



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:02 PM GMT

PDB ID : 1N77
Title : Crystal structure of Thermus thermophilus glutamyl-tRNA synthetase complexed with tRNA(Glu) and ATP.
Authors : Sekine, S.; Nureki, O.; Dubois, D.Y.; Bernier, S.; Chenevert, R.; Lapointe, J.; Vassilyev, D.G.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2002-11-13
Resolution : 2.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

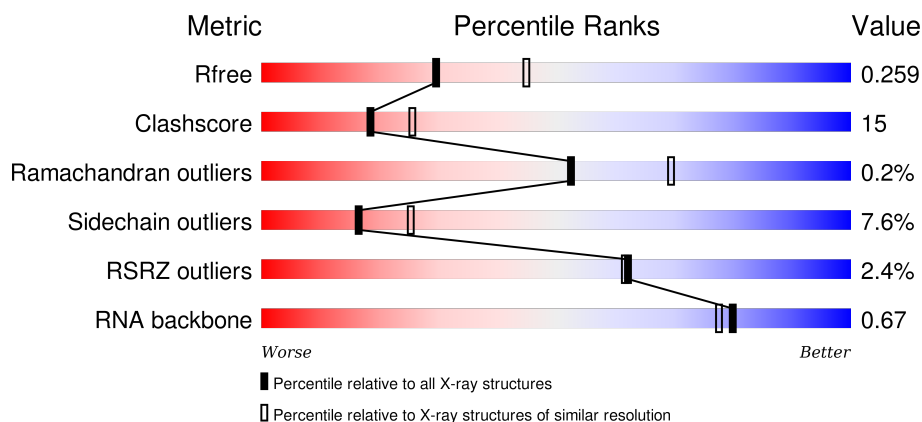
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)
RNA backbone	2183	1073 (2.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	C	75	<div> <div>52%</div> <div>29%</div> <div>13%</div> <div>5%</div> </div>
1	D	75	<div>8%</div> <div>44%</div> <div>35%</div> <div>16%</div> <div>5%</div>
2	A	468	<div>69%</div> <div>28%</div> <div>.</div>
2	B	468	<div>3%</div> <div>63%</div> <div>33%</div> <div>.</div>

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	ATP	A	600	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11391 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 tRNA(Glu).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	75	Total	C	N	O	P	0	0	0
			1597	711	284	527	75			
1	D	75	Total	C	N	O	P	0	0	0
			1597	711	284	527	75			

- Molecule 2 is a protein called Glutamyl-tRNA synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	468	Total	C	N	O	S	0	0	0
			3814	2443	676	687	8			
2	B	468	Total	C	N	O	S	0	0	0
			3814	2443	676	687	8			

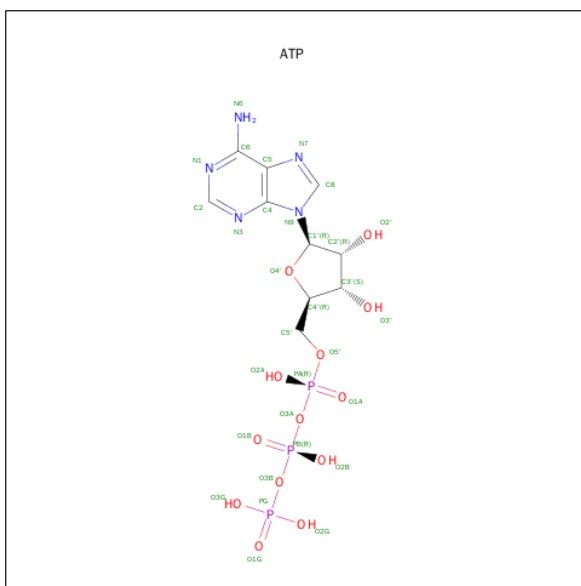
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	74	GLY	ALA	SEE REMARK 999	UNP P27000
A	75	GLY	ALA	SEE REMARK 999	UNP P27000
A	77	HIS	THR	SEE REMARK 999	UNP P27000
B	74	GLY	ALA	SEE REMARK 999	UNP P27000
B	75	GLY	ALA	SEE REMARK 999	UNP P27000
B	77	HIS	THR	SEE REMARK 999	UNP P27000

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula:

$$\text{C}_{10}\text{H}_{16}\text{N}_5\text{O}_{13}\text{P}_3).$$


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

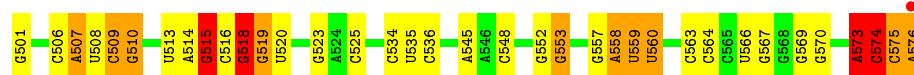
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	175	Total O 175 175	0	0
5	B	185	Total O 185 185	0	0
5	C	70	Total O 70 70	0	0
5	D	75	Total O 75 75	0	0

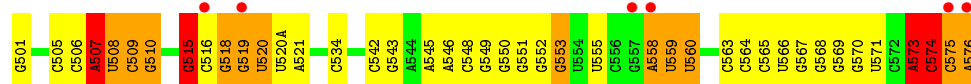
3 Residue-property plots

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.

