



# Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 12:45 AM GMT

PDB ID : 3AMT  
Title : Crystal structure of the TiaS-tRNA(Ile2)-ATP complex  
Authors : Numata, T.; Osawa, T.  
Deposited on : 2010-08-23  
Resolution : 2.90 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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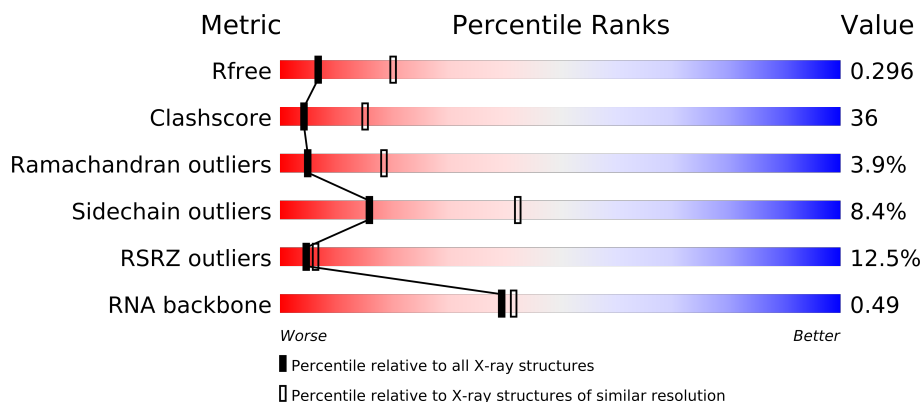
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.90 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))
$R_{free}$	66092	1053 (2.90-2.90)
Clashscore	79885	1326 (2.90-2.90)
Ramachandran outliers	78287	1290 (2.90-2.90)
Sidechain outliers	78261	1292 (2.90-2.90)
RSRZ outliers	66119	1054 (2.90-2.90)
RNA backbone	1838	1055 (3.40-2.40)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	440	
2	B	78	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5044 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Putative uncharacterized protein.

Mol	Chain	Residues	Atoms								ZeroOcc	AltConf	Trace
1	A	417	Total	C	N	O	P	S	Se		0	0	0
			3348	2127	579	625	1	5	11				

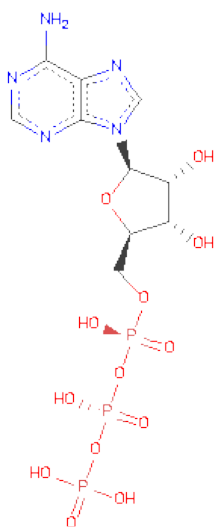
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MSE	-	EXPRESSION TAG	UNP O28025
A	-18	GLY	-	EXPRESSION TAG	UNP O28025
A	-17	SER	-	EXPRESSION TAG	UNP O28025
A	-16	SER	-	EXPRESSION TAG	UNP O28025
A	-15	HIS	-	EXPRESSION TAG	UNP O28025
A	-14	HIS	-	EXPRESSION TAG	UNP O28025
A	-13	HIS	-	EXPRESSION TAG	UNP O28025
A	-12	HIS	-	EXPRESSION TAG	UNP O28025
A	-11	HIS	-	EXPRESSION TAG	UNP O28025
A	-10	HIS	-	EXPRESSION TAG	UNP O28025
A	-9	SER	-	EXPRESSION TAG	UNP O28025
A	-8	SER	-	EXPRESSION TAG	UNP O28025
A	-7	GLY	-	EXPRESSION TAG	UNP O28025
A	-6	LEU	-	EXPRESSION TAG	UNP O28025
A	-5	VAL	-	EXPRESSION TAG	UNP O28025
A	-4	PRO	-	EXPRESSION TAG	UNP O28025
A	-3	ARG	-	EXPRESSION TAG	UNP O28025
A	-2	GLY	-	EXPRESSION TAG	UNP O28025
A	-1	SER	-	EXPRESSION TAG	UNP O28025
A	0	HIS	-	EXPRESSION TAG	UNP O28025

- Molecule 2 is a RNA chain called RNA (78-MER).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	78	Total	C	N	O	P		0	0	0
			1665	741	301	545	78				

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula:  $C_{10}H_{16}N_5O_{13}P_3$ ).

