



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:49 PM UTC

PDB ID : 3AMT / pdb_00003amt
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 X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

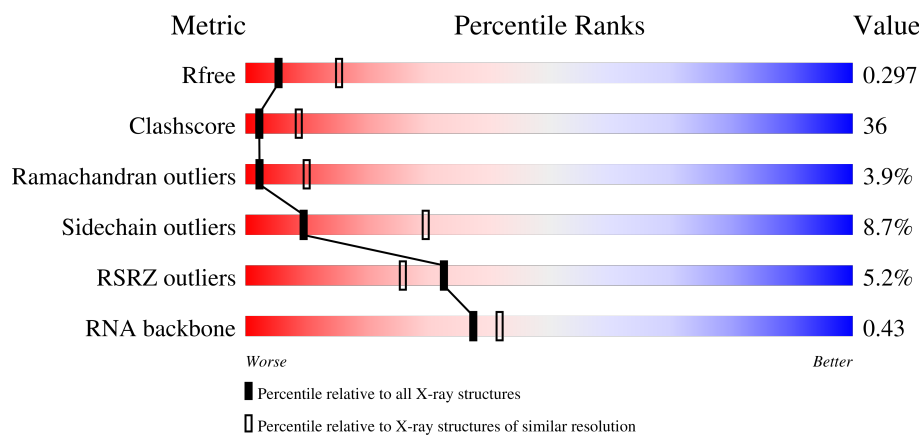
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)
RNA backbone	3983	1120 (3.10-2.70)

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 $\leq 5\%$. 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	440	
2	B	78	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	TPO	A	18	-	-	X	-
3	ATP	A	421	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5044 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 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			

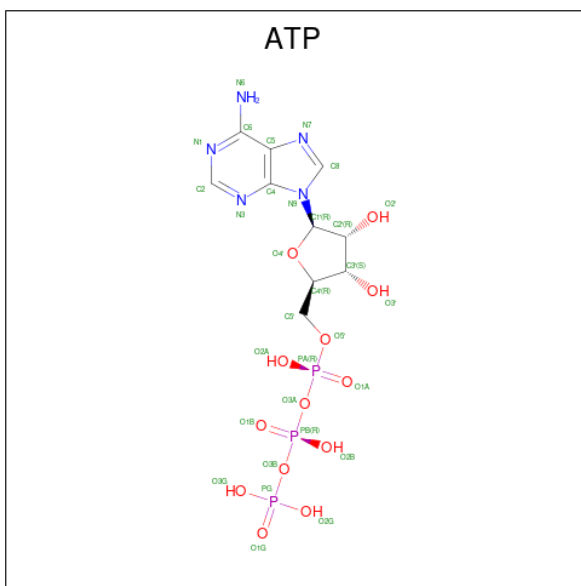
There are 20 discrepancies between the modelled and reference sequences:

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 (CCD ID: ATP) (formula: $\text{C}_{10}\text{H}_{16}\text{N}_5\text{O}_{13}\text{P}_3$).