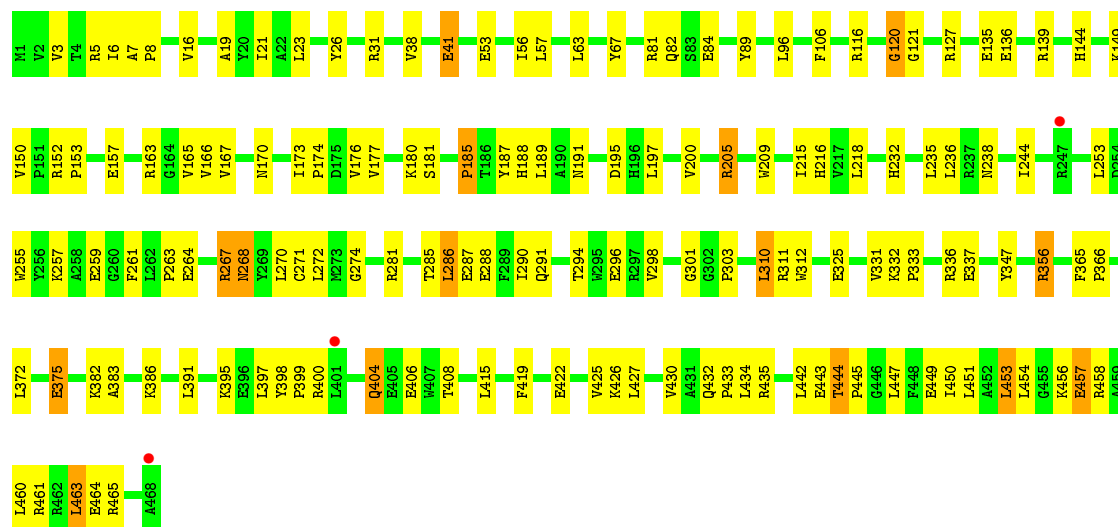
• Molecule 1: tRNA(Glu)



• Molecule 1: tRNA(Glu)

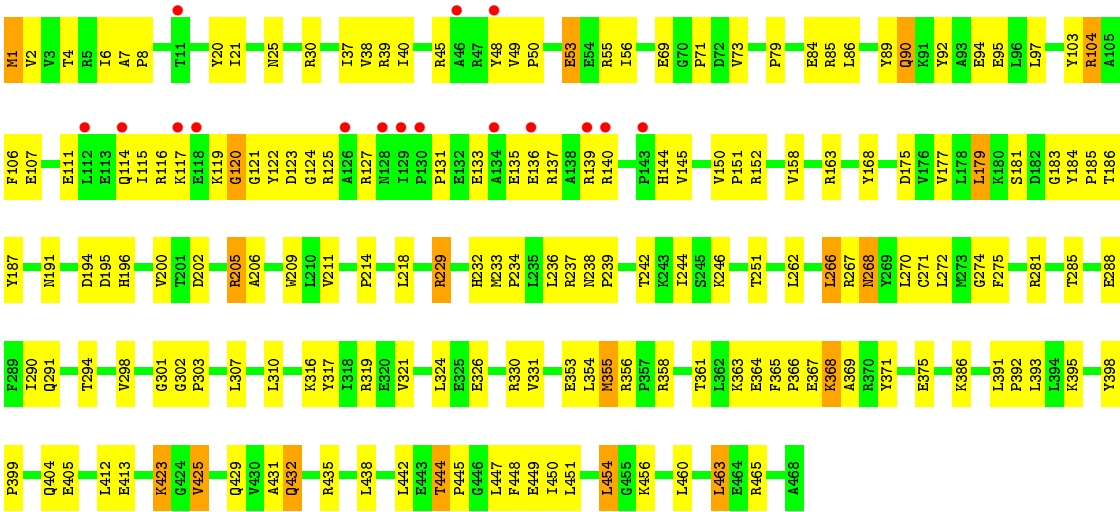


• Molecule 2: Glutamyl-tRNA synthetase



• Molecule 2: Glutamyl-tRNA synthetase





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	110.69Å 219.12Å 135.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.40 – 2.40 49.40 – 2.38	Depositor EDS
% Data completeness (in resolution range)	91.4 (49.40-2.40) 90.1 (49.40-2.38)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 2.37Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.213 , 0.257 0.213 , 0.259	Depositor DCC
R_{free} test set	2972 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	31.7	Xtriage
Anisotropy	0.262	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 46.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 59840 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11391	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.93% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: MG, 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	C	0.40	1/1782 (0.1%)	0.83	7/2774 (0.3%)
1	D	0.41	1/1782 (0.1%)	0.84	10/2774 (0.4%)
2	A	0.39	0/3910	0.62	0/5293
2	B	0.38	0/3910	0.60	0/5293
All	All	0.39	2/11384 (0.0%)	0.70	17/16134 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
1	D	1	1
All	All	1	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	501	G	OP3-P	-7.07	1.52	1.61
1	D	501	G	OP3-P	-6.92	1.52	1.61

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	574	C	C2'-C3'-O3'	8.50	128.20	109.50
1	D	574	C	C2'-C3'-O3'	7.52	126.04	109.50
1	C	518	G	C2'-C3'-O3'	7.40	125.78	109.50
1	C	519	G	C2'-C3'-O3'	7.32	125.60	109.50
1	D	515	G	C2'-C3'-O3'	7.30	125.57	109.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	518	G	C2'-C3'-O3'	7.25	125.45	109.50
1	D	507	A	C2'-C3'-O3'	7.24	125.42	109.50
1	C	573	A	C2'-C3'-O3'	7.06	125.04	109.50
1	D	519	G	C2'-C3'-O3'	7.01	124.92	109.50
1	D	573	A	C2'-C3'-O3'	6.90	124.74	113.70
1	D	515	G	C4'-C3'-O3'	6.00	124.99	113.00
1	C	515	G	C4'-C3'-O3'	5.97	124.93	113.00
1	D	515	G	C4'-C3'-C2'	5.79	108.39	102.60
1	D	507	A	C4'-C3'-C2'	5.55	108.16	102.60
1	C	574	C	C4'-C3'-C2'	5.20	107.80	102.60
1	D	573	A	C4'-C3'-C2'	5.16	107.76	102.60
1	C	501	G	OP1-P-OP2	-5.03	112.05	119.60

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	D	515	G	C3'

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	515	G	Sidechain
1	D	515	G	Sidechain

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1597	0	813	28	0
1	D	1597	0	813	43	0
2	A	3814	0	3818	125	0
2	B	3814	0	3818	137	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	31	0	12	4	0
4	B	31	0	12	4	0
5	A	175	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	185	0	0	8	0
5	C	70	0	0	0	0
5	D	75	0	0	6	0
All	All	11391	0	9286	312	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 15.

