



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 27, 2016 – 03:50 PM EDT

PDB ID : 5ITU
Title : Crystal Structure of Human NEIL1(242K) bound to duplex DNA containing THF
Authors : Zhu, C.; Lu, L.; Zhang, J.; Yue, Z.; Song, J.; Zong, S.; Liu, M.; Stovicek, O.; Gao, Y.; Yi, C.
Deposited on : 2016-03-17
Resolution : 2.41 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	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20027939
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)	:	rb-20027939

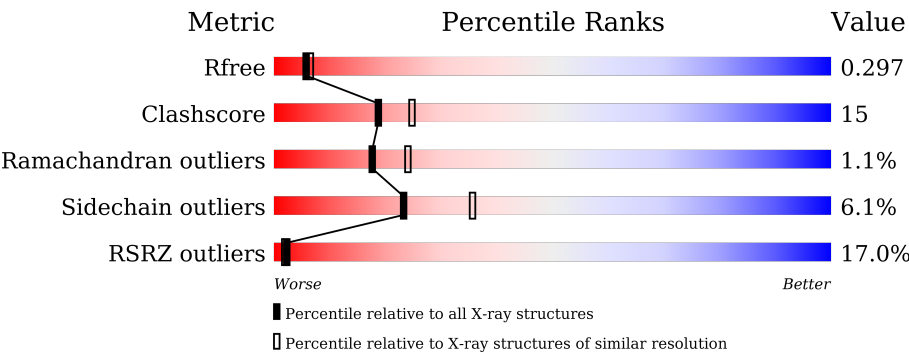
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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.



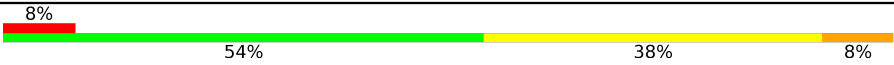


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3386 (2.44-2.40)
Clashscore	102246	3897 (2.44-2.40)
Ramachandran outliers	100387	3837 (2.44-2.40)
Sidechain outliers	100360	3838 (2.44-2.40)
RSRZ outliers	91569	3396 (2.44-2.40)

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. 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	400	<div><div></div><div><div></div><div>53%</div><div>11%</div><div>.</div><div>33%</div></div></div>
1	B	400	<div><div></div><div><div></div><div>50%</div><div>13%</div><div>.</div><div>36%</div></div></div>
1	C	400	<div><div>32%</div><div><div></div><div>49%</div><div>17%</div><div>.</div><div>32%</div></div></div>
2	D	13	<div><div>8%</div><div><div></div><div>46%</div><div>54%</div></div></div>
2	F	13	<div><div></div><div><div></div><div>38%</div><div>62%</div></div></div>
2	H	13	<div><div>15%</div><div><div></div><div>31%</div><div>69%</div></div></div>

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Mol	Chain	Length	Quality of chain
3	E	13	
3	G	13	
3	I	13	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8023 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 Endonuclease 8-like 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	0	1	0
			2120	1355	382	372	11			
1	B	255	Total	C	N	O	S	0	0	0
			2033	1302	371	350	10			
1	C	271	Total	C	N	O	S	0	0	0
			2143	1373	382	378	10			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	391	ALA	-	expression tag	UNP Q96FI4
A	392	ALA	-	expression tag	UNP Q96FI4
A	393	LEU	-	expression tag	UNP Q96FI4
A	394	GLY	-	expression tag	UNP Q96FI4
A	395	HIS	-	expression tag	UNP Q96FI4
A	396	HIS	-	expression tag	UNP Q96FI4
A	397	HIS	-	expression tag	UNP Q96FI4
A	398	HIS	-	expression tag	UNP Q96FI4
A	399	HIS	-	expression tag	UNP Q96FI4
A	400	HIS	-	expression tag	UNP Q96FI4
B	391	ALA	-	expression tag	UNP Q96FI4
B	392	ALA	-	expression tag	UNP Q96FI4
B	393	LEU	-	expression tag	UNP Q96FI4
B	394	GLY	-	expression tag	UNP Q96FI4
B	395	HIS	-	expression tag	UNP Q96FI4
B	396	HIS	-	expression tag	UNP Q96FI4
B	397	HIS	-	expression tag	UNP Q96FI4
B	398	HIS	-	expression tag	UNP Q96FI4
B	399	HIS	-	expression tag	UNP Q96FI4
B	400	HIS	-	expression tag	UNP Q96FI4
C	391	ALA	-	expression tag	UNP Q96FI4
C	392	ALA	-	expression tag	UNP Q96FI4
C	393	LEU	-	expression tag	UNP Q96FI4

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Chain	Residue	Modelled	Actual	Comment	Reference
C	394	GLY	-	expression tag	UNP Q96FI4
C	395	HIS	-	expression tag	UNP Q96FI4
C	396	HIS	-	expression tag	UNP Q96FI4
C	397	HIS	-	expression tag	UNP Q96FI4
C	398	HIS	-	expression tag	UNP Q96FI4
C	399	HIS	-	expression tag	UNP Q96FI4
C	400	HIS	-	expression tag	UNP Q96FI4

