



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:47 PM GMT

PDB ID : 1QLN  
Title : STRUCTURE OF A TRANSCRIBING T7 RNA POLYMERASE INITIATION COMPLEX  
Authors : Cheetham, G.M.T.; Steitz, T.A.  
Deposited on : 1999-09-01  
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](mailto: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.

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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 : 1.7 (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

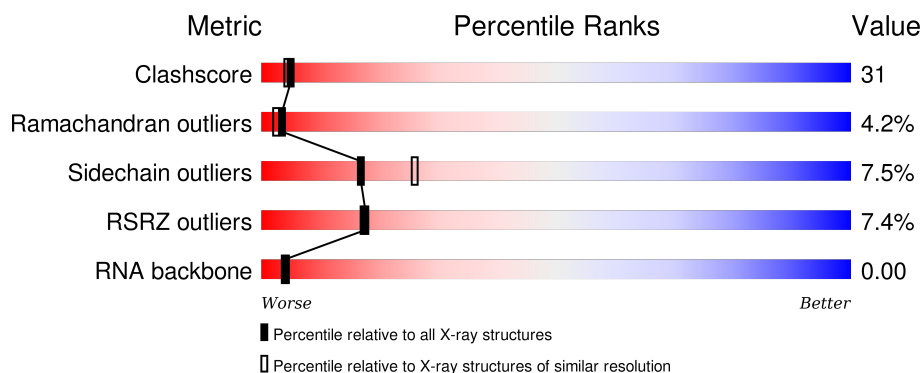
# 1 Overall quality at a glance

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))
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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	883	
2	N	17	
3	R	3	
4	T	22	

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7992 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 BACTERIOPHAGE T7 RNA POLYMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	862	Total	C	N	O	S	0	0	0
			6696	4269	1167	1224	36			

- Molecule 2 is a DNA chain called DNA (5-D(P\*TP\*AP\*AP\*TP\*AP\*CP\*GP\*AP\*CP\*TP\*CP\*AP\*CP\*TP\*A)-3).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	N	14	Total	C	N	O	P	0	0	0
			283	136	50	83	14			

- Molecule 3 is a RNA chain called RNA (5- R(PPP\*GP\*GP\*G)-3).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	R	3	Total	C	N	O	P	0	0	0
			78	30	15	28	5			

- Molecule 4 is a DNA chain called DNA (5- D (P\*CP\*TP\*CP\*CP\*CP\*TP\*AP\*TP\*AP\*GP\*TP\*GP\*AP\*GP\*TP\*CP\*GP\*TP\* AP\*TP\*TP\*A)-3).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	T	22	Total	C	N	O	P	0	0	0
			448	215	76	135	22			

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	429	Total	O	0	0
			429	429		
5	N	21	Total	O	0	0
			21	21		
5	R	2	Total	O	0	0
			2	2		

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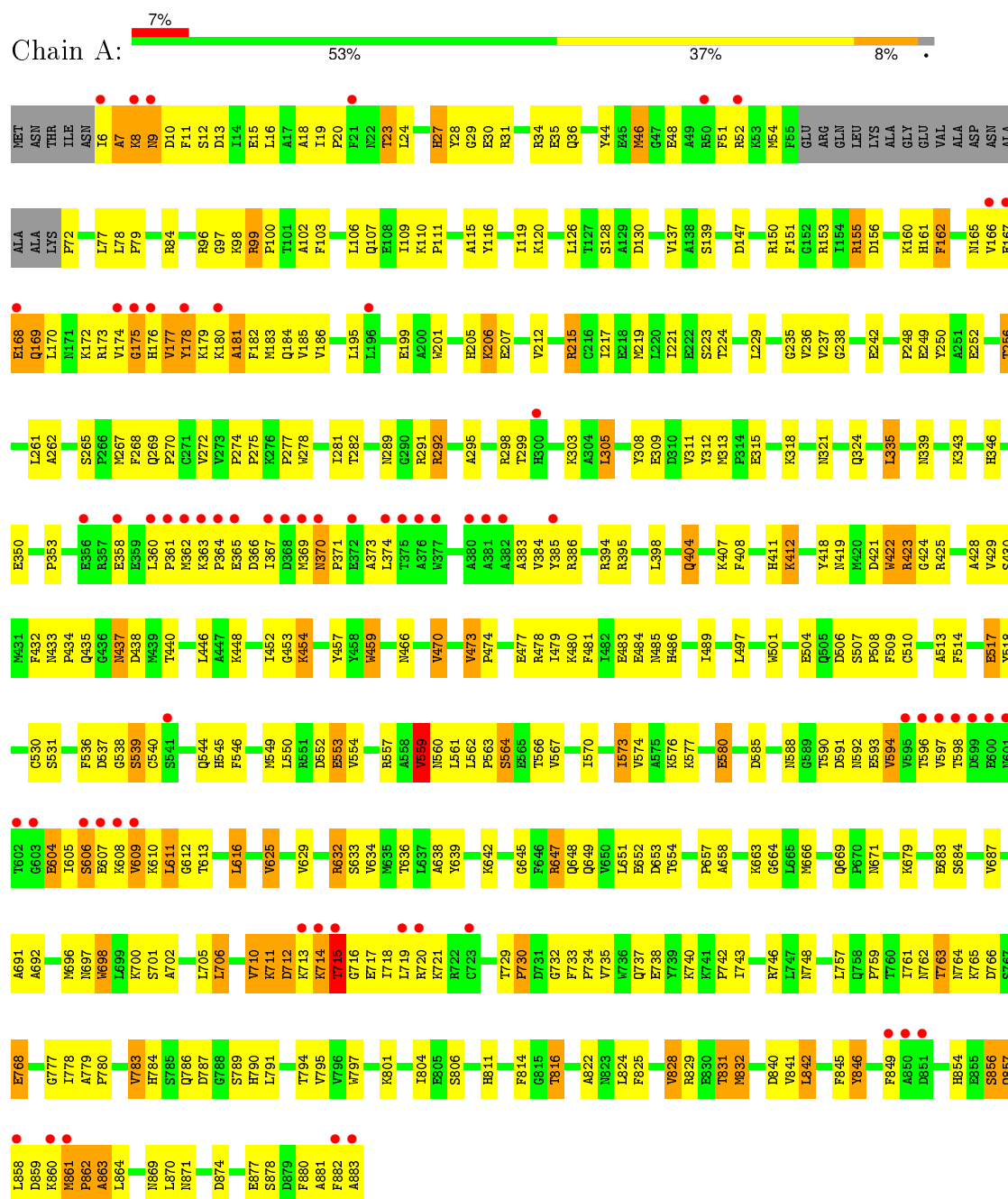
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	T	35	Total	O	0	0
			35	35		

