Full wwPDB X-ray Structure Validation Report (i)

Aug 20, 2023 – 11:52 AM EDT

PDB ID	:	2NRE
Title	:	Crystal structure of pseudoudirinde synthase TruA in complex with leucyl tRNA
Authors	:	Hur, S.; Stroud, R.M.
Deposited on	:	2006-11-01
Resolution	:	4.00 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 (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

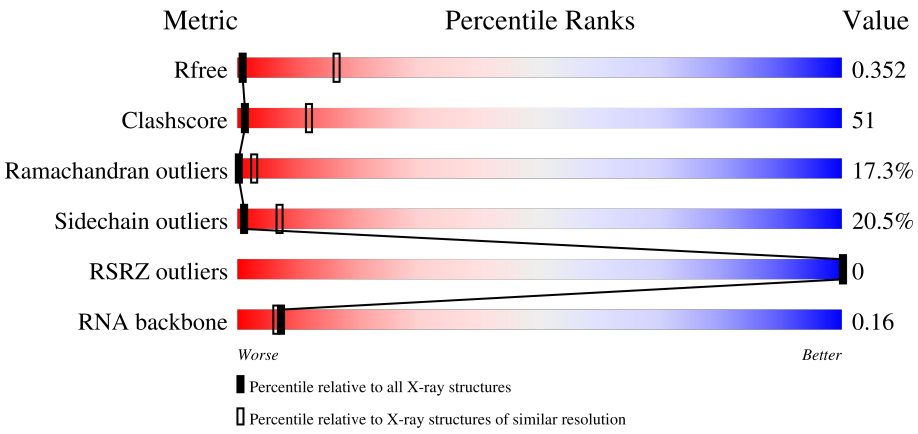
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 4.00 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	1087 (4.30-3.70)
Clashscore	141614	1148 (4.30-3.70)
Ramachandran outliers	138981	1108 (4.30-3.70)
Sidechain outliers	138945	1099 (4.30-3.70)
RSRZ outliers	127900	1028 (4.34-3.66)
RNA backbone	3102	1048 (5.00-3.00)

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 of 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 <=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	F	87	10% 26% 23% 5% 36%
2	A	270	20% 54% 20% . .

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3159 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 RNA chain called leucyl tRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	F	56	1196	533	212	395	56	0	0	0

- Molecule 2 is a protein called tRNA pseudouridine synthase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	A	259	1961	1251	356	347	7	0	0	0

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

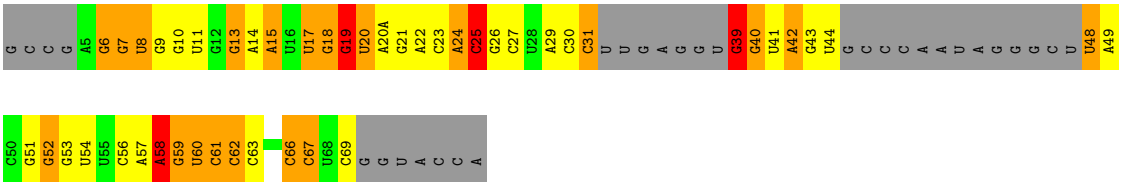
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	K		
3	F	2	2	2	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