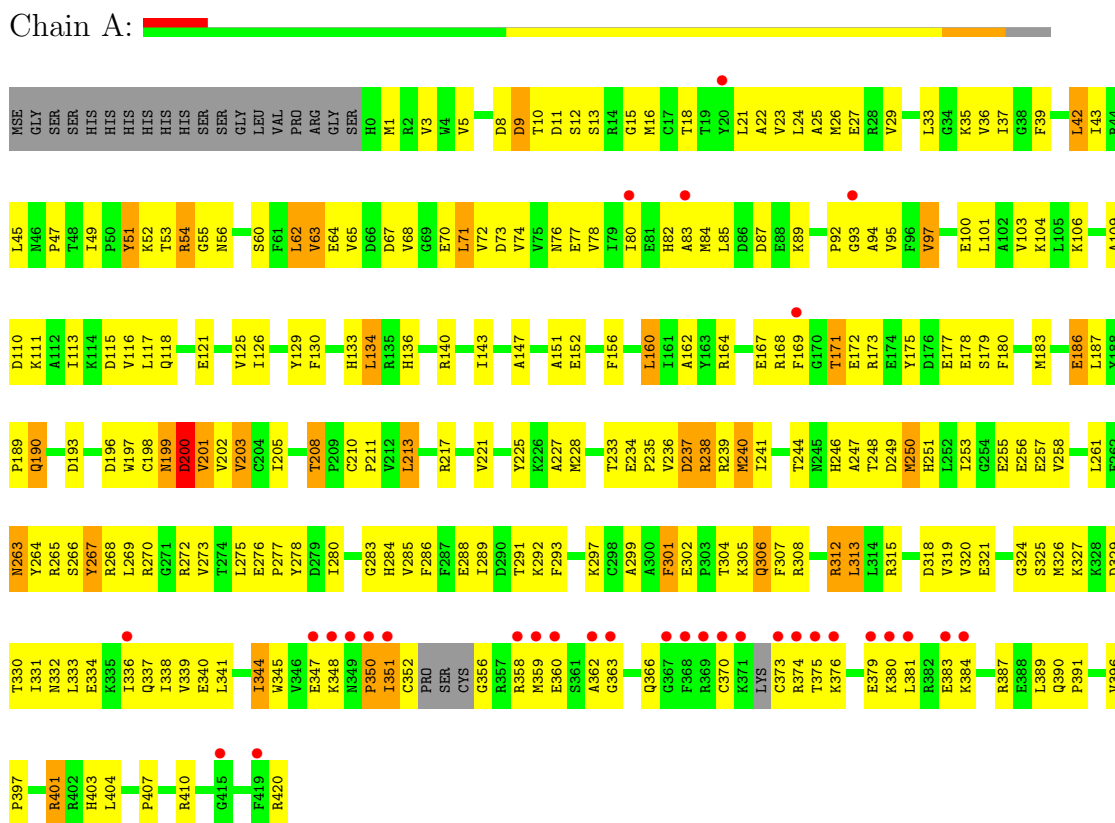


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
3	A	1	31	10	5	13	3	0	0

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

#### • Molecule 1: Putative uncharacterized protein



#### • Molecule 2: RNA (78-MER)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	131.10Å 131.10Å 86.62Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	47.48 – 2.90 47.48 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.5 (47.48-2.90) 99.8 (47.48-2.90)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.93 (at 2.91Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.240 , 0.286 0.243 , 0.296	Depositor DCC
$R_{free}$ test set	968 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	62.4	Xtriage
Anisotropy	0.409	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 45.1	EDS
Estimated twinning fraction	0.037 for -h,-k,l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 36617 reflections	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	5044	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

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

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	0.41	0/3392	0.69	0/4559
2	B	0.46	1/1859 (0.1%)	0.78	1/2895 (0.0%)
All	All	0.43	1/5251 (0.0%)	0.72	1/7454 (0.0%)

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
2	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	G	OP3-P	-6.90	1.52	1.61

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	G	OP1-P-OP2	-5.02	112.08	119.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	41	G	Sidechain

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3348	0	3304	251	0
2	B	1665	0	848	87	0
3	A	31	0	12	13	0
All	All	5044	0	4164	330	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 36.

All (330) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:111:LYS:HD3	1:A:117:LEU:HD21	1.32	1.12
1:A:26:MSE:HE3	1:A:36:VAL:HG11	1.22	1.11
1:A:341:LEU:HD21	1:A:389:LEU:HD12	1.30	1.06
1:A:362:ALA:O	1:A:366:GLN:HB2	1.63	0.99
1:A:33:LEU:HD11	1:A:74:VAL:HG21	1.55	0.89
2:B:4:C:H42	2:B:69:G:H1	1.20	0.88
1:A:11:ASP:CG	1:A:18:TPO:O3P	2.12	0.88
1:A:92:PRO:HG2	1:A:136:HIS:HB2	1.55	0.85
1:A:370:CYS:CB	1:A:375:THR:HG22	2.08	0.84
2:B:35:A:O2'	2:B:36:U:OP1	1.93	0.84
2:B:17:C:H2'	2:B:17:C:O2	1.77	0.82
1:A:83:ALA:O	1:A:85:LEU:HD12	1.79	0.81
1:A:11:ASP:OD2	1:A:18:TPO:O3P	1.98	0.81
1:A:12:SER:HB3	1:A:84:MSE:HG3	1.61	0.81
1:A:401:ARG:NH1	1:A:407:PRO:HB3	1.96	0.81
2:B:52:G:H1	2:B:62:C:H42	1.25	0.81
1:A:299:ALA:HB3	1:A:332:ASN:HD22	1.47	0.80
1:A:321:GLU:HB2	1:A:337:GLN:HB3	1.64	0.79
1:A:253:ILE:O	1:A:267:TYR:HB3	1.83	0.79
1:A:350:PRO:HB2	1:A:359:MSE:HE3	1.65	0.78
1:A:175:TYR:HB2	1:A:233:THR:HG22	1.63	0.78
1:A:12:SER:HB3	1:A:15:GLY:HA3	1.65	0.77
1:A:304:THR:HG22	1:A:307:PHE:CB	2.16	0.76
1:A:113:ILE:CD1	1:A:250:MSE:HE2	2.16	0.75