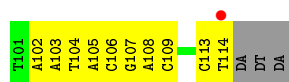
### 3 Residue-property plots

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: BACTERIOPHAGE T7 RNA POLYMERASE



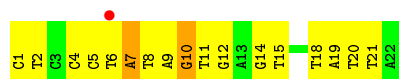
- Molecule 2: DNA (5-D(P\*TP\*AP\*AP\*TP\*AP\*CP\*GP\*AP\*CP\*TP\*CP\*AP\*CP\*TP\*A)-3)



- Molecule 3: RNA (5- R(PPP\*GP\*GP\*G)-3)



- Molecule 4: DNA (5- D (P\*CP\*TP\*CP\*CP\*CP\*TP\*AP\*TP\*AP\*GP\*TP\*GP\*AP\*GP\*TP\*C P\*GP\*TP\* AP\*TP\*TP\*A)-3)



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	221.20 Å 73.60 Å 80.90 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.40 39.79 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.8 (40.00-2.40) 97.5 (39.79-2.40)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.16 (at 2.39 Å)	Xtriage
Refinement program	CNS 0.5	Depositor
R, $R_{free}$	0.223 , 0.267 0.228 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	56.2	Xtriage
Anisotropy	0.225	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 47.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 51475 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	7992	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.93% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup> Theoretical 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: GTP

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	0.58	0/6851	0.78	4/9281 (0.0%)
2	N	0.83	0/316	0.96	0/484
3	R	0.79	0/51	0.75	0/78
4	T	0.79	0/500	0.93	0/769
All	All	0.61	0/7718	0.80	4/10612 (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
4	T	0	2

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	559	VAL	N-CA-C	-5.83	95.26	111.00
1	A	715	THR	N-CA-C	-5.71	95.60	111.00
1	A	346	HIS	N-CA-C	-5.52	96.10	111.00
1	A	181	ALA	N-CA-C	-5.41	96.41	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	T	10	DG	Sidechain
4	T	7	DA	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	A	6696	0	6561	416	0
2	N	283	0	159	13	0
3	R	78	0	34	2	0
4	T	448	0	251	42	0
5	A	429	0	0	17	0
5	N	21	0	0	0	0
5	R	2	0	0	0	0
5	T	35	0	0	2	0
All	All	7992	0	7005	451	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 31.

