



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 02:53 pm BST

PDB ID : 1QZH
Title : Crystal structure of Pot1 (protection of telomere)- ssDNA complex
Authors : Lei, M.; Podell, E.R.; Baumann, P.; Cech, T.R.
Deposited on : 2003-09-16
Resolution : 2.40 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

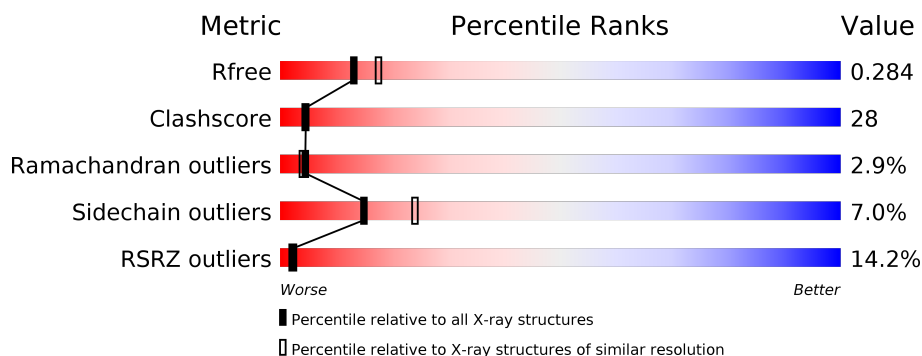
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (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 ≥ 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	G	6	<div> <div>33%</div> <div>67%</div> </div>
1	H	6	<div> <div>33%</div> <div>67%</div> </div>
1	I	6	<div> <div>100%</div> </div>
1	J	6	<div> <div>100%</div> </div>
1	K	6	<div> <div>100%</div> </div>
1	L	6	<div> <div>33%</div> <div>67%</div> </div>

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Mol	Chain	Length	Quality of chain
2	A	187	<div><div></div><div>11%</div><div>42%</div><div>44%</div><div>5%</div><div>9%</div></div>
2	B	187	<div><div></div><div>7%</div><div>54%</div><div>30%</div><div>6%</div><div>9%</div></div>
2	C	187	<div><div></div><div>11%</div><div>56%</div><div>30%</div><div>5%</div><div>9%</div></div>
2	D	187	<div><div></div><div>16%</div><div>53%</div><div>33%</div><div>5%</div><div>9%</div></div>
2	E	187	<div><div></div><div>18%</div><div>52%</div><div>35%</div><div></div><div>9%</div></div>
2	F	187	<div><div></div><div>18%</div><div>60%</div><div>26%</div><div></div><div>9%</div></div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 9033 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 DNA chain called telomeric single-stranded DNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	G	6	Total	C	N	O	P	0	0	0
			121	59	22	35	5			
1	H	6	Total	C	N	O	P	0	0	0
			121	59	22	35	5			
1	I	6	Total	C	N	O	P	0	0	0
			121	59	22	35	5			
1	J	6	Total	C	N	O	P	0	0	0
			121	59	22	35	5			
1	K	6	Total	C	N	O	P	0	0	0
			121	59	22	35	5			
1	L	6	Total	C	N	O	P	0	0	0
			121	59	22	35	5			

- Molecule 2 is a protein called Protection of telomeres protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	170	Total	C	N	O	S	0	0	0
			1365	868	234	259	4			
2	B	170	Total	C	N	O	S	0	0	0
			1365	868	234	259	4			
2	C	170	Total	C	N	O	S	0	0	0
			1351	859	229	259	4			
2	D	170	Total	C	N	O	S	0	0	0
			1351	859	229	259	4			
2	E	170	Total	C	N	O	S	0	0	0
			1351	859	229	259	4			
2	F	170	Total	C	N	O	S	0	0	0
			1351	859	229	259	4			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	CLONING ARTIFACT	UNP O13988

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Chain	Residue	Modelled	Actual	Comment	Reference
A	0	PRO	-	CLONING ARTIFACT	UNP O13988
A	1	GLY	MET	CLONING ARTIFACT	UNP O13988
B	-1	GLY	-	CLONING ARTIFACT	UNP O13988
B	0	PRO	-	CLONING ARTIFACT	UNP O13988
B	1	GLY	MET	CLONING ARTIFACT	UNP O13988
C	-1	GLY	-	CLONING ARTIFACT	UNP O13988
C	0	PRO	-	CLONING ARTIFACT	UNP O13988
C	1	GLY	MET	CLONING ARTIFACT	UNP O13988
D	-1	GLY	-	CLONING ARTIFACT	UNP O13988
D	0	PRO	-	CLONING ARTIFACT	UNP O13988
D	1	GLY	MET	CLONING ARTIFACT	UNP O13988
E	-1	GLY	-	CLONING ARTIFACT	UNP O13988
E	0	PRO	-	CLONING ARTIFACT	UNP O13988
E	1	GLY	MET	CLONING ARTIFACT	UNP O13988
F	-1	GLY	-	CLONING ARTIFACT	UNP O13988
F	0	PRO	-	CLONING ARTIFACT	UNP O13988
F	1	GLY	MET	CLONING ARTIFACT	UNP O13988

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	H	2	Total O 2 2	0	0
3	I	3	Total O 3 3	0	0
3	J	3	Total O 3 3	0	0
3	K	1	Total O 1 1	0	0
3	L	2	Total O 2 2	0	0
3	A	43	Total O 43 43	0	0
3	B	35	Total O 35 35	0	0
3	C	28	Total O 28 28	0	0
3	D	14	Total O 14 14	0	0
3	E	21	Total O 21 21	0	0
3	F	21	Total O 21 21	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 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: telomeric single-stranded DNA



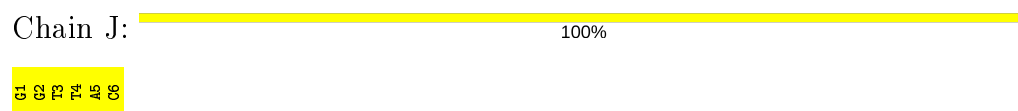
- Molecule 1: telomeric single-stranded DNA



- Molecule 1: telomeric single-stranded DNA



- Molecule 1: telomeric single-stranded DNA



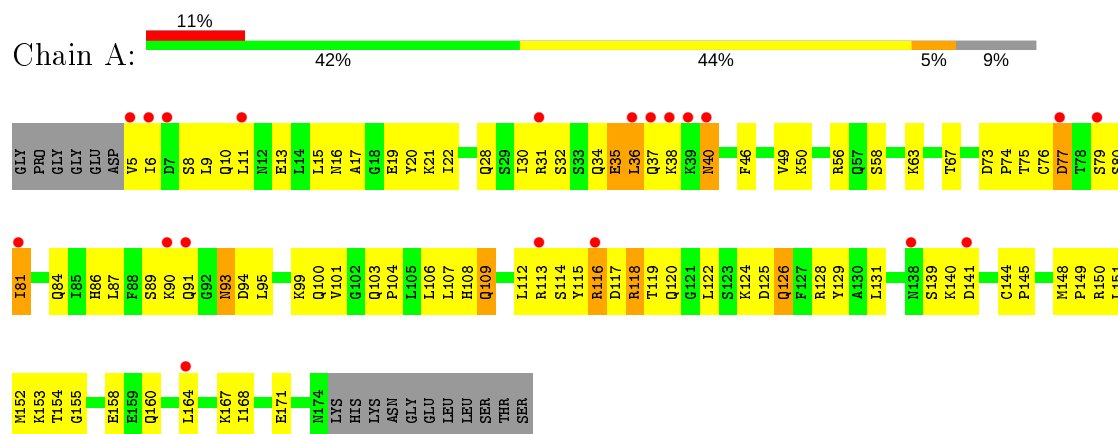
- Molecule 1: telomeric single-stranded DNA



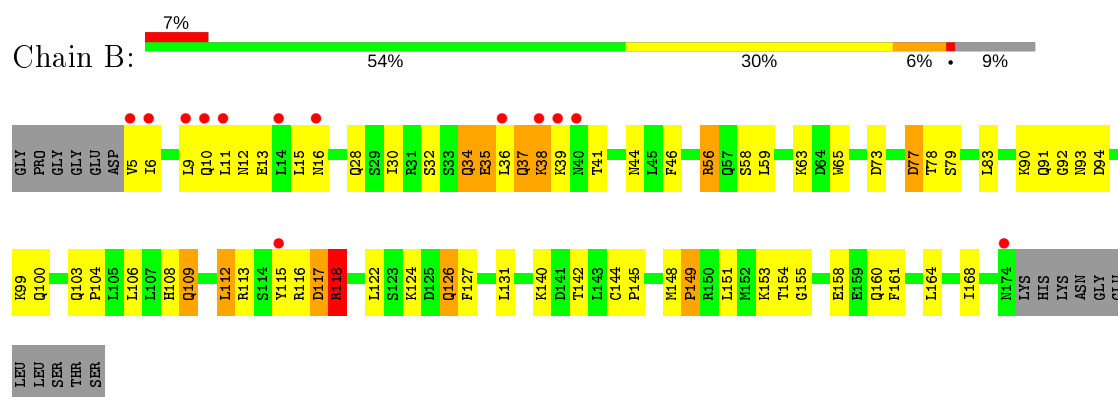
- Molecule 1: telomeric single-stranded DNA