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:9:ASP:HB2	3:A:421:ATP:H3'	1.67	0.75
1:A:70:GLU:O	1:A:74:VAL:HG23	1.87	0.75
1:A:9:ASP:N	1:A:18:TPO:O2P	2.20	0.74
1:A:140:ARG:HG2	3:A:421:ATP:O3'	1.87	0.74
1:A:319:VAL:HG12	1:A:339:VAL:HB	1.69	0.73
1:A:101:LEU:HD21	1:A:129:TYR:HB3	1.70	0.72
1:A:351:ILE:HA	1:A:358:ARG:HA	1.70	0.72
1:A:111:LYS:HD3	1:A:117:LEU:CD2	2.16	0.71
1:A:67:ASP:OD2	1:A:70:GLU:HG3	1.90	0.71
2:B:63:C:H2'	2:B:64:U:C6	2.26	0.71
2:B:21(B):C:H1'	2:B:48:C:H41	1.56	0.71
2:B:47:U:O2'	2:B:48:C:OP2	2.08	0.70
1:A:111:LYS:HE2	1:A:115:ASP:OD2	1.92	0.70
1:A:115:ASP:CG	1:A:116:VAL:H	1.95	0.70
2:B:16:C:O2'	2:B:60:U:H1'	1.91	0.70
1:A:304:THR:HG22	1:A:307:PHE:HB2	1.73	0.69
2:B:52:G:H2'	2:B:53:G:H8	1.58	0.68
1:A:347:GLU:OE1	1:A:380:LYS:HD2	1.93	0.68
2:B:38:A:O2'	2:B:39:C:OP2	2.12	0.68
1:A:73:ASP:O	1:A:77:GLU:HG3	1.94	0.67
1:A:18:TPO:O3P	3:A:421:ATP:O1A	2.13	0.67
1:A:18:TPO:O3P	3:A:421:ATP:O3B	2.12	0.67
1:A:68:VAL:O	1:A:72:VAL:HG23	1.95	0.67
1:A:324:GLY:HA2	1:A:334:GLU:HG3	1.78	0.66
1:A:12:SER:CB	1:A:84:MSE:HG3	2.26	0.66
1:A:187:LEU:O	1:A:190:GLN:HB2	1.95	0.66
2:B:21(A):U:H5'	2:B:21(B):C:OP2	1.96	0.66
1:A:92:PRO:HG2	1:A:136:HIS:CB	2.26	0.66
1:A:167:GLU:CD	1:A:167:GLU:H	1.99	0.66
1:A:95:VAL:HG12	1:A:97:VAL:HG12	1.78	0.65
1:A:240:MSE:HE2	1:A:241:ILE:H	1.61	0.65
1:A:29:VAL:HG11	1:A:63:VAL:HG21	1.79	0.65
1:A:186:GLU:OE1	1:A:186:GLU:HA	1.96	0.65
2:B:3:G:H2'	2:B:4:C:H5''	1.79	0.65
1:A:173:ARG:HH11	1:A:173:ARG:HG3	1.61	0.65
1:A:22:ALA:O	1:A:26:MSE:HG3	1.96	0.65
2:B:4:C:N4	2:B:69:G:H1	1.93	0.64
1:A:403:HIS:CE1	2:B:34:C:H2'	2.32	0.63
1:A:389:LEU:HD13	1:A:390:GLN:N	2.13	0.63
2:B:35:A:HO2'	2:B:36:U:P	2.21	0.63
2:B:50:A:H2'	2:B:51:G:H8	1.64	0.63
1:A:193:ASP:OD1	1:A:401:ARG:HD2	1.98	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:9:ASP:CA	1:A:18:TPO:O2P	2.46	0.63
1:A:45:LEU:HB2	1:A:56:ASN:ND2	2.13	0.63
1:A:312:ARG:HG3	1:A:312:ARG:HH11	1.62	0.63
1:A:280:ILE:HG12	1:A:284:HIS:HB2	1.80	0.63
1:A:95:VAL:CG1	1:A:97:VAL:HG12	2.29	0.63
1:A:168:ARG:O	1:A:171:THR:HG23	1.99	0.62
2:B:17:C:C2'	2:B:17:C:O2	2.47	0.62
2:B:15:G:H5''	2:B:16:C:OP2	1.99	0.62
1:A:16:MSE:HE1	1:A:82:HIS:HB3	1.81	0.61
1:A:350:PRO:CB	1:A:359:MSE:HE3	2.29	0.61
1:A:54:ARG:NH2	2:B:33:U:O3'	2.33	0.61
1:A:101:LEU:HD21	1:A:129:TYR:CB	2.31	0.61
1:A:200:ASP:O	1:A:201:VAL:HG23	2.00	0.61
2:B:16:C:HO2'	2:B:60:U:H1'	1.64	0.61
1:A:313:LEU:HB3	1:A:389:LEU:HB2	1.81	0.60
2:B:64:U:H2'	2:B:65:C:H6	1.65	0.60
2:B:17:C:H5''	2:B:18:A:H4'	1.82	0.60
1:A:160:LEU:HD11	1:A:228:MSE:HE3	1.84	0.60
2:B:8:U:H1'	2:B:48:C:O2'	2.02	0.60