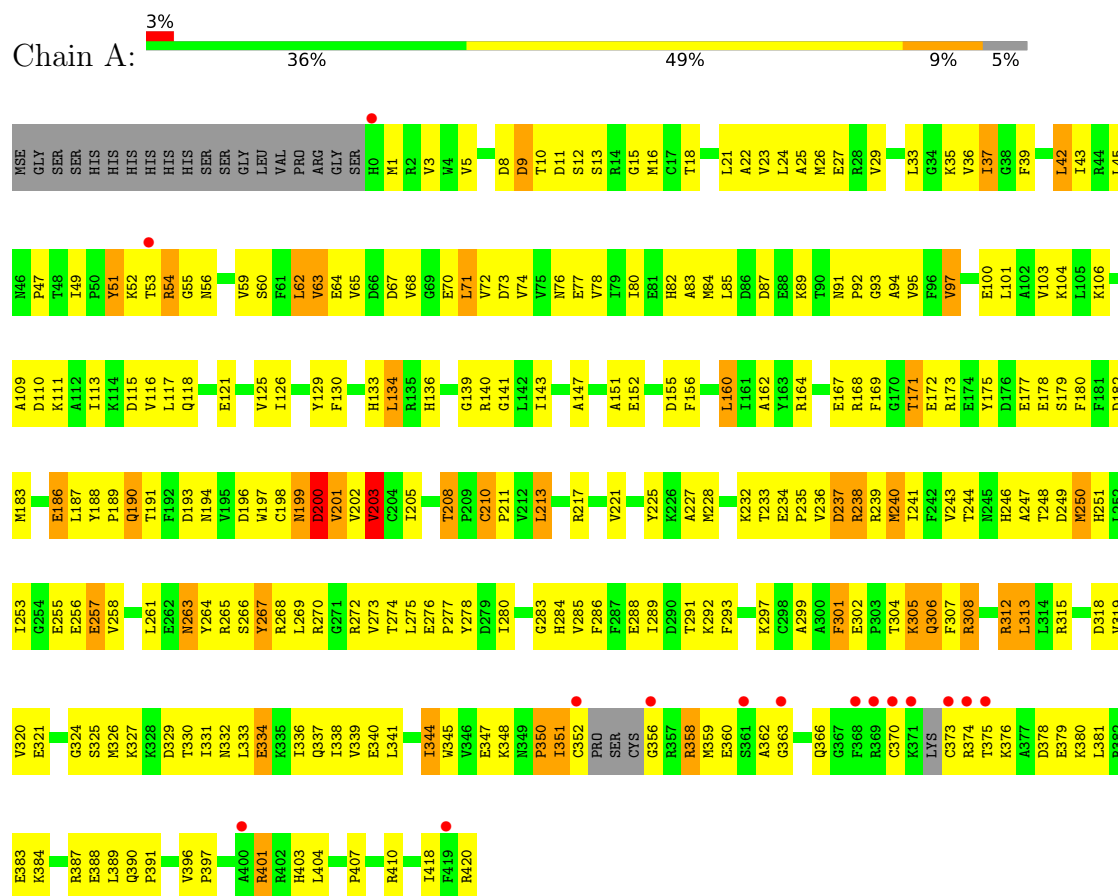


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

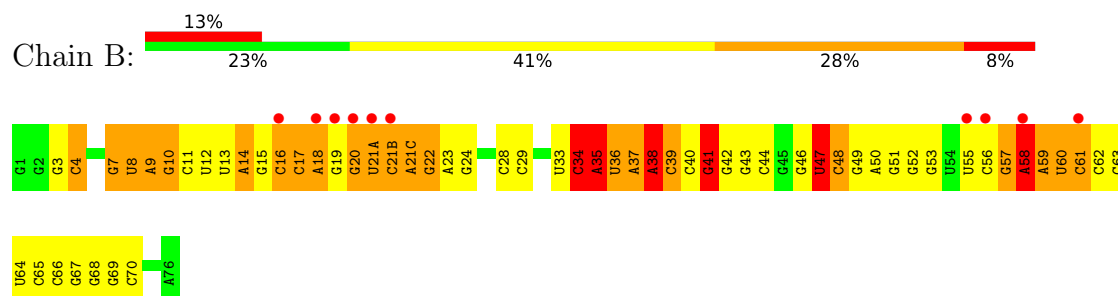
3 Residue-property plots

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: 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, α , β , γ	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) 99.7 (47.48-2.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.91Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.240 , 0.286 0.246 , 0.297	Depositor DCC
R_{free} test set	1844 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	62.4	Xtriage
Anisotropy	0.409	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 52.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.037 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5044	wwPDB-VP
Average B, all atoms (Å ²)	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.*

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

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	0.47	0/3392	1.05	20/4559 (0.4%)
2	B	0.37	0/1859	0.80	5/2895 (0.2%)
All	All	0.44	0/5251	0.96	25/7454 (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
2	B	0	1