All (451) 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:713:LYS:HD3	1:A:714:LYS:H	1.05	1.11
1:A:710:VAL:HG11	1:A:720:ARG:H	1.03	1.10
1:A:710:VAL:HG11	1:A:720:ARG:N	1.65	1.08
4:T:4:DC:H2'	4:T:5:DC:O4'	1.54	1.08
1:A:298:ARG:HD2	4:T:6:DT:H5'	1.35	1.07
1:A:710:VAL:HG21	1:A:719:LEU:HB3	1.40	1.01
1:A:663:LYS:HG2	1:A:664:GLY:H	1.21	1.01
1:A:714:LYS:HA	1:A:714:LYS:HE2	1.40	1.00
1:A:560:ASN:O	1:A:881:ALA:HB2	1.62	0.99
4:T:6:DT:H4'	4:T:7:DA:H5''	1.41	0.99
1:A:155:ARG:HG3	1:A:155:ARG:HH11	1.29	0.96
1:A:713:LYS:HD3	1:A:714:LYS:N	1.81	0.94
4:T:6:DT:H4'	4:T:7:DA:C5'	1.99	0.91
1:A:298:ARG:HD3	4:T:6:DT:H2'	1.51	0.91
1:A:412:LYS:HE2	1:A:412:LYS:HA	1.53	0.90
1:A:645:GLY:HA3	4:T:2:DT:OP1	1.71	0.89
1:A:369:MET:C	1:A:371:PRO:HD2	1.92	0.88
4:T:20:DT:H2''	4:T:21:DT:H5'	1.56	0.87
1:A:710:VAL:CG1	1:A:720:ARG:H	1.86	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:729:THR:HG22	1:A:733:PHE:O	1.76	0.86
1:A:278:TRP:H	1:A:321:ASN:HD21	1.22	0.86
1:A:206:LYS:HE2	4:T:9:DA:N1	1.90	0.85
1:A:729:THR:HG21	1:A:733:PHE:HB3	1.60	0.83
1:A:446:LEU:HD22	1:A:806:SER:HB3	1.61	0.83
1:A:363:LYS:HD3	1:A:367:ILE:CB	2.08	0.82
1:A:281:ILE:HG23	1:A:282:THR:HG23	1.61	0.82
1:A:763:THR:CG2	1:A:765:LYS:H	1.92	0.82
1:A:763:THR:HG23	1:A:765:LYS:H	1.44	0.81
1:A:30:GLU:OE1	1:A:173:ARG:HD2	1.79	0.81
1:A:632:ARG:HH11	1:A:632:ARG:HB2	1.46	0.80
1:A:715:THR:O	1:A:717:GLU:N	2.15	0.80
1:A:730:PRO:CD	1:A:786:GLN:HE22	1.93	0.80
1:A:363:LYS:NZ	1:A:374:LEU:HD11	1.96	0.80
4:T:5:DC:H3'	4:T:7:DA:OP1	1.81	0.80
1:A:715:THR:C	1:A:717:GLU:H	1.85	0.79
1:A:553:GLU:CD	1:A:553:GLU:H	1.86	0.79
1:A:806:SER:O	1:A:816:THR:HG23	1.84	0.78
1:A:269:GLN:HE22	1:A:407:LYS:NZ	1.82	0.77
1:A:729:THR:CG2	1:A:733:PHE:HB3	2.15	0.76
1:A:109:ILE:HD12	1:A:109:ILE:H	1.52	0.75
1:A:155:ARG:HG3	1:A:155:ARG:NH1	1.99	0.75
1:A:861:MET:H	1:A:862:PRO:HD2	1.50	0.75
1:A:663:LYS:HG2	1:A:664:GLY:N	1.99	0.75
1:A:794:THR:OG1	1:A:831:THR:HG21	1.86	0.74
1:A:858:LEU:O	1:A:858:LEU:HG	1.86	0.74
1:A:398:LEU:C	1:A:398:LEU:HD23	2.08	0.74
2:N:103:DA:H2''	2:N:104:DT:H5'	1.69	0.73
1:A:298:ARG:CD	4:T:6:DT:H5'	2.18	0.73
1:A:324:GLN:HE21	1:A:418:TYR:H	1.35	0.71
2:N:108:DA:H2''	2:N:109:DC:C5'	2.20	0.71
4:T:19:DA:H1'	4:T:20:DT:H5''	1.72	0.71
1:A:170:LEU:O	1:A:174:VAL:HG23	1.90	0.71
1:A:176:HIS:O	1:A:177:VAL:O	2.10	0.70
1:A:335:LEU:HD22	1:A:339:ASN:ND2	2.06	0.70
1:A:647:ARG:HH11	1:A:671:ASN:HD21	1.40	0.69
1:A:743:ILE:O	1:A:763:THR:HB	1.93	0.69
1:A:846:TYR:HA	1:A:849:PHE:CE1	2.27	0.69
1:A:486:HIS:HD2	1:A:518:TYR:OH	1.76	0.69
1:A:229:LEU:HD11	1:A:242:GLU:HG3	1.75	0.69
1:A:710:VAL:O	1:A:710:VAL:HG13	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:299:THR:H	4:T:6:DT:H72	1.59	0.68
1:A:281:ILE:HD11	1:A:305:LEU:O	1.94	0.68
1:A:713:LYS:HA	1:A:713:LYS:HE2	1.76	0.68
1:A:156:ASP:OD1	1:A:160:LYS:HE2	1.94	0.68
1:A:72:PRO:HB3	1:A:262:ALA:HB3	1.76	0.67
1:A:252:GLU:O	1:A:256:THR:HB	1.94	0.67
1:A:269:GLN:HE22	1:A:407:LYS:HZ3	1.43	0.66
1:A:438:ASP:OD1	1:A:507:SER:HB3	1.96	0.66
