



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 08:25 pm GMT

PDB ID : 3WC1
Title : Crystal structure of *C. albicans* tRNA(His) guanylyltransferase (Thg1) with a G-1 deleted tRNA(His)
Authors : Nakamura, A.; Nemoto, T.; Sonoda, T.; Yamashita, K.; Tanaka, I.; Yao, M.
Deposited on : 2013-05-24
Resolution : 4.18 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

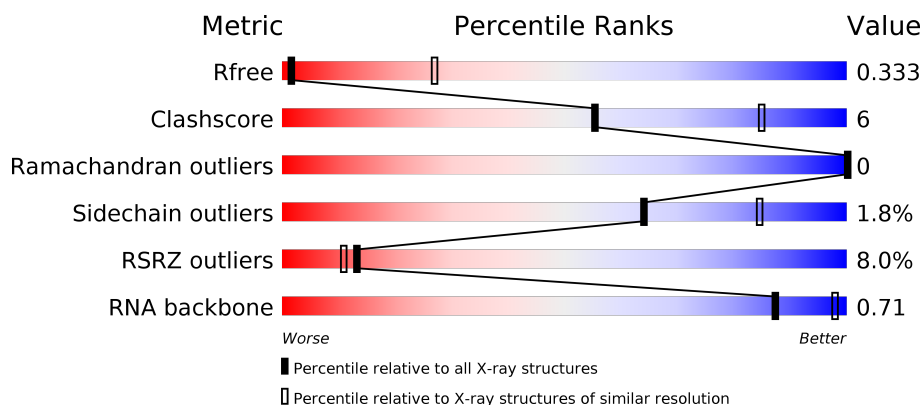
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.18 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}	100719	1167 (4.76-3.60)
Clashscore	112137	1024 (4.70-3.66)
Ramachandran outliers	110173	1022 (4.72-3.62)
Sidechain outliers	110143	1007 (4.72-3.62)
RSRZ outliers	101464	1179 (4.76-3.60)
RNA backbone	2435	1026 (5.04-2.90)

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	271	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>13%</div> <div>12%</div> </div> </div>
1	B	271	<div> <div>7%</div> <div> <div></div> <div>74%</div> <div>13%</div> <div>12%</div> </div> </div>
1	C	271	<div> <div>3%</div> <div> <div></div> <div>68%</div> <div>20%</div> <div>12%</div> </div> </div>
1	D	271	<div> <div></div> <div> <div></div> <div>71%</div> <div>16%</div> <div>12%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
2	P	75	<div><div><div>9%</div><div>72%</div><div>23%</div><div></div><div></div></div></div>
2	Q	75	<div><div><div>63%</div><div>57%</div><div>37%</div><div></div><div></div></div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11230 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 Likely histidyl tRNA-specific guanylyltransferase.

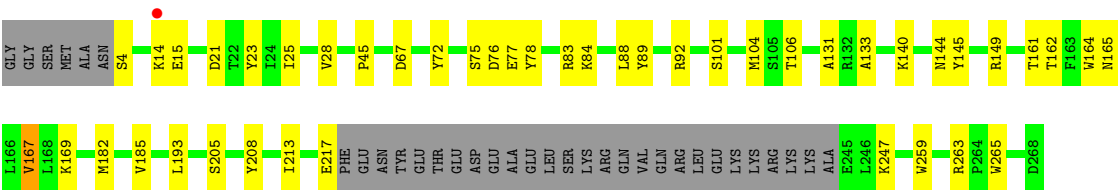
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	238	Total	C	N	O	S	0	0	0
			2032	1320	331	369	12			
1	B	238	Total	C	N	O	S	0	0	0
			2032	1320	331	369	12			
1	C	238	Total	C	N	O	S	0	0	0
			2032	1320	331	369	12			
1	D	238	Total	C	N	O	S	0	0	0
			2032	1320	331	369	12			

There are 12 discrepancies between the modelled and reference sequences:

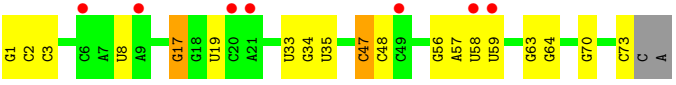
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
A	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
A	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
B	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
B	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
B	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
C	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
C	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
C	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
D	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
D	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
D	0	SER	-	EXPRESSION TAG	UNP Q5AFK5

- Molecule 2 is a RNA chain called 75-mer tRNA.