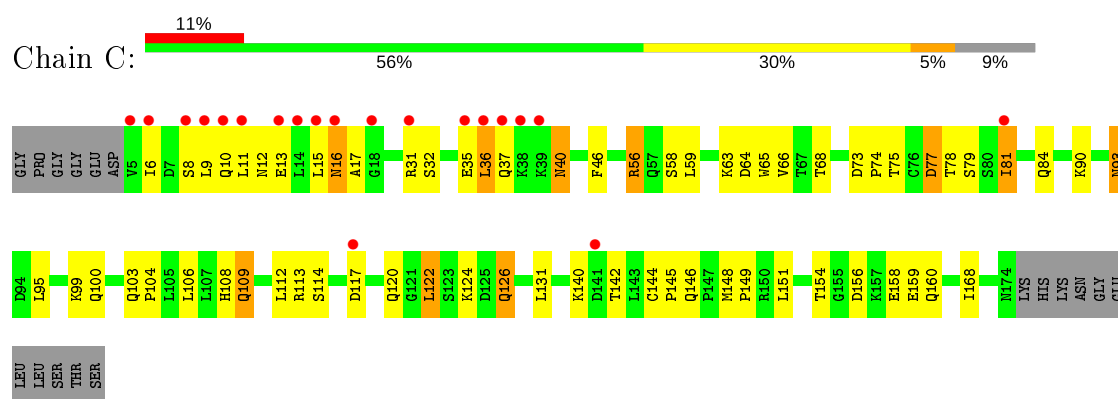
- Molecule 2: Protection of telomeres protein 1



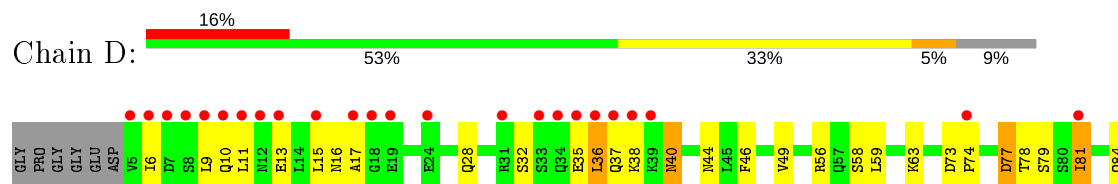
- Molecule 2: Protection of telomeres protein 1

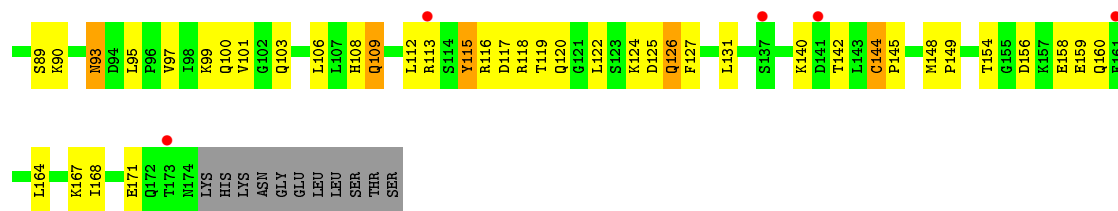


- Molecule 2: Protection of telomeres protein 1

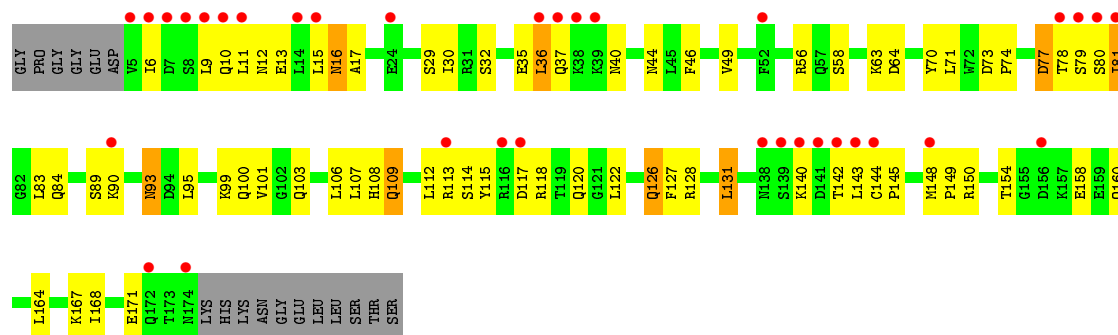


- Molecule 2: Protection of telomeres protein 1

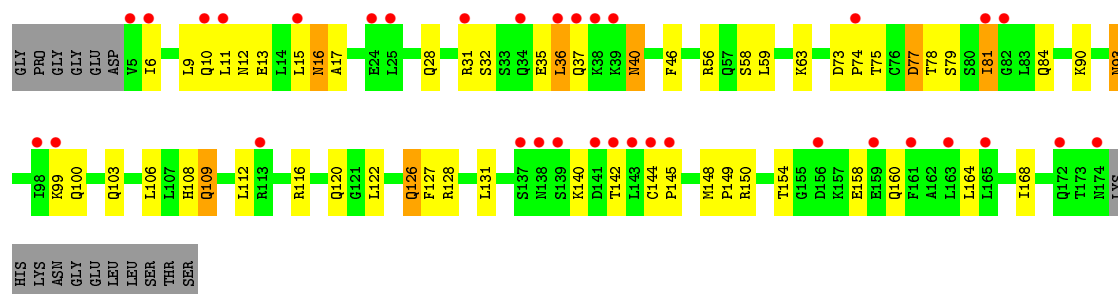




• Molecule 2: Protection of telomeres protein 1



• Molecule 2: Protection of telomeres protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	131.61 Å 76.33 Å 140.13 Å 90.00° 115.21° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 39.75 – 2.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.40) 76.9 (39.75-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.39 Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.249 , 0.283 0.257 , 0.284	Depositor DCC
R_{free} test set	2180 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	56.7	Xtriage
Anisotropy	0.429	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.056 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9033	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.86% 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	G	0.24	0/135	0.72	0/207
1	H	0.29	0/135	0.76	0/207
1	I	0.25	0/135	0.70	0/207
1	J	0.26	0/135	0.74	0/207
1	K	0.42	0/135	0.68	0/207
1	L	0.23	0/135	0.74	0/207
2	A	0.45	0/1394	0.68	0/1890
2	B	0.49	0/1394	0.75	0/1890
2	C	0.37	0/1380	0.62	0/1875
2	D	0.37	0/1380	0.69	2/1875 (0.1%)
2	E	0.35	0/1380	0.62	0/1875
2	F	0.34	0/1380	0.62	0/1875
All	All	0.39	0/9118	0.67	2/12522 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	115	TYR	CB-CG-CD1	6.37	124.82	121.00
2	D	115	TYR	CB-CG-CD2	-6.23	117.26	121.00

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	G	121	0	70	10	0
1	H	121	0	70	7	0
1	I	121	0	70	5	0
1	J	121	0	70	9	0
1	K	121	0	70	10	0
1	L	121	0	70	5	0
2	A	1365	0	1347	116	0
2	B	1365	0	1347	88	0
2	C	1351	0	1314	81	0
2	D	1351	0	1314	66	0
2	E	1351	0	1314	65	0
2	F	1351	0	1314	66	0
3	A	43	0	0	27	0
3	B	35	0	0	14	0
3	C	28	0	0	20	0
3	D	14	0	0	8	0
3	E	21	0	0	13	0
3	F	21	0	0	12	0
3	H	2	0	0	0	0
3	I	3	0	0	1	0
3	J	3	0	0	1	0
3	K	1	0	0	0	0
3	L	2	0	0	0	0
All	All	9033	0	8370	489	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 28.