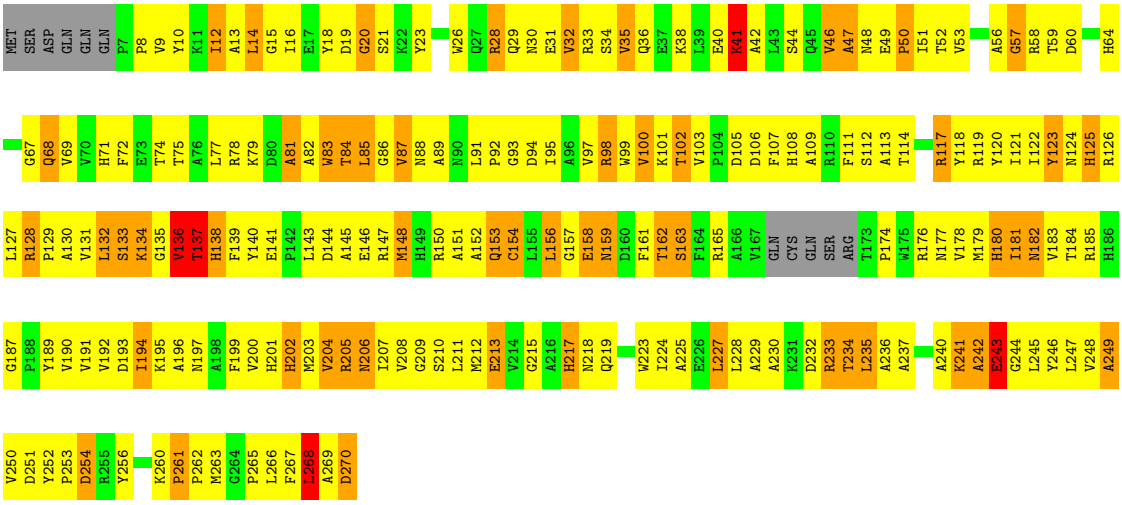
- Molecule 1: leucyl tRNA

Chain F: 10% 26% 23% 5% 36%



- Molecule 2: tRNA pseudouridine synthase A

Chain A: 20% 54% 20%



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	80.36Å 80.36Å 205.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 4.00 41.32 – 3.50	Depositor EDS
% Data completeness (in resolution range)	89.2 (20.00-4.00) 84.7 (41.32-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.17	Depositor
$< I/\sigma(I) >$ [1]	2.35 (at 3.48Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.259 , 0.357 0.266 , 0.352	Depositor DCC
R_{free} test set	501 reflections (5.35%)	wwPDB-VP
Wilson B-factor (Å^2)	74.3	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 33.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	0.059 for -h,-k,l	Xtriage
F_o,F_c correlation	0.81	EDS
Total number of atoms	3159	wwPDB-VP
Average B, all atoms (Å^2)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.50% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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)

Bond lengths and bond angles in the following residue types are not validated in this section: K

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	F	0.72	0/1334	1.35	11/2073 (0.5%)
2	A	0.53	0/2014	0.69	0/2746
All	All	0.61	0/3348	1.03	11/4819 (0.2%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	F	25	C	P-O3'-C3'	-14.72	102.03	119.70
1	F	58	A	P-O3'-C3'	8.75	130.20	119.70
1	F	26	G	OP1-P-OP2	-8.53	106.81	119.60
1	F	26	G	P-O5'-C5'	-8.29	107.64	120.90
1	F	25	C	O3'-P-O5'	7.50	118.25	104.00
1	F	39	G	OP1-P-OP2	-7.08	108.98	119.60
1	F	48	U	OP1-P-OP2	-6.73	109.50	119.60
1	F	39	G	P-O3'-C3'	6.00	126.90	119.70
1	F	19	G	C3'-C2'-C1'	5.49	105.89	101.50
1	F	13	G	C3'-C2'-C1'	5.47	105.88	101.50
1	F	8	U	C3'-C2'-C1'	5.22	105.67	101.50

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	F	1196	0	605	41	0
2	A	1961	0	1870	242	0
3	F	2	0	0	0	0
All	All	3159	0	2475	283	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 51.