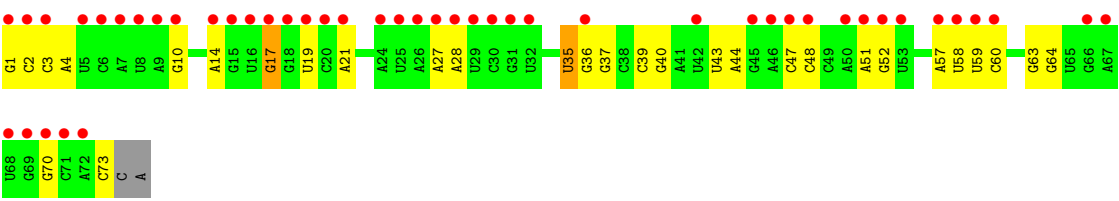
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	73	Total	C	N	O	P	0	0	0
			1551	693	268	517	73			
2	Q	73	Total	C	N	O	P	0	0	0
			1551	693	268	517	73			



• Molecule 2: 75-mer tRNA



• Molecule 2: 75-mer tRNA



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	96.53Å 96.53Å 300.39Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.95 – 4.18 48.79 – 4.18	Depositor EDS
% Data completeness (in resolution range)	98.8 (42.95-4.18) 98.9 (48.79-4.18)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 4.14Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.292 , 0.335 0.290 , 0.333	Depositor DCC
R_{free} test set	875 reflections (6.99%)	DCC
Wilson B-factor (Å ²)	174.3	Xtriage
Anisotropy	0.369	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 135.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.059 for -h,-k,l	Xtriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	11230	wwPDB-VP
Average B, all atoms (Å ²)	135.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% 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	A	0.25	0/2089	0.43	0/2819
1	B	0.25	0/2089	0.43	0/2819
1	C	0.25	0/2089	0.47	0/2819
1	D	0.26	0/2089	0.44	0/2819
2	P	0.39	1/1732 (0.1%)	0.81	0/2694
2	Q	0.40	1/1732 (0.1%)	0.84	0/2694
All	All	0.30	2/11820 (0.0%)	0.59	0/16664

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	Q	1	G	OP3-P	-10.60	1.48	1.61
2	P	1	G	OP3-P	-10.35	1.48	1.61

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2032	0	1975	20	0
1	B	2032	0	1975	22	0
1	C	2032	0	1975	31	0
1	D	2032	0	1975	26	0
2	P	1551	0	781	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	Q	1551	0	781	23	0
All	All	11230	0	9462	119	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 6.