1:A:292:LYS:HG2	1:A:292:LYS:O	2.02	0.60
1:A:11:ASP:OD1	3:A:421:ATP:O2G	2.19	0.60
2:B:21(C):A:O2'	2:B:22:G:OP1	2.16	0.60
1:A:247:ALA:HB1	1:A:407:PRO:HD2	1.83	0.60
2:B:14:A:H5'	2:B:14:A:H8	1.67	0.60
1:A:26:MSE:CE	1:A:36:VAL:HG11	2.15	0.59
1:A:407:PRO:HG2	1:A:410:ARG:HG3	1.84	0.59
1:A:113:ILE:HD12	1:A:250:MSE:HE2	1.85	0.59
1:A:42:LEU:HB3	1:A:244:THR:HG22	1.82	0.59
1:A:401:ARG:HH11	1:A:407:PRO:HB3	1.66	0.59
1:A:39:PHE:CE2	1:A:241:ILE:HB	2.38	0.59
1:A:228:MSE:HE1	1:A:239:ARG:HD2	1.85	0.59
1:A:234:GLU:HB2	1:A:235:PRO:HD2	1.84	0.59
1:A:10:THR:HG22	1:A:18:TPO:HG22	1.85	0.58
1:A:10:THR:CG2	1:A:18:TPO:HG22	2.34	0.58
1:A:51:TYR:CD1	1:A:140:ARG:HD3	2.39	0.58
1:A:53:THR:OG1	3:A:421:ATP:O3B	2.21	0.58
1:A:315:ARG:HG2	1:A:387:ARG:NH2	2.18	0.58
1:A:352:CYS:O	1:A:356:GLY:N	2.37	0.57
1:A:272:ARG:O	1:A:289:ILE:HA	2.02	0.57
1:A:52:LYS:HA	3:A:421:ATP:H5'2	1.85	0.57
1:A:237:ASP:CG	1:A:238:ARG:H	2.08	0.57
1:A:197:TRP:O	1:A:199:ASN:N	2.35	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:270:ARG:HH12	1:A:339:VAL:HG11	1.69	0.57
1:A:63:VAL:HG13	1:A:65:VAL:HG13	1.86	0.57
1:A:196:ASP:CG	1:A:199:ASN:HB2	2.25	0.57
1:A:29:VAL:HG11	1:A:63:VAL:CG2	2.34	0.57
1:A:332:ASN:HB2	2:B:36:U:O2'	2.04	0.56
1:A:272:ARG:HA	1:A:318:ASP:O	2.05	0.56
2:B:17:C:OP2	2:B:19:G:H5'	2.05	0.56
1:A:348:LYS:HB2	1:A:381:LEU:HB2	1.87	0.56
1:A:312:ARG:HG3	1:A:312:ARG:NH1	2.18	0.56
1:A:338:ILE:HB	1:A:391:PRO:HA	1.87	0.56
1:A:3:VAL:HG13	1:A:63:VAL:HG12	1.85	0.56
1:A:315:ARG:HG3	1:A:344:ILE:CD1	2.35	0.56
1:A:253:ILE:H	1:A:267:TYR:HA	1.71	0.56
1:A:264:TYR:HE1	1:A:327:LYS:HZ1	1.52	0.56
1:A:11:ASP:OD2	1:A:53:THR:HG21	2.06	0.56
1:A:94:ALA:HB3	1:A:134:LEU:HB2	1.88	0.56
1:A:173:ARG:NH1	1:A:173:ARG:HG3	2.20	0.55
1:A:115:ASP:CG	1:A:116:VAL:N	2.59	0.55
1:A:180:PHE:CZ	1:A:205:ILE:HD11	2.41	0.55
2:B:43:G:H2'	2:B:44:C:C6	2.40	0.55
2:B:65:C:O2'	2:B:66:C:H5'	2.06	0.55
1:A:263:ASN:O	1:A:264:TYR:HB2	2.07	0.55
1:A:24:LEU:HD12	1:A:82:HIS:ND1	2.21	0.55
2:B:36:U:H4'	2:B:37:A:O5'	2.07	0.55
1:A:228:MSE:CE	1:A:239:ARG:HD2	2.37	0.55
1:A:297:LYS:O	1:A:330:THR:HG23	2.06	0.55
2:B:13:U:O2'	2:B:14:A:H5''	2.07	0.55
1:A:33:LEU:HD11	1:A:74:VAL:CG2	2.35	0.54
1:A:396:VAL:HB	1:A:397:PRO:HD2	1.87	0.54
1:A:115:ASP:HB2	1:A:265:ARG:NH2	2.23	0.54
1:A:109:ALA:O	1:A:113:ILE:HG13	2.07	0.54
1:A:345:TRP:HA	1:A:383:GLU:O	2.08	0.54
1:A:65:VAL:HG11	1:A:71:LEU:HG	1.89	0.54
1:A:304:THR:HG21	1:A:307:PHE:CD2	2.42	0.54
2:B:55:U:N3	2:B:57:G:H5''	2.22	0.54
1:A:162:ALA:HB1	1:A:236:VAL:HG13	1.89	0.53
1:A:304:THR:HG22	1:A:307:PHE:HB3	1.90	0.53
1:A:304:THR:HG23	1:A:396:VAL:HG12	1.90	0.53