All (283) 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:A:123:TYR:HB3	2:A:138:HIS:HB2	1.15	1.15
2:A:261:PRO:HB2	2:A:262:PRO:HD2	1.33	1.06
2:A:88:ASN:HA	2:A:91:LEU:HB2	1.39	1.05
2:A:83:TRP:O	2:A:87:VAL:HG21	1.57	1.03
2:A:152:ALA:HB2	2:A:211:LEU:HD11	1.42	1.01
2:A:77:LEU:C	2:A:78:ARG:HD3	1.88	0.94
2:A:77:LEU:O	2:A:78:ARG:HD3	1.69	0.93
2:A:82:ALA:O	2:A:87:VAL:HG23	1.69	0.92
2:A:248:VAL:HB	2:A:263:MET:HB2	1.53	0.90
2:A:157:GLY:O	2:A:158:GLU:HB2	1.72	0.90
2:A:88:ASN:HD22	2:A:97:VAL:H	1.21	0.87
2:A:213:GLU:OE1	2:A:213:GLU:HA	1.72	0.86
2:A:12:ILE:HG22	2:A:102:THR:HA	1.57	0.86
2:A:185:ARG:O	2:A:185:ARG:HG2	1.77	0.83
2:A:108:HIS:O	2:A:112:SER:HB2	1.79	0.82
2:A:180:HIS:ND1	2:A:180:HIS:O	2.12	0.82
2:A:32:VAL:HG23	2:A:33:ARG:H	1.43	0.82
2:A:261:PRO:CB	2:A:262:PRO:HD2	2.11	0.80
2:A:261:PRO:HB2	2:A:262:PRO:CD	2.12	0.79
2:A:123:TYR:CB	2:A:138:HIS:HB2	2.06	0.79
2:A:205:ARG:HB3	2:A:205:ARG:HH11	1.48	0.79
1:F:15:A:H2'	1:F:59:G:O6	1.84	0.78
1:F:18:G:N2	1:F:57:A:H2'	2.00	0.77
2:A:83:TRP:O	2:A:87:VAL:CG2	2.32	0.77
2:A:200:VAL:HG22	2:A:203:MET:SD	2.26	0.76
2:A:101:LYS:O	2:A:103:VAL:HG13	1.85	0.76
2:A:227:LEU:HA	2:A:230:ALA:HB3	1.68	0.76
2:A:144:ASP:O	2:A:148:MET:HG2	1.86	0.76
2:A:16:ILE:HD11	2:A:68:GLN:HB3	1.68	0.74
2:A:49:GLU:OE2	2:A:75:THR:HG22	1.87	0.74
2:A:82:ALA:O	2:A:87:VAL:CG2	2.36	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:84:THR:HG23	2:A:85:LEU:H	1.53	0.74
2:A:120:TYR:HE1	2:A:245:LEU:HD12	1.53	0.73
2:A:181:ILE:O	2:A:182:ASN:HB2	1.88	0.73
2:A:23:TYR:HE1	2:A:33:ARG:NH2	1.87	0.73
2:A:114:THR:HG22	2:A:251:ASP:O	1.89	0.72
2:A:108:HIS:CG	2:A:111:PHE:HB2	2.25	0.72
2:A:12:ILE:HD13	2:A:12:ILE:H	1.54	0.72
2:A:123:TYR:HB3	2:A:138:HIS:CB	2.08	0.72
1:F:21:G:H5"	1:F:22:A:H5'	1.71	0.71
2:A:136:VAL:N	2:A:243:GLU:O	2.22	0.71
2:A:12:ILE:HD13	2:A:12:ILE:N	2.05	0.71
2:A:107:PHE:HB2	2:A:256:TYR:CE1	2.26	0.71
2:A:147:ARG:O	2:A:148:MET:C	2.30	0.70
2:A:163:SER:HB3	2:A:232:ASP:H	1.55	0.70
2:A:240:ALA:O	2:A:241:LYS:O	2.09	0.70
2:A:16:ILE:C	2:A:16:ILE:HD12	2.11	0.70
2:A:119:ARG:NH2	2:A:270:ASP:OXT	2.22	0.70
2:A:15:GLY:HA3	2:A:99:TRP:NE1	2.06	0.70
2:A:148:MET:SD	2:A:215:GLY:HA3	2.32	0.69
1:F:14:A:N6	1:F:21:G:H1'	2.08	0.69
2:A:49:GLU:CD	2:A:75:THR:HG22	2.12	0.69
2:A:185:ARG:HG3	2:A:187:GLY:O	1.93	0.69
2:A:13:ALA:HA	2:A:71:HIS:HA	1.75	0.69
2:A:146:GLU:O	2:A:150:ARG:HB2	1.93	0.68
2:A:205:ARG:HH11	2:A:205:ARG:CB	2.06	0.68
2:A:135:GLY:C	2:A:136:VAL:HG22	2.15	0.67
1:F:41:U:H2'	1:F:42:A:C8	2.29	0.67
2:A:88:ASN:CA	2:A:91:LEU:HB2	2.22	0.67
2:A:240:ALA:O	2:A:241:LYS:C	2.33	0.67
2:A:162:THR:HG22	2:A:174:PRO:HB2	1.77	0.67
2:A:147:ARG:HA	2:A:150:ARG:HB3	1.77	0.66
2:A:208:VAL:HG12	2:A:212:MET:HG3	1.77	0.66
2:A:152:ALA:CB	2:A:211:LEU:HD11	2.21	0.66
2:A:32:VAL:HG23	2:A:33:ARG:N	2.11	0.65
2:A:79:LYS:HG3	2:A:79:LYS:O	1.96	0.65
2:A:120:TYR:CE1	2:A:245:LEU:HD12	2.31	0.65
2:A:14:LEU:HD21	2:A:72:PHE:HE1	1.62	0.65
1:F:39:G:H2'	1:F:40:G:C8	2.31	0.64
2:A:28:ARG:HG3	2:A:36:GLN:NE2	2.12	0.64