All (119) 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:D:4:SER:N	2:Q:10:G:HO2'	1.61	0.96
1:C:204:GLU:O	1:C:209:LYS:NZ	2.23	0.70
1:C:140:LYS:O	1:C:144:ASN:ND2	2.26	0.67
1:B:217:GLU:OE2	1:B:263:ARG:NH1	2.27	0.67
1:D:259:TRP:O	1:D:263:ARG:NH1	2.28	0.66
1:D:164:TRP:HA	1:D:167:VAL:HG13	1.79	0.65
1:D:15:GLU:OE2	1:D:92:ARG:NH1	2.30	0.64
1:A:4:SER:HB3	1:A:7:GLU:HG3	1.79	0.64
1:A:140:LYS:O	1:A:144:ASN:ND2	2.28	0.63
1:A:263:ARG:HG2	1:A:265:TRP:CZ2	2.35	0.62
1:D:140:LYS:O	1:D:144:ASN:ND2	2.32	0.62
1:D:88:LEU:HG	1:D:89:TYR:HD2	1.66	0.60
1:B:28:VAL:HG21	1:B:104:MET:HG3	1.85	0.58
1:C:217:GLU:HG3	1:C:247:LYS:HD2	1.85	0.58
1:B:4:SER:HB2	2:P:73:C:H2'	1.86	0.57
1:C:67:ASP:OD1	1:C:83:ARG:NH1	2.34	0.57
2:P:17:G:H4'	2:P:59:U:H3	1.69	0.57
1:C:259:TRP:O	1:C:263:ARG:NH1	2.37	0.57
1:C:263:ARG:HG2	1:C:265:TRP:CZ2	2.40	0.56
1:C:191:GLU:HB2	2:Q:37:G:O2'	2.06	0.56
1:B:217:GLU:HG3	1:B:247:LYS:HD2	1.88	0.56
1:B:67:ASP:OD1	1:B:83:ARG:NH1	2.34	0.56
1:C:209:LYS:HD2	2:Q:36:G:O6	2.05	0.56
1:D:165:ASN:O	1:D:169:LYS:HB2	2.07	0.55
1:A:28:VAL:HG21	1:A:104:MET:HG3	1.88	0.54
1:A:162:THR:HA	1:A:193:LEU:HD21	1.90	0.53
1:B:166:LEU:HD23	1:B:170:LEU:HD13	1.90	0.53
1:B:72:TYR:HD1	1:B:213:ILE:HG13	1.75	0.52
1:A:217:GLU:HG3	1:A:247:LYS:HD2	1.92	0.51
2:Q:43:U:H2'	2:Q:44:A:H8	1.75	0.51
1:B:84:LYS:O	1:B:91:ARG:NH2	2.44	0.51
1:C:4:SER:HB3	1:C:7:GLU:HG3	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:263:ARG:HG2	1:D:265:TRP:CZ2	2.46	0.50
2:Q:3:C:H2'	2:Q:4:A:H8	1.77	0.50
1:D:23:TYR:CE1	1:D:84:LYS:HG3	2.46	0.50
1:A:6:TYR:HB3	1:B:148:TRP:HB2	1.94	0.50
1:B:75:SER:OG	1:B:76:ASP:N	2.45	0.49
1:B:4:SER:HB3	1:B:7:GLU:HG3	1.94	0.49
1:A:16:ASN:OD1	1:D:14:LYS:HE2	2.13	0.49
1:C:47:ASP:O	1:C:51:LEU:HD23	2.13	0.49
1:A:173:THR:HG22	1:A:175:GLN:H	1.78	0.49
1:D:217:GLU:HG3	1:D:247:LYS:HD2	1.95	0.48
1:C:172:MET:HE2	1:C:176:GLN:HB3	1.96	0.48
1:D:28:VAL:HG21	1:D:104:MET:HG3	1.94	0.48
1:C:28:VAL:HG21	1:C:104:MET:HG3	1.96	0.47
2:P:33:U:O2'	2:P:34:G:O5'	2.32	0.47
2:Q:14:A:H61	2:Q:21:A:H2	1.61	0.47
2:Q:51:A:H2'	2:Q:52:G:H8	1.80	0.47
1:C:75:SER:OG	1:C:76:ASP:N	2.48	0.46
1:A:75:SER:OG	1:A:76:ASP:N	2.47	0.46
1:A:25:ILE:HD13	1:A:145:TYR:CD2	2.50	0.46
1:B:33:PHE:HA	1:B:36:PHE:HB3	1.97	0.46
2:P:17:G:O2'	2:P:56:G:N1	2.45	0.46
1:A:185:VAL:HG22	1:A:188:ASP:OD2	2.16	0.46
1:B:45:PRO:HG3	1:B:161:THR:HA	1.97	0.46
2:Q:63:G:H2'	2:Q:64:G:H8	1.81	0.45
1:A:205:SER:HB3	1:A:208:TYR:CD2	2.52	0.45
1:C:200:ASN:OD1	2:Q:35:U:H1'	2.17	0.45
1:A:89:TYR:HE2	1:A:99:LEU:HD22	1.80	0.45
2:Q:39:C:H2'	2:Q:40:G:H8	1.81	0.45
1:D:75:SER:OG	1:D:76:ASP:N	2.50	0.45
1:C:88:LEU:HG	1:C:89:TYR:CD2	2.52	0.45
1:C:173:THR:HG23	1:C:174:PRO:HD2	1.99	0.45