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	334	GLU	N-CA-C	-9.24	102.58	114.04
1	A	257	GLU	N-CA-C	7.63	121.42	112.72
1	A	191	THR	N-CA-C	-7.36	97.70	109.76
1	A	305	LYS	CB-CA-C	-6.96	108.53	116.54
2	B	58	A	C2'-C3'-O3'	6.93	119.89	109.50
2	B	38	A	C2'-C3'-O3'	6.87	119.81	109.50
1	A	37	ILE	N-CA-C	6.83	118.30	107.98
1	A	182	ASP	N-CA-C	-6.57	104.12	111.28
1	A	274	THR	N-CA-C	-6.32	104.28	112.68
1	A	91	ASN	CA-C-N	6.26	126.23	120.03
1	A	91	ASN	C-N-CA	6.26	126.23	120.03
1	A	203	VAL	N-CA-C	5.88	116.66	110.72
2	B	47	U	C2'-C3'-O3'	5.76	118.14	109.50
1	A	210	CYS	CA-C-N	5.76	127.04	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	210	CYS	C-N-CA	5.76	127.04	119.84
1	A	210	CYS	N-CA-C	5.68	117.67	109.48
1	A	388	GLU	N-CA-C	-5.58	106.14	113.12
1	A	59	VAL	N-CA-C	5.55	116.73	108.46
1	A	358	ARG	N-CA-C	5.48	117.61	110.43
1	A	378	ASP	N-CA-C	5.41	119.84	112.04
2	B	34	C	N1-C1'-C2'	5.35	120.03	112.00
1	A	243	VAL	N-CA-C	-5.20	100.39	107.99
2	B	35	A	C2'-C3'-O3'	5.16	117.23	109.50
1	A	418	ILE	N-CA-C	5.13	115.67	108.89
1	A	155	ASP	N-CA-C	-5.04	101.62	109.24

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	41	G	Sidechain

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	3348	0	3304	256	0
2	B	1665	0	848	85	0
3	A	31	0	12	13	0
All	All	5044	0	4164	333	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 36.