All (312) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:506:C:O2'	1:D:507:A:H5'	1.72	0.89
2:B:454:LEU:H	2:B:454:LEU:HD22	1.38	0.89
1:D:576:A:H3'	2:B:209:TRP:HH2	1.39	0.86
2:A:205:ARG:HD3	2:A:232:HIS:CE1	2.10	0.85
2:B:454:LEU:N	2:B:454:LEU:HD22	1.91	0.85
2:B:205:ARG:HD3	2:B:232:HIS:CE1	2.11	0.84
2:A:294:THR:HG22	2:A:296:GLU:H	1.41	0.84
1:C:574:C:H5'	2:A:177:VAL:HG11	1.60	0.82
2:B:4:THR:HB	2:B:25:ASN:HD22	1.43	0.81
2:A:136:GLU:HG2	2:A:139:ARG:HH21	1.46	0.81
1:C:576:A:H5''	2:A:187:TYR:HB2	1.62	0.81
2:B:116:ARG:HA	2:B:121:GLY:H	1.46	0.79
2:A:460:LEU:O	2:A:464:GLU:HG3	1.84	0.77
2:B:432:GLN:HA	2:B:432:GLN:HE21	1.48	0.77
2:A:365:PHE:HB3	2:A:366:PRO:HD3	1.67	0.76
2:B:262:LEU:HD22	2:B:330:ARG:HH11	1.50	0.75
2:B:365:PHE:HB3	2:B:366:PRO:HD3	1.67	0.75
2:B:435:ARG:HD3	2:B:444:THR:HG22	1.69	0.74
2:A:435:ARG:HD3	2:A:444:THR:HG22	1.69	0.73
1:D:576:A:H5''	2:B:187:TYR:HB2	1.69	0.73
2:A:267:ARG:HD3	5:A:1690:HOH:O	1.88	0.73
1:C:574:C:H5'	2:A:177:VAL:CG1	2.18	0.73
5:D:1313:HOH:O	2:B:316:LYS:HE2	1.88	0.72
2:B:354:LEU:HD22	2:B:449:GLU:HG3	1.70	0.72
2:A:285:THR:HG23	2:A:288:GLU:OE1	1.90	0.72
2:B:432:GLN:HA	2:B:432:GLN:NE2	2.05	0.71
2:B:454:LEU:CD2	2:B:454:LEU:H	2.04	0.70
2:A:267:ARG:HD2	2:A:286:LEU:HG	1.74	0.70
2:A:375:GLU:HG3	2:A:465:ARG:HH22	1.57	0.70
1:D:573:A:H5'	5:D:1271:HOH:O	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:506:C:O2'	1:C:507:A:H5'	1.92	0.69
2:A:157:GLU:HG2	2:A:167:VAL:HG22	1.73	0.69
2:B:361:THR:OG1	2:B:364:GLU:HG3	1.93	0.69
2:A:205:ARG:HD3	2:A:232:HIS:HE1	1.53	0.69
1:D:576:A:H3'	2:B:209:TRP:CH2	2.25	0.67
2:A:188:HIS:HD2	2:A:216:HIS:HE1	1.41	0.67
4:B:601:ATP:H2'	5:B:1456:HOH:O	1.94	0.67
2:B:326:GLU:HG3	5:B:1790:HOH:O	1.94	0.66
2:B:413:GLU:HG3	2:B:448:PHE:CZ	2.31	0.66
2:A:257:LYS:HE3	5:A:1017:HOH:O	1.95	0.66
2:B:465:ARG:HH11	2:B:465:ARG:HB3	1.61	0.66
1:D:576:A:C5'	2:B:187:TYR:HB2	2.25	0.65
2:A:445:PRO:HG2	2:A:450:ILE:HD11	1.77	0.65
2:B:444:THR:HG23	2:B:445:PRO:O	1.97	0.65
2:A:444:THR:HG23	2:A:445:PRO:O	1.97	0.65
2:A:461:ARG:NH1	2:A:465:ARG:HD3	2.11	0.64
1:D:570:G:H21	2:B:211:VAL:HG21	1.63	0.64
2:B:432:GLN:CA	2:B:432:GLN:HE21	2.06	0.64
2:B:205:ARG:HD3	2:B:232:HIS:HE1	1.60	0.63
2:B:6:ILE:CD1	2:B:21:ILE:HG22	2.29	0.62
2:B:25:ASN:ND2	5:B:1786:HOH:O	2.31	0.62
2:A:255:TRP:CZ2	2:A:259:GLU:HG3	2.35	0.62
2:B:122:TYR:CE2	2:B:124:GLY:HA2	2.34	0.62
2:B:391:LEU:N	2:B:392:PRO:HD2	2.15	0.62
1:D:574:C:C5'	2:B:177:VAL:HG11	2.30	0.61
2:B:454:LEU:CD2	2:B:454:LEU:N	2.62	0.61
2:A:400:ARG:HH22	2:A:422:GLU:CD	2.03	0.61
1:D:573:A:H2'	5:D:1136:HOH:O	1.99	0.61
5:D:1024:HOH:O	2:B:211:VAL:HG23	2.00	0.61
2:A:453:LEU:HB3	2:A:454:LEU:HD22	1.81	0.61
2:A:6:ILE:HD11	2:A:38:VAL:HG22	1.81	0.61
2:A:188:HIS:HD2	2:A:216:HIS:CE1	2.19	0.60