All (489) 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)
2:B:115:TYR:HD1	2:B:118:ARG:NH2	1.41	1.19
2:A:28:GLN:HA	3:A:202:HOH:O	1.56	1.02
1:J:5:DA:H3'	1:J:6:DC:H5''	1.40	1.00
1:K:3:DT:H2''	1:K:4:DT:C5'	1.92	0.99
1:K:3:DT:H2''	1:K:4:DT:H5''	1.39	0.99
2:B:115:TYR:CD1	2:B:118:ARG:NH2	2.29	0.97
2:B:56:ARG:HG2	2:B:56:ARG:HH11	1.26	0.97
2:A:115:TYR:O	2:A:116:ARG:O	1.85	0.94
2:E:144:CYS:HA	3:E:200:HOH:O	1.67	0.94
2:C:81:ILE:H	2:C:81:ILE:HD12	1.33	0.93
2:F:81:ILE:H	2:F:81:ILE:HD12	1.33	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:9:LEU:HD12	3:B:186:HOH:O	1.70	0.91
2:E:81:ILE:H	2:E:81:ILE:HD12	1.37	0.90
2:B:30:ILE:HA	3:B:217:HOH:O	1.72	0.90
2:F:140:LYS:HB2	2:F:140:LYS:NZ	1.88	0.89
2:C:35:GLU:HB2	3:C:188:HOH:O	1.76	0.86
2:A:38:LYS:HD2	2:C:79:SER:CB	2.06	0.85
2:D:81:ILE:H	2:D:81:ILE:HD12	1.41	0.84
2:E:100:GLN:H	2:E:103:GLN:HE21	1.24	0.84
2:C:151:LEU:HG	3:C:212:HOH:O	1.77	0.83
2:A:31:ARG:HH21	2:C:114:SER:HB2	1.43	0.82
2:D:125:ASP:HB3	3:D:195:HOH:O	1.81	0.81
2:A:17:ALA:HA	3:A:217:HOH:O	1.78	0.81
2:F:99:LYS:H	2:F:103:GLN:HE22	1.29	0.81
2:B:15:LEU:HD21	3:B:214:HOH:O	1.81	0.81
2:C:8:SER:HB2	3:C:195:HOH:O	1.80	0.81
1:G:4:DT:H2''	1:G:5:DA:H5''	1.61	0.81
2:A:8:SER:HA	3:A:222:HOH:O	1.78	0.80
2:B:115:TYR:HD1	2:B:118:ARG:HH21	0.83	0.79
2:A:31:ARG:HD2	2:C:117:ASP:OD2	1.82	0.79
2:D:36:LEU:HD22	3:D:188:HOH:O	1.83	0.79
2:D:100:GLN:H	2:D:103:GLN:HE21	1.31	0.78
2:E:143:LEU:HD13	3:E:199:HOH:O	1.81	0.78
2:C:93:ASN:HB2	3:C:189:HOH:O	1.83	0.78
2:A:99:LYS:H	2:A:103:GLN:HE22	1.29	0.78
2:F:140:LYS:HB2	2:F:140:LYS:HZ2	1.48	0.77
2:A:31:ARG:HG3	2:C:36:LEU:HD13	1.65	0.77
2:B:99:LYS:H	2:B:103:GLN:HE22	1.32	0.77
2:C:6:ILE:HD12	3:C:196:HOH:O	1.85	0.77
2:E:99:LYS:H	2:E:103:GLN:HE22	1.32	0.77
1:K:3:DT:C2'	1:K:4:DT:H5''	2.15	0.77
2:B:34:GLN:O	2:B:35:GLU:O	2.04	0.76
3:I:149:HOH:O	2:C:66:VAL:HG21	1.85	0.76
1:L:5:DA:H3'	1:L:6:DC:H5''	1.67	0.75
2:B:65:TRP:CG	2:B:92:GLY:HA3	2.21	0.75
1:H:5:DA:H3'	1:H:6:DC:H5''	1.68	0.75
2:A:148:MET:HG2	3:A:205:HOH:O	1.87	0.75
2:A:81:ILE:HD12	2:A:81:ILE:H	1.50	0.75
1:K:5:DA:H3'	1:K:6:DC:H5''	1.69	0.74
2:A:21:LYS:NZ	2:B:108:HIS:HE1	1.86	0.74
2:A:19:GLU:HA	3:A:202:HOH:O	1.89	0.73
2:D:99:LYS:H	2:D:103:GLN:HE22	1.36	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:144:CYS:HA	3:F:192:HOH:O	1.89	0.73
2:A:36:LEU:HD21	2:A:118:ARG:C	2.09	0.72
2:B:77:ASP:C	2:B:79:SER:H	1.92	0.72
2:D:40:ASN:N	2:D:40:ASN:HD22	1.88	0.72
1:K:1:DG:H2"	1:K:2:DG:C8	2.25	0.72
2:A:155:GLY:HA2	2:B:153:LYS:HD3	1.72	0.71
2:E:131:LEU:HD22	3:E:199:HOH:O	1.88	0.71
2:A:31:ARG:HD2	2:C:117:ASP:CG	2.10	0.71
2:E:29:SER:HA	3:E:204:HOH:O	1.91	0.70
1:K:3:DT:H2"	1:K:4:DT:H5'	1.73	0.70
2:A:126:GLN:CD	2:A:126:GLN:H	1.95	0.69
2:C:37:GLN:HA	3:C:193:HOH:O	1.90	0.69
2:F:100:GLN:H	2:F:103:GLN:HE21	1.38	0.69
2:B:5:VAL:HA	3:B:211:HOH:O	1.91	0.69
1:I:5:DA:OP1	2:C:59:LEU:N	2.22	0.69
2:F:40:ASN:HD22	2:F:40:ASN:N	1.89	0.69
2:A:50:LYS:HG2	3:A:226:HOH:O	1.92	0.68
2:E:154:THR:HB	2:E:158:GLU:HG3	1.73	0.68
1:J:5:DA:C3'	1:J:6:DC:H5"	2.21	0.68
2:B:154:THR:HB	2:B:158:GLU:HG3	1.75	0.68
2:F:140:LYS:CB	2:F:140:LYS:NZ	2.57	0.68
2:A:100:GLN:H	2:A:103:GLN:HE21	1.41	0.68
2:B:12:ASN:HB2	3:B:186:HOH:O	1.94	0.68
2:D:115:TYR:O	2:D:116:ARG:C	2.31	0.67
2:A:20:TYR:HE1	2:A:22:ILE:CD1	2.06	0.67
2:D:140:LYS:NZ	2:D:140:LYS:HB2	2.10	0.67
2:C:154:THR:HB	2:C:158:GLU:HG3	1.76	0.67
1:L:5:DA:OP1	2:F:59:LEU:N	2.19	0.67
1:G:4:DT:C2'	1:G:5:DA:H5"	2.25	0.67
2:A:40:ASN:HD22	2:A:40:ASN:N	1.93	0.66
2:D:119:THR:HG23	3:D:194:HOH:O	1.94	0.66
2:C:11:LEU:HA	3:C:196:HOH:O	1.95	0.66
1:J:5:DA:OP1	2:D:59:LEU:N	2.27	0.66
2:E:117:ASP:HA	3:E:202:HOH:O	1.96	0.66
2:B:116:ARG:O	2:B:117:ASP:O	2.14	0.65
2:A:31:ARG:HH21	2:C:114:SER:CB	2.09	0.65
1:H:1:DG:N7	2:B:124:LYS:NZ	2.44	0.65
2:C:40:ASN:N	2:C:40:ASN:HD22	1.93	0.65
2:A:21:LYS:NZ	2:B:108:HIS:CE1	2.64	0.65
2:A:76:CYS:HA	3:A:206:HOH:O	1.96	0.65
2:B:37:GLN:NE2	2:B:112:LEU:HD23	2.10	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:126:GLN:H	2:E:126:GLN:CD	1.99	0.65
2:A:148:MET:N	2:A:149:PRO:HD3	2.11	0.65
2:C:9:LEU:O	2:C:13:GLU:HG3	1.97	0.65
2:B:148:MET:N	2:B:149:PRO:HD3	2.12	0.65
1:I:5:DA:H3'	1:I:6:DC:H5''	1.78	0.64
2:A:140:LYS:NZ	2:A:140:LYS:HB2	2.12	0.64
2:F:154:THR:HB	2:F:158:GLU:HG3	1.80	0.64
2:D:9:LEU:O	2:D:13:GLU:HG3	1.98	0.64
2:B:116:ARG:O	2:B:117:ASP:C	2.35	0.64
2:B:34:GLN:C	2:B:35:GLU:O	2.32	0.64
2:D:154:THR:HB	2:D:158:GLU:HG3	1.80	0.64
1:J:5:DA:H3'	1:J:6:DC:C5'	2.23	0.64
2:F:58:SER:HB2	2:F:63:LYS:HA	1.78	0.64
2:C:140:LYS:HB2	2:C:140:LYS:NZ	2.13	0.64
2:A:84:GLN:OE1	2:A:118:ARG:NH1	2.31	0.63
2:F:108:HIS:O	2:F:109:GLN:HB2	1.98	0.63
2:B:90:LYS:NZ	2:F:90:LYS:HB3	2.13	0.63
2:B:65:TRP:CD1	2:B:92:GLY:HA3	2.34	0.63
2:E:71:LEU:HD23	3:E:187:HOH:O	1.99	0.63
2:B:160:GLN:O	2:B:164:LEU:HD13	1.99	0.63
2:A:86:HIS:HD2	3:A:199:HOH:O	1.81	0.63
2:A:108:HIS:O	2:A:109:GLN:HB2	1.99	0.62
2:B:140:LYS:HB2	2:B:140:LYS:NZ	2.13	0.62
2:E:40:ASN:N	2:E:40:ASN:HD22	1.97	0.62
2:E:80:SER:HA	3:E:186:HOH:O	1.99	0.62
2:A:89:SER:HB2	2:A:95:LEU:HD21	1.80	0.62
2:C:99:LYS:H	2:C:103:GLN:HE22	1.45	0.62
2:B:77:ASP:O	2:B:79:SER:N	2.33	0.62
2:B:77:ASP:C	2:B:79:SER:N	2.53	0.62
2:C:160:GLN:HB3	3:C:210:HOH:O	1.99	0.62
2:E:11:LEU:HD21	2:E:74:PRO:HG2	1.82	0.61