4:T:18:DT:H2''	4:T:19:DA:C8	2.30	0.66
1:A:545:HIS:O	1:A:549:MET:HG2	1.95	0.66
1:A:169:GLN:HG2	1:A:172:LYS:HE3	1.76	0.66
1:A:761:ILE:O	1:A:761:ILE:HG13	1.94	0.65
1:A:608:LYS:O	1:A:609:VAL:HB	1.96	0.65
1:A:109:ILE:HD12	1:A:109:ILE:N	2.11	0.65
1:A:647:ARG:NH1	1:A:671:ASN:HD21	1.95	0.65
1:A:801:LYS:HD3	1:A:801:LYS:O	1.96	0.65
1:A:221:ILE:O	1:A:224:THR:O	2.15	0.64
1:A:19:ILE:O	1:A:23:THR:HG23	1.98	0.64
1:A:480:LYS:O	1:A:484:GLU:HG3	1.97	0.64
1:A:363:LYS:HZ2	1:A:374:LEU:HD11	1.61	0.64
1:A:854:HIS:CB	1:A:858:LEU:HD13	2.27	0.64
1:A:829:ARG:NH1	1:A:878:SER:O	2.31	0.64
1:A:632:ARG:HB2	1:A:632:ARG:NH1	2.12	0.64
1:A:19:ILE:HG22	1:A:195:LEU:HD21	1.79	0.64
1:A:638:ALA:O	1:A:780:PRO:HB3	1.98	0.64
1:A:437:ASN:C	1:A:437:ASN:HD22	2.00	0.63
1:A:679:LYS:O	1:A:683:GLU:HB2	1.99	0.63
1:A:730:PRO:HD3	1:A:786:GLN:HE22	1.62	0.63
2:N:108:DA:H2''	2:N:109:DC:H5'	1.81	0.63
1:A:177:VAL:O	1:A:179:LYS:HG3	1.98	0.62
1:A:371:PRO:HG2	1:A:373:ALA:CB	2.30	0.62
1:A:452:ILE:HD11	1:A:457:TYR:HA	1.82	0.62
1:A:102:ALA:HB3	1:A:215:ARG:HG2	1.82	0.61
1:A:738:GLU:O	1:A:738:GLU:HG3	2.00	0.61
1:A:394:ARG:HG2	1:A:394:ARG:HH11	1.64	0.61
1:A:308:TYR:HA	1:A:311:VAL:CG2	2.30	0.61
1:A:308:TYR:HA	1:A:311:VAL:HG23	1.82	0.61
1:A:804:ILE:HG23	1:A:816:THR:HG21	1.83	0.61
1:A:517:GLU:HG3	1:A:530:CYS:SG	2.41	0.61
1:A:715:THR:C	1:A:717:GLU:N	2.51	0.61
1:A:19:ILE:HG22	1:A:20:PRO:N	2.16	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:T:1:DC:H4'	4:T:1:DC:OP1	2.00	0.61
1:A:182:PHE:CE2	1:A:448:LYS:HB3	2.36	0.61
1:A:278:TRP:H	1:A:321:ASN:ND2	1.98	0.61
1:A:710:VAL:CG2	1:A:719:LEU:HB3	2.25	0.60
1:A:545:HIS:CD2	1:A:787:ASP:HB3	2.35	0.60
1:A:647:ARG:HG3	1:A:648:GLN:N	2.15	0.60
3:R:1:GTP:O1G	3:R:1:GTP:O3A	2.13	0.60
4:T:5:DC:H5'	4:T:7:DA:H4'	1.84	0.60
1:A:710:VAL:HG12	1:A:720:ARG:O	2.02	0.60
2:N:108:DA:H2''	2:N:109:DC:H5''	1.83	0.60
1:A:120:LYS:HE2	1:A:265:SER:O	2.00	0.60
1:A:185:VAL:HG22	1:A:186:VAL:N	2.17	0.60
1:A:383:ALA:HA	1:A:386:ARG:NH1	2.17	0.60
1:A:18:ALA:HA	1:A:161:HIS:CD2	2.37	0.60
1:A:713:LYS:O	1:A:714:LYS:C	2.39	0.60
1:A:861:MET:N	1:A:862:PRO:HD2	2.10	0.59
1:A:364:PRO:C	1:A:366:ASP:H	2.05	0.59
1:A:778:ILE:HG23	1:A:779:ALA:N	2.17	0.59
1:A:358:GLU:OE1	1:A:384:VAL:HG13	2.02	0.59
1:A:97:GLY:HA2	4:T:20:DT:O2	2.02	0.59
1:A:593:GLU:O	1:A:594:VAL:O	2.20	0.59
1:A:562:LEU:HD21	1:A:870:LEU:CD1	2.32	0.59
1:A:177:VAL:O	1:A:179:LYS:N	2.31	0.59
1:A:596:THR:O	1:A:598:THR:N	2.36	0.59
1:A:34:ARG:HH22	1:A:165:ASN:HD22	1.51	0.59
1:A:369:MET:C	1:A:371:PRO:CD	2.67	0.59
1:A:12:SER:HB3	1:A:15:GLU:HG3	1.85	0.59
1:A:501:TRP:HA	1:A:504:GLU:OE1	2.02	0.58
1:A:701:SER:HB3	1:A:861:MET:HG2	1.85	0.58
1:A:199:GLU:OE1	1:A:199:GLU:HA	2.03	0.58
1:A:278:TRP:HE1	1:A:324:GLN:HE22	1.50	0.58
1:A:485:ASN:O	1:A:489:ILE:HG13	2.03	0.58
1:A:298:ARG:HH11	4:T:6:DT:C2'	2.15	0.58
1:A:36:GLN:NE2	1:A:272:VAL:H	2.01	0.58
1:A:272:VAL:HG12	1:A:272:VAL:O	2.04	0.58
4:T:14:DG:H2''	4:T:15:DT:H5'	1.86	0.57
1:A:72:PRO:HB3	1:A:262:ALA:CB	2.34	0.57