2:B:435:ARG:HD3	2:B:444:THR:CG2	2.32	0.60
1:D:574:C:H5'	2:B:177:VAL:HG11	1.82	0.60
2:A:5:ARG:HH21	2:A:191:ASN:ND2	2.00	0.60
1:C:509:C:H5'	1:C:510:G:OP2	2.01	0.59
2:A:287:GLU:O	2:A:291:GLN:HG3	2.03	0.59
2:B:398:TYR:HB3	2:B:399:PRO:HD3	1.84	0.59
1:D:569:G:O2'	1:D:570:G:H5'	2.01	0.59
2:B:90:GLN:O	2:B:94:GLU:HG3	2.03	0.59
2:A:21:ILE:HG13	4:A:600:ATP:H1'	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:238:ASN:HB3	2:B:239:PRO:HD2	1.85	0.58
2:A:205:ARG:HG3	2:A:209:TRP:CD1	2.39	0.58
1:D:569:G:H5''	2:B:242:THR:HA	1.86	0.58
2:B:106:PHE:CD2	2:B:144:HIS:HB3	2.39	0.58
2:A:53:GLU:O	2:A:56:ILE:HG22	2.03	0.58
1:D:574:C:O2'	1:D:575:C:OP1	2.20	0.57
2:B:6:ILE:HG23	2:B:8:PRO:HD3	1.85	0.57
2:B:158:VAL:HG21	2:B:214:PRO:HG3	1.86	0.57
2:A:383:ALA:HA	2:A:442:LEU:HD13	1.87	0.57
2:A:435:ARG:HD3	2:A:444:THR:CG2	2.34	0.57
2:A:435:ARG:HH11	2:A:444:THR:HG22	1.68	0.57
2:A:445:PRO:CG	2:A:450:ILE:HD11	2.34	0.57
1:D:573:A:H3'	5:D:1161:HOH:O	2.04	0.57
2:A:454:LEU:HD22	2:A:454:LEU:N	2.18	0.57
2:A:333:PRO:HA	2:A:336:ARG:NH1	2.19	0.57
1:C:563:C:H2'	1:C:564:C:C6	2.40	0.57
2:B:425:VAL:HG22	2:B:429:GLN:OE1	2.06	0.56
2:A:116:ARG:O	2:A:120:GLY:HA2	2.05	0.56
2:B:116:ARG:O	2:B:120:GLY:HA2	2.06	0.56
2:A:152:ARG:HH11	2:A:152:ARG:HG2	1.70	0.56
2:A:31:ARG:HD3	2:A:290:ILE:O	2.06	0.55
2:A:116:ARG:HA	2:A:121:GLY:H	1.70	0.55
2:B:363:LYS:O	2:B:367:GLU:HG3	2.06	0.55
2:B:37:ILE:HG12	2:B:69:GLU:HB2	1.88	0.55
2:B:79:PRO:HB3	2:B:84:GLU:OE1	2.07	0.55
2:A:404:GLN:HG2	2:A:415:LEU:HD22	1.89	0.55
2:A:8:PRO:HA	4:A:600:ATP:O2A	2.07	0.55
2:A:7:ALA:HB1	2:A:41:GLU:CD	2.26	0.55
2:B:111:GLU:O	2:B:115:ILE:HD13	2.07	0.55
2:A:427:LEU:O	2:A:430:VAL:HG12	2.07	0.55
2:B:111:GLU:O	2:B:114:GLN:HB2	2.07	0.54
2:A:274:GLY:O	2:A:298:VAL:HA	2.07	0.54
2:B:136:GLU:HA	2:B:139:ARG:HD2	1.88	0.54
2:A:301:GLY:O	2:A:303:PRO:HD3	2.08	0.54
2:B:413:GLU:HG3	2:B:448:PHE:CE2	2.43	0.53
2:A:5:ARG:HH21	2:A:191:ASN:HD21	1.55	0.53
2:B:133:GLU:O	2:B:137:ARG:HG3	2.09	0.53
2:B:445:PRO:HG2	2:B:450:ILE:HD11	1.90	0.53
2:A:435:ARG:HH11	2:A:444:THR:CG2	2.21	0.53
2:A:5:ARG:NH2	2:A:191:ASN:ND2	2.57	0.53
1:C:563:C:H2'	1:C:564:C:H6	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:358:ARG:O	2:B:368:LYS:HE2	2.09	0.53
2:B:391:LEU:O	2:B:395:LYS:HG3	2.08	0.53
1:D:563:C:H2'	1:D:564:C:C6	2.44	0.53
2:B:237:ARG:NH1	2:B:302:GLY:HA3	2.23	0.53
1:D:558:A:H4'	1:D:559:U:OP1	2.09	0.53
2:A:375:GLU:HG3	2:A:465:ARG:NH2	2.24	0.52
2:A:170:ASN:HA	2:A:173:ILE:HD12	1.91	0.52
2:B:331:VAL:HG11	2:B:365:PHE:CD2	2.44	0.52
2:B:244:ILE:HG12	2:B:251:THR:HG22	1.91	0.52
2:A:6:ILE:C	2:A:8:PRO:HD3	2.30	0.52
2:B:425:VAL:HG22	2:B:429:GLN:CD	2.29	0.52
1:D:555:U:H5''	5:D:1491:HOH:O	2.10	0.52
2:B:454:LEU:HG	5:B:1127:HOH:O	2.08	0.52