2:A:44:SER:HB3	2:A:51:ILE:HB	1.80	0.64
2:A:206:ASN:HD21	2:A:233:ARG:NH1	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:57:A:O2'	1:F:58:A:H5'	1.98	0.64
2:A:92:PRO:O	2:A:94:ASP:N	2.31	0.63
2:A:252:TYR:HB3	2:A:253:PRO:CD	2.28	0.63
2:A:233:ARG:O	2:A:235:LEU:N	2.30	0.63
2:A:205:ARG:HB3	2:A:205:ARG:NH1	2.12	0.62
2:A:64:HIS:CB	2:A:246:TYR:CE2	2.83	0.62
2:A:47:ALA:O	2:A:49:GLU:N	2.31	0.62
2:A:217:HIS:C	2:A:219:GLN:H	2.02	0.61
2:A:125:HIS:O	2:A:126:ARG:CB	2.48	0.61
1:F:14:A:H61	1:F:21:G:H1'	1.65	0.61
2:A:125:HIS:O	2:A:126:ARG:HB3	2.01	0.61
2:A:181:ILE:HG23	2:A:182:ASN:N	2.15	0.61
2:A:234:THR:O	2:A:234:THR:HG22	2.00	0.61
2:A:88:ASN:HA	2:A:91:LEU:CB	2.23	0.61
2:A:49:GLU:OE1	2:A:75:THR:HG22	2.00	0.60
2:A:151:ALA:O	2:A:224:ILE:HG21	2.01	0.60
2:A:210:SER:OG	2:A:236:ALA:HB1	2.01	0.60
2:A:147:ARG:O	2:A:150:ARG:N	2.33	0.60
2:A:44:SER:HB2	2:A:51:ILE:H	1.66	0.60
2:A:133:SER:O	2:A:134:LYS:C	2.39	0.60
1:F:30:C:H2'	1:F:31:C:H5	1.66	0.60
2:A:64:HIS:HB2	2:A:246:TYR:CD2	2.37	0.60
2:A:77:LEU:O	2:A:78:ARG:CD	2.48	0.60
2:A:108:HIS:ND1	2:A:109:ALA:N	2.50	0.59
2:A:122:ILE:O	2:A:189:TYR:HA	2.02	0.59
2:A:267:PHE:O	2:A:268:LEU:C	2.41	0.59
2:A:165:ARG:HH12	2:A:174:PRO:HB3	1.67	0.59
2:A:16:ILE:HD12	2:A:16:ILE:O	2.02	0.59
2:A:98:ARG:NH2	2:A:262:PRO:O	2.35	0.59
2:A:12:ILE:N	2:A:12:ILE:CD1	2.66	0.58
2:A:86:GLY:O	2:A:89:ALA:HB3	2.03	0.58
1:F:17:U:H4'	1:F:17:U:OP2	1.96	0.58
1:F:53:G:H2'	1:F:54:U:C6	2.39	0.58
2:A:47:ALA:C	2:A:49:GLU:H	2.07	0.58
2:A:185:ARG:O	2:A:185:ARG:CG	2.45	0.58
2:A:248:VAL:O	2:A:249:ALA:HB2	2.03	0.58
1:F:53:G:H2'	1:F:54:U:H6	1.66	0.57
1:F:39:G:H2'	1:F:40:G:H8	1.69	0.57
1:F:30:C:H2'	1:F:31:C:C5	2.39	0.57
2:A:203:MET:O	2:A:204:VAL:C	2.42	0.57
2:A:196:ALA:HB3	2:A:199:PHE:CZ	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:123:TYR:HD2	2:A:129:PRO:HD3	1.69	0.56
2:A:152:ALA:O	2:A:154:CYS:N	2.37	0.56
2:A:161:PHE:C	2:A:163:SER:N	2.59	0.56
2:A:265:PRO:O	2:A:266:LEU:HD12	2.06	0.56
1:F:48:U:C6	1:F:59:G:H5'	2.41	0.56
2:A:161:PHE:C	2:A:163:SER:H	2.09	0.56
2:A:180:HIS:ND1	2:A:180:HIS:C	2.59	0.56
2:A:208:VAL:CG1	2:A:212:MET:HG3	2.36	0.56
1:F:51:G:H2'	1:F:52:G:O4'	2.06	0.56
2:A:35:VAL:O	2:A:36:GLN:C	2.44	0.55
2:A:15:GLY:HA3	2:A:99:TRP:CE2	2.41	0.55
2:A:64:HIS:HB2	2:A:246:TYR:CE2	2.42	0.55
2:A:108:HIS:N	2:A:112:SER:OG	2.36	0.55
2:A:210:SER:HA	2:A:237:ALA:HB3	1.88	0.55
2:A:35:VAL:O	2:A:38:LYS:N	2.39	0.55
2:A:121:ILE:O	2:A:122:ILE:HG23	2.06	0.55
2:A:203:MET:O	2:A:206:ASN:N	2.40	0.55
2:A:49:GLU:OE1	2:A:75:THR:CG2	2.55	0.55
2:A:57:GLY:O	2:A:59:THR:N	2.40	0.55
2:A:254:ASP:OD1	2:A:254:ASP:N	2.39	0.55
2:A:250:VAL:O	2:A:260:LYS:HE3	2.06	0.54
2:A:146:GLU:O	2:A:150:ARG:CB	2.54	0.54
2:A:139:PHE:HE2	2:A:141:GLU:HG3	1.72	0.54
2:A:202:HIS:HB3	2:A:205:ARG:HD3	1.89	0.54
1:F:29:A:H2'	1:F:30:C:O4'	2.07	0.54
2:A:83:TRP:C	2:A:87:VAL:CG2	2.76	0.54
2:A:88:ASN:HA	2:A:91:LEU:HD12	1.88	0.54
1:F:18:G:H22	1:F:57:A:H2'	1.73	0.54
2:A:135:GLY:H	2:A:243:GLU:HB2	1.73	0.54