2:B:21(C):A:N6	2:B:46:G:H2'	2.24	0.53
1:A:54:ARG:NH1	2:B:34:C:OP1	2.41	0.53
1:A:10:THR:N	1:A:18:TPO:O2P	2.42	0.53
2:B:52:G:H2'	2:B:53:G:C8	2.42	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:29:C:N3	2:B:41:G:O6	2.41	0.53
1:A:318:ASP:OD1	1:A:341:LEU:HA	2.08	0.52
2:B:19:G:O2'	2:B:20:G:H5'	2.09	0.52
2:B:58:A:H4'	2:B:59:A:OP1	2.09	0.52
1:A:11:ASP:OD1	1:A:53:THR:OG1	2.16	0.52
1:A:160:LEU:HD11	1:A:228:MSE:HB2	1.91	0.52
1:A:286:PHE:CZ	2:B:38:A:N3	2.78	0.52
2:B:21(C):A:H2'	2:B:46:G:O6	2.10	0.52
2:B:58:A:O2'	2:B:59:A:H3'	2.09	0.52
1:A:164:ARG:HG2	1:A:213:LEU:HG	1.91	0.52
1:A:80:ILE:O	1:A:80:ILE:HG22	2.09	0.52
1:A:275:LEU:HD23	1:A:276:GLU:O	2.09	0.52
1:A:106:LYS:HG3	1:A:152:GLU:HB2	1.91	0.52
1:A:338:ILE:HG21	1:A:389:LEU:HD11	1.92	0.52
1:A:286:PHE:HD2	1:A:299:ALA:HA	1.73	0.52
1:A:270:ARG:HA	1:A:320:VAL:O	2.11	0.51
1:A:93:GLY:HA2	1:A:134:LEU:O	2.10	0.51
2:B:23:A:H2'	2:B:24:G:C8	2.46	0.51
1:A:126:ILE:HG21	1:A:133:HIS:CD2	2.45	0.51
1:A:113:ILE:HD13	1:A:250:MSE:HE2	1.92	0.51
1:A:12:SER:OG	1:A:84:MSE:HE2	2.10	0.51
1:A:246:HIS:O	1:A:247:ALA:HB3	2.11	0.51
1:A:326:MSE:HA	1:A:330:THR:O	2.11	0.51
1:A:187:LEU:CD1	1:A:227:ALA:HB2	2.42	0.50
1:A:240:MSE:HE2	1:A:240:MSE:HA	1.93	0.50
1:A:197:TRP:C	1:A:199:ASN:H	2.15	0.50
1:A:190:GLN:HA	1:A:190:GLN:OE1	2.10	0.50
1:A:301:PHE:HE2	1:A:332:ASN:ND2	2.09	0.50
2:B:13:U:C2'	2:B:14:A:H5''	2.41	0.50
1:A:255:GLU:HB3	1:A:269:LEU:HD22	1.94	0.50
1:A:263:ASN:HD21	1:A:329:ASP:H	1.60	0.50
1:A:273:VAL:HA	1:A:288:GLU:O	2.12	0.50
1:A:8:ASP:OD2	3:A:421:ATP:O2A	2.30	0.50
1:A:291:THR:C	1:A:293:PHE:H	2.15	0.50
2:B:17:C:P	2:B:60:U:O2'	2.69	0.49
2:B:7:G:H1	2:B:66:C:H42	1.60	0.49
2:B:67:G:C6	2:B:68:G:N7	2.80	0.49
1:A:255:GLU:HB3	1:A:269:LEU:CD2	2.43	0.49
1:A:248:THR:O	1:A:249:ASP:HB2	2.12	0.49
2:B:63:C:H2'	2:B:64:U:H6	1.74	0.49
1:A:356:GLY:N	1:A:373:CYS:CB	2.76	0.49
1:A:103:VAL:HG13	1:A:104:LYS:N	2.28	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:266:SER:HA	1:A:325:SER:HA	1.95	0.49
1:A:101:LEU:HD12	1:A:104:LYS:HD2	1.94	0.48
1:A:1:MSE:O	1:A:64:GLU:HA	2.12	0.48
1:A:11:ASP:OD2	3:A:421:ATP:O1A	2.30	0.48
1:A:332:ASN:CB	2:B:36:U:O2'	2.61	0.48
1:A:87:ASP:OD2	1:A:89:LYS:HB2	2.13	0.48
2:B:17:C:H5''	2:B:18:A:C4'	2.42	0.48
1:A:12:SER:CB	1:A:84:MSE:HE2	2.44	0.48
1:A:268:ARG:O	1:A:269:LEU:HD23	2.14	0.48
1:A:341:LEU:CD2	1:A:389:LEU:HD12	2.23	0.48
1:A:18:TPO:O1P	3:A:421:ATP:O1B	2.31	0.48
3:A:421:ATP:O2A	3:A:421:ATP:O1B	2.32	0.48
1:A:261:LEU:HD11	1:A:269:LEU:HD11	1.95	0.48
2:B:50:A:O5'	2:B:50:A:H8	1.97	0.48
2:B:17:C:C4'	2:B:18:A:H4'	2.44	0.47
1:A:169:PHE:O	1:A:171:THR:N	2.44	0.47