2:E:9:LEU:O	2:E:13:GLU:HG3	2.00	0.61
2:F:126:GLN:CD	2:F:126:GLN:H	2.03	0.61
2:B:9:LEU:O	2:B:13:GLU:HG3	2.01	0.61
2:B:144:CYS:HB2	2:B:145:PRO:HD2	1.81	0.61
2:A:9:LEU:O	2:A:13:GLU:HG3	2.01	0.61
2:C:77:ASP:C	2:C:79:SER:H	2.04	0.61
2:E:100:GLN:H	2:E:103:GLN:NE2	1.96	0.60
2:B:148:MET:N	2:B:149:PRO:CD	2.63	0.60
2:F:9:LEU:O	2:F:13:GLU:HG3	2.01	0.60
2:B:56:ARG:CG	2:B:56:ARG:HH11	2.05	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:36:LEU:HD23	2:D:119:THR:OG1	2.01	0.60
2:A:120:GLN:CD	3:A:199:HOH:O	2.39	0.59
2:A:11:LEU:HD21	2:A:74:PRO:HG2	1.84	0.59
2:D:126:GLN:H	2:D:126:GLN:CD	2.05	0.59
2:A:31:ARG:HG3	2:C:36:LEU:CD1	2.32	0.59
2:B:90:LYS:HZ1	2:F:90:LYS:HB3	1.66	0.59
2:E:77:ASP:C	2:E:79:SER:H	2.05	0.59
2:A:148:MET:N	2:A:149:PRO:CD	2.64	0.59
2:C:11:LEU:HD21	2:C:74:PRO:HG2	1.84	0.59
2:B:9:LEU:HA	3:B:186:HOH:O	2.03	0.59
2:C:148:MET:HG2	3:C:208:HOH:O	2.02	0.59
1:H:5:DA:H5"	1:H:5:DA:H8	1.68	0.59
2:F:77:ASP:C	2:F:79:SER:H	2.05	0.58
1:J:3:DT:H2"	1:J:4:DT:O5'	2.03	0.58
2:A:109:GLN:HG3	3:A:219:HOH:O	2.03	0.58
2:A:139:SER:HB3	3:A:191:HOH:O	2.04	0.58
2:C:108:HIS:O	2:C:109:GLN:HB2	2.04	0.58
2:A:20:TYR:CE1	2:A:22:ILE:HD12	2.39	0.58
2:B:154:THR:HB	2:B:158:GLU:CG	2.33	0.58
2:C:146:GLN:HB3	3:C:212:HOH:O	2.03	0.58
2:F:10:GLN:HA	2:F:13:GLU:OE2	2.04	0.58
2:D:77:ASP:C	2:D:79:SER:H	2.07	0.58
2:A:114:SER:HB3	3:A:227:HOH:O	2.04	0.57
2:B:37:GLN:O	2:B:39:LYS:N	2.37	0.57
2:E:154:THR:HB	2:E:158:GLU:CG	2.33	0.57
2:F:35:GLU:HB2	3:F:187:HOH:O	2.05	0.57
2:B:37:GLN:OE1	2:B:41:THR:HG21	2.05	0.57
2:C:154:THR:HB	2:C:158:GLU:CG	2.34	0.57
2:B:56:ARG:NH1	2:B:56:ARG:HG2	2.06	0.56
2:D:36:LEU:HD21	2:D:119:THR:N	2.20	0.56
1:L:2:DG:C4	2:F:122:LEU:HD23	2.39	0.56
2:F:93:ASN:HD22	2:F:93:ASN:C	2.08	0.56
2:A:10:GLN:HA	2:A:13:GLU:OE2	2.05	0.56
2:C:37:GLN:HG3	3:C:206:HOH:O	2.05	0.56
2:F:116:ARG:HD3	3:F:188:HOH:O	2.04	0.56
2:C:126:GLN:H	2:C:126:GLN:CD	2.08	0.56
2:F:11:LEU:HD21	2:F:74:PRO:HG2	1.86	0.56
2:E:84:GLN:NE2	2:E:120:GLN:HE21	2.03	0.56
2:A:36:LEU:HD21	2:A:118:ARG:CA	2.36	0.56
2:B:32:SER:HA	2:B:35:GLU:HG3	1.87	0.56
2:C:100:GLN:H	2:C:103:GLN:HE21	1.51	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:37:GLN:C	2:B:39:LYS:N	2.57	0.56
2:F:93:ASN:ND2	3:F:193:HOH:O	2.38	0.56
2:E:144:CYS:HB2	2:E:145:PRO:HD2	1.87	0.56
2:B:126:GLN:H	2:B:126:GLN:CD	2.08	0.56
2:A:144:CYS:HB2	2:A:145:PRO:HD2	1.88	0.56
2:A:35:GLU:O	2:A:37:GLN:HG2	2.06	0.55
2:E:108:HIS:O	2:E:109:GLN:HB2	2.06	0.55
2:A:20:TYR:CE1	2:A:22:ILE:CD1	2.88	0.55
2:A:31:ARG:NH2	2:C:114:SER:HB2	2.16	0.55
2:D:10:GLN:HA	2:D:13:GLU:OE2	2.06	0.55
1:J:2:DG:C4	2:D:122:LEU:HD23	2.41	0.55
2:D:93:ASN:HD22	2:D:93:ASN:C	2.09	0.55
2:F:99:LYS:HG2	3:F:200:HOH:O	2.06	0.55
2:A:36:LEU:CD2	2:A:118:ARG:C	2.75	0.55
2:E:36:LEU:HD22	3:E:202:HOH:O	2.06	0.55
2:D:115:TYR:C	2:D:117:ASP:N	2.60	0.54
2:B:91:GLN:C	2:B:93:ASN:H	2.11	0.54
2:F:154:THR:HB	2:F:158:GLU:CG	2.38	0.54
2:E:100:GLN:O	2:E:103:GLN:HG3	2.07	0.54
2:D:108:HIS:O	2:D:109:GLN:HB2	2.07	0.54
2:A:77:ASP:C	2:A:79:SER:H	2.10	0.54
2:B:100:GLN:H	2:B:103:GLN:HE21	1.54	0.54
2:C:77:ASP:C	2:C:79:SER:N	2.62	0.54
2:B:37:GLN:O	2:B:39:LYS:O	2.26	0.53
2:D:44:ASN:OD1	2:D:108:HIS:HD2	1.91	0.53
2:B:11:LEU:HD13	2:B:168:ILE:HD12	1.89	0.53
2:E:77:ASP:C	2:E:79:SER:N	2.62	0.53
2:A:21:LYS:HZ3	2:B:108:HIS:CE1	2.26	0.53
2:E:115:TYR:HD1	3:E:191:HOH:O	1.90	0.53
1:H:2:DG:C4	2:B:122:LEU:HD23	2.43	0.53
1:I:3:DT:H2"	1:I:4:DT:O5'	2.09	0.53
2:A:32:SER:HA	2:A:35:GLU:HG3	1.91	0.53
2:B:12:ASN:O	2:B:16:ASN:ND2	2.42	0.53
2:A:140:LYS:HZ2	2:A:140:LYS:HB2	1.73	0.53
2:A:20:TYR:HE1	2:A:22:ILE:HD12	1.73	0.53
1:G:4:DT:H3'	1:G:5:DA:H5'	1.90	0.53
2:F:12:ASN:O	2:F:16:ASN:ND2	2.42	0.53
2:B:99:LYS:H	2:B:103:GLN:NE2	2.03	0.53
1:H:1:DG:H2"	1:H:2:DG:C8	2.43	0.53
2:D:113:ARG:N	3:D:194:HOH:O	2.42	0.53
2:F:109:GLN:HG2	3:F:204:HOH:O	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:144:CYS:HB2	2:C:145:PRO:HD2	1.90	0.53
2:C:10:GLN:HA	2:C:13:GLU:OE2	2.09	0.52
2:C:140:LYS:HB2	2:C:140:LYS:HZ2	1.74	0.52
2:F:32:SER:HB2	3:F:201:HOH:O	2.09	0.52
2:E:114:SER:HB3	3:E:202:HOH:O	2.08	0.52
2:F:100:GLN:NE2	2:F:142:THR:HG22	2.23	0.52
2:C:77:ASP:O	2:C:79:SER:N	2.42	0.52
2:B:34:GLN:O	2:B:35:GLU:C	2.48	0.52
2:C:56:ARG:NH2	3:C:191:HOH:O	2.43	0.52
2:B:90:LYS:HZ1	2:F:90:LYS:HD3	1.75	0.52
1:G:4:DT:C3'	1:G:5:DA:C5'	2.88	0.52
2:D:154:THR:HB	2:D:158:GLU:CG	2.40	0.52
2:D:77:ASP:C	2:D:79:SER:N	2.63	0.52
2:E:32:SER:HA	2:E:35:GLU:HG3	1.91	0.51
2:A:34:GLN:O	2:A:36:LEU:N	2.37	0.51
2:E:140:LYS:NZ	2:E:140:LYS:HB2	2.24	0.51
2:F:40:ASN:ND2	2:F:40:ASN:N	2.57	0.51
2:F:77:ASP:C	2:F:79:SER:N	2.63	0.51
2:A:115:TYR:O	2:A:116:ARG:C	2.46	0.51
2:E:89:SER:HB2	2:E:95:LEU:HD21	1.92	0.51
2:C:11:LEU:HD13	2:C:168:ILE:HD12	1.92	0.51
2:C:31:ARG:HB2	2:C:75:THR:OG1	2.10	0.51
1:I:1:DG:H2''	1:I:2:DG:C8	2.45	0.51
2:B:6:ILE:HA	2:B:10:GLN:OE1	2.10	0.51
2:D:115:TYR:O	2:D:117:ASP:N	2.44	0.51
2:F:160:GLN:O	2:F:164:LEU:HD13	2.11	0.51
2:A:149:PRO:HD2	3:A:189:HOH:O	2.11	0.51
2:D:140:LYS:HB2	2:D:140:LYS:HZ2	1.76	0.51
1:G:5:DA:C8	1:G:5:DA:H5'	2.46	0.51
2:D:97:VAL:CG1	2:F:145:PRO:HG3	2.41	0.50
2:F:32:SER:HA	2:F:35:GLU:HG3	1.92	0.50
1:I:1:DG:N7	2:C:124:LYS:NZ	2.52	0.50
2:A:73:ASP:OD1	2:A:73:ASP:C	2.49	0.50