All (333) 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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:11:ASP:CG	1:A:18:TPO:O3P	2.12	0.92
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: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
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:115:ASP:CG	1:A:116:VAL:H	1.95	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:167:GLU:CD	1:A:167:GLU:H	1.99	0.70
1:A:111:LYS:HE2	1:A:115:ASP:OD2	1.92	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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
1:A:173:ARG:HG3	1:A:173:ARG:HH11	1.61	0.65
2:B:3:G:H2'	2:B:4:C:H5''	1.79	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
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:HH11	1:A:312:ARG:HG3	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:O2	2:B:17:C:C2'	2.47	0.62
2:B:15:G:H5''	2:B:16:C:OP2	1.99	0.62
1:A:196:ASP:CG	1:A:199:ASN:HB2	2.25	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
1:A:237:ASP:CG	1:A:238:ARG:H	2.08	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
1:A:160:LEU:HD11	1:A:228:MSE:HE3	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:17:C:H5'	2:B:18:A:H4'	1.82	0.60
1:A:292:LYS:O	1:A:292:LYS:HG2	2.02	0.60
2:B:8:U:H1'	2:B:48:C:O2'	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:42:LEU:HB3	1:A:244:THR:HG22	1.82	0.59
1:A:113:ILE:HD12	1:A:250:MSE:HE2	1.85	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:197:TRP:O	1:A:199:ASN:N	2.35	0.57
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: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
1:A:312:ARG:HG3	1:A:312:ARG:NH1	2.18	0.56
1:A:348:LYS:HB2	1:A:381:LEU:HB2	1.87	0.56
2:B:17:C:OP2	2:B:19:G:H5'	2.05	0.56
1:A:3:VAL:HG13	1:A:63:VAL:HG12	1.85	0.56
1:A:338:ILE:HB	1:A:391:PRO:HA	1.87	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: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:HG3	1:A:173:ARG:NH1	2.20	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:65:C:O2'	2:B:66:C:H5'	2.06	0.55
1:A:197:TRP:C	1:A:199:ASN:H	2.15	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
1:A:115:ASP:CG	1:A:116:VAL:N	2.59	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:109:ALA:O	1:A:113:ILE:HG13	2.07	0.54
1:A:115:ASP:HB2	1:A:265:ARG:NH2	2.23	0.54
1:A:291:THR:C	1:A:293:PHE:H	2.15	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
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:160:LEU:HD11	1:A:228:MSE:HB2	1.91	0.52
1:A:80:ILE:HG22	1:A:80:ILE:O	2.09	0.52
1:A:164:ARG:HG2	1:A:213:LEU:HG	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: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:286:PHE:HD2	1:A:299:ALA:HA	1.73	0.52
1:A:338:ILE:HG21	1:A:389:LEU:HD11	1.92	0.52
1:A:237:ASP:CG	1:A:238:ARG:N	2.68	0.51
1:A:93:GLY:HA2	1:A:134:LEU:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:ILE:HG21	1:A:133:HIS:CD2	2.45	0.51
1:A:270:ARG:HA	1:A:320:VAL:O	2.11	0.51
2:B:23:A:H2'	2:B:24:G:C8	2.46	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:190:GLN:OE1	1:A:190:GLN:HA	2.10	0.50
1:A:189:PRO:O	1:A:190:GLN:C	2.54	0.50
1:A:275:LEU:HD23	1:A:275:LEU:C	2.36	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:100:GLU:CD	1:A:100:GLU:H	2.19	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
2:B:17:C:P	2:B:60:U:O2'	2.69	0.49
1:A:291:THR:C	1:A:293:PHE:N	2.70	0.49
2:B:7:G:H1	2:B:66:C:H42	1.60	0.49
1:A:255:GLU:HB3	1:A:269:LEU:CD2	2.43	0.49
2:B:67:G:C6	2:B:68:G:N7	2.80	0.49
1:A:248:THR:O	1:A:249:ASP:HB2	2.12	0.49
1:A:356:GLY:N	1:A:373:CYS:CB	2.76	0.49
2:B:63:C:H2'	2:B:64:U:H6	1.74	0.49
1:A:103:VAL:HG13	1:A:104:LYS:N	2.28	0.49
1:A:266:SER:HA	1:A:325:SER:HA	1.95	0.49
1:A:1:MSE:O	1:A:64:GLU:HA	2.12	0.48
1:A:101:LEU:HD12	1:A:104:LYS:HD2	1.94	0.48
1:A:347:GLU:C	1:A:348:LYS:HG3	2.38	0.48
1:A:11:ASP:OD2	3:A:421:ATP:O1A	2.30	0.48
1:A:87:ASP:OD2	1:A:89:LYS:HB2	2.13	0.48
1:A:332:ASN:CB	2:B:36:U:O2'	2.61	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:129:TYR:O	1:A:130:PHE:C	2.57	0.48
1:A:268:ARG:O	1:A:269:LEU:HD23	2.14	0.48
1:A:18:TPO:O1P	3:A:421:ATP:O1B	2.31	0.48
1:A:261:LEU:HD11	1:A:269:LEU:HD11	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:LEU:CD2	1:A:389:LEU:HD12	2.23	0.48