2:B:131:PRO:O	2:B:135:GLU:HG2	2.10	0.52
2:B:21:ILE:HD13	4:B:601:ATP:N3	2.26	0.51
2:A:391:LEU:O	2:A:395:LYS:HG2	2.10	0.51
2:A:406:GLU:HG2	2:A:408:THR:HG23	1.93	0.51
2:A:163:ARG:NH2	2:A:232:HIS:O	2.43	0.51
1:D:552:G:H5'	1:D:553:G:OP2	2.11	0.51
2:A:268:ASN:C	2:A:268:ASN:HD22	2.12	0.51
2:B:431:ALA:CB	2:B:447:LEU:HD22	2.41	0.51
2:A:152:ARG:NH1	2:A:152:ARG:HG2	2.26	0.51
2:A:26:TYR:CE1	2:A:290:ILE:HD11	2.45	0.51
2:A:41:GLU:HG3	2:A:82:GLN:OE1	2.11	0.51
1:D:549:G:N2	1:D:566:U:H1'	2.26	0.51
1:D:563:C:H2'	1:D:564:C:H6	1.75	0.51
2:B:274:GLY:O	2:B:298:VAL:HA	2.10	0.51
1:C:558:A:O2'	1:C:560:U:H5	1.93	0.51
2:A:81:ARG:HD2	2:A:84:GLU:OE2	2.11	0.51
2:B:431:ALA:HB1	2:B:447:LEU:HD22	1.92	0.50
1:D:509:C:O2'	1:D:545:A:H2'	2.10	0.50
2:A:444:THR:OG1	2:A:445:PRO:HD2	2.11	0.50
1:C:552:G:H5'	1:C:553:G:OP2	2.11	0.50
2:A:144:HIS:O	2:A:181:SER:HA	2.10	0.50
1:C:518:G:O2'	1:C:557:G:N2	2.33	0.50
1:C:574:C:C4'	2:A:177:VAL:HG21	2.41	0.50
2:A:347:TYR:CE1	2:A:453:LEU:HD22	2.46	0.50
2:B:179:LEU:HD21	2:B:183:GLY:O	2.12	0.50
1:D:565:C:O2'	1:D:566:U:H5'	2.12	0.50
2:A:81:ARG:NH1	2:A:84:GLU:OE2	2.38	0.50
2:B:152:ARG:HH11	2:B:152:ARG:HG2	1.77	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:558:A:O2'	1:D:560:U:H5	1.94	0.50
2:B:244:ILE:HG22	4:B:601:ATP:HN62	1.77	0.49
2:A:106:PHE:CD1	2:A:106:PHE:N	2.80	0.49
2:A:136:GLU:HG2	2:A:139:ARG:NH2	2.22	0.49
1:C:569:G:O2'	1:C:570:G:H5'	2.12	0.49
1:D:506:C:C2'	1:D:507:A:H5'	2.42	0.49
2:A:386:LYS:HE2	2:A:432:GLN:HG2	1.95	0.49
2:A:347:TYR:CE1	2:A:453:LEU:CD2	2.96	0.49
2:B:438:LEU:HD11	2:B:451:LEU:HD13	1.94	0.48
2:B:355:MET:HE2	2:B:369:ALA:HB2	1.95	0.48
1:C:575:C:OP2	2:A:180:LYS:HE2	2.13	0.48
2:B:301:GLY:O	2:B:303:PRO:HD3	2.13	0.48
1:D:509:C:H5'	1:D:510:G:OP2	2.14	0.48
2:A:383:ALA:HA	2:A:442:LEU:CD1	2.44	0.48
1:D:550:G:H2'	1:D:551:G:C8	2.49	0.48
2:B:145:VAL:HG12	2:B:181:SER:N	2.28	0.48
1:D:508:U:O5'	1:D:508:U:H6	1.97	0.48
2:B:30:ARG:HB3	2:B:290:ILE:HD12	1.96	0.48
1:C:534:C:O2	2:A:435:ARG:NH2	2.43	0.48
2:B:211:VAL:O	2:B:214:PRO:HD2	2.13	0.48
2:A:332:LYS:HB2	2:A:333:PRO:HD3	1.95	0.48
2:B:393:LEU:HD21	2:B:423:LYS:HG2	1.96	0.47
2:A:89:TYR:CD2	2:A:185:PRO:HG3	2.49	0.47
2:B:73:VAL:O	2:B:73:VAL:HG12	2.15	0.47
2:B:246:LYS:HG2	5:B:1420:HOH:O	2.14	0.47
2:A:23:LEU:HD21	2:A:271:CYS:SG	2.55	0.47
2:B:444:THR:OG1	2:B:445:PRO:HD2	2.14	0.47
1:C:576:A:H2'	2:A:209:TRP:CZ2	2.50	0.47
2:A:7:ALA:HB1	2:A:41:GLU:OE2	2.14	0.47
1:C:536:C:H4'	2:A:443:GLU:HG2	1.96	0.47
2:A:281:ARG:HE	2:A:281:ARG:HB3	1.47	0.47
2:B:45:ARG:HD3	2:B:184:TYR:CE2	2.49	0.47
2:A:398:TYR:N	2:A:399:PRO:HD2	2.30	0.46
1:C:573:A:O2'	1:C:574:C:H3'	2.15	0.46
2:B:89:TYR:CD2	2:B:185:PRO:HG3	2.51	0.46
1:C:558:A:O2'	1:C:559:U:P	2.74	0.46
1:D:573:A:O2'	1:D:574:C:H3'	2.16	0.46