2:E:10:GLN:HA	2:E:13:GLU:OE2	2.11	0.50
1:H:5:DA:H5''	1:H:5:DA:C8	2.47	0.50
2:D:11:LEU:HD21	2:D:74:PRO:HG2	1.94	0.50
2:C:100:GLN:NE2	2:C:142:THR:HG22	2.27	0.50
2:E:77:ASP:O	2:E:79:SER:N	2.45	0.50
2:A:126:GLN:NE2	3:A:219:HOH:O	2.45	0.50
2:B:161:PHE:HZ	3:B:214:HOH:O	1.93	0.50
2:C:106:LEU:HD23	2:C:106:LEU:C	2.32	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:160:GLN:O	2:A:164:LEU:HD13	2.12	0.50
2:A:31:ARG:HB2	2:A:75:THR:HB	1.94	0.50
2:D:100:GLN:H	2:D:103:GLN:NE2	2.04	0.49
2:E:93:ASN:HD22	2:E:93:ASN:C	2.14	0.49
2:B:30:ILE:HG12	2:B:73:ASP:OD2	2.11	0.49
1:J:5:DA:H5"	1:J:5:DA:H8	1.77	0.49
2:D:35:GLU:O	2:D:37:GLN:HG2	2.12	0.49
2:D:40:ASN:N	2:D:40:ASN:ND2	2.56	0.49
2:A:19:GLU:HG2	3:A:202:HOH:O	2.13	0.49
2:E:106:LEU:HD23	2:E:106:LEU:C	2.33	0.49
2:E:100:GLN:NE2	2:E:142:THR:HG22	2.28	0.49
2:A:100:GLN:H	2:A:103:GLN:NE2	2.07	0.49
2:A:125:ASP:HB3	3:A:214:HOH:O	2.12	0.49
2:F:140:LYS:CB	2:F:140:LYS:HZ3	2.24	0.49
2:F:77:ASP:O	2:F:79:SER:N	2.46	0.49
2:B:106:LEU:C	2:B:106:LEU:HD23	2.32	0.49
2:C:84:GLN:NE2	2:C:120:GLN:HE21	2.09	0.49
2:C:32:SER:HA	2:C:35:GLU:HG3	1.95	0.49
2:D:156:ASP:HA	2:D:159:GLU:HB3	1.95	0.49
2:A:6:ILE:HA	2:A:10:GLN:OE1	2.13	0.48
2:A:99:LYS:H	2:A:103:GLN:NE2	2.04	0.48
2:B:100:GLN:NE2	2:B:142:THR:HG22	2.28	0.48
2:A:141:ASP:HB2	3:A:209:HOH:O	2.13	0.48
2:A:154:THR:HB	2:A:158:GLU:CG	2.43	0.48
2:C:40:ASN:N	2:C:40:ASN:ND2	2.61	0.48
2:B:11:LEU:HD23	2:B:11:LEU:C	2.34	0.48
2:F:6:ILE:HA	2:F:10:GLN:OE1	2.13	0.48
2:F:100:GLN:H	2:F:103:GLN:NE2	2.06	0.48
2:D:167:LYS:O	2:D:171:GLU:HG3	2.13	0.48
1:G:2:DG:O6	2:A:124:LYS:HD2	2.13	0.48
2:A:38:LYS:HB3	3:C:207:HOH:O	2.13	0.48
2:C:158:GLU:H	2:C:158:GLU:CD	2.17	0.48
2:A:15:LEU:C	2:A:17:ALA:H	2.17	0.48
2:C:35:GLU:O	2:C:37:GLN:HG2	2.14	0.48
2:A:20:TYR:HE1	2:A:22:ILE:HD11	1.78	0.47
2:E:99:LYS:H	2:E:103:GLN:NE2	2.08	0.47
2:B:10:GLN:HA	2:B:13:GLU:OE2	2.14	0.47
2:C:113:ARG:HE	2:C:122:LEU:HD22	1.78	0.47
2:E:6:ILE:HA	2:E:10:GLN:OE1	2.14	0.47
2:D:148:MET:N	2:D:149:PRO:CD	2.77	0.47
2:E:148:MET:N	2:E:149:PRO:CD	2.77	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:40:ASN:ND2	2:E:40:ASN:N	2.62	0.47
2:A:77:ASP:C	2:A:79:SER:N	2.67	0.47
2:D:89:SER:HB2	2:D:95:LEU:HD21	1.96	0.47
2:A:21:LYS:HZ3	2:B:108:HIS:HE1	1.59	0.47
2:D:160:GLN:O	2:D:164:LEU:HD13	2.14	0.47
2:D:58:SER:HB2	2:D:63:LYS:HA	1.96	0.47
2:A:153:LYS:HD2	2:B:155:GLY:HA2	1.96	0.47
2:C:73:ASP:C	2:C:73:ASP:OD1	2.53	0.47
2:F:144:CYS:HB2	2:F:145:PRO:HD2	1.97	0.47
2:F:35:GLU:CB	3:F:187:HOH:O	2.60	0.47
1:L:5:DA:C3'	1:L:6:DC:H5''	2.41	0.47
2:A:6:ILE:HG13	2:A:164:LEU:HD21	1.96	0.47
2:D:77:ASP:O	2:D:79:SER:N	2.48	0.47
2:D:84:GLN:NE2	2:D:120:GLN:HE21	2.13	0.47
2:E:35:GLU:O	2:E:37:GLN:HG2	2.15	0.47
2:B:9:LEU:HD23	3:B:196:HOH:O	2.15	0.47
2:D:113:ARG:HB2	3:D:189:HOH:O	2.14	0.47
2:E:81:ILE:HD12	2:E:81:ILE:N	2.19	0.46
2:F:37:GLN:HG2	3:F:187:HOH:O	2.15	0.46
2:F:73:ASP:C	2:F:73:ASP:OD1	2.53	0.46
2:C:6:ILE:HA	2:C:10:GLN:OE1	2.15	0.46
2:E:73:ASP:OD1	2:E:73:ASP:C	2.54	0.46
2:F:37:GLN:CG	3:F:187:HOH:O	2.62	0.46
2:C:81:ILE:N	2:C:81:ILE:HD12	2.16	0.46
2:D:100:GLN:NE2	2:D:142:THR:HG22	2.29	0.46
2:A:6:ILE:HG12	3:A:204:HOH:O	2.14	0.46
2:D:15:LEU:C	2:D:17:ALA:H	2.19	0.46
2:A:128:ARG:HD3	2:A:150:ARG:HA	1.98	0.46
2:A:158:GLU:CD	2:A:158:GLU:H	2.19	0.46
2:E:126:GLN:N	2:E:126:GLN:CD	2.68	0.46
2:F:35:GLU:O	2:F:37:GLN:HG2	2.16	0.46
2:A:31:ARG:HA	2:C:117:ASP:OD1	2.15	0.46
2:C:12:ASN:O	2:C:16:ASN:ND2	2.48	0.46
2:C:11:LEU:HD13	2:C:168:ILE:CD1	2.45	0.46
2:C:93:ASN:C	2:C:93:ASN:HD22	2.17	0.46
2:B:77:ASP:HB2	3:B:207:HOH:O	2.15	0.46
2:D:32:SER:HA	2:D:35:GLU:HG3	1.96	0.46
2:F:128:ARG:HD3	2:F:150:ARG:HA	1.97	0.45
1:G:2:DG:C4	2:A:122:LEU:HD23	2.51	0.45
2:C:148:MET:N	2:C:149:PRO:CD	2.79	0.45
2:E:58:SER:HB2	2:E:63:LYS:HA	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:97:VAL:HG11	2:F:145:PRO:CG	2.45	0.45
1:J:1:DG:H5'	3:E:205:HOH:O	2.15	0.45
2:A:35:GLU:O	2:A:36:LEU:C	2.55	0.45
2:A:91:GLN:HG3	2:A:94:ASP:OD2	2.17	0.45
2:B:79:SER:CB	3:B:207:HOH:O	2.64	0.45
2:C:58:SER:HB2	2:C:63:LYS:HA	1.97	0.45
2:A:31:ARG:HD3	2:A:75:THR:O	2.16	0.45
2:E:160:GLN:O	2:E:164:LEU:HD13	2.17	0.45
2:A:58:SER:HB2	2:A:63:LYS:HA	1.97	0.45
2:D:144:CYS:HB2	2:D:145:PRO:HD2	1.98	0.45
2:E:70:TYR:HA	2:E:83:LEU:O	2.17	0.45
1:G:4:DT:C2'	1:G:5:DA:C5'	2.93	0.45
2:A:30:ILE:HG12	2:A:73:ASP:OD2	2.17	0.45
2:D:158:GLU:CD	2:D:158:GLU:H	2.20	0.45
2:F:106:LEU:C	2:F:106:LEU:HD23	2.37	0.45
2:B:94:ASP:OD1	2:B:94:ASP:C	2.55	0.45
2:C:113:ARG:HE	2:C:122:LEU:CD2	2.30	0.45
2:D:6:ILE:HA	2:D:10:GLN:OE1	2.16	0.45
2:E:15:LEU:HD22	3:E:204:HOH:O	2.16	0.45
2:F:35:GLU:O	2:F:36:LEU:C	2.55	0.45
2:A:40:ASN:ND2	2:A:40:ASN:N	2.61	0.45
2:D:35:GLU:O	2:D:36:LEU:C	2.55	0.45
2:A:115:TYR:N	3:A:227:HOH:O	2.49	0.44
2:A:106:LEU:HD22	2:A:152:MET:CE	2.47	0.44
2:B:108:HIS:O	2:B:109:GLN:HB2	2.18	0.44
1:H:5:DA:OP1	2:B:59:LEU:N	2.39	0.44
2:B:10:GLN:O	2:B:13:GLU:HB2	2.17	0.44
1:J:2:DG:O6	2:D:124:LYS:HD2	2.17	0.44
1:L:1:DG:H2''	1:L:2:DG:C8	2.52	0.44
2:A:106:LEU:C	2:A:106:LEU:HD23	2.38	0.44
2:C:11:LEU:CD2	2:C:74:PRO:HG2	2.48	0.44
2:E:158:GLU:H	2:E:158:GLU:CD	2.21	0.44
1:K:5:DA:H8	1:K:5:DA:H5''	1.82	0.44
2:C:37:GLN:NE2	3:C:193:HOH:O	2.50	0.44
2:E:113:ARG:HE	2:E:122:LEU:HD22	1.82	0.44
2:E:36:LEU:HD21	2:E:118:ARG:CA	2.47	0.44