2:A:241:LYS:NZ	2:A:241:LYS:HB3	2.22	0.54
2:A:107:PHE:HB2	2:A:256:TYR:HE1	1.71	0.53
2:A:28:ARG:HG3	2:A:36:GLN:HE21	1.72	0.53
2:A:36:GLN:O	2:A:40:GLU:N	2.38	0.53
2:A:183:VAL:HG22	2:A:192:VAL:HG13	1.90	0.53
2:A:136:VAL:O	2:A:137:THR:O	2.28	0.52
2:A:213:GLU:OE1	2:A:213:GLU:CA	2.51	0.52
2:A:196:ALA:HB3	2:A:199:PHE:CE2	2.45	0.52
2:A:261:PRO:CB	2:A:262:PRO:CD	2.83	0.52
2:A:165:ARG:NH1	2:A:174:PRO:HB3	2.24	0.51
2:A:156:LEU:HA	2:A:178:VAL:HB	1.91	0.51
2:A:46:VAL:HG12	2:A:46:VAL:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:108:HIS:O	2:A:112:SER:CB	2.56	0.51
2:A:180:HIS:O	2:A:194:ILE:HA	2.11	0.51
2:A:224:ILE:CG2	2:A:225:ALA:N	2.73	0.51
2:A:44:SER:CB	2:A:51:ILE:HB	2.40	0.50
1:F:19:G:HO2'	1:F:20:U:P	2.34	0.50
1:F:24:A:H3'	1:F:25:C:H5"	1.93	0.50
2:A:117:ARG:HA	2:A:194:ILE:O	2.10	0.50
2:A:20:GLY:O	2:A:21:SER:HB3	2.09	0.50
2:A:100:VAL:CG2	2:A:101:LYS:N	2.75	0.50
2:A:158:GLU:O	2:A:159:ASN:O	2.29	0.50
2:A:83:TRP:C	2:A:87:VAL:HG21	2.29	0.49
2:A:83:TRP:HA	2:A:83:TRP:CE3	2.47	0.49
2:A:101:LYS:O	2:A:103:VAL:CG1	2.59	0.49
2:A:124:ASN:HD21	2:A:143:LEU:H	1.59	0.49
2:A:18:TYR:CE1	2:A:26:TRP:HZ3	2.30	0.49
2:A:87:VAL:HB	2:A:97:VAL:HG21	1.92	0.49
1:F:7:G:C2	1:F:67:C:O2	2.65	0.49
2:A:108:HIS:CB	2:A:111:PHE:HB2	2.42	0.49
2:A:161:PHE:O	2:A:163:SER:N	2.46	0.49
2:A:98:ARG:NH2	2:A:261:PRO:HG2	2.28	0.49
2:A:159:ASN:HB2	2:A:228:LEU:HD21	1.95	0.48
1:F:66:C:H2'	1:F:67:C:O5'	2.13	0.48
2:A:83:TRP:HA	2:A:83:TRP:HE3	1.77	0.48
1:F:53:G:O2'	1:F:54:U:H5'	2.13	0.48
2:A:145:ALA:O	2:A:146:GLU:C	2.52	0.48
2:A:23:TYR:HE2	2:A:95:ILE:HD11	1.79	0.48
2:A:41:LYS:O	2:A:42:ALA:C	2.52	0.48
2:A:108:HIS:HB3	2:A:111:PHE:HB2	1.95	0.48
2:A:108:HIS:CD2	2:A:111:PHE:HB2	2.49	0.48
1:F:18:G:H1'	1:F:58:A:C2	2.48	0.48
1:F:48:U:C5	1:F:59:G:H5'	2.49	0.48
2:A:14:LEU:HD21	2:A:72:PHE:CE1	2.46	0.48
2:A:135:GLY:C	2:A:136:VAL:CG2	2.81	0.48
2:A:28:ARG:HH21	2:A:40:GLU:HG2	1.79	0.48
2:A:148:MET:SD	2:A:215:GLY:CA	3.02	0.48
2:A:217:HIS:O	2:A:218:ASN:HB2	2.14	0.48
2:A:203:MET:O	2:A:207:ILE:N	2.41	0.47
1:F:25:C:H4'	1:F:25:C:OP1	2.14	0.47
2:A:217:HIS:C	2:A:219:GLN:N	2.69	0.47
2:A:252:TYR:HB3	2:A:253:PRO:HD2	1.96	0.46
2:A:81:ALA:HA	2:A:84:THR:HG23	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:201:HIS:O	2:A:203:MET:N	2.49	0.46
2:A:208:VAL:O	2:A:212:MET:N	2.32	0.46
1:F:51:G:N2	1:F:63:C:N3	2.47	0.46
2:A:18:TYR:OH	2:A:59:THR:OG1	2.33	0.46
2:A:67:GLY:O	2:A:69:VAL:HG23	2.15	0.46
2:A:162:THR:HG22	2:A:174:PRO:CB	2.43	0.46
2:A:203:MET:O	2:A:205:ARG:N	2.49	0.46
2:A:94:ASP:N	2:A:94:ASP:OD2	2.47	0.45
2:A:233:ARG:C	2:A:235:LEU:H	2.16	0.45
2:A:32:VAL:CG2	2:A:33:ARG:N	2.77	0.45
2:A:10:TYR:N	2:A:74:THR:O	2.34	0.45
2:A:47:ALA:HA	2:A:78:ARG:NH2	2.31	0.45
2:A:51:ILE:CG2	2:A:52:THR:N	2.79	0.45
2:A:204:VAL:O	2:A:208:VAL:HG23	2.16	0.45
2:A:100:VAL:HG22	2:A:101:LYS:N	2.30	0.45
2:A:163:SER:HB3	2:A:232:ASP:N	2.28	0.45
2:A:217:HIS:O	2:A:219:GLN:N	2.49	0.45
1:F:15:A:H2	1:F:20(A):A:H1'	1.82	0.45