2:B:28:C:O2'	2:B:29:C:H5'	2.14	0.47
2:B:50:A:H2'	2:B:51:G:C8	2.45	0.47
1:A:331:ILE:O	1:A:331:ILE:HG22	2.13	0.47
1:A:275:LEU:HD13	1:A:288:GLU:OE2	2.15	0.47
1:A:237:ASP:CG	1:A:238:ARG:N	2.68	0.47
1:A:251:HIS:H	1:A:251:HIS:CD2	2.31	0.47
2:B:7:G:H1	2:B:66:C:N4	2.12	0.47
1:A:304:THR:CG2	1:A:396:VAL:HG12	2.45	0.47
1:A:3:VAL:CG1	1:A:63:VAL:HG12	2.45	0.47
2:B:7:G:O2'	2:B:8:U:OP1	2.29	0.47
2:B:14:A:H2'	2:B:15:G:O4'	2.14	0.46
2:B:36:U:H5''	2:B:37:A:OP1	2.15	0.46
1:A:285:VAL:HG23	1:A:308:ARG:HG2	1.97	0.46
1:A:5:VAL:O	1:A:60:SER:HA	2.15	0.46
1:A:305:LYS:O	1:A:308:ARG:HD3	2.15	0.46
1:A:27:GLU:OE2	1:A:238:ARG:HD3	2.16	0.46
1:A:35:LYS:HG2	1:A:36:VAL:N	2.30	0.46
1:A:18:TPO:P	3:A:421:ATP:O1G	2.74	0.46
2:B:17:C:H5''	2:B:18:A:O3'	2.16	0.46
1:A:160:LEU:CD1	1:A:228:MSE:HB2	2.46	0.46
1:A:76:ASN:HB2	1:A:134:LEU:HD21	1.98	0.46
1:A:285:VAL:CG2	1:A:308:ARG:HG2	2.45	0.46
1:A:275:LEU:HD23	1:A:275:LEU:C	2.36	0.45
1:A:318:ASP:OD1	1:A:341:LEU:HD23	2.16	0.45
2:B:8:U:C1'	2:B:48:C:O2'	2.64	0.45
1:A:189:PRO:O	1:A:190:GLN:C	2.54	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:11:C:H2'	2:B:11:C:O2	2.16	0.45
1:A:100:GLU:H	1:A:100:GLU:CD	2.19	0.45
2:B:62:C:H2'	2:B:63:C:H6	1.81	0.45
1:A:304:THR:HG21	1:A:307:PHE:HD2	1.81	0.45
1:A:51:TYR:CD2	1:A:51:TYR:N	2.84	0.45
2:B:58:A:HO2'	2:B:59:A:P	2.39	0.45
2:B:9:A:H5'	2:B:10:G:OP2	2.16	0.45
1:A:12:SER:HB3	1:A:15:GLY:CA	2.43	0.45
1:A:221:VAL:O	1:A:225:TYR:HB2	2.15	0.45
1:A:37:ILE:HB	1:A:62:LEU:HD13	1.98	0.45
2:B:50:A:C5	2:B:51:G:N7	2.84	0.45
1:A:275:LEU:HB3	1:A:288:GLU:HB2	1.97	0.45
1:A:291:THR:C	1:A:293:PHE:N	2.70	0.45
2:B:21(B):C:O2	2:B:47:U:O2	2.35	0.45
1:A:180:PHE:CE2	1:A:205:ILE:HD11	2.52	0.45
1:A:180:PHE:O	1:A:183:MSE:HB3	2.17	0.45
1:A:217:ARG:NH2	1:A:403:HIS:ND1	2.66	0.44
2:B:20:G:N2	2:B:56:C:O2	2.45	0.44
1:A:302:GLU:C	1:A:304:THR:H	2.20	0.44
1:A:21:LEU:O	1:A:21:LEU:HD23	2.17	0.44
1:A:169:PHE:O	1:A:171:THR:HG22	2.16	0.44
1:A:327:LYS:HE2	1:A:327:LYS:HB3	1.85	0.44
1:A:384:LYS:HG3	1:A:384:LYS:O	2.16	0.44
2:B:38:A:O2'	2:B:39:C:P	2.75	0.44
2:B:43:G:H2'	2:B:44:C:O4'	2.17	0.44
1:A:121:GLU:O	1:A:125:VAL:HG23	2.17	0.44
2:B:68:G:H2'	2:B:69:G:O4'	2.17	0.44
1:A:263:ASN:HA	1:A:263:ASN:HD22	1.62	0.44
1:A:25:ALA:HA	1:A:78:VAL:HG21	2.00	0.44
1:A:53:THR:O	1:A:54:ARG:HB2	2.17	0.44
2:B:18:A:H2'	2:B:19:G:OP1	2.17	0.44
1:A:351:ILE:O	1:A:359:MSE:HE2	2.18	0.44
1:A:347:GLU:C	1:A:348:LYS:HG3	2.38	0.44
1:A:283:GLY:HA3	1:A:308:ARG:HH21	1.83	0.44
1:A:54:ARG:NH1	1:A:208:THR:CG2	2.81	0.43
1:A:351:ILE:O	1:A:359:MSE:HG3	2.18	0.43
1:A:304:THR:CG2	1:A:307:PHE:HB3	2.48	0.43