2:F:158:GLU:H	2:F:158:GLU:CD	2.20	0.44
2:A:11:LEU:C	2:A:11:LEU:HD23	2.38	0.44
2:B:100:GLN:H	2:B:103:GLN:NE2	2.14	0.44
2:B:56:ARG:NH1	2:B:56:ARG:CG	2.70	0.44
2:B:65:TRP:CB	2:B:92:GLY:HA3	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:15:LEU:C	2:C:17:ALA:H	2.21	0.44
2:A:140:LYS:CB	2:A:140:LYS:NZ	2.80	0.44
2:F:99:LYS:H	2:F:103:GLN:NE2	2.06	0.44
2:A:109:GLN:CG	3:A:219:HOH:O	2.64	0.44
2:B:58:SER:HB2	2:B:63:LYS:HA	2.00	0.44
2:C:65:TRP:CE2	3:C:189:HOH:O	2.68	0.44
2:A:21:LYS:HZ1	2:B:108:HIS:CE1	2.36	0.43
2:D:36:LEU:HD21	2:D:118:ARG:C	2.39	0.43
2:D:140:LYS:NZ	2:D:140:LYS:CB	2.79	0.43
2:F:84:GLN:NE2	2:F:120:GLN:HE21	2.16	0.43
2:D:106:LEU:HD23	2:D:106:LEU:C	2.37	0.43
2:D:11:LEU:HD13	2:D:168:ILE:HD12	1.99	0.43
2:F:93:ASN:ND2	2:F:93:ASN:C	2.71	0.43
2:E:15:LEU:C	2:E:17:ALA:H	2.22	0.43
2:E:11:LEU:CD2	2:E:74:PRO:HG2	2.48	0.43
2:B:103:GLN:HA	2:B:104:PRO:HD3	1.87	0.43
2:B:140:LYS:CB	2:B:140:LYS:NZ	2.80	0.43
2:E:12:ASN:O	2:E:16:ASN:ND2	2.52	0.43
2:C:77:ASP:HB2	3:C:207:HOH:O	2.19	0.43
2:E:131:LEU:HD13	3:E:199:HOH:O	2.18	0.43
1:G:1:DG:H2"	1:G:2:DG:C8	2.54	0.43
2:A:106:LEU:HD23	2:A:107:LEU:N	2.34	0.43
2:A:129:TYR:CA	2:A:152:MET:SD	3.07	0.43
2:A:21:LYS:CE	2:B:108:HIS:HE1	2.31	0.43
2:C:65:TRP:CD1	3:C:189:HOH:O	2.70	0.43
2:D:11:LEU:HD13	2:D:168:ILE:CD1	2.49	0.43
2:A:11:LEU:HD13	2:A:168:ILE:HD12	2.01	0.42
2:B:113:ARG:HD3	3:B:210:HOH:O	2.19	0.42
2:D:28:GLN:NE2	3:D:192:HOH:O	2.50	0.42
2:F:100:GLN:NE2	3:F:200:HOH:O	2.40	0.42
2:F:99:LYS:N	2:F:103:GLN:HE22	2.08	0.42
2:A:80:SER:O	3:A:226:HOH:O	2.21	0.42
2:F:15:LEU:C	2:F:17:ALA:H	2.22	0.42
2:F:28:GLN:NE2	3:F:201:HOH:O	2.51	0.42
2:B:144:CYS:CB	2:B:145:PRO:HD2	2.48	0.42
1:K:2:DG:C4	2:E:122:LEU:HD23	2.54	0.42
2:E:93:ASN:HD22	2:E:95:LEU:H	1.67	0.42
2:B:28:GLN:HG2	3:B:208:HOH:O	2.19	0.42
2:C:103:GLN:HA	2:C:104:PRO:HD3	1.80	0.42
2:E:128:ARG:HD3	2:E:150:ARG:HA	2.00	0.42
2:C:35:GLU:O	2:C:36:LEU:C	2.58	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:68:THR:HB	3:C:191:HOH:O	2.18	0.42
2:C:93:ASN:N	3:C:189:HOH:O	2.51	0.42
2:E:167:LYS:O	2:E:171:GLU:HG3	2.19	0.42
2:C:108:HIS:O	2:C:109:GLN:CB	2.66	0.42
2:E:64:ASP:OD1	2:E:90:LYS:HG2	2.20	0.42
2:F:11:LEU:HD13	2:F:168:ILE:HD12	2.00	0.42
1:G:4:DT:C3'	1:G:5:DA:H5'	2.50	0.42
2:A:113:ARG:HD3	3:A:216:HOH:O	2.19	0.42
2:A:20:TYR:CE1	2:A:22:ILE:HD11	2.55	0.42
2:C:100:GLN:O	2:C:103:GLN:HG3	2.20	0.42
2:C:37:GLN:HG2	3:C:188:HOH:O	2.20	0.42
2:A:38:LYS:CD	2:C:79:SER:CB	2.89	0.42
2:A:155:GLY:CA	2:B:153:LYS:HD3	2.45	0.42
2:C:93:ASN:HD22	2:C:95:LEU:H	1.67	0.42
2:D:44:ASN:OD1	2:D:108:HIS:CD2	2.73	0.42
2:D:90:LYS:NZ	3:D:186:HOH:O	2.52	0.42
2:B:37:GLN:O	2:B:38:LYS:C	2.58	0.42
2:E:49:VAL:O	2:E:101:VAL:HG13	2.20	0.41
2:A:11:LEU:HD12	2:A:164:LEU:HB3	2.01	0.41
2:B:126:GLN:N	2:B:126:GLN:CD	2.73	0.41
2:C:64:ASP:OD1	2:C:90:LYS:HG2	2.20	0.41
2:B:37:GLN:C	2:B:39:LYS:H	2.21	0.41
3:J:153:HOH:O	2:D:90:LYS:HE2	2.20	0.41
2:A:108:HIS:O	2:A:109:GLN:CB	2.66	0.41
2:A:90:LYS:HE2	3:A:218:HOH:O	2.19	0.41
2:F:31:ARG:HB2	2:F:75:THR:OG1	2.21	0.41
2:E:44:ASN:OD1	2:E:108:HIS:HD2	2.03	0.41
1:K:5:DA:C3'	1:K:6:DC:H5''	2.47	0.41
2:D:38:LYS:HD3	3:D:196:HOH:O	2.19	0.41
2:A:37:GLN:HA	3:A:211:HOH:O	2.21	0.41
2:A:5:VAL:N	3:A:204:HOH:O	2.53	0.41
2:C:100:GLN:H	2:C:103:GLN:NE2	2.18	0.41
2:A:31:ARG:NH2	2:C:114:SER:CB	2.78	0.41
2:C:31:ARG:HB2	2:C:75:THR:CB	2.50	0.41
2:E:30:ILE:HG12	2:E:73:ASP:OD2	2.21	0.41
2:F:148:MET:N	2:F:149:PRO:CD	2.83	0.41
2:F:81:ILE:HD12	2:F:81:ILE:N	2.16	0.41
2:A:118:ARG:HG3	2:A:119:THR:N	2.36	0.41
2:D:11:LEU:HD12	2:D:164:LEU:HB3	2.03	0.41
2:A:67:THR:HG22	2:A:87:LEU:HB2	2.01	0.41
2:D:73:ASP:C	2:D:73:ASP:OD1	2.59	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:106:LEU:HD23	2:E:107:LEU:N	2.36	0.41
2:F:108:HIS:O	2:F:109:GLN:CB	2.66	0.41
2:A:154:THR:HB	2:A:158:GLU:HG3	2.01	0.41
2:A:167:LYS:O	2:A:171:GLU:HG3	2.21	0.41
2:A:90:LYS:HB2	3:A:210:HOH:O	2.21	0.41
2:B:11:LEU:HD12	2:B:164:LEU:HB3	2.03	0.41
2:D:49:VAL:O	2:D:101:VAL:HG13	2.20	0.41
2:E:11:LEU:HD13	2:E:168:ILE:HD12	2.01	0.41
2:A:103:GLN:HA	2:A:104:PRO:HD3	1.83	0.40
2:A:36:LEU:HD21	2:A:118:ARG:HA	2.03	0.40
2:A:91:GLN:HG2	3:A:210:HOH:O	2.20	0.40
2:D:97:VAL:HG12	2:F:145:PRO:HG3	2.02	0.40
2:B:83:LEU:HD22	3:B:217:HOH:O	2.21	0.40
2:C:140:LYS:CB	2:C:140:LYS:NZ	2.82	0.40
2:C:156:ASP:HA	2:C:159:GLU:HB3	2.02	0.40
2:D:11:LEU:HD23	2:D:11:LEU:C	2.41	0.40
2:F:11:LEU:CD2	2:F:74:PRO:HG2	2.50	0.40
2:A:49:VAL:O	2:A:101:VAL:HG13	2.21	0.40
2:A:93:ASN:C	2:A:93:ASN:HD22	2.23	0.40
2:B:44:ASN:OD1	2:B:108:HIS:HD2	2.04	0.40
2:E:35:GLU:O	2:E:36:LEU:C	2.59	0.40
2:B:160:GLN:NE2	3:B:211:HOH:O	2.54	0.40
2:B:32:SER:O	2:B:35:GLU:HG3	2.22	0.40
2:D:99:LYS:H	2:D:103:GLN:NE2	2.13	0.40
1:K:5:DA:H5"	1:K:5:DA:C8	2.56	0.40
2:E:36:LEU:HD21	2:E:118:ARG:C	2.41	0.40
2:F:81:ILE:CD1	2:F:81:ILE:H	2.11	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	
2	A	168/187 (90%)	145 (86%)	17 (10%)	6 (4%)	3	3
2	B	168/187 (90%)	147 (88%)	14 (8%)	7 (4%)	3	2
2	C	168/187 (90%)	147 (88%)	17 (10%)	4 (2%)	6	6
2	D	168/187 (90%)	150 (89%)	14 (8%)	4 (2%)	6	6
2	E	168/187 (90%)	148 (88%)	16 (10%)	4 (2%)	6	6
2	F	168/187 (90%)	148 (88%)	16 (10%)	4 (2%)	6	6
All	All	1008/1122 (90%)	885 (88%)	94 (9%)	29 (3%)	4	4