1:F:58:A:H4'	1:F:59:G:OP1	2.17	0.45
2:A:119:ARG:HA	2:A:192:VAL:O	2.17	0.45
2:A:16:ILE:C	2:A:16:ILE:CD1	2.81	0.45
2:A:77:LEU:C	2:A:78:ARG:CD	2.75	0.45
2:A:13:ALA:H	2:A:103:VAL:CG1	2.30	0.45
2:A:47:ALA:C	2:A:49:GLU:N	2.70	0.45
2:A:148:MET:HB3	2:A:183:VAL:HG11	1.98	0.45
2:A:179:MET:SD	2:A:195:LYS:O	2.75	0.45
2:A:44:SER:CB	2:A:51:ILE:H	2.28	0.44
2:A:242:ALA:O	2:A:244:GLY:N	2.40	0.44
2:A:120:TYR:O	2:A:191:VAL:HA	2.17	0.44
2:A:267:PHE:O	2:A:269:ALA:N	2.50	0.44
2:A:182:ASN:HB3	2:A:193:ASP:OD2	2.17	0.44
1:F:29:A:C2	1:F:42:A:C2	3.05	0.44
2:A:137:THR:HG22	2:A:243:GLU:HG3	2.00	0.44
2:A:266:LEU:HB3	2:A:267:PHE:H	1.45	0.44
2:A:40:GLU:OE2	2:A:53:VAL:HG22	2.18	0.44
2:A:130:ALA:O	2:A:132:LEU:N	2.51	0.43
2:A:113:ALA:HA	2:A:252:TYR:CE2	2.53	0.43
2:A:161:PHE:HE1	2:A:176:ARG:O	2.00	0.43
2:A:265:PRO:C	2:A:266:LEU:HD12	2.38	0.43
2:A:127:LEU:O	2:A:128:ARG:C	2.56	0.43
2:A:81:ALA:HA	2:A:84:THR:CG2	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:119:ARG:HG2	2:A:193:ASP:HB3	2.01	0.43
2:A:181:ILE:CD1	2:A:182:ASN:H	2.32	0.43
1:F:22:A:N7	1:F:23:C:C4	2.87	0.43
2:A:35:VAL:HG12	2:A:36:GLN:N	2.33	0.43
2:A:252:TYR:CB	2:A:253:PRO:CD	2.95	0.42
2:A:153:GLN:OE1	2:A:181:ILE:HG22	2.19	0.42
2:A:15:GLY:HA2	2:A:69:VAL:HG22	2.01	0.42
2:A:49:GLU:O	2:A:50:PRO:C	2.57	0.42
2:A:14:LEU:N	2:A:14:LEU:HD23	2.35	0.42
2:A:177:ASN:O	2:A:197:ASN:N	2.41	0.42
1:F:14:A:H2'	1:F:14:A:N3	2.35	0.42
1:F:22:A:C8	1:F:23:C:C5	3.08	0.42
2:A:64:HIS:CB	2:A:246:TYR:CD2	3.02	0.42
2:A:200:VAL:O	2:A:203:MET:HB2	2.19	0.42
2:A:268:LEU:HA	2:A:268:LEU:HD22	1.73	0.42
2:A:15:GLY:O	2:A:98:ARG:HB2	2.20	0.42
2:A:136:VAL:C	2:A:137:THR:HG23	2.40	0.42
2:A:91:LEU:HA	2:A:92:PRO:HD3	1.78	0.42
2:A:163:SER:HA	2:A:232:ASP:C	2.40	0.42
2:A:209:GLY:C	2:A:211:LEU:H	2.23	0.42
2:A:227:LEU:HD22	2:A:236:ALA:HB2	2.00	0.42
2:A:232:ASP:OD1	2:A:233:ARG:N	2.53	0.42
1:F:51:G:N2	1:F:63:C:C2	2.85	0.41
2:A:64:HIS:CG	2:A:246:TYR:CE2	3.08	0.41
1:F:58:A:C4	1:F:61:C:C5	3.08	0.41
2:A:46:VAL:HG11	2:A:82:ALA:HB1	2.02	0.41
2:A:148:MET:O	2:A:211:LEU:HD21	2.21	0.41
2:A:91:LEU:HA	2:A:91:LEU:HD23	1.73	0.41
2:A:15:GLY:CA	2:A:99:TRP:CE2	3.03	0.41
2:A:196:ALA:O	2:A:197:ASN:C	2.57	0.41
2:A:208:VAL:HG12	2:A:212:MET:HE2	2.03	0.41
1:F:52:G:H1	1:F:62:C:H42	1.67	0.41
1:F:18:G:H21	1:F:57:A:H2'	1.79	0.41
1:F:62:C:H2'	1:F:63:C:H6	1.86	0.41
2:A:23:TYR:CE1	2:A:33:ARG:NH2	2.77	0.41
2:A:92:PRO:C	2:A:94:ASP:H	2.23	0.41
2:A:118:TYR:CZ	2:A:247:LEU:HD12	2.56	0.41
2:A:229:ALA:O	2:A:230:ALA:C	2.59	0.41
1:F:59:G:H2'	1:F:60:U:O4'	2.21	0.41
2:A:42:ALA:O	2:A:46:VAL:HG23	2.21	0.41
2:A:56:ALA:HB2	2:A:109:ALA:HB1	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:202:HIS:CB	2:A:205:ARG:HD3	2.51	0.41
1:F:40:G:C2	1:F:41:U:N3	2.90	0.40
2:A:193:ASP:OD2	2:A:193:ASP:N	2.55	0.40
2:A:26:TRP:CZ3	2:A:68:GLN:HG3	2.56	0.40
1:F:6:G:H1	1:F:67:C:H42	1.69	0.40
2:A:205:ARG:H	2:A:205:ARG:HG3	1.47	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
2	A	255/270 (94%)	161 (63%)	50 (20%)	44 (17%)	0 3