2:A:19:ALA:HB1	2:A:63:LEU:CD1	2.46	0.46
2:A:5:ARG:NH1	5:A:1521:HOH:O	2.46	0.46
2:A:195:ASP:HB3	2:A:200:VAL:HG22	1.98	0.46
2:A:294:THR:CG2	2:A:296:GLU:HB2	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:107:GLU:N	2:B:107:GLU:OE2	2.39	0.46
2:A:445:PRO:HG2	2:A:450:ILE:CG1	2.46	0.46
2:A:127:ARG:HH11	2:A:149:LYS:HE2	1.82	0.46
1:C:509:C:O2'	1:C:545:A:C2'	2.64	0.45
2:A:106:PHE:CD2	2:A:144:HIS:HB3	2.52	0.45
2:B:294:THR:HG23	2:B:294:THR:O	2.16	0.45
2:A:176:VAL:HG11	2:A:215:ILE:HD13	1.99	0.45
2:A:375:GLU:CG	2:A:465:ARG:HH12	2.29	0.45
2:A:235:LEU:HD13	4:A:600:ATP:C2	2.51	0.45
2:B:268:ASN:C	2:B:268:ASN:HD22	2.17	0.45
2:A:165:VAL:HG12	2:A:166:VAL:N	2.32	0.45
1:D:570:G:H2'	1:D:571:U:C6	2.51	0.45
2:A:191:ASN:HA	5:A:1685:HOH:O	2.15	0.45
2:B:136:GLU:O	2:B:140:ARG:HG2	2.17	0.45
1:D:558:A:O2'	1:D:559:U:P	2.73	0.45
2:A:325:GLU:H	2:A:325:GLU:CD	2.19	0.45
2:A:432:GLN:N	2:A:433:PRO:HD2	2.32	0.45
2:B:423:LYS:HB2	2:B:423:LYS:HZ2	1.82	0.45
2:B:85:ARG:NH1	2:B:194:ASP:OD2	2.37	0.45
2:A:3:VAL:HB	2:A:200:VAL:HA	1.98	0.45
1:C:566:U:O2'	1:C:567:G:H5'	2.17	0.45
2:A:5:ARG:NH2	2:A:191:ASN:HD21	2.12	0.45
2:B:206:ALA:HB3	4:B:601:ATP:O2'	2.17	0.45
2:A:372:LEU:HD22	2:A:453:LEU:HD13	1.99	0.45
2:A:337:GLU:O	2:A:337:GLU:HG2	2.16	0.45
1:D:534:C:O2	2:B:435:ARG:NH2	2.40	0.44
2:B:386:LYS:HE2	2:B:432:GLN:HG2	1.99	0.44
2:B:371:TYR:HH	2:B:444:THR:HG1	1.63	0.44
1:D:508:U:H2'	1:D:546:A:N3	2.32	0.44
2:B:55:ARG:HD3	5:B:1204:HOH:O	2.17	0.44
2:B:103:TYR:CE2	2:B:127:ARG:HG2	2.52	0.44
2:B:354:LEU:CD2	2:B:449:GLU:HG3	2.41	0.44
2:A:434:LEU:HD23	2:A:447:LEU:HD21	1.99	0.44
1:C:509:C:H5	1:C:523:G:O6	2.00	0.44
2:A:261:PHE:CE2	2:A:310:LEU:HD13	2.53	0.44
2:B:324:LEU:HD11	2:B:353:GLU:HB2	2.00	0.44
2:B:444:THR:HG23	2:B:445:PRO:N	2.32	0.44
2:B:115:ILE:HD12	2:B:115:ILE:N	2.33	0.44
2:B:40:ILE:HG12	2:B:71:PRO:HG2	1.98	0.44
1:D:505:C:H4'	2:B:163:ARG:HD3	2.00	0.44
1:D:542:C:O2'	1:D:543:G:H5'	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:285:THR:OG1	2:B:288:GLU:HG3	2.18	0.44
2:A:135:GLU:O	2:A:139:ARG:HG3	2.18	0.43
2:B:116:ARG:HA	2:B:121:GLY:N	2.24	0.43
1:D:550:G:H2'	1:D:551:G:H8	1.82	0.43
2:B:53:GLU:O	2:B:56:ILE:HG22	2.19	0.43
2:B:463:LEU:HD23	2:B:463:LEU:O	2.18	0.43
2:B:317:TYR:HA	2:B:321:VAL:HG23	2.01	0.43
2:A:445:PRO:HG2	2:A:450:ILE:CD1	2.44	0.43
2:A:56:ILE:HG23	2:A:57:LEU:N	2.33	0.43
2:B:262:LEU:HD22	2:B:330:ARG:NH1	2.26	0.43
2:A:397:LEU:HD13	2:A:419:PHE:CD2	2.53	0.43
2:A:187:TYR:CD2	2:A:209:TRP:CZ3	3.06	0.43
2:A:255:TRP:CH2	2:A:259:GLU:HG3	2.53	0.43
2:B:115:ILE:HG13	2:B:125:ARG:NH2	2.34	0.43
2:B:39:ARG:NH2	2:B:195:ASP:OD2	2.34	0.43
1:D:520:U:H2'	1:D:521:A:H5'	2.01	0.43
2:B:432:GLN:HE22	2:B:435:ARG:HH21	1.65	0.42
2:B:450:ILE:HA	2:B:450:ILE:HD13	1.84	0.42