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	116	ARG
2	A	117	ASP
2	B	35	GLU
2	B	117	ASP
2	B	118	ARG
2	E	36	LEU
2	A	16	ASN
2	A	36	LEU
2	A	109	GLN
2	B	38	LYS
2	B	78	THR
2	C	36	LEU
2	C	109	GLN
2	D	36	LEU
2	D	109	GLN
2	F	36	LEU
2	B	109	GLN
2	C	78	THR
2	D	16	ASN
2	E	16	ASN
2	F	78	THR
2	F	109	GLN
2	C	16	ASN
2	E	78	THR
2	E	109	GLN
2	F	16	ASN
2	A	35	GLU
2	D	78	THR
2	B	149	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	
2	A	153/171 (90%)	142 (93%)	11 (7%)	14	23
2	B	153/171 (90%)	141 (92%)	12 (8%)	12	19
2	C	150/171 (88%)	140 (93%)	10 (7%)	16	26
2	D	150/171 (88%)	139 (93%)	11 (7%)	14	22
2	E	150/171 (88%)	141 (94%)	9 (6%)	19	31
2	F	150/171 (88%)	140 (93%)	10 (7%)	16	26
All	All	906/1026 (88%)	843 (93%)	63 (7%)	15	24

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	40	ASN
2	A	46	PHE
2	A	56	ARG
2	A	77	ASP
2	A	81	ILE
2	A	93	ASN
2	A	112	LEU
2	A	118	ARG
2	A	126	GLN
2	A	131	LEU
2	A	151	LEU
2	B	34	GLN
2	B	36	LEU
2	B	37	GLN
2	B	46	PHE
2	B	56	ARG
2	B	77	ASP
2	B	112	LEU
2	B	118	ARG
2	B	126	GLN
2	B	127	PHE
2	B	131	LEU