3:A:421:ATP:O2A	3:A:421:ATP:O1B	2.32	0.48
2:B:50:A:O5'	2:B:50:A:H8	1.97	0.48
1:A:302:GLU:C	1:A:304:THR:H	2.20	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:156:PHE:C	1:A:156:PHE:CD1	2.93	0.47
1:A:331:ILE:HG22	1:A:331:ILE:O	2.13	0.47
1:A:275:LEU:HD13	1:A:288:GLU:OE2	2.15	0.47
1:A:251:HIS:CD2	1:A:251:HIS:H	2.31	0.47
2:B:7:G:H1	2:B:66:C:N4	2.12	0.47
1:A:3:VAL:CG1	1:A:63:VAL:HG12	2.45	0.47
1:A:304:THR:CG2	1:A:396: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:5:VAL:O	1:A:60:SER:HA	2.15	0.46
1:A:285:VAL:HG23	1:A:308:ARG:HG2	1.97	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:18:TPO:P	3:A:421:ATP:O1G	2.74	0.46
1:A:35:LYS:HG2	1:A:36:VAL:N	2.30	0.46
2:B:17:C:H5''	2:B:18:A:O3'	2.16	0.46
1:A:76:ASN:HB2	1:A:134:LEU:HD21	1.98	0.46
1:A:160:LEU:CD1	1:A:228:MSE:HB2	2.46	0.46
1:A:285:VAL:CG2	1:A:308:ARG:HG2	2.45	0.46
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
2:B:11:C:H2'	2:B:11:C:O2	2.16	0.45
1:A:304:THR:HG21	1:A:307:PHE:HD2	1.81	0.45
2:B:62:C:H2'	2:B:63:C:H6	1.81	0.45
1:A:51:TYR:CD2	1:A:51:TYR:N	2.84	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:37:ILE:HB	1:A:62:LEU:HD13	1.98	0.45
1:A:221:VAL:O	1:A:225:TYR:HB2	2.15	0.45
1:A:275:LEU:HB3	1:A:288:GLU:HB2	1.97	0.45
2:B:50:A:C5	2:B:51:G:N7	2.84	0.45
1:A:180:PHE:O	1:A:183:MSE:HB3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:PHE:CE2	1:A:205:ILE:HD11	2.52	0.45
2:B:21(B):C:O2	2:B:47:U:O2	2.35	0.45
1:A:21:LEU:HD23	1:A:21:LEU:O	2.17	0.44
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:169:PHE:O	1:A:171:THR:HG22	2.16	0.44
1:A:302:GLU:C	1:A:304:THR:N	2.74	0.44
1:A:327:LYS:HE2	1:A:327:LYS:HB3	1.85	0.44
1:A:384:LYS:O	1:A:384:LYS:HG3	2.16	0.44
1:A:121:GLU:O	1:A:125:VAL:HG23	2.17	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:25:ALA:HA	1:A:78:VAL:HG21	2.00	0.44
1:A:263:ASN:HD22	1:A:263:ASN:HA	1.62	0.44
2:B:68:G:H2'	2:B:69:G:O4'	2.17	0.44
2:B:58:A:O2'	2:B:59:A:O5'	2.35	0.44
1:A:232:LYS:C	1:A:233:THR:HG23	2.43	0.44
1:A:53:THR:O	1:A:54:ARG:HB2	2.17	0.44
1:A:283:GLY:HA3	1:A:308:ARG:HH21	1.83	0.44
1:A:351:ILE:O	1:A:359:MSE:HE2	2.18	0.44
2:B:18:A:H2'	2:B:19:G:OP1	2.17	0.44
1:A:54:ARG:NH1	1:A:208:THR:CG2	2.81	0.43
1:A:106:LYS:O	1:A:110:ASP:HB2	2.18	0.43
1:A:297:LYS:C	1:A:330:THR:HG23	2.42	0.43
1:A:304:THR:CG2	1:A:307:PHE:HB3	2.48	0.43
1:A:351:ILE:O	1:A:359:MSE:HG3	2.18	0.43
1:A:23:VAL:O	1:A:27:GLU:HG3	2.19	0.43
1:A:16:MSE:CE	1:A:82:HIS:HB3	2.47	0.43
1:A:85:LEU:HD12	1:A:85:LEU:H	1.84	0.43
1:A:167:GLU:CD	1:A:167:GLU:N	2.74	0.43
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
1:A:139:GLY:C	1:A:141:GLY:N	2.76	0.43
1:A:348:LYS:O	1:A:381:LEU:HB2	2.18	0.43
2:B:66:C:H2'	2:B:67:G:O4'	2.18	0.43
1:A:54:ARG:NH1	1:A:208:THR:HG23	2.34	0.43
2:B:3:G:C2'	2:B:4:C:H5''	2.45	0.43
1:A:200:ASP:O	1:A:200:ASP:OD1	2.37	0.42
1:A:338:ILE:HG21	1:A:389:LEU:CD1	2.48	0.42
2:B:17:C:C5'	2:B:18:A:H4'	2.46	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:319:VAL:HB	1:A:340:GLU:HB3	2.01	0.42
1:A:341:LEU:HD12	1:A:391:PRO:HD3	2.02	0.42
1:A:80:ILE:O	1:A:80:ILE:CG2	2.67	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: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:200:ASP:O	1:A:200:ASP:CG	2.64	0.41
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
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:188:TYR:CD2	1:A:188:TYR:C	2.98	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
2:B:37:A:H4'	2:B:38:A:OP2	2.19	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
1:A:194:ASN:C	1:A:203:VAL:HG23	2.46	0.41
1:A:315:ARG:CG	1:A:387:ARG:NH2	2.84	0.41
2:B:4:C:H6	2:B:4:C:C5'	2.34	0.41
1:A:16:MSE:HE3	1:A:82:HIS:O	2.21	0.41
1:A:109:ALA:HB2	1:A:147:ALA:HA	2.03	0.41
1:A:331:ILE:HG22	1:A:333:LEU:HD23	2.03	0.41
1:A:375:THR:OG1	1:A:376:LYS:N	2.54	0.41
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:43:ILE:HG21	1:A:151:ALA:CB	2.51	0.40
1:A:197:TRP:HD1	1:A:420:ARG:OXT	2.04	0.40
1:A:284:HIS:CE1	1:A:302:GLU:H	2.40	0.40
2:B:11:C:C2	2:B:12:U:C5	3.09	0.40
1:A:264:TYR:CE1	1:A:327:LYS:HE3	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:20:G:H21	2:B:57:G:H1'	1.86	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 [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	410/440 (93%)	347 (85%)	47 (12%)	16 (4%)	2	10