1:A:298:ARG:HH11	4:T:6:DT:H2'	1.68	0.57
1:A:423:ARG:NH2	1:A:784:HIS:ND1	2.51	0.57
1:A:174:VAL:HG12	1:A:177:VAL:HG22	1.86	0.57
1:A:115:ALA:O	1:A:119:ILE:HG12	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:828:VAL:O	1:A:831:THR:HG22	2.05	0.57
1:A:311:VAL:O	1:A:312:TYR:HB3	2.03	0.57
1:A:365:GLU:C	1:A:367:ILE:H	2.08	0.57
1:A:593:GLU:HB3	1:A:610:LYS:CB	2.35	0.57
1:A:99:ARG:HD2	1:A:99:ARG:N	2.19	0.57
1:A:371:PRO:HG2	1:A:373:ALA:HB3	1.88	0.56
1:A:806:SER:O	1:A:816:THR:CG2	2.52	0.56
1:A:24:LEU:O	1:A:27:HIS:O	2.23	0.56
4:T:14:DG:H1'	4:T:15:DT:H5''	1.85	0.56
1:A:103:PHE:CZ	1:A:107:GLN:HG3	2.40	0.56
1:A:363:LYS:HZ3	1:A:374:LEU:HD11	1.68	0.56
1:A:147:ASP:OD1	1:A:292:ARG:HD2	2.04	0.56
1:A:816:THR:OG1	1:A:824:LEU:HD22	2.05	0.56
1:A:281:ILE:CG2	1:A:282:THR:HG23	2.34	0.56
1:A:169:GLN:HG2	1:A:172:LYS:CE	2.34	0.56
1:A:34:ARG:HH22	1:A:165:ASN:ND2	2.03	0.56
1:A:184:GLN:HE21	1:A:185:VAL:HG12	1.70	0.56
1:A:364:PRO:O	1:A:366:ASP:N	2.34	0.56
1:A:109:ILE:CD1	1:A:109:ILE:H	2.18	0.55
1:A:362:MET:CE	1:A:363:LYS:H	2.19	0.55
1:A:473:VAL:HG21	1:A:477:GLU:OE1	2.07	0.55
1:A:179:LYS:O	1:A:181:ALA:N	2.34	0.55
1:A:470:VAL:HG12	1:A:470:VAL:O	2.06	0.55
1:A:29:GLY:CA	1:A:175:GLY:HA3	2.37	0.55
1:A:859:ASP:O	1:A:860:LYS:HB2	2.05	0.55
4:T:4:DC:H2'	4:T:5:DC:C4'	2.37	0.55
1:A:714:LYS:HA	1:A:714:LYS:CE	2.26	0.55
2:N:108:DA:C2'	2:N:109:DC:H5''	2.36	0.55
1:A:147:ASP:O	1:A:150:ARG:O	2.25	0.55
1:A:421:ASP:HB3	4:T:6:DT:OP2	2.06	0.54
1:A:371:PRO:C	1:A:373:ALA:H	2.09	0.54
1:A:642:LYS:NZ	1:A:697:ASN:HD21	2.06	0.54
1:A:421:ASP:OD2	1:A:423:ARG:NH1	2.40	0.54
1:A:730:PRO:CD	1:A:786:GLN:NE2	2.68	0.54
1:A:201:TRP:NE1	4:T:6:DT:H73	2.23	0.54
1:A:763:THR:HG22	1:A:765:LYS:H	1.72	0.54
1:A:298:ARG:HH11	4:T:6:DT:H3'	1.73	0.54
1:A:269:GLN:HE22	1:A:407:LYS:HZ2	1.56	0.54
1:A:454:LYS:C	1:A:454:LYS:HE3	2.28	0.54
1:A:361:PRO:HB3	5:A:2057:HOH:O	2.08	0.54
1:A:178:TYR:HE1	1:A:412:LYS:HG2	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:862:PRO:O	1:A:863:ALA:HB3	2.07	0.53
1:A:713:LYS:CD	1:A:714:LYS:H	1.98	0.53
1:A:363:LYS:CG	1:A:364:PRO:HD2	2.39	0.53
1:A:715:THR:O	1:A:715:THR:HG22	2.08	0.53
1:A:713:LYS:CD	1:A:714:LYS:N	2.66	0.53
1:A:27:HIS:O	1:A:28:TYR:HB2	2.09	0.53
1:A:856:SER:O	1:A:857:GLN:C	2.47	0.53
1:A:398:LEU:HD23	1:A:398:LEU:O	2.07	0.53
1:A:235:GLY:HA2	2:N:113:DC:O2	2.09	0.53
1:A:23:THR:O	1:A:27:HIS:HD2	1.92	0.53
1:A:705:LEU:HD11	1:A:861:MET:HB2	1.92	0.52
1:A:669:GLN:NE2	1:A:669:GLN:HA	2.23	0.52
1:A:265:SER:HB2	5:A:2015:HOH:O	2.09	0.52
1:A:398:LEU:C	1:A:398:LEU:CD2	2.78	0.52
1:A:150:ARG:O	1:A:151:PHE:HB2	2.10	0.52
1:A:9:ASN:O	1:A:11:PHE:N	2.43	0.52
1:A:871:ASN:HB3	1:A:874:ASP:OD2	2.10	0.52
1:A:608:LYS:O	1:A:609:VAL:CB	2.58	0.52
1:A:696:MET:O	1:A:700:LYS:HG3	2.10	0.52
1:A:369:MET:O	1:A:371:PRO:HD2	2.08	0.51
1:A:367:ILE:O	1:A:374:LEU:HD22	2.11	0.51
1:A:295:ALA:HB1	1:A:305:LEU:HD11	1.92	0.51
1:A:652:GLU:O	1:A:657:PRO:HD3	2.10	0.51
1:A:651:LEU:O	1:A:651:LEU:HD23	2.10	0.51
1:A:557:ARG:O	1:A:559:VAL:O	2.27	0.51