All (44) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	28	ARG
2	A	29	GLN
2	A	31	GLU
2	A	48	ASN
2	A	87	VAL
2	A	133	SER
2	A	137	THR
2	A	153	GLN
2	A	158	GLU
2	A	159	ASN
2	A	202	HIS
2	A	204	VAL
2	A	234	THR
2	A	241	LYS
2	A	20	GLY

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Mol	Chain	Res	Type
2	A	30	ASN
2	A	41	LYS
2	A	47	ALA
2	A	58	ARG
2	A	93	GLY
2	A	132	LEU
2	A	136	VAL
2	A	148	MET
2	A	154	CYS
2	A	182	ASN
2	A	223	TRP
2	A	243	GLU
2	A	268	LEU
2	A	32	VAL
2	A	105	ASP
2	A	128	ARG
2	A	134	LYS
2	A	217	HIS
2	A	242	ALA
2	A	57	GLY
2	A	249	ALA
2	A	35	VAL
2	A	50	PRO
2	A	81	ALA
2	A	131	VAL
2	A	140	TYR
2	A	8	PRO
2	A	261	PRO
2	A	46	VAL

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
2	A	190/223 (85%)	151 (80%)	39 (20%)	1 7

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	9	VAL
2	A	12	ILE
2	A	14	LEU
2	A	19	ASP
2	A	34	SER
2	A	41	LYS
2	A	60	ASP
2	A	68	GLN
2	A	83	TRP
2	A	84	THR
2	A	85	LEU
2	A	98	ARG
2	A	100	VAL
2	A	102	THR
2	A	106	ASP
2	A	117	ARG
2	A	123	TYR
2	A	125	HIS
2	A	136	VAL
2	A	137	THR
2	A	138	HIS
2	A	156	LEU
2	A	162	THR
2	A	163	SER
2	A	180	HIS
2	A	181	ILE
2	A	184	THR
2	A	190	VAL
2	A	194	ILE
2	A	205	ARG
2	A	206	ASN
2	A	213	GLU
2	A	227	LEU
2	A	233	ARG
2	A	235	LEU
2	A	243	GLU
2	A	254	ASP
2	A	268	LEU
2	A	270	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	36	GLN
2	A	68	GLN
2	A	88	ASN
2	A	201	HIS