- Molecule 2 is a DNA chain called DNA (5'-D(*CP*GP*TP*CP*CP*AP*CP*GP*TP*CP*TP*AP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	13	Total	C	N	O	P	0	0	0
			249	120	41	76	12			
2	F	13	Total	C	N	O	P	0	0	0
			249	120	41	76	12			
2	H	13	Total	C	N	O	P	0	0	0
			249	120	41	76	12			

- Molecule 3 is a DNA chain called DNA (5'-D(*TP*AP*GP*AP*CP*CP*TP*GP*GP*AP*CP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	13	Total	C	N	O	P	0	0	0
			267	127	53	75	12			
3	G	13	Total	C	N	O	P	0	0	0
			267	127	53	75	12			
3	I	13	Total	C	N	O	P	0	0	0
			267	127	53	75	12			

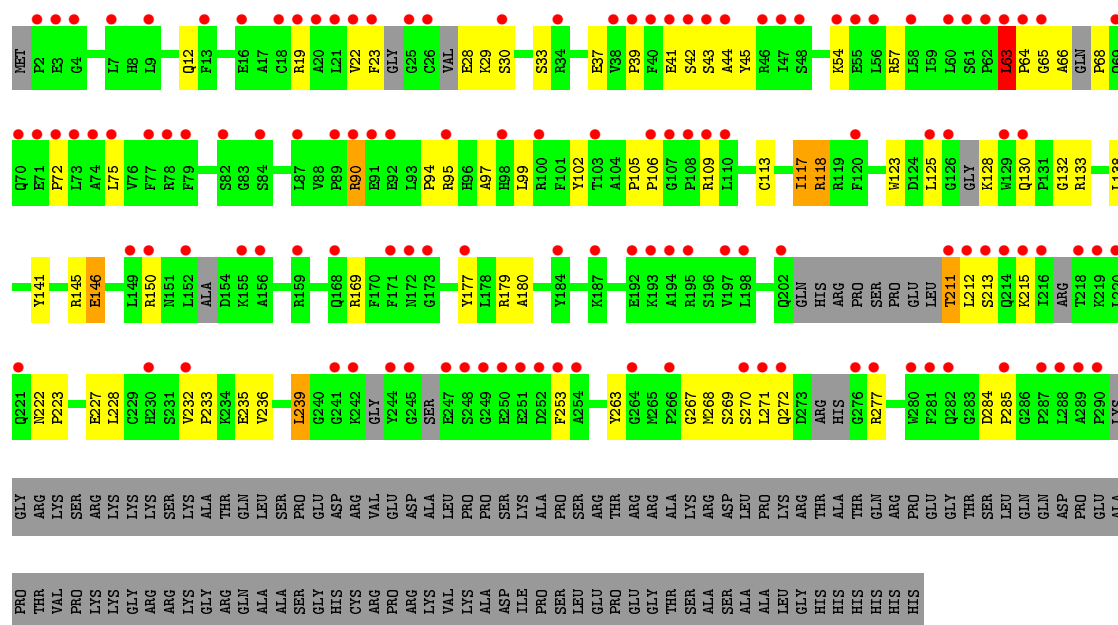
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	97	Total	O	0	0
			97	97		
4	B	58	Total	O	0	0
			58	58		
4	D	6	Total	O	0	0
			6	6		
4	E	8	Total	O	0	0
			8	8		

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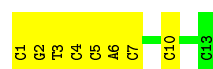
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	F	6	Total	O	0	0
			6	6		
4	G	4	Total	O	0	0
			4	4		



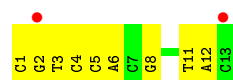
- Molecule 2: DNA (5'-D(*CP*GP*TP*CP*CP*AP*CP*GP*TP*CP*TP*AP*C)-3')



- Molecule 2: DNA (5'-D(*CP*GP*TP*CP*CP*AP*CP*GP*TP*CP*TP*AP*C)-3')



- Molecule 2: DNA (5'-D(*CP*GP*TP*CP*CP*AP*CP*GP*TP*CP*TP*AP*C)-3')



- Molecule 3: DNA (5'-D(*TP*AP*GP*AP*CP*CP*TP*GP*GP*AP*CP*GP*G)-3')

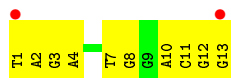


- Molecule 3: DNA (5'-D(*TP*AP*GP*AP*CP*CP*TP*GP*GP*AP*CP*GP*G)-3')





- Molecule 3: DNA (5'-D(*TP*AP*GP*AP*CP*CP*TP*GP*GP*AP*CP*GP*G)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	73.70Å 109.29Å 171.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	92.23 – 2.41 39.54 – 2.41	Depositor EDS
% Data completeness (in resolution range)	98.8 (92.23-2.41) 98.8 (39.54-2.41)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.89 (at 2.42Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.210 , 0.260 0.258 , 0.297	Depositor DCC
R_{free} test set	2722 reflections (5.08%)	DCC
Wilson B-factor (Å ²)	56.4	Xtriage
Anisotropy	0.347	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8023	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% 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