1:A:422:TRP:C	1:A:422:TRP:CD1	2.84	0.51
1:A:433:ASN:HB2	1:A:434:PRO:CD	2.40	0.51
1:A:362:MET:HE2	1:A:363:LYS:H	1.75	0.51
1:A:459:TRP:CZ3	1:A:822:ALA:HB2	2.46	0.51
1:A:466:ASN:OD1	1:A:478:ARG:NH1	2.44	0.51
1:A:609:VAL:O	1:A:609:VAL:HG13	2.11	0.51
1:A:35:GLU:OE2	1:A:411:HIS:HE1	1.92	0.51
1:A:607:GLU:O	1:A:609:VAL:N	2.40	0.51
1:A:573:ILE:CD1	1:A:577:LYS:HE2	2.41	0.51
1:A:371:PRO:C	1:A:373:ALA:N	2.64	0.51
1:A:861:MET:HB3	1:A:862:PRO:CD	2.41	0.51
2:N:107:DG:H2''	2:N:108:DA:H5'	1.92	0.51
1:A:248:PRO:O	1:A:252:GLU:HG3	2.10	0.51
1:A:669:GLN:HE21	1:A:669:GLN:HA	1.75	0.51
1:A:437:ASN:ND2	1:A:437:ASN:C	2.64	0.51
1:A:576:LYS:O	1:A:580:GLU:HG2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:VAL:HG13	1:A:173:ARG:HD3	1.92	0.51
1:A:609:VAL:HG21	1:A:669:GLN:HG3	1.91	0.51
1:A:8:LYS:O	1:A:9:ASN:CB	2.59	0.51
1:A:473:VAL:HG22	1:A:474:PRO:HD2	1.92	0.50
1:A:185:VAL:CG2	1:A:277:PRO:HG3	2.41	0.50
2:N:106:DC:H1'	2:N:107:DG:C8	2.46	0.50
1:A:563:PRO:HB3	1:A:877:GLU:O	2.11	0.50
1:A:538:GLY:HA3	1:A:883:ALA:HB3	1.94	0.50
1:A:605:ILE:O	1:A:606:SER:O	2.30	0.50
1:A:562:LEU:HD21	1:A:870:LEU:HD12	1.94	0.50
1:A:814:PHE:HE1	1:A:883:ALA:HB2	1.77	0.50
4:T:5:DC:C5'	4:T:7:DA:H4'	2.42	0.50
1:A:840:ASP:HA	5:A:2399:HOH:O	2.12	0.50
1:A:712:ASP:CG	1:A:713:LYS:N	2.64	0.50
1:A:564:SER:HB3	1:A:566:THR:O	2.11	0.50
1:A:370:ASN:N	1:A:371:PRO:CD	2.75	0.50
1:A:831:THR:CG2	1:A:832:MET:N	2.75	0.50
1:A:692:ALA:O	1:A:696:MET:HG3	2.11	0.50
1:A:78:LEU:HB3	1:A:79:PRO:CD	2.42	0.50
1:A:553:GLU:OE1	1:A:869:ASN:N	2.45	0.49
1:A:861:MET:O	1:A:862:PRO:O	2.30	0.49
1:A:452:ILE:HG23	1:A:453:GLY:N	2.27	0.49
1:A:120:LYS:NZ	1:A:267:MET:SD	2.85	0.49
1:A:178:TYR:HE1	1:A:412:LYS:CG	2.26	0.49
1:A:31:ARG:HD3	1:A:176:HIS:CE1	2.47	0.49
1:A:632:ARG:HH11	1:A:632:ARG:CB	2.20	0.49
1:A:562:LEU:HD21	1:A:870:LEU:HD11	1.95	0.49
1:A:173:ARG:C	1:A:175:GLY:N	2.65	0.49
1:A:412:LYS:HE2	1:A:412:LYS:CA	2.35	0.49
1:A:175:GLY:HA2	5:A:2070:HOH:O	2.13	0.49
1:A:19:ILE:HG22	1:A:195:LEU:CD2	2.43	0.49
1:A:6:ILE:O	1:A:7:ALA:C	2.50	0.49
1:A:706:LEU:HD21	1:A:849:PHE:HB2	1.93	0.49
1:A:437:ASN:ND2	1:A:440:THR:H	2.11	0.48
1:A:358:GLU:OE1	1:A:384:VAL:HG22	2.14	0.48
1:A:737:GLN:HE22	1:A:777:GLY:C	2.16	0.48
1:A:585:ASP:O	1:A:588:ASN:O	2.31	0.48
2:N:102:DA:H1'	2:N:103:DA:C8	2.48	0.48
1:A:270:PRO:HD2	1:A:408:PHE:CE2	2.48	0.48
4:T:7:DA:H2''	4:T:8:DT:O5'	2.14	0.48
1:A:369:MET:CB	1:A:371:PRO:HD2	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:ILE:HD12	1:A:309:GLU:N	2.28	0.48
1:A:236:VAL:HG12	1:A:236:VAL:O	2.12	0.48
1:A:364:PRO:C	1:A:366:ASP:N	2.67	0.48
1:A:730:PRO:CG	1:A:786:GLN:HE22	2.26	0.48
1:A:446:LEU:HB2	1:A:531:SER:O	2.14	0.48
1:A:861:MET:HE3	1:A:862:PRO:HD3	1.95	0.48
1:A:27:HIS:O	1:A:28:TYR:CB	2.60	0.48
1:A:433:ASN:HB2	1:A:434:PRO:HD3	1.96	0.48
1:A:459:TRP:HZ3	1:A:822:ALA:HB2	1.78	0.48
1:A:236:VAL:O	1:A:238:GLY:N	2.47	0.48
1:A:36:GLN:HE22	1:A:272:VAL:H	1.62	0.47
1:A:404:GLN:HA	1:A:404:GLN:HE21	1.78	0.47
1:A:861:MET:SD	1:A:862:PRO:N	2.87	0.47