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Mol	Chain	Res	Type
2	B	151	LEU
2	C	40	ASN
2	C	46	PHE
2	C	56	ARG
2	C	77	ASP
2	C	81	ILE
2	C	93	ASN
2	C	112	LEU
2	C	122	LEU
2	C	126	GLN
2	C	131	LEU
2	D	40	ASN
2	D	46	PHE
2	D	56	ARG
2	D	77	ASP
2	D	81	ILE
2	D	93	ASN
2	D	112	LEU
2	D	126	GLN
2	D	127	PHE
2	D	131	LEU
2	D	144	CYS
2	E	46	PHE
2	E	56	ARG
2	E	77	ASP
2	E	81	ILE
2	E	93	ASN
2	E	112	LEU
2	E	126	GLN
2	E	127	PHE
2	E	131	LEU
2	F	40	ASN
2	F	46	PHE
2	F	56	ARG
2	F	77	ASP
2	F	81	ILE
2	F	93	ASN
2	F	112	LEU
2	F	126	GLN
2	F	127	PHE
2	F	131	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such

sidechains are listed below:

Mol	Chain	Res	Type
2	A	40	ASN
2	A	93	ASN
2	A	100	GLN
2	A	103	GLN
2	A	108	HIS
2	B	12	ASN
2	B	16	ASN
2	B	40	ASN
2	B	93	ASN
2	B	100	GLN
2	B	103	GLN
2	B	108	HIS
2	C	16	ASN
2	C	40	ASN
2	C	84	GLN
2	C	93	ASN
2	C	100	GLN
2	C	103	GLN
2	C	108	HIS
2	D	16	ASN
2	D	40	ASN
2	D	84	GLN
2	D	93	ASN
2	D	100	GLN
2	D	103	GLN
2	D	108	HIS
2	E	12	ASN
2	E	16	ASN
2	E	40	ASN
2	E	84	GLN
2	E	93	ASN
2	E	100	GLN
2	E	103	GLN
2	E	108	HIS
2	F	12	ASN
2	F	16	ASN
2	F	40	ASN
2	F	84	GLN
2	F	93	ASN
2	F	100	GLN
2	F	103	GLN
2	F	108	HIS

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	G	6/6 (100%)	0.44	0 100 100	49, 52, 68, 81	0
1	H	6/6 (100%)	0.52	0 100 100	52, 62, 64, 78	0
1	I	6/6 (100%)	0.39	0 100 100	49, 54, 66, 80	0
1	J	6/6 (100%)	0.20	0 100 100	55, 61, 66, 70	0
1	K	6/6 (100%)	0.27	0 100 100	56, 61, 73, 87	0
1	L	6/6 (100%)	0.32	0 100 100	57, 60, 73, 77	0
2	A	170/187 (90%)	0.55	20 (11%) 4 4	24, 55, 98, 117	0
2	B	170/187 (90%)	0.37	13 (7%) 13 12	24, 51, 95, 111	0
2	C	170/187 (90%)	0.92	20 (11%) 4 4	36, 67, 133, 147	0
2	D	170/187 (90%)	1.06	29 (17%) 1 1	36, 70, 127, 148	0
2	E	170/187 (90%)	1.23	34 (20%) 1 0	43, 69, 119, 144	0
2	F	170/187 (90%)	1.29	34 (20%) 1 0	40, 81, 122, 133	0
All	All	1056/1158 (91%)	0.89	150 (14%) 2 2	24, 65, 117, 148	0

All (150) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	142	THR	14.3
2	D	38	LYS	11.0
2	E	5	VAL	11.0
2	E	38	LYS	10.8
2	D	5	VAL	10.4
2	F	142	THR	10.2
2	F	5	VAL	9.9
2	C	5	VAL	9.9
2	B	5	VAL	9.0
2	D	9	LEU	8.9
2	C	9	LEU	8.7

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Mol	Chain	Res	Type	RSRZ
2	E	6	ILE	8.5
2	F	156	ASP	8.1
2	C	15	LEU	7.4
2	D	36	LEU	7.4
2	C	37	GLN	7.3
2	A	38	LYS	7.3
2	E	143	LEU	7.3
2	F	99	LYS	7.2
2	C	6	ILE	7.0
2	D	11	LEU	7.0
2	E	11	LEU	6.8
2	F	38	LYS	6.8
2	E	39	LYS	6.7
2	D	74	PRO	6.7
2	C	38	LYS	6.5
2	D	6	ILE	6.4
2	D	39	LYS	6.3
2	D	7	ASP	6.2
2	F	11	LEU	6.0
2	E	36	LEU	5.8
2	C	10	GLN	5.7
2	F	144	CYS	5.5
2	B	38	LYS	5.5
2	F	6	ILE	5.4
2	F	36	LEU	5.3
2	F	39	LYS	5.3
2	E	140	LYS	5.3
2	A	11	LEU	5.3
2	E	8	SER	5.3
2	C	13	GLU	5.2
2	F	143	LEU	5.0
2	D	15	LEU	5.0
2	E	138	ASN	5.0
2	F	81	ILE	5.0
2	D	8	SER	5.0
2	C	36	LEU	4.9
2	E	9	LEU	4.8
2	B	36	LEU	4.7
2	E	116	ARG	4.6
2	C	39	LYS	4.4
2	E	79	SER	4.3
2	E	10	GLN	4.3

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Mol	Chain	Res	Type	RSRZ
2	E	144	CYS	4.2
2	A	5	VAL	4.2
2	E	139	SER	4.1
2	F	34	GLN	4.1
2	F	141	ASP	4.1
2	C	11	LEU	4.1
2	B	174	ASN	4.1
2	C	81	ILE	4.1
2	C	117	ASP	4.0
2	F	138	ASN	4.0
2	E	81	ILE	3.9
2	A	90	LYS	3.9
2	F	24	GLU	3.9
2	F	165	LEU	3.8
2	C	16	ASN	3.8
2	D	31	ARG	3.8
2	D	18	GLY	3.8
2	A	39	LYS	3.7
2	A	7	ASP	3.6
2	A	31	ARG	3.6
2	E	15	LEU	3.6
2	E	14	LEU	3.5
2	F	137	SER	3.5
2	C	141	ASP	3.4
2	D	37	GLN	3.4
2	A	116	ARG	3.4
2	A	91	GLN	3.4
2	D	10	GLN	3.4
2	F	161	PHE	3.3
2	E	78	THR	3.3
2	D	13	GLU	3.3
2	F	37	GLN	3.3
2	C	18	GLY	3.2
2	E	80	SER	3.2
2	B	9	LEU	3.2
2	B	39	LYS	3.2
2	E	113	ARG	3.1
2	E	141	ASP	3.0
2	A	37	GLN	3.0
2	B	16	ASN	2.9
2	F	25	LEU	2.9
2	D	173	THR	2.9

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Mol	Chain	Res	Type	RSRZ
2	F	31	ARG	2.9
2	F	98	ILE	2.8
2	E	37	GLN	2.8
2	F	82	GLY	2.8
2	E	90	LYS	2.8
2	F	159	GLU	2.8
2	D	12	ASN	2.7
2	A	141	ASP	2.7
2	E	7	ASP	2.7
2	A	36	LEU	2.7
2	F	145	PRO	2.7
2	A	6	ILE	2.6
2	B	6	ILE	2.6
2	E	52	PHE	2.6
2	A	113	ARG	2.5
2	D	113	ARG	2.5
2	F	139	SER	2.5
2	E	148	MET	2.5
2	D	19	GLU	2.5
2	B	115	TYR	2.4
2	C	31	ARG	2.4
2	B	10	GLN	2.4
2	C	8	SER	2.4
2	B	14	LEU	2.4
2	D	33	SER	2.3
2	B	40	ASN	2.3
2	F	15	LEU	2.3
2	D	34	GLN	2.3
2	A	79	SER	2.3
2	B	11	LEU	2.3
2	F	174	ASN	2.3
2	A	81	ILE	2.3
2	D	17	ALA	2.3
2	D	24	GLU	2.3
2	F	172	GLN	2.2
2	A	164	LEU	2.2
2	D	81	ILE	2.2
2	F	74	PRO	2.2
2	D	35	GLU	2.2
2	D	141	ASP	2.2
2	A	138	ASN	2.2
2	F	163	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
2	A	77	ASP	2.2
2	E	117	ASP	2.2
2	E	156	ASP	2.2
2	F	113	ARG	2.2
2	E	172	GLN	2.1
2	E	174	ASN	2.1
2	C	14	LEU	2.0
2	C	35	GLU	2.0
2	E	24	GLU	2.0
2	F	10	GLN	2.0
2	A	40	ASN	2.0
2	D	137	SER	2.0
2	D	161	PHE	2.0

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 carbohydrates 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.