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	1.06	2/2181 (0.1%)	1.24	16/2949 (0.5%)
1	B	0.94	1/2089 (0.0%)	1.15	15/2826 (0.5%)
1	C	0.53	0/2192	0.77	3/2952 (0.1%)
2	D	0.60	0/277	0.90	0/424
2	F	0.65	0/277	0.99	0/424
2	H	0.43	0/277	0.85	0/424
3	E	0.50	0/300	0.93	1/462 (0.2%)
3	G	0.58	0/300	0.92	0/462
3	I	0.33	0/300	0.77	0/462
All	All	0.81	3/8193 (0.0%)	1.03	35/11385 (0.3%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	31	SER	CB-OG	-5.64	1.34	1.42
1	B	68	PRO	N-CD	5.14	1.55	1.47
1	A	45	TYR	CE2-CZ	5.04	1.45	1.38

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	133	ARG	NE-CZ-NH2	-18.24	111.18	120.30
1	B	133	ARG	NE-CZ-NH2	-16.44	112.08	120.30
1	A	133	ARG	NE-CZ-NH1	13.52	127.06	120.30
1	B	119	ARG	NE-CZ-NH2	-12.41	114.09	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	128	LYS	N-CA-C	11.17	141.15	111.00
1	C	128	LYS	N-CA-CB	-10.94	90.90	110.60
1	A	179	ARG	NE-CZ-NH2	-10.36	115.12	120.30
1	B	100	ARG	NE-CZ-NH2	-9.18	115.71	120.30
1	B	119	ARG	NE-CZ-NH1	8.51	124.56	120.30
1	A	179	ARG	NE-CZ-NH1	8.30	124.45	120.30
1	A	119	ARG	NE-CZ-NH2	-7.83	116.39	120.30
3	E	7	DT	O5'-P-OP2	-7.54	98.91	105.70
1	A	119	ARG	NE-CZ-NH1	7.01	123.80	120.30
1	B	57	ARG	NE-CZ-NH1	6.82	123.71	120.30
1	A	167	ASP	CB-CG-OD1	6.71	124.33	118.30
1	A	57	ARG	NE-CZ-NH1	6.66	123.63	120.30
1	A	67	GLN	C-N-CD	-6.62	106.03	120.60
1	A	277	ARG	NE-CZ-NH2	-6.59	117.01	120.30
1	A	90	ARG	NE-CZ-NH1	6.50	123.55	120.30
1	B	118	ARG	NE-CZ-NH1	-6.42	117.09	120.30
1	A	133	ARG	CG-CD-NE	-6.36	98.45	111.80
1	A	118	ARG	NE-CZ-NH2	-6.06	117.27	120.30
1	B	277	ARG	NE-CZ-NH1	6.04	123.32	120.30
1	B	277	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	A	57	ARG	NE-CZ-NH2	-5.94	117.33	120.30
1	B	133	ARG	CG-CD-NE	-5.86	99.49	111.80
1	C	63	LEU	C-N-CD	5.86	140.70	128.40
1	B	100	ARG	NE-CZ-NH1	5.67	123.14	120.30
1	B	133	ARG	NE-CZ-NH1	5.60	123.10	120.30
1	B	288	LEU	CA-CB-CG	5.54	128.05	115.30
1	A	273	ASP	CB-CG-OD2	-5.46	113.39	118.30
1	B	118	ARG	NE-CZ-NH2	5.46	123.03	120.30
1	A	150	ARG	NE-CZ-NH1	5.30	122.95	120.30
1	B	57	ARG	NE-CZ-NH2	-5.26	117.67	120.30
1	B	140	GLU	OE1-CD-OE2	-5.05	117.24	123.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	248	SER	Peptide
1	A	67	GLN	Mainchain,Peptide

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	2120	0	2091	46	1
1	B	2033	0	2026	39	1
1	C	2143	0	2109	75	0
2	D	249	0	143	6	0
2	F	249	0	143	14	0
2	H	249	0	143	27	0
3	E	267	0	147	19	0
3	G	267	0	147	12	0
3	I	267	0	147	44	0
4	A	97	0	0	4	1
4	B	58	0	0	1	0
4	D	6	0	0	0	0
4	E	8	0	0	0	0
4	F	6	0	0	0	0
4	G	4	0	0	0	0
All	All	8023	0	7096	229	2

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 15.