1:A:711:LYS:HG3	5:A:2350:HOH:O	2.12	0.47
1:A:430:SER:O	1:A:433:ASN:HB3	2.14	0.47
1:A:360:LEU:HD11	1:A:385:TYR:OH	2.14	0.47
1:A:370:ASN:OD1	1:A:370:ASN:O	2.33	0.47
1:A:84:ARG:HB2	1:A:223:SER:HB3	1.96	0.47
1:A:298:ARG:HH11	4:T:6:DT:C3'	2.27	0.47
1:A:663:LYS:HE3	1:A:666:MET:HE1	1.97	0.47
1:A:536:PHE:HB3	1:A:882:PHE:HB3	1.97	0.47
1:A:215:ARG:O	1:A:219:MET:HG3	2.15	0.47
1:A:473:VAL:CG2	1:A:477:GLU:OE1	2.63	0.47
1:A:651:LEU:HD23	1:A:651:LEU:C	2.35	0.47
1:A:268:PHE:O	1:A:430:SER:HB2	2.15	0.47
1:A:790:HIS:CD2	1:A:831:THR:HG23	2.50	0.47
1:A:537:ASP:N	1:A:882:PHE:HD2	2.12	0.47
1:A:44:TYR:HB2	1:A:155:ARG:HH12	1.80	0.47
1:A:610:LYS:O	1:A:611:LEU:C	2.53	0.47
1:A:611:LEU:HB3	1:A:616:LEU:HD13	1.96	0.47
1:A:404:GLN:HG2	1:A:432:PHE:HB3	1.96	0.47
1:A:48:GLU:OE2	1:A:52:ARG:HB2	2.14	0.46
1:A:710:VAL:HG11	1:A:719:LEU:N	2.30	0.46
1:A:882:PHE:CD1	1:A:882:PHE:N	2.80	0.46
1:A:881:ALA:C	1:A:882:PHE:O	2.52	0.46
1:A:281:ILE:HG23	1:A:282:THR:CG2	2.40	0.46
3:R:1:GTP:O6	4:T:5:DC:O2	2.33	0.46
1:A:394:ARG:NH1	1:A:394:ARG:HG2	2.30	0.46
1:A:116:TYR:CZ	1:A:120:LYS:HD2	2.51	0.46
1:A:343:LYS:HD3	5:A:2174:HOH:O	2.14	0.46
1:A:139:SER:OG	1:A:206:LYS:HE3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:831:THR:HG23	1:A:832:MET:N	2.30	0.46
1:A:311:VAL:O	1:A:312:TYR:CB	2.64	0.46
1:A:479:ILE:O	1:A:483:GLU:HG3	2.16	0.46
4:T:2:DT:OP2	4:T:2:DT:H4'	2.16	0.46
1:A:363:LYS:HG3	1:A:364:PRO:HD2	1.98	0.46
1:A:353:PRO:O	1:A:395:ARG:NH1	2.49	0.46
1:A:732:GLY:O	1:A:734:PRO:HD3	2.16	0.46
1:A:654:THR:O	1:A:658:ALA:HB2	2.16	0.45
1:A:606:SER:C	1:A:608:LYS:N	2.67	0.45
1:A:729:THR:HG23	1:A:733:PHE:H	1.80	0.45
1:A:452:ILE:HG23	1:A:453:GLY:H	1.80	0.45
1:A:791:LEU:O	1:A:795:VAL:HG23	2.17	0.45
1:A:570:ILE:O	1:A:574:VAL:HG23	2.17	0.45
1:A:729:THR:OG1	1:A:789:SER:HB3	2.17	0.45
1:A:77:LEU:HD11	1:A:250:TYR:CE1	2.51	0.45
1:A:663:LYS:HE3	1:A:666:MET:CE	2.46	0.45
1:A:763:THR:HG23	1:A:764:ASN:N	2.31	0.45
1:A:185:VAL:CG2	1:A:186:VAL:N	2.79	0.45
1:A:324:GLN:NE2	1:A:418:TYR:H	2.11	0.45
1:A:281:ILE:CD1	1:A:305:LEU:HD22	2.47	0.45
1:A:510:CYS:O	1:A:513:ALA:N	2.50	0.45
1:A:606:SER:O	1:A:607:GLU:C	2.54	0.45
1:A:649:GLN:O	1:A:653:ASP:HB2	2.17	0.45
1:A:422:TRP:O	1:A:422:TRP:CD1	2.70	0.45
1:A:710:VAL:CG1	1:A:720:ARG:N	2.55	0.45
1:A:746:ARG:HA	1:A:759:PRO:O	2.17	0.45
1:A:717:GLU:CG	1:A:718:ILE:N	2.80	0.44
1:A:215:ARG:H	1:A:215:ARG:HD2	1.82	0.44
1:A:315:GLU:OE2	1:A:318:LYS:HE2	2.16	0.44
1:A:205:HIS:O	1:A:206:LYS:C	2.55	0.44
1:A:768:GLU:HB2	5:A:2146:HOH:O	2.17	0.44
1:A:714:LYS:CA	1:A:714:LYS:HE2	2.28	0.44
1:A:797:TRP:CH2	1:A:801:LYS:HG3	2.53	0.44
1:A:590:THR:HG22	1:A:591:ASP:N	2.32	0.44
1:A:786:GLN:NE2	1:A:841:VAL:HG11	2.32	0.44
1:A:567:VAL:HG22	1:A:880:PHE:CD2	2.53	0.44
1:A:205:HIS:HB3	1:A:207:GLU:OE1	2.18	0.44
1:A:96:ARG:HD2	2:N:103:DA:H1'	1.99	0.44
1:A:23:THR:HG21	1:A:195:LEU:HD11	1.99	0.44
1:A:592:ASN:HA	1:A:612:GLY:H	1.81	0.44
1:A:778:ILE:HG23	1:A:779:ALA:H	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:T:2:DT:OP2	4:T:2:DT:C4'	2.66	0.43
