47	12, 25, 50, 76	0
1	C	183/190 (96%)	0.38	8 (4%) 38 47	15, 35, 56, 73	0
1	D	183/190 (96%)	0.95	29 (15%) 3 4	18, 36, 80, 90	0
1	E	188/190 (98%)	0.86	30 (15%) 3 4	17, 39, 83, 95	0
All	All	924/950 (97%)	0.52	87 (9%) 11 16	12, 34, 71, 95	0

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	114	LEU	11.1
1	D	113	GLU	10.0
1	B	114	LEU	7.9
1	D	49	PHE	7.4
1	D	116	GLY	7.1
1	D	115	THR	7.0
1	E	130	ALA	7.0
1	E	113	GLU	6.8
1	E	115	THR	6.5
1	D	130	ALA	6.5
1	E	132	GLY	6.1
1	E	129	THR	5.5
1	A	1	GLY	5.4
1	C	7	ALA	5.3
1	E	117	LEU	5.1
1	E	127	GLU	5.1
1	B	115	THR	4.9
1	E	112	ASP	4.9
1	A	3	HIS	4.9
1	D	117	LEU	4.8

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Mol	Chain	Res	Type	RSRZ
1	D	129	THR	4.7
1	D	75	GLY	4.7
1	E	126	ALA	4.7
1	D	48	GLU	4.6
1	E	108	THR	4.4
1	E	131	GLY	4.3
1	C	114	LEU	4.3
1	A	5	ALA	4.3
1	E	189	SER	4.3
1	D	51	ARG	4.3
1	B	113	GLU	4.2
1	A	155	ASP	4.0
1	D	13	VAL	4.0
1	E	110	GLN	4.0
1	D	15	ASN	4.0
1	D	131	GLY	3.9
1	E	118	SER	3.7
1	D	14	LYS	3.7
1	D	134	ASP	3.7
1	C	153	GLY	3.6
1	D	50	GLY	3.6
1	C	189	SER	3.5
1	A	114	LEU	3.3
1	C	8	GLU	3.2
1	D	45	LEU	3.2
1	D	47	ASP	3.2
1	C	188	GLU	3.0
1	B	111	THR	2.9
1	D	74	GLY	2.9
1	E	111	THR	2.9
1	D	53	LYS	2.8
1	A	12	GLN	2.8
1	B	112	ASP	2.8
1	E	169	GLN	2.8
1	D	114	LEU	2.8
1	E	133	LYS	2.8
1	D	7	ALA	2.8
1	D	127	GLU	2.8
1	E	2	SER	2.7
1	E	102	ILE	2.6
1	B	188	GLU	2.6
1	E	145	GLN	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	19	ILE	2.6
1	E	116	GLY	2.5
1	A	113	GLU	2.4
1	C	130	ALA	2.4
1	C	145	GLN	2.4
1	E	109	ARG	2.4
1	D	132	GLY	2.4
1	E	187	GLN	2.4
1	A	126	ALA	2.4
1	E	123	LEU	2.4
1	D	16	LYS	2.3
1	E	3	HIS	2.3
1	E	84	HIS	2.3
1	B	8	GLU	2.2
1	E	188	GLU	2.2
1	B	110	GLN	2.2
1	A	153	GLY	2.2
1	D	128	ARG	2.2
1	E	128	ARG	2.2
1	D	112	ASP	2.1
1	A	4	MET	2.1
1	A	150	LEU	2.1
1	E	107	ILE	2.1
1	A	6	ALA	2.0
1	D	110	GLN	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.