All (229) 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:C:118:ARG:NH2	2:H:6:DA:C2	1.83	1.45
1:C:118:ARG:NH2	2:H:6:DA:N1	1.74	1.35
1:C:117:ILE:CD1	3:I:8:DG:H5'	1.60	1.30
2:H:2:DG:H1	3:I:11:DC:N4	1.43	1.14
3:E:1:DT:H5''	3:I:13:DG:H2''	1.30	1.09
1:C:117:ILE:HD12	3:I:8:DG:H5'	1.31	1.06
3:E:3:DG:H2''	3:E:4:DA:H5'	1.38	1.04
2:H:2:DG:N1	3:I:11:DC:N4	2.02	1.04
3:E:1:DT:C6	3:I:13:DG:H2'	1.94	1.02
2:H:12:DA:N6	3:I:1:DT:O4	1.94	1.00
1:A:250:GLU:O	1:A:251:GLU:HG2	1.61	0.98
1:C:118:ARG:CZ	3:I:7:DT:O2	2.10	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:PHE:O	1:B:288:LEU:HD22	1.64	0.96
2:F:1:DC:H2''	2:F:2:DG:OP2	1.65	0.95
1:A:266:PRO:O	4:A:501:HOH:O	1.82	0.95
1:C:118:ARG:HH22	2:H:6:DA:H2	1.11	0.94
2:H:12:DA:N1	3:I:1:DT:N3	2.14	0.94
1:C:29:LYS:HG2	1:C:30:SER:H	1.30	0.93
3:I:12:DG:H2''	3:I:13:DG:H5''	1.49	0.93
1:C:117:ILE:HD11	3:I:8:DG:H5'	1.48	0.92
4:A:550:HOH:O	1:B:143:GLN:HG2	1.67	0.92
1:A:200:ALA:H	1:A:201:LEU:HD23	1.32	0.92
1:B:104:ALA:HB1	1:B:105:PRO:HD2	1.51	0.91
1:C:63:LEU:HD22	1:C:64:PRO:HD2	1.50	0.91
2:H:2:DG:C6	3:I:11:DC:N4	2.37	0.90
3:E:3:DG:H2''	3:E:4:DA:C5'	2.02	0.90
1:A:251:GLU:HG3	1:A:252:ASP:H	1.36	0.90
3:E:1:DT:C5'	3:I:13:DG:H2''	2.03	0.88
1:C:23:PHE:O	1:C:44:ALA:HA	1.74	0.87
1:C:117:ILE:CD1	3:I:8:DG:C5'	2.52	0.87
3:E:1:DT:H6	3:I:13:DG:H2'	1.40	0.86
1:C:118:ARG:NH2	2:H:6:DA:H2	1.58	0.85
1:C:41:GLU:O	1:C:42:SER:OG	1.94	0.84
1:B:78:ARG:NH1	2:F:10:DC:OP1	2.10	0.84
1:C:118:ARG:NH2	3:I:7:DT:O2	2.12	0.83
3:G:12:DG:H2''	3:G:13:DG:C8	2.14	0.83
3:I:12:DG:H2''	3:I:13:DG:C5'	2.08	0.83
2:H:2:DG:H1	3:I:11:DC:H42	0.90	0.82
1:A:249:GLY:O	1:A:250:GLU:HG2	1.80	0.82
1:B:118:ARG:HD2	3:G:6:DC:O2	1.80	0.81
1:A:250:GLU:HA	1:A:253:PHE:HB3	1.61	0.81
1:A:250:GLU:O	1:A:251:GLU:CG	2.29	0.80
1:C:118:ARG:CZ	2:H:6:DA:C2	2.64	0.80
1:A:274:ARG:CZ	1:A:275:HIS:HB2	2.11	0.80
1:B:191:PHE:O	1:B:288:LEU:CD2	2.31	0.79
3:I:11:DC:H2''	3:I:12:DG:C8	2.18	0.78
2:H:2:DG:O6	3:I:11:DC:N4	2.17	0.78
2:H:3:DT:O4	3:I:10:DA:N6	2.18	0.77
1:A:130:GLN:OE1	1:A:133:ARG:HD2	1.83	0.76
3:E:2:DA:H2''	3:E:3:DG:O5'	1.85	0.75
1:C:277:ARG:NH2	2:H:8:DG:H2'	2.02	0.74
1:C:29:LYS:HG2	1:C:30:SER:N	2.01	0.74
3:E:1:DT:H5''	3:I:13:DG:C2'	2.17	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:GLN:HG2	1:B:238:GLN:OE1	1.89	0.72
3:I:3:DG:H2'	3:I:4:DA:C8	2.25	0.72
3:I:10:DA:H1'	3:I:11:DC:H5'	1.72	0.71
1:A:250:GLU:HG3	1:A:251:GLU:H	1.55	0.70
1:C:28:GLU:O	1:C:99:LEU:HD12	1.90	0.70
1:C:146:GLU:O	1:C:150:ARG:HB2	1.92	0.70
2:F:2:DG:H1	3:G:11:DC:H42	1.40	0.70
3:I:1:DT:H2''	3:I:2:DA:C8	2.28	0.69
1:C:130:GLN:HE22	1:C:133:ARG:HG3	1.56	0.69
3:I:3:DG:H2''	3:I:4:DA:H5'	1.75	0.69
2:F:3:DT:H1'	2:F:4:DC:H5''	1.75	0.68
2:H:4:DC:H2'	2:H:5:DC:C6	2.28	0.68
2:F:2:DG:H2''	2:F:3:DT:H5'	1.76	0.67
3:G:1:DT:H4'	3:G:2:DA:O5'	1.94	0.67
1:C:277:ARG:HH22	2:H:8:DG:H2'	1.60	0.67
3:I:10:DA:H2''	3:I:11:DC:H5'	1.76	0.66
1:A:105:PRO:O	1:A:106:PRO:C	2.34	0.66
1:A:251:GLU:HG3	1:A:252:ASP:N	2.07	0.66
1:C:117:ILE:HD12	3:I:8:DG:C5'	2.17	0.66
3:E:3:DG:H2'	3:E:4:DA:C8	2.31	0.66
3:E:3:DG:H4'	3:E:4:DA:OP1	1.96	0.65
1:A:200:ALA:N	1:A:201:LEU:HD23	2.07	0.64
1:A:250:GLU:HG3	1:A:251:GLU:N	2.12	0.64
1:C:29:LYS:CG	1:C:30:SER:H	2.07	0.64
1:C:23:PHE:HB2	1:C:45:TYR:CZ	2.32	0.64