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
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%)	325 (91%)	31 (9%)	9 30

All (31) 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
1	A	267	TYR
1	A	306	GLN
1	A	308	ARG
1	A	312	ARG
1	A	313	LEU
1	A	344	ILE
1	A	379	GLU

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Mol	Chain	Res	Type
1	A	401	ARG

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

Mol	Chain	Res	Type
1	A	76	ASN
1	A	118	GLN
1	A	133	HIS
1	A	199	ASN
1	A	251	HIS
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
2	B	22	G
2	B	34	C
2	B	35	A
2	B	36	U
2	B	37	A
2	B	38	A

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Mol	Chain	Res	Type
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	8,10,11	1.08	0	10,14,16	1.04	0

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	-	2/9/11/13	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	18	TPO	O-C-CA-CB
1	A	18	TPO	CB-OG1-P-O2P

There are no ring outliers.

1 monomer is involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	18	TPO	11	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	-	32,33,33	1.23	5 (15%)	48,52,52	1.91	11 (22%)

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	-	-	1/22/38/38	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	421	ATP	C5-N7	-3.14	1.33	1.39
3	A	421	ATP	C8-N9	-2.29	1.33	1.37
3	A	421	ATP	C4-N9	-2.06	1.33	1.37
3	A	421	ATP	PG-O2G	2.03	1.62	1.54
3	A	421	ATP	PG-O3G	2.01	1.62	1.54

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	421	ATP	C5-C4-N3	-5.35	119.35	126.72
3	A	421	ATP	N3-C2-N1	-4.24	122.17	128.58
3	A	421	ATP	O2B-PB-O3A	-4.16	96.02	107.27
3	A	421	ATP	O3A-PB-O1B	-3.81	99.23	110.70
3	A	421	ATP	N3-C4-N9	3.70	133.47	127.17
3	A	421	ATP	C2-N3-C4	3.42	120.17	111.83
3	A	421	ATP	O2B-PB-O1B	3.03	126.54	112.44
3	A	421	ATP	N9-C8-N7	-2.82	109.94	113.94
3	A	421	ATP	C4-C5-N7	-2.64	107.57	110.58
3	A	421	ATP	C5-N7-C8	2.61	107.55	103.45
3	A	421	ATP	O2B-PB-O3B	2.42	113.81	107.27

There are no chirality outliers.

All (1) torsion outliers are listed below:

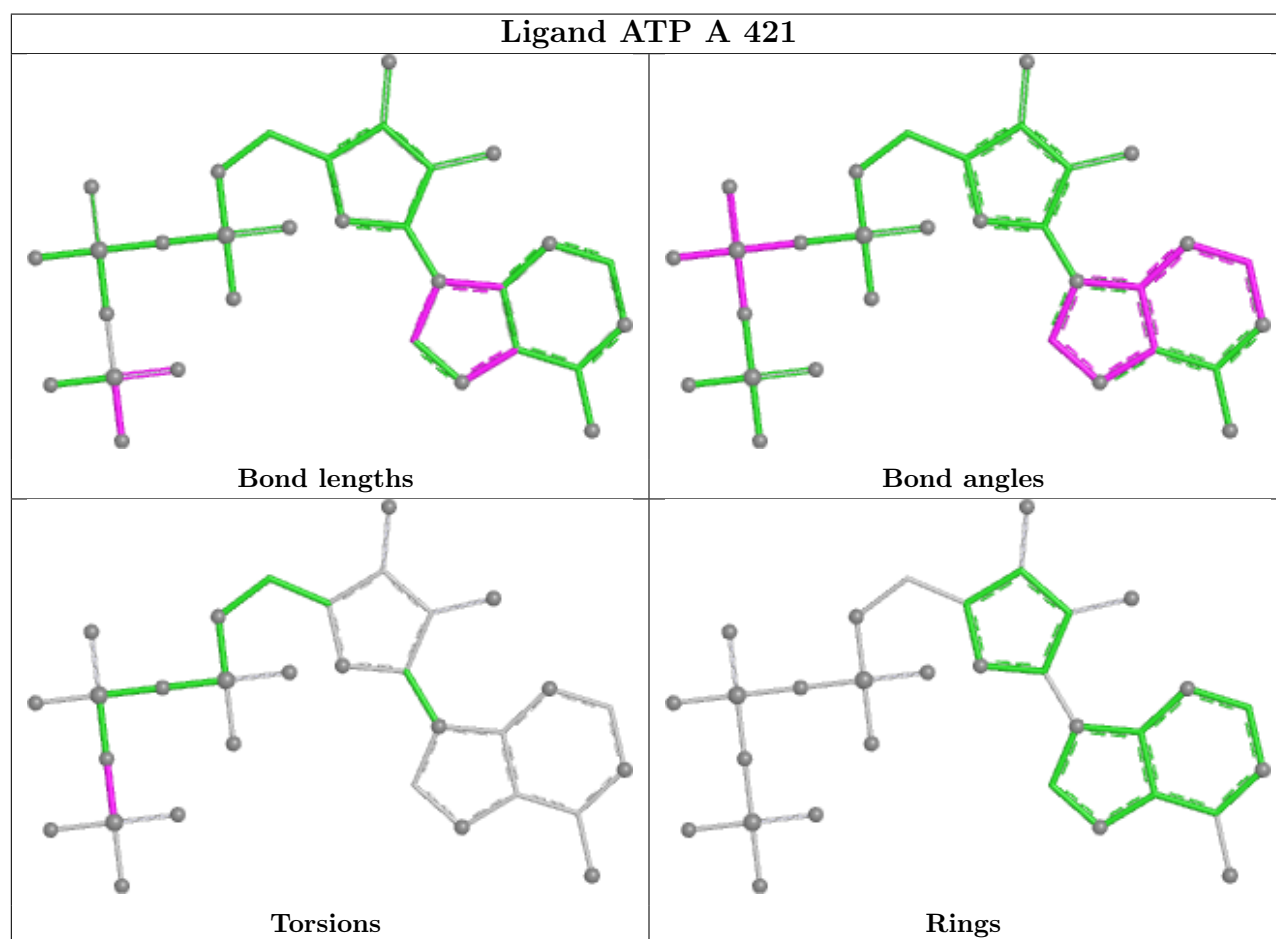
Mol	Chain	Res	Type	Atoms
3	A	421	ATP	PB-O3B-PG-O2G

There are no ring outliers.

1 monomer is involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	421	ATP	13	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	405/440 (92%)	0.27	15 (3%) 45 37	34, 59, 91, 114	0
2	B	78/78 (100%)	0.80	10 (12%) 7 6	51, 80, 135, 143	0
All	All	483/518 (93%)	0.36	25 (5%) 33 26	34, 62, 112, 143	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	373	CYS	7.2
1	A	352	CYS	5.4
1	A	356	GLY	4.9
1	A	370	CYS	3.8
2	B	18	A	3.0
1	A	375	THR	3.0
1	A	369	ARG	2.9
2	B	21(B)	C	2.9
2	B	61	C	2.8
2	B	55	U	2.7
2	B	19	G	2.6
1	A	419	PHE	2.6
1	A	400	ALA	2.5
2	B	16	C	2.5
2	B	20	G	2.5
2	B	58	A	2.4
2	B	21(A)	U	2.4
1	A	0	HIS	2.3
1	A	363	GLY	2.3
1	A	374	ARG	2.2
1	A	371	LYS	2.2
2	B	56	C	2.1
1	A	53	THR	2.1
1	A	368	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	361	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	TPO	A	18	11/12	0.97	0.07	38,42,53,59	0