2:A:150:VAL:O	2:A:152:ARG:HG3	2.19	0.42
1:D:558:A:O2'	1:D:559:U:OP1	2.35	0.42
2:A:261:PHE:CD2	2:A:310:LEU:HD13	2.53	0.42
2:B:404:GLN:NE2	2:B:404:GLN:HA	2.34	0.42
2:A:456:LYS:HG2	2:A:457:GLU:OE1	2.18	0.42
1:D:576:A:H5'	2:B:187:TYR:HB2	2.01	0.42
2:B:196:HIS:HD2	5:B:1115:HOH:O	2.01	0.42
2:B:127:ARG:HH21	2:B:175:ASP:CG	2.23	0.42
2:A:294:THR:HG22	2:A:296:GLU:HB2	2.00	0.42
2:B:319:ARG:HD2	5:B:1045:HOH:O	2.19	0.42
2:B:465:ARG:NH1	2:B:465:ARG:HB3	2.29	0.42
2:B:123:ASP:OD1	2:B:125:ARG:HB3	2.20	0.42
2:A:311:ARG:HB3	2:A:311:ARG:HH11	1.83	0.42
2:A:463:LEU:HD23	2:A:463:LEU:O	2.20	0.42
1:D:574:C:C4'	2:B:177:VAL:HG11	2.50	0.42
2:B:1:MET:HB2	2:B:2:VAL:H	1.57	0.42
2:B:177:VAL:HG13	2:B:186:THR:HG21	2.02	0.41
2:B:20:TYR:CD2	2:B:21:ILE:HD12	2.55	0.41
1:C:525:C:H5''	2:A:312:TRP:CE2	2.55	0.41
2:B:233:MET:HA	2:B:234:PRO:HD3	1.93	0.41
2:B:326:GLU:O	2:B:330:ARG:HG3	2.20	0.41
1:D:564:C:O2'	1:D:565:C:H5'	2.20	0.41
2:A:19:ALA:HB1	2:A:63:LEU:HD11	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:263:PRO:HG2	2:A:264:GLU:OE1	2.20	0.41
2:B:202:ASP:OD2	2:B:229:ARG:NH1	2.54	0.41
2:B:92:TYR:O	2:B:95:GLU:HB3	2.20	0.41
2:B:266:LEU:O	2:B:270:LEU:HD13	2.20	0.41
2:B:45:ARG:O	2:B:48:TYR:HB3	2.20	0.41
1:D:567:G:H2'	1:D:568:G:H8	1.85	0.41
2:A:174:PRO:HG2	2:A:176:VAL:HG13	2.03	0.41
2:B:195:ASP:HB3	2:B:200:VAL:CG2	2.50	0.41
2:B:195:ASP:CG	2:B:200:VAL:HG21	2.41	0.41
2:B:456:LYS:O	2:B:460:LEU:HG	2.21	0.41
1:C:573:A:H5''	1:C:573:A:H8	1.86	0.41
2:A:7:ALA:HB1	2:A:41:GLU:OE1	2.20	0.41
2:B:111:GLU:O	2:B:115:ILE:CD1	2.69	0.41
1:C:535:U:C4'	2:A:432:GLN:HE22	2.33	0.41
2:B:7:ALA:HA	2:B:39:ARG:O	2.21	0.41
2:A:356:ARG:NH2	5:A:1432:HOH:O	2.53	0.41
2:B:271:CYS:O	2:B:275:PHE:HB3	2.21	0.41
1:C:513:U:O2'	1:C:514:A:H5'	2.20	0.41
1:C:576:A:H3'	2:A:209:TRP:HH2	1.86	0.41
2:A:382:LYS:HG3	5:A:1356:HOH:O	2.20	0.41
2:B:375:GLU:HG3	2:B:465:ARG:HH21	1.85	0.40
2:B:168:TYR:CD2	2:B:214:PRO:HG3	2.55	0.40
2:B:391:LEU:N	2:B:392:PRO:CD	2.84	0.40
1:D:520(A):U:O5'	1:D:520(A):U:H6	2.03	0.40
2:A:16:VAL:HA	2:A:253:LEU:HD21	2.03	0.40
2:B:150:VAL:HA	2:B:151:PRO:HD3	1.91	0.40
1:C:558:A:O2'	1:C:559:U:OP1	2.30	0.40
1:D:509:C:O2'	1:D:545:A:C2'	2.70	0.40
2:A:425:VAL:HG22	2:A:426:LYS:N	2.37	0.40
2:B:97:LEU:HD11	2:B:104:ARG:HG2	2.04	0.40
1:C:575:C:H2'	1:C:576:A:O4'	2.22	0.40
2:B:6:ILE:HG22	2:B:38:VAL:HA	2.03	0.40
2:A:454:LEU:CD2	2:A:454:LEU:N	2.82	0.40
4:A:600:ATP:O1G	4:A:600:ATP:O1B	2.40	0.40
2:B:49:VAL:HA	2:B:50:PRO:HD3	1.91	0.40
2:A:238:ASN:CG	2:A:244:ILE:HA	2.42	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	
2	A	466/468 (100%)	450 (97%)	15 (3%)	1 (0%)	52	69
2	B	466/468 (100%)	449 (96%)	16 (3%)	1 (0%)	52	69
All	All	932/936 (100%)	899 (96%)	31 (3%)	2 (0%)	52	69