1:A:88:VAL:HG21	1:A:94:PRO:HD3	1.81	0.63
1:A:117:ILE:HD11	3:E:7:DT:H1'	1.81	0.62
1:C:118:ARG:NH2	3:I:7:DT:C2	2.67	0.62
1:C:105:PRO:HB2	1:C:106:PRO:HD3	1.80	0.62
3:G:2:DA:C8	3:G:2:DA:H5''	2.35	0.62
1:A:142:GLN:HG2	1:B:238:GLN:CD	2.19	0.62
1:C:130:GLN:OE1	1:C:133:ARG:HD2	1.99	0.62
3:E:1:DT:C7	3:I:13:DG:H3'	2.30	0.62
1:A:78:ARG:NH1	2:D:10:DC:OP1	2.30	0.62
1:B:160:PRO:HG3	1:B:288:LEU:HD23	1.82	0.61
3:I:10:DA:H1'	3:I:11:DC:C5'	2.30	0.61
1:A:274:ARG:NH1	1:A:275:HIS:HB2	2.14	0.61
1:B:65:GLY:O	1:B:67:GLN:NE2	2.32	0.60
1:B:46:ARG:HH11	1:B:46:ARG:HG2	1.67	0.60
1:C:23:PHE:HB2	1:C:45:TYR:CE1	2.37	0.59
3:I:10:DA:C2'	3:I:11:DC:H5'	2.32	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:3:DG:H2''	3:I:4:DA:C5'	2.31	0.59
1:C:118:ARG:NH1	3:I:7:DT:O2	2.35	0.59
3:I:10:DA:C1'	3:I:11:DC:H5'	2.31	0.59
1:C:212:LEU:HA	1:C:215:LYS:HB2	1.85	0.59
3:I:2:DA:H2''	3:I:3:DG:C8	2.37	0.59
1:C:90:ARG:HD3	1:C:102:TYR:CE1	2.38	0.58
3:G:11:DC:C2'	3:G:12:DG:O5'	2.51	0.58
1:B:130:GLN:OE1	1:B:133:ARG:HD2	2.04	0.57
1:C:130:GLN:NE2	1:C:133:ARG:CD	2.67	0.57
1:C:277:ARG:NH2	2:H:8:DG:H8	2.02	0.57
2:D:11:DT:H4'	2:D:12:DA:OP1	2.05	0.57
1:C:41:GLU:C	1:C:42:SER:HG	2.01	0.57
1:C:45:TYR:CD1	1:C:45:TYR:N	2.73	0.56
1:C:65:GLY:O	1:C:66:ALA:HB3	2.06	0.56
1:A:159:ARG:HB3	1:A:160:PRO:HD2	1.87	0.56
1:B:88:VAL:HB	1:B:89:PRO:HD2	1.88	0.56
2:H:11:DT:H2''	2:H:12:DA:C8	2.41	0.56
1:C:130:GLN:NE2	1:C:133:ARG:HG3	2.21	0.56
1:C:118:ARG:CZ	3:I:7:DT:C2	2.89	0.55
1:B:190:PRO:HG3	1:B:281:PHE:CG	2.41	0.55
1:B:46:ARG:NE	1:B:63:LEU:HD21	2.22	0.54
1:C:37:GLU:O	1:C:39:PRO:HD3	2.07	0.54
1:C:44:ALA:C	1:C:45:TYR:HD1	2.10	0.54
1:A:117:ILE:HD13	3:E:8:DG:H5'	1.90	0.54
3:I:12:DG:H2''	3:I:13:DG:O5'	2.06	0.54
1:C:130:GLN:CD	1:C:133:ARG:HD2	2.28	0.54
3:G:12:DG:C2'	3:G:13:DG:C8	2.89	0.54
1:C:118:ARG:CZ	2:H:6:DA:H2	2.12	0.53
3:E:1:DT:H72	3:I:13:DG:H3'	1.90	0.53
1:C:130:GLN:NE2	1:C:133:ARG:HD2	2.24	0.53
1:A:117:ILE:HD12	3:E:7:DT:H4'	1.90	0.52
1:A:201:LEU:N	1:A:201:LEU:CD2	2.72	0.52
1:B:159:ARG:HB3	1:B:160:PRO:HD2	1.92	0.52
1:B:67:GLN:HA	1:B:68:PRO:C	2.30	0.52
2:H:2:DG:H2''	2:H:3:DT:OP1	2.09	0.52
2:F:1:DC:C2'	2:F:2:DG:OP2	2.50	0.51
1:C:94:PRO:HD2	1:C:97:ALA:HB2	1.93	0.51
3:G:11:DC:H2'	3:G:12:DG:C8	2.46	0.51
1:A:8:HIS:HE1	4:A:567:HOH:O	1.93	0.51
1:B:21:LEU:HD12	1:B:21:LEU:N	2.26	0.51
2:F:2:DG:H1	3:G:11:DC:N4	2.06	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:ALA:O	1:A:201:LEU:HB2	2.12	0.50
1:C:211:THR:O	1:C:215:LYS:HB2	2.11	0.50
1:B:67:GLN:HA	1:B:68:PRO:O	2.12	0.50
1:A:200:ALA:H	1:A:201:LEU:CD2	2.14	0.50
1:A:56:LEU:C	1:A:56:LEU:HD23	2.31	0.50
1:B:56:LEU:HD23	1:B:56:LEU:C	2.31	0.50
1:C:236:VAL:HA	1:C:239:LEU:HD23	1.91	0.50
1:C:23:PHE:CB	1:C:45:TYR:CZ	2.96	0.49
2:D:11:DT:H1'	2:D:12:DA:C8	2.47	0.49
1:B:118:ARG:CD	3:G:6:DC:O2	2.56	0.49
1:C:22:VAL:CG1	1:C:44:ALA:HB1	2.43	0.48
2:H:12:DA:N1	3:I:1:DT:C4	2.79	0.48
1:A:105:PRO:O	1:A:107:GLY:N	2.45	0.48
3:G:9:DG:H2''	3:G:10:DA:OP2	2.13	0.48
1:C:37:GLU:O	1:C:125:LEU:HD21	2.14	0.48
2:F:5:DC:H2''	2:F:6:DA:C8	2.48	0.48
1:C:99:LEU:HD22	1:C:123:TRP:CE2	2.49	0.48
1:B:35:ASN:HB2	1:B:123:TRP:CZ2	2.49	0.48
3:I:3:DG:H2'	3:I:4:DA:H8	1.72	0.48
1:A:117:ILE:HD12	3:E:7:DT:C4'	2.44	0.48