2:Q:2:C:H2'	2:Q:3:C:H6	1.82	0.45
1:D:4:SER:N	2:Q:10:G:O2'	2.38	0.44
1:D:25:ILE:HD13	1:D:145:TYR:CD2	2.53	0.44
1:A:4:SER:N	2:Q:73:C:O2	2.50	0.44
2:Q:3:C:H2'	2:Q:4:A:C8	2.53	0.44
2:Q:57:A:N6	2:Q:60:C:C2	2.86	0.44
2:Q:63:G:H2'	2:Q:64:G:C8	2.52	0.44
1:B:19:LEU:HA	1:B:20:PRO:HD3	1.87	0.44
1:C:101:SER:OG	1:D:101:SER:OG	2.34	0.44
1:C:173:THR:HG22	1:C:175:GLN:H	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:205:SER:HB3	1:D:208:TYR:CD2	2.52	0.43
1:B:22:THR:HG21	1:B:135:LEU:HB3	2.00	0.43
1:A:27:ARG:NH2	1:A:132:ARG:HD2	2.33	0.43
1:D:28:VAL:HG12	1:D:131:ALA:HB2	2.00	0.43
2:Q:27:A:H2'	2:Q:28:A:H8	1.83	0.43
1:C:94:MET:HE1	1:D:133:ALA:H	1.83	0.43
1:D:77:GLU:OE1	1:D:149:ARG:NH2	2.47	0.43
1:A:88:LEU:HG	1:A:89:TYR:CD2	2.53	0.43
1:C:72:TYR:HD1	1:C:213:ILE:HG13	1.84	0.43
1:A:33:PHE:HA	1:A:36:PHE:HB3	2.01	0.43
1:C:77:GLU:CD	1:C:149:ARG:HH21	2.22	0.43
1:B:145:TYR:O	1:B:148:TRP:HB3	2.19	0.42
1:B:162:THR:HA	1:B:193:LEU:HD21	2.00	0.42
2:Q:17:G:H4'	2:Q:59:U:N3	2.34	0.42
2:Q:43:U:H2'	2:Q:44:A:C8	2.54	0.42
1:A:89:TYR:CE2	1:A:99:LEU:HD22	2.54	0.42
2:P:2:C:H2'	2:P:3:C:C6	2.54	0.42
2:P:56:G:H2'	2:P:57:A:H5'	2.00	0.42
2:P:63:G:H2'	2:P:64:G:C8	2.54	0.42
1:C:19:LEU:HA	1:C:20:PRO:HD3	1.81	0.42
2:Q:2:C:H2'	2:Q:3:C:C6	2.55	0.42
1:D:67:ASP:OD1	1:D:83:ARG:NH2	2.50	0.42
2:Q:39:C:H2'	2:Q:40:G:C8	2.55	0.42
1:C:27:ARG:NH2	1:C:132:ARG:HD2	2.35	0.42
1:A:7:GLU:HG2	1:B:148:TRP:HE1	1.85	0.41
1:C:89:TYR:CE2	1:C:99:LEU:HD22	2.55	0.41
1:D:162:THR:HA	1:D:193:LEU:HD21	2.03	0.41
2:P:8:U:O4'	2:P:47:C:O2'	2.31	0.41
1:C:161:THR:HG1	1:C:201:TYR:HE1	1.67	0.41
1:C:25:ILE:HD13	1:C:145:TYR:CD2	2.55	0.41
1:B:263:ARG:HG2	1:B:265:TRP:CZ2	2.56	0.41
1:D:21:ASP:O	1:D:84:LYS:HE2	2.20	0.41
2:P:17:G:H4'	2:P:59:U:N3	2.34	0.41
1:C:154:HIS:CE1	2:Q:36:G:N1	2.88	0.41
2:Q:51:A:H2'	2:Q:52:G:C8	2.55	0.41
2:P:63:G:H2'	2:P:64:G:H8	1.84	0.41
1:B:165:ASN:O	1:B:169:LYS:HB2	2.20	0.41
1:B:99:LEU:HA	1:B:99:LEU:HD23	1.90	0.41
1:C:48:LEU:HB2	1:C:208:TYR:OH	2.21	0.41
1:C:127:PRO:O	1:D:106:THR:HG21	2.21	0.41
1:D:72:TYR:HD1	1:D:213:ILE:HG13	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:179:GLN:HA	1:C:182:MET:HG3	2.02	0.40
1:D:45:PRO:HG3	1:D:161:THR:HG22	2.02	0.40
1:C:18:LEU:HB3	1:C:135:LEU:HD21	2.03	0.40
2:P:17:G:C5	2:P:57:A:C6	3.09	0.40
1:B:28:VAL:HA	1:B:130:ASP:O	2.21	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	234/271 (86%)	232 (99%)	2 (1%)	0	100	100
1	B	234/271 (86%)	234 (100%)	0	0	100	100
1	C	234/271 (86%)	234 (100%)	0	0	100	100
1	D	234/271 (86%)	234 (100%)	0	0	100	100
All	All	936/1084 (86%)	934 (100%)	2 (0%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	224/252 (89%)	221 (99%)	3 (1%)	73	87