1:A:106:LYS:O	1:A:110:ASP:HB2	2.18	0.43
1:A:23:VAL:O	1:A:27:GLU:HG3	2.19	0.43
1:A:129:TYR:O	1:A:130:PHE:C	2.57	0.43
1:A:85:LEU:HD12	1:A:85:LEU:H	1.84	0.43
1:A:16:MSE:CE	1:A:82:HIS:HB3	2.47	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:277:PRO:O	1:A:278:TYR:HB3	2.18	0.43
1:A:269:LEU:O	1:A:321:GLU:HA	2.18	0.43
1:A:360:GLU:OE2	2:B:70:C:N4	2.50	0.43
2:B:66:C:H2'	2:B:67:G:O4'	2.18	0.43
1:A:348:LYS:O	1:A:381:LEU:HB2	2.18	0.43
2:B:3:G:C2'	2:B:4:C:H5''	2.45	0.43
1:A:54:ARG:NH1	1:A:208:THR:HG23	2.34	0.43
1:A:200:ASP:OD1	1:A:200:ASP:O	2.37	0.42
1:A:338:ILE:HG21	1:A:389:LEU:CD1	2.48	0.42
1:A:156:PHE:C	1:A:156:PHE:CD1	2.93	0.42
2:B:17:C:C5'	2:B:18:A:H4'	2.46	0.42
1:A:304:THR:CG2	1:A:307:PHE:CB	2.91	0.42
1:A:49:ILE:O	1:A:52:LYS:HG2	2.20	0.42
2:B:60:U:H5''	2:B:61:C:OP2	2.18	0.42
1:A:313:LEU:N	1:A:313:LEU:CD2	2.82	0.42
2:B:66:C:O2'	2:B:67:G:H5'	2.20	0.42
1:A:210:CYS:HA	1:A:211:PRO:HD3	1.84	0.42
1:A:341:LEU:HD12	1:A:391:PRO:HD3	2.02	0.42
1:A:319:VAL:HB	1:A:340:GLU:HB3	2.01	0.42
1:A:304:THR:HG23	1:A:396:VAL:CG1	2.50	0.42
1:A:319:VAL:CG1	1:A:339:VAL:HB	2.47	0.42
1:A:80:ILE:CG2	1:A:80:ILE:O	2.67	0.42
1:A:203:VAL:HB	2:B:34:C:C6	2.55	0.42
1:A:286:PHE:CZ	2:B:38:A:C2	3.08	0.42
1:A:257:GLU:O	1:A:258:VAL:HG13	2.20	0.41
1:A:27:GLU:OE1	1:A:238:ARG:NH1	2.54	0.41
2:B:37:A:H4'	2:B:38:A:OP2	2.19	0.41
1:A:228:MSE:HE1	1:A:239:ARG:CD	2.50	0.41
1:A:330:THR:HG22	1:A:331:ILE:N	2.36	0.41
1:A:118:GLN:O	1:A:121:GLU:N	2.53	0.41
1:A:177:GLU:O	1:A:179:SER:N	2.53	0.41
1:A:286:PHE:CD2	1:A:299:ALA:HA	2.53	0.41
1:A:307:PHE:CZ	1:A:336:ILE:HG22	2.56	0.41
2:B:43:G:H2'	2:B:44:C:H6	1.85	0.41
2:B:65:C:O2	2:B:66:C:C6	2.74	0.41
2:B:17:C:H4'	2:B:18:A:H4'	2.02	0.41
2:B:4:C:H6	2:B:4:C:C5'	2.34	0.41
1:A:315:ARG:CG	1:A:387:ARG:NH2	2.84	0.41
1:A:375:THR:OG1	1:A:376:LYS:N	2.54	0.41
1:A:109:ALA:HB2	1:A:147:ALA:HA	2.03	0.41
1:A:16:MSE:HE3	1:A:82:HIS:O	2.21	0.41
1:A:331:ILE:HG22	1:A:333:LEU:HD23	2.03	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:47:PRO:HD3	1:A:404:LEU:HD12	2.03	0.41
2:B:40:C:H2'	2:B:41:G:O4'	2.21	0.40
1:A:284:HIS:CE1	1:A:302:GLU:H	2.40	0.40
1:A:197:TRP:HD1	1:A:420:ARG:OXT	2.04	0.40
2:B:11:C:C2	2:B:12:U:C5	3.09	0.40
1:A:43:ILE:HG21	1:A:151:ALA:CB	2.51	0.40
1:A:264:TYR:CE1	1:A:327:LYS:HE3	2.55	0.40
2:B:20:G:H21	2:B:57:G:H1'	1.86	0.40
1:A:302:GLU:C	1:A:304:THR:N	2.74	0.40
2:B:58:A:O2'	2:B:59:A:O5'	2.35	0.40
1:A:9:ASP:HA	3:A:421:ATP:O1A	2.21	0.40
1:A:45:LEU:HB2	1:A:56:ASN:HD21	1.82	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	
1	A	410/440 (93%)	347 (85%)	47 (12%)	16 (4%)	5	18