1:C:235:GLU:O	1:C:239:LEU:HB3	2.14	0.47
1:A:81:MET:CE	2:D:6:DA:H2'	2.44	0.47
2:H:11:DT:H2''	2:H:12:DA:H8	1.80	0.47
1:B:189:PRO:HA	1:B:190:PRO:HD2	1.68	0.47
2:D:9:DT:H2''	2:D:10:DC:H5''	1.96	0.46
2:F:3:DT:H2''	2:F:4:DC:H5''	1.96	0.46
1:C:141:TYR:OH	1:C:145:ARG:NH1	2.46	0.46
1:A:35:ASN:HB3	1:A:123:TRP:CE2	2.50	0.46
1:C:44:ALA:O	1:C:45:TYR:HB3	2.16	0.46
1:C:284:ASP:HA	1:C:285:PRO:HD3	1.80	0.46
1:A:4:GLY:N	1:A:5:PRO:CD	2.78	0.46
1:C:130:GLN:HE22	1:C:133:ARG:CG	2.25	0.45
1:C:177:TYR:HB3	1:C:263:TYR:CD2	2.51	0.45
2:H:12:DA:C6	3:I:1:DT:O4	2.68	0.45
1:A:169:ARG:NE	4:A:508:HOH:O	2.48	0.45
1:B:70:GLN:NE2	4:B:503:HOH:O	2.45	0.45
1:C:132:GLY:O	1:C:169:ARG:HG2	2.16	0.45
2:H:3:DT:H2'	2:H:3:DT:O2	2.16	0.45
2:H:1:DC:H2'	2:H:2:DG:C8	2.51	0.45
1:A:117:ILE:CD1	3:E:7:DT:H1'	2.45	0.45
1:A:249:GLY:O	1:A:250:GLU:CG	2.60	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:130:GLN:HB3	1:B:133:ARG:CD	2.45	0.45
2:F:3:DT:C2'	2:F:4:DC:H5''	2.46	0.45
2:H:4:DC:H2'	2:H:5:DC:C5	2.52	0.45
3:E:3:DG:H2''	3:E:4:DA:O5'	2.13	0.45
1:B:101:PHE:O	1:B:111:ALA:HA	2.16	0.44
1:C:99:LEU:HD22	1:C:123:TRP:CZ2	2.52	0.44
1:C:44:ALA:C	1:C:45:TYR:CD1	2.90	0.44
1:C:54:LYS:HB2	1:C:133:ARG:NH2	2.31	0.44
1:A:93:LEU:HD13	1:A:100:ARG:CD	2.48	0.44
2:D:2:DG:H2'	2:D:3:DT:H72	1.99	0.44
3:E:1:DT:C5'	3:I:13:DG:C2'	2.87	0.44
1:C:23:PHE:CB	1:C:45:TYR:CE1	3.00	0.44
1:C:177:TYR:HB3	1:C:263:TYR:CE2	2.53	0.44
1:C:179:ARG:HG3	1:C:180:ALA:N	2.33	0.44
1:B:20:ALA:C	1:B:21:LEU:HD12	2.38	0.44
1:C:268:MET:HA	1:C:268:MET:CE	2.47	0.44
1:A:222:ASN:OD1	1:A:222:ASN:N	2.51	0.43
1:B:3:GLU:OE2	1:B:177:TYR:CD2	2.71	0.43
1:C:130:GLN:NE2	1:C:133:ARG:CG	2.80	0.43
1:B:21:LEU:CD1	1:B:21:LEU:N	2.81	0.43
1:C:57:ARG:HA	1:C:75:LEU:O	2.18	0.43
1:B:232:VAL:HB	1:B:233:PRO:CD	2.48	0.43
1:A:116:ASP:OD1	1:A:116:ASP:C	2.56	0.43
1:A:230:HIS:O	1:A:234:LYS:HB2	2.19	0.43
1:A:250:GLU:CG	1:A:251:GLU:H	2.23	0.43
1:B:17:ALA:HB1	1:B:87:LEU:HD22	2.01	0.42
1:C:99:LEU:HB2	1:C:123:TRP:CH2	2.55	0.42
1:B:95:ARG:HB3	1:B:95:ARG:HE	1.31	0.42
1:A:191:PHE:HB2	1:A:285:PRO:O	2.20	0.42
1:C:118:ARG:NH2	3:I:7:DT:H3	2.18	0.42
1:B:26:CYS:SG	1:B:41:GLU:HG3	2.60	0.42
1:A:199:GLU:O	1:A:200:ALA:HB2	2.20	0.41
1:A:35:ASN:HB2	1:A:123:TRP:CZ2	2.55	0.41
2:F:4:DC:H2''	2:F:5:DC:O4'	2.21	0.41
1:A:200:ALA:C	1:A:201:LEU:HD23	2.41	0.41
1:C:223:PRO:HA	1:C:227:GLU:OE1	2.21	0.41
1:B:163:GLU:OE2	1:B:274:ARG:HB2	2.21	0.41
1:C:228:LEU:O	1:C:232:VAL:HB	2.20	0.41
2:F:3:DT:H2''	2:F:4:DC:C5'	2.51	0.41
1:B:197:VAL:HG11	1:B:225:LEU:HD22	2.02	0.41
1:B:225:LEU:HD12	1:B:225:LEU:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:232:VAL:HB	1:C:233:PRO:HD3	2.03	0.41
1:C:267:GLY:O	1:C:268:MET:HE2	2.21	0.41
1:C:29:LYS:NZ	1:C:33:SER:O	2.54	0.41
2:F:2:DG:H2''	2:F:3:DT:C5'	2.48	0.41
1:C:270:SER:O	1:C:271:LEU:HD23	2.21	0.41
3:G:6:DC:H2'	3:G:7:DT:C6	2.56	0.41
1:B:154:ASP:O	1:B:155:LYS:C	2.59	0.41
2:H:3:DT:C4	2:H:4:DC:C4	3.09	0.40
1:A:67:GLN:C	1:A:68:PRO:O	2.60	0.40
1:B:46:ARG:HH11	1:B:46:ARG:CG	2.29	0.40
1:B:2:PRO:N	2:F:7:DC:H2''	2.36	0.40
1:C:179:ARG:HG3	1:C:180:ALA:H	1.87	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:501:HOH:O	4:A:503:HOH:O[3_554]	1.99	0.21
1:A:284:ASP:OD2	1:B:95:ARG:NH2[3_454]	2.09	0.11