6.3 Carbohydrates [i](#)

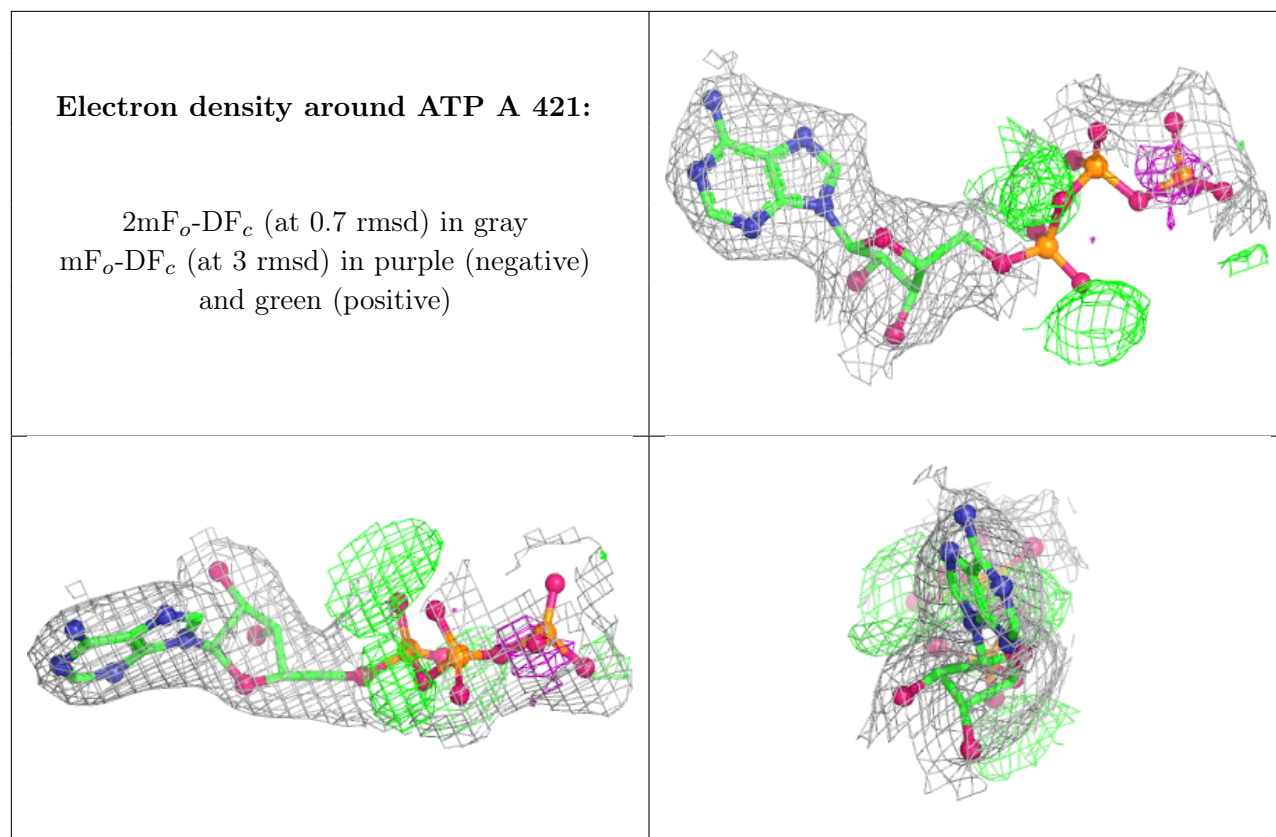
There are no oligosaccharides in this entry.

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	ATP	A	421	31/31	0.94	0.09	38,42,56,58	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.