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	224/252 (89%)	220 (98%)	4 (2%)	64	84
1	C	224/252 (89%)	219 (98%)	5 (2%)	57	80
1	D	224/252 (89%)	220 (98%)	4 (2%)	64	84
All	All	896/1008 (89%)	880 (98%)	16 (2%)	64	84

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	TYR
1	A	158	LEU
1	A	267	LYS
1	B	78	TYR
1	B	170	LEU
1	B	173	THR
1	B	267	LYS
1	C	5	LYS
1	C	78	TYR
1	C	154	HIS
1	C	184	THR
1	C	190	ASN
1	D	78	TYR
1	D	167	VAL
1	D	182	MET
1	D	185	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	154	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	P	72/75 (96%)	7 (9%)	0
2	Q	72/75 (96%)	7 (9%)	0
All	All	144/150 (96%)	14 (9%)	0

All (14) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	P	17	G
2	P	19	U
2	P	35	U
2	P	47	C
2	P	48	C
2	P	58	U
2	P	70	G
2	Q	17	G
2	Q	19	U
2	Q	35	U
2	Q	47	C
2	Q	48	C
2	Q	58	U
2	Q	70	G

There are no RNA pucker outliers to report.

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	238/271 (87%)	0.08	5 (2%) 64 54	78, 118, 148, 157	0
1	B	238/271 (87%)	0.33	19 (7%) 13 11	27, 140, 274, 298	0
1	C	238/271 (87%)	0.12	9 (3%) 41 32	17, 130, 235, 262	0
1	D	238/271 (87%)	-0.12	1 (0%) 92 88	16, 69, 106, 132	0
2	P	73/75 (97%)	0.88	7 (9%) 9 8	96, 104, 110, 111	0
2	Q	73/75 (97%)	3.06	47 (64%) 0 1	255, 262, 270, 272	0
All	All	1098/1234 (88%)	0.35	88 (8%) 13 11	16, 111, 264, 298	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	Q	26	A	12.4
1	B	182	MET	10.2
2	Q	59	U	8.0
2	Q	1	G	7.9
1	B	180	ARG	7.8
1	B	181	LEU	7.5
2	Q	57	A	7.3
2	Q	30	C	7.2
2	Q	6	C	6.6
2	Q	58	U	6.5
2	Q	36	G	6.5
2	Q	25	U	6.1
2	Q	47	C	5.6
1	B	162	THR	5.6
2	Q	31	G	5.2
2	Q	60	C	5.0
2	Q	15	G	4.8
1	A	201	TYR	4.8
1	B	183	GLY	4.7

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Mol	Chain	Res	Type	RSRZ
2	Q	19	U	4.7
2	Q	7	A	4.6
1	C	172	MET	4.6
2	Q	50	A	4.5
1	A	121	LEU	4.3
2	Q	29	U	4.3
1	C	186	ALA	4.2
1	B	189	LYS	4.2
1	B	159	TYR	4.2
2	Q	20	C	4.2
1	C	127	PRO	4.1
2	Q	9	A	4.0
1	B	184	THR	4.0
2	Q	51	A	4.0
2	Q	48	C	4.0
2	Q	2	C	3.9
2	Q	27	A	3.8
1	B	178	GLU	3.8
1	B	158	LEU	3.8
2	Q	21	A	3.7
2	P	58	U	3.6
2	Q	8	U	3.5
2	Q	10	G	3.5
1	B	156	ASN	3.4
1	B	196	GLU	3.3
2	Q	14	A	3.2
1	B	163	PHE	3.1
2	Q	69	G	3.1
1	C	201	TYR	3.0
2	Q	16	U	3.0
2	Q	3	C	2.9
2	Q	53	U	2.9
2	Q	66	G	2.9
1	A	180	ARG	2.8
2	Q	70	G	2.8
2	P	20	C	2.8
1	A	159	TYR	2.7
2	Q	17	G	2.7
2	Q	5	U	2.7
1	B	165	ASN	2.7
2	Q	45	G	2.6
2	Q	32	U	2.6

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Mol	Chain	Res	Type	RSRZ
2	Q	42	U	2.6
2	P	59	U	2.6
2	Q	18	G	2.6
2	Q	67	A	2.6
2	P	49	C	2.6
2	Q	46	A	2.6
2	Q	52	G	2.6
2	P	9	A	2.5
2	Q	72	A	2.5
2	Q	28	A	2.5
1	C	188	ASP	2.4
1	B	73	GLY	2.4
2	Q	24	A	2.4
2	P	6	C	2.4
1	B	179	GLN	2.2
1	B	14	LYS	2.2
1	C	76	ASP	2.2
1	C	154	HIS	2.2
1	B	195	LYS	2.2
2	Q	68	U	2.1
1	C	32	GLY	2.1
1	B	169	LYS	2.1
1	C	253	ILE	2.1
1	D	14	LYS	2.1
1	A	209	LYS	2.1
2	P	21	A	2.1
2	Q	71	C	2.1

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.