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	263/400 (66%)	249 (95%)	8 (3%)	6 (2%)	8	8
1	B	249/400 (62%)	232 (93%)	16 (6%)	1 (0%)	39	54
1	C	249/400 (62%)	217 (87%)	31 (12%)	1 (0%)	39	54
All	All	761/1200 (63%)	698 (92%)	55 (7%)	8 (1%)	17	24

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	68	PRO
1	A	105	PRO
1	A	200	ALA
1	A	250	GLU
1	A	251	GLU
1	A	106	PRO
1	C	72	PRO
1	B	106	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	222/335 (66%)	211 (95%)	11 (5%)	30	47
1	B	214/335 (64%)	205 (96%)	9 (4%)	36	55
1	C	225/335 (67%)	205 (91%)	20 (9%)	12	17
All	All	661/1005 (66%)	621 (94%)	40 (6%)	23	35

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	SER
1	A	70	GLN
1	A	105	PRO
1	A	117	ILE
1	A	124	ASP
1	A	155	LYS
1	A	201	LEU
1	A	222	ASN
1	A	234	LYS
1	A	257	ARG
1	A	274	ARG
1	B	90	ARG
1	B	95	ARG
1	B	100	ARG
1	B	144	PHE

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Mol	Chain	Res	Type
1	B	193	LYS
1	B	238	GLN
1	B	257	ARG
1	B	265	MET
1	B	288	LEU
1	C	12	GLN
1	C	19	ARG
1	C	43	SER
1	C	63	LEU
1	C	68	PRO
1	C	90	ARG
1	C	95	ARG
1	C	109	ARG
1	C	113	CYS
1	C	117	ILE
1	C	118	ARG
1	C	138	LEU
1	C	146	GLU
1	C	211	THR
1	C	213	SER
1	C	222	ASN
1	C	239	LEU
1	C	253	PHE
1	C	269	SER
1	C	272	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	139	GLN
1	B	12	GLN
1	B	275	HIS
1	C	151	ASN

5.3.3 RNA ⓘ

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	268/400 (67%)	0.28	3 (1%) 82 82	2, 9, 34, 60	0
1	B	255/400 (63%)	0.53	11 (4%) 39 39	4, 17, 41, 52	0
1	C	271/400 (67%)	2.16	127 (46%) 0 0	38, 62, 111, 169	0
2	D	13/13 (100%)	0.08	1 (7%) 16 15	11, 30, 65, 67	0
2	F	13/13 (100%)	0.13	0 100 100	17, 33, 46, 59	0
2	H	13/13 (100%)	0.78	2 (15%) 3 3	27, 39, 59, 62	0
3	E	13/13 (100%)	-0.26	1 (7%) 16 15	16, 31, 48, 50	0
3	G	13/13 (100%)	0.33	1 (7%) 16 15	12, 33, 44, 54	0
3	I	13/13 (100%)	0.78	2 (15%) 3 3	33, 48, 70, 70	0
All	All	872/1278 (68%)	0.94	148 (16%) 2 2	2, 25, 84, 169	0