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	54	ARG
1	A	198	CYS
1	A	200	ASP
1	A	237	ASP
1	A	350	PRO
1	A	178	GLU
1	A	363	GLY
1	A	9	ASP
1	A	190	GLN
1	A	199	ASN
1	A	351	ILE

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Mol	Chain	Res	Type
1	A	374	ARG
1	A	55	GLY
1	A	306	GLN
1	A	301	PHE
1	A	143	ILE

### 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	356/372 (96%)	326 (92%)	30 (8%)	16	42

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	SER
1	A	42	LEU
1	A	51	TYR
1	A	62	LEU
1	A	63	VAL
1	A	71	LEU
1	A	97	VAL
1	A	134	LEU
1	A	160	LEU
1	A	171	THR
1	A	172	GLU
1	A	186	GLU
1	A	200	ASP
1	A	201	VAL
1	A	202	VAL
1	A	203	VAL
1	A	208	THR
1	A	213	LEU
1	A	238	ARG
1	A	240	MSE
1	A	250	MSE
1	A	256	GLU
1	A	263	ASN

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Mol	Chain	Res	Type
1	A	267	TYR
1	A	306	GLN
1	A	312	ARG
1	A	313	LEU
1	A	344	ILE
1	A	379	GLU
1	A	401	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	118	GLN
1	A	133	HIS
1	A	199	ASN
1	A	251	HIS
1	A	263	ASN
1	A	309	ASN
1	A	332	ASN
1	A	337	GLN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	77/78 (98%)	26 (33%)	12 (15%)

All (26) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	4	C
2	B	8	U
2	B	9	A
2	B	10	G
2	B	14	A
2	B	16	C
2	B	17	C
2	B	18	A
2	B	20	G
2	B	21(A)	U
2	B	21(B)	C
2	B	21(C)	A

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Mol	Chain	Res	Type
2	B	22	G
2	B	34	C
2	B	35	A
2	B	36	U
2	B	37	A
2	B	38	A
2	B	39	C
2	B	41	G
2	B	42	G
2	B	48	C
2	B	49	G
2	B	57	G
2	B	59	A
2	B	61	C

All (12) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	7	G
2	B	9	A
2	B	21(B)	C
2	B	21(C)	A
2	B	35	A
2	B	36	U
2	B	37	A
2	B	38	A
2	B	47	U
2	B	48	C
2	B	58	A
2	B	60	U

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	TPO	A	18	1	10,10,11	5.28	2 (20%)	12,14,16	1.23	1 (8%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	TPO	A	18	1	-	0/9/11/13	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	18	TPO	O-C	16.15	1.22	1.11
1	A	18	TPO	CA-C	2.93	1.53	1.48

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	18	TPO	OG1-P-O1P	3.11	115.48	106.79

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	ATP	A	421	-	33,33,33	1.12	3 (9%)	52,52,52	2.04	9 (17%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ATP	A	421	-	-	0/22/38/38	0/1/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	421	ATP	C4-N9	-2.97	1.33	1.37
3	A	421	ATP	PG-O2G	2.08	1.62	1.54
3	A	421	ATP	PG-O3G	2.07	1.62	1.54

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	421	ATP	N3-C2-N1	-7.82	122.17	128.71
3	A	421	ATP	PB-O3B-PG	-5.98	114.14	131.68
3	A	421	ATP	PA-O3A-PB	-4.95	117.17	131.68
3	A	421	ATP	N3-C4-N9	4.45	133.47	125.43
3	A	421	ATP	C5-C4-N3	-2.92	119.35	125.70
3	A	421	ATP	O4'-C1'-C2'	-2.79	102.49	106.77
3	A	421	ATP	O2B-PB-O1B	2.56	126.54	112.21
3	A	421	ATP	C4-C5-N7	-2.28	107.57	109.52
3	A	421	ATP	C2-N3-C4	2.17	120.17	114.01

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	417/440 (94%)	0.63	32 (7%) 13 16	34, 59, 92, 114	0
2	B	78/78 (100%)	1.89	30 (38%) 1 0	51, 80, 135, 143	0
All	All	495/518 (95%)	0.83	62 (12%) 5 6	34, 62, 112, 143	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	21(B)	C	9.3
1	A	370	CYS	6.7
1	A	369	ARG	6.1
1	A	381	LEU	6.0
2	B	58	A	5.8
1	A	373	CYS	5.5
2	B	62	C	5.0
1	A	374	ARG	5.0
2	B	51	G	4.9
1	A	371	LYS	4.6
2	B	67	G	4.4
1	A	362	ALA	4.3
2	B	61	C	4.0
1	A	348	LYS	4.0
1	A	350	PRO	3.9
1	A	375	THR	3.9
2	B	16	C	3.8
1	A	351	ILE	3.7
2	B	49	G	3.7
1	A	368	PHE	3.6
1	A	376	LYS	3.6
1	A	349	ASN	3.4
1	A	419	PHE	3.4
2	B	65	C	3.4

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Mol	Chain	Res	Type	RSRZ
2	B	7	G	3.3
1	A	359	MSE	3.3
2	B	19	G	3.3
1	A	336	ILE	3.3
1	A	169	PHE	3.2
2	B	55	U	3.2
1	A	384	LYS	3.2
2	B	68	G	3.1
1	A	383	GLU	3.1
1	A	20	TYR	3.0
2	B	66	C	3.0
1	A	80	ILE	3.0
2	B	17	C	2.8
2	B	60	U	2.7
1	A	380	LYS	2.7
2	B	69	G	2.7
2	B	48	C	2.6
2	B	38	A	2.6
2	B	47	U	2.6
2	B	5	C	2.6
2	B	56	C	2.6
2	B	18	A	2.6
2	B	57	G	2.6
2	B	50	A	2.4
1	A	358	ARG	2.4
2	B	53	G	2.3
1	A	360	GLU	2.3
1	A	363	GLY	2.3
2	B	21(A)	U	2.3
1	A	415	GLY	2.2
2	B	76	A	2.2
2	B	59	A	2.2
1	A	347	GLU	2.2
1	A	367	GLY	2.1
1	A	83	ALA	2.1
2	B	52	G	2.1
1	A	93	GLY	2.0
1	A	379	GLU	2.0



## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
1	TPO	A	18	11/12	0.17	-	38,42,53,59	0

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	ATP	A	421	31/31	0.20	-	38,42,56,58	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.