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	120	GLY
2	B	120	GLY

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	
2	A	393/393 (100%)	366 (93%)	27 (7%)	19	30
2	B	393/393 (100%)	360 (92%)	33 (8%)	14	20
All	All	786/786 (100%)	726 (92%)	60 (8%)	16	25

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	41	GLU
2	A	67	TYR
2	A	96	LEU
2	A	153	PRO

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Mol	Chain	Res	Type
2	A	185	PRO
2	A	189	LEU
2	A	197	LEU
2	A	205	ARG
2	A	218	LEU
2	A	236	LEU
2	A	267	ARG
2	A	268	ASN
2	A	270	LEU
2	A	272	LEU
2	A	286	LEU
2	A	310	LEU
2	A	331	VAL
2	A	356	ARG
2	A	375	GLU
2	A	404	GLN
2	A	444	THR
2	A	449	GLU
2	A	451	LEU
2	A	453	LEU
2	A	457	GLU
2	A	458	ARG
2	A	463	LEU
2	B	1	MET
2	B	53	GLU
2	B	86	LEU
2	B	90	GLN
2	B	104	ARG
2	B	117	LYS
2	B	119	LYS
2	B	179	LEU
2	B	191	ASN
2	B	205	ARG
2	B	218	LEU
2	B	229	ARG
2	B	236	LEU
2	B	266	LEU
2	B	267	ARG
2	B	268	ASN
2	B	272	LEU
2	B	281	ARG
2	B	291	GLN

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Mol	Chain	Res	Type
2	B	307	LEU
2	B	310	LEU
2	B	355	MET
2	B	356	ARG
2	B	368	LYS
2	B	405	GLU
2	B	412	LEU
2	B	423	LYS
2	B	425	VAL
2	B	432	GLN
2	B	442	LEU
2	B	444	THR
2	B	454	LEU
2	B	463	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	90	GLN
2	A	188	HIS
2	A	191	ASN
2	A	216	HIS
2	A	232	HIS
2	A	384	GLN
2	A	429	GLN
2	B	25	ASN
2	B	171	GLN
2	B	188	HIS
2	B	196	HIS
2	B	232	HIS
2	B	404	GLN
2	B	432	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	C	74/75 (98%)	16 (21%)	9 (12%)
1	D	74/75 (98%)	14 (18%)	9 (12%)
All	All	148/150 (98%)	30 (20%)	18 (12%)

All (30) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	C	507	A
1	C	508	U
1	C	510	G
1	C	515	G
1	C	516	C
1	C	518	G
1	C	519	G
1	C	520	U
1	C	548	C
1	C	553	G
1	C	558	A
1	C	559	U
1	C	573	A
1	C	574	C
1	C	575	C
1	C	576	A
1	D	508	U
1	D	510	G
1	D	516	C
1	D	518	G
1	D	519	G
1	D	520	U
1	D	548	C
1	D	553	G
1	D	558	A
1	D	559	U
1	D	573	A
1	D	574	C
1	D	575	C
1	D	576	A

All (18) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	C	507	A
1	C	509	C
1	C	515	G
1	C	518	G
1	C	519	G
1	C	558	A
1	C	560	U
1	C	573	A

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Mol	Chain	Res	Type
1	C	574	C
1	D	507	A
1	D	509	C
1	D	515	G
1	D	518	G
1	D	519	G
1	D	558	A
1	D	560	U
1	D	573	A
1	D	574	C

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
4	ATP	A	600	-	24,33,33	0.66	0	31,52,52	1.39	5 (16%)
4	ATP	B	601	-	24,33,33	0.66	0	31,52,52	1.37	5 (16%)

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
4	ATP	A	600	-	-	0/18/38/38	0/3/3/3
4	ATP	B	601	-	-	0/18/38/38	0/3/3/3

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	600	ATP	PA-O3A-PB	-4.07	121.31	132.73
4	B	601	ATP	PA-O3A-PB	-3.97	121.57	132.73
4	B	601	ATP	PB-O3B-PG	-3.55	120.75	132.67
4	A	600	ATP	PB-O3B-PG	-3.24	121.80	132.67
4	B	601	ATP	C2'-C3'-C4'	2.15	107.03	102.61
4	B	601	ATP	O3'-C3'-C2'	2.29	119.28	111.83
4	A	600	ATP	C2'-C3'-C4'	2.31	107.35	102.61
4	B	601	ATP	O3'-C3'-C4'	2.47	118.46	111.05
4	A	600	ATP	O3'-C3'-C4'	2.62	118.91	111.05
4	A	600	ATP	O3'-C3'-C2'	2.80	120.93	111.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	600	ATP	4	0
4	B	601	ATP	4	0

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	C	75/75 (100%)	-0.32	1 (1%) 79 79	27, 48, 85, 104	0
1	D	75/75 (100%)	-0.07	6 (8%) 15 15	25, 49, 96, 113	0
2	A	468/468 (100%)	-0.38	3 (0%) 90 90	12, 27, 58, 73	0
2	B	468/468 (100%)	-0.27	16 (3%) 49 49	12, 30, 79, 93	0
All	All	1086/1086 (100%)	-0.31	26 (2%) 62 61	12, 31, 80, 113	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	576	A	4.8
2	B	129	ILE	4.2
2	B	48	TYR	3.9
2	B	11	THR	3.7
2	B	130	PRO	3.6
2	B	140	ARG	3.6
2	B	126	ALA	3.5
1	D	575	C	3.5
1	D	516	C	3.4
2	B	128	ASN	3.3
2	B	114	GLN	3.3
2	B	118	GLU	2.9
2	B	143	PRO	2.8
2	B	46	ALA	2.8
2	B	139	ARG	2.8
1	C	576	A	2.7
2	A	401	LEU	2.7
2	B	117	LYS	2.6
2	B	136	GLU	2.5
2	A	247	ARG	2.4
2	B	134	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	519	G	2.3
2	A	468	ALA	2.1
2	B	112	LEU	2.1
1	D	558	A	2.1
1	D	557	G	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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, 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	LLDF	B-factors(Å ²)	Q<0.9
4	ATP	A	600	31/31	0.71	0.30	3.18	76,86,104,105	0
4	ATP	B	601	31/31	0.80	0.29	1.47	84,94,110,111	0
3	MG	C	700	1/1	0.95	0.27	-	38,38,38,38	0
3	MG	D	701	1/1	0.97	0.27	-	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.