All (148) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	245	GLY	11.8
1	C	211	THR	9.5
1	C	271	LEU	7.9
1	C	71	GLU	7.4
1	C	60	LEU	7.0
1	C	125	LEU	6.8
1	C	220	LEU	6.8
1	C	23	PHE	6.7
1	C	22	VAL	6.6
1	C	156	ALA	5.8
1	C	155	LYS	5.8
1	C	26	CYS	5.7
1	C	202	GLN	5.6
1	C	110	LEU	5.5
1	C	73	LEU	5.4

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Mol	Chain	Res	Type	RSRZ
1	C	25	GLY	5.4
1	C	92	GLU	5.3
1	C	18	CYS	5.2
1	C	149	LEU	5.2
1	C	40	PHE	5.2
1	C	44	ALA	5.0
1	C	290	PRO	5.0
1	C	242	LYS	5.0
1	C	272	GLN	4.8
1	C	75	LEU	4.8
1	C	19	ARG	4.8
1	C	109	ARG	4.8
1	C	152	LEU	4.5
1	C	48	SER	4.5
1	C	47	ILE	4.5
1	C	65	GLY	4.4
1	C	70	GLN	4.4
1	C	216	ILE	4.3
1	C	108	PRO	4.3
1	C	214	GLN	4.2
1	C	250	GLU	4.2
1	C	248	SER	4.1
1	C	46	ARG	4.1
1	C	87	LEU	4.1
1	C	159	ARG	4.0
1	C	126	GLY	3.9
1	C	289	ALA	3.9
1	C	13	PHE	3.9
1	C	90	ARG	3.9
1	C	91	GLU	3.9
1	C	107	GLY	3.8
1	C	2	PRO	3.8
1	C	38	VAL	3.8
1	C	282	GLN	3.8
1	C	64	PRO	3.8
1	C	4	GLY	3.8
1	C	247	GLU	3.7
1	C	61	SER	3.6
1	C	129	TRP	3.6
1	C	244	TYR	3.6
1	C	89	PRO	3.5
1	C	78	ARG	3.5

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Mol	Chain	Res	Type	RSRZ
1	C	215	LYS	3.4
1	C	100	ARG	3.4
1	C	288	LEU	3.4
1	C	55	GLU	3.4
1	C	63	LEU	3.4
1	C	221	GLN	3.3
1	C	276	GLY	3.3
1	C	218	THR	3.3
2	H	13	DC	3.3
3	I	1	DT	3.2
1	C	21	LEU	3.2
1	C	213	SER	3.2
1	C	270	SER	3.2
1	C	280	TRP	3.1
1	C	193	LYS	3.1
1	C	69	GLN	3.1
1	C	277	ARG	3.1
1	C	150	ARG	3.0
1	C	198	LEU	2.9
1	C	281	PHE	2.9
1	C	287	PRO	2.9
1	C	219	LYS	2.9
1	C	79	PHE	2.8
1	C	39	PRO	2.8
1	C	173	GLY	2.7
1	C	177	TYR	2.7
3	I	13	DG	2.7
2	H	2	DG	2.7
1	C	82	SER	2.7
1	B	266	PRO	2.7
1	C	54	LYS	2.7
1	C	56	LEU	2.7
1	C	187	LYS	2.7
1	C	120	PHE	2.6
1	C	95	ARG	2.6
1	C	62	PRO	2.6
1	C	251	GLU	2.6
1	C	172	ASN	2.6
1	B	79	PHE	2.5
1	C	197	VAL	2.5
1	C	30	SER	2.5
1	C	253	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	7	LEU	2.4
1	C	58	LEU	2.4
1	C	41	GLU	2.4
1	B	121	GLY	2.4
1	C	249	GLY	2.4
1	C	230	HIS	2.4
1	C	43	SER	2.4
1	B	173	GLY	2.4
1	C	184	TYR	2.4
1	C	72	PRO	2.4
1	C	266	PRO	2.3
1	C	74	ALA	2.3
1	C	241	GLY	2.3
1	C	106	PRO	2.3
1	B	283	GLY	2.3
1	C	252	ASP	2.3
1	C	254	ALA	2.2
1	C	264	GLY	2.2
1	C	232	VAL	2.2
1	B	51	ALA	2.2
3	G	13	DG	2.2
1	C	195	ARG	2.2
1	C	20	ALA	2.2
1	C	16	GLU	2.2
1	C	77	PHE	2.2
1	C	42	SER	2.2
1	C	168	GLN	2.2
1	A	249	GLY	2.2
1	A	290	PRO	2.2
1	B	63	LEU	2.2
1	A	233	PRO	2.2
1	B	65	GLY	2.1
1	C	9	LEU	2.1
1	B	255	ALA	2.1
1	C	98	HIS	2.1
1	C	171	PHE	2.1
2	D	13	DC	2.1
1	B	233	PRO	2.1
1	C	84	SER	2.1
1	C	212	LEU	2.1
1	C	3	GLU	2.1
1	B	137	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	34	ARG	2.1
1	C	192	GLU	2.0
3	E	1	DT	2.0
1	C	130	GLN	2.0
1	C	194	ALA	2.0
1	C	103	THR	2.0
1	C	285	PRO	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.