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	F	54/87 (62%)	31 (57%)	8 (14%)

All (31) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	F	6	G
1	F	7	G
1	F	8	U
1	F	9	G
1	F	10	G
1	F	11	U
1	F	13	G
1	F	15	A
1	F	17	U
1	F	18	G
1	F	19	G
1	F	20	U
1	F	24	A
1	F	25	C
1	F	27	C
1	F	31	C
1	F	40	G
1	F	42	A
1	F	43	G
1	F	44	U
1	F	49	A
1	F	52	G
1	F	56	C
1	F	58	A
1	F	59	G
1	F	60	U
1	F	61	C
1	F	62	C
1	F	66	C

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Mol	Chain	Res	Type
1	F	67	C
1	F	69	C

All (8) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	F	8	U
1	F	15	A
1	F	18	G
1	F	19	G
1	F	24	A
1	F	39	G
1	F	58	A
1	F	60	U

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 (i)

6.1 Protein, DNA and RNA chains (i)

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	F	56/87 (64%)	0.04	0	100	100	74, 101, 145, 150	0
2	A	259/270 (95%)	-0.41	0	100	100	16, 39, 59, 66	0
All	All	315/357 (88%)	-0.33	0	100	100	16, 42, 123, 150	0

There are no RSRZ outliers to report.

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 monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	K	F	78	1/1	0.84	0.14	85,85,85,85	0
3	K	F	77	1/1	0.86	0.14	82,82,82,82	0

6.5 Other polymers (i)

There are no such residues in this entry.