1:A:242:GLU:HG2	1:A:757:LEU:CD1	2.48	0.43
2:N:104:DT:H2''	2:N:105:DA:C8	2.53	0.43
1:A:742:PRO:HD2	5:A:2364:HOH:O	2.18	0.43
1:A:698:TRP:CZ2	1:A:864:LEU:HD23	2.53	0.43
1:A:669:GLN:HE21	1:A:669:GLN:CA	2.32	0.43
1:A:825:PHE:O	1:A:829:ARG:HG2	2.18	0.43
1:A:538:GLY:CA	1:A:883:ALA:HB3	2.47	0.43
1:A:249:GLU:CD	1:A:249:GLU:H	2.21	0.43
1:A:702:ALA:O	1:A:706:LEU:HD12	2.19	0.43
1:A:590:THR:HB	1:A:613:THR:H	1.84	0.43
1:A:546:PHE:CZ	1:A:783:VAL:HG13	2.54	0.43
1:A:299:THR:H	4:T:6:DT:C7	2.29	0.43
1:A:849:PHE:HE2	1:A:861:MET:HE2	1.83	0.43
1:A:604:GLU:C	1:A:606:SER:H	2.21	0.43
1:A:165:ASN:HA	1:A:165:ASN:HD22	1.69	0.43
1:A:269:GLN:NE2	1:A:407:LYS:HZ3	2.13	0.43
1:A:609:VAL:HA	5:A:2303:HOH:O	2.19	0.43
1:A:183:MET:HG3	5:A:2081:HOH:O	2.19	0.43
1:A:738:GLU:O	1:A:740:LYS:HG3	2.19	0.43
1:A:609:VAL:HG11	1:A:669:GLN:OE1	2.19	0.43
1:A:766:ASP:HB2	5:A:2361:HOH:O	2.18	0.43
1:A:497:LEU:HA	1:A:497:LEU:HD23	1.82	0.42
1:A:552:ASP:HB2	1:A:691:ALA:HB2	2.01	0.42
4:T:2:DT:H72	5:T:2007:HOH:O	2.18	0.42
1:A:748:ASN:HB3	2:N:106:DC:H2''	2.00	0.42
1:A:550:LEU:HD21	1:A:842:LEU:CD2	2.48	0.42
1:A:98:LYS:HD3	5:T:2031:HOH:O	2.19	0.42
1:A:168:GLU:HG3	5:A:2068:HOH:O	2.19	0.42
1:A:711:LYS:O	1:A:712:ASP:CB	2.66	0.42
1:A:27:HIS:CD2	1:A:27:HIS:N	2.87	0.42
1:A:470:VAL:HG21	1:A:481:PHE:CG	2.54	0.42
1:A:636:THR:HG23	5:A:2322:HOH:O	2.19	0.42
1:A:371:PRO:HG2	1:A:373:ALA:HB2	2.01	0.42
1:A:647:ARG:NH1	1:A:671:ASN:ND2	2.65	0.42
1:A:28:TYR:HB3	1:A:183:MET:HE1	2.00	0.42
1:A:779:ALA:N	1:A:780:PRO:HD2	2.34	0.42
1:A:473:VAL:HG13	1:A:477:GLU:HB2	2.01	0.42
1:A:162:PHE:C	1:A:162:PHE:CD1	2.93	0.42
1:A:126:LEU:HD23	1:A:126:LEU:HA	1.81	0.42
1:A:849:PHE:CD2	1:A:849:PHE:O	2.72	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:604:GLU:C	1:A:606:SER:N	2.72	0.42
1:A:215:ARG:HD2	1:A:215:ARG:N	2.34	0.42
1:A:684:SER:O	1:A:687:VAL:HG22	2.18	0.42
1:A:51:PHE:O	1:A:54:MET:HG3	2.20	0.42
1:A:155:ARG:CG	1:A:155:ARG:NH1	2.70	0.42
1:A:30:GLU:HG3	1:A:31:ARG:N	2.33	0.42
1:A:438:ASP:OD2	1:A:509:PHE:HB2	2.19	0.42
1:A:130:ASP:HB2	5:A:2051:HOH:O	2.20	0.42
4:T:10:DG:H2''	4:T:11:DT:O5'	2.18	0.42
1:A:298:ARG:HD3	4:T:6:DT:C2'	2.35	0.42
1:A:560:ASN:C	1:A:881:ALA:HB2	2.33	0.42
1:A:639:TYR:CD2	4:T:2:DT:O2	2.73	0.42
1:A:825:PHE:CE2	1:A:829:ARG:NH2	2.88	0.42
1:A:206:LYS:HG3	1:A:762:ASN:O	2.20	0.42
1:A:242:GLU:HG2	1:A:757:LEU:HD11	2.02	0.42
1:A:634:VAL:HG22	1:A:634:VAL:O	2.20	0.42
1:A:166:VAL:HG23	1:A:167:GLU:OE2	2.20	0.41
4:T:14:DG:H1'	4:T:15:DT:C5'	2.48	0.41
1:A:573:ILE:HD12	1:A:577:LYS:HE2	2.01	0.41
4:T:11:DT:H2''	4:T:12:DG:C8	2.55	0.41
1:A:429:VAL:O	1:A:429:VAL:CG1	2.68	0.41
1:A:424:GLY:O	1:A:425:ARG:C	2.59	0.41
1:A:537:ASP:H	1:A:882:PHE:HD2	1.67	0.41
1:A:106:LEU:HD11	1:A:212:VAL:HG13	2.03	0.41
4:T:6:DT:O5'	4:T:7:DA:OP1	2.38	0.41
1:A:206:LYS:HA	1:A:206:LYS:HD3	1.82	0.41
1:A:29:GLY:HA3	1:A:175:GLY:HA3	2.01	0.41
1:A:881:ALA:O	1:A:882:PHE:C	2.58	0.41
1:A:634:VAL:N	5:A:2322:HOH:O	2.53	0.41
1:A:46:MET:HG2	1:A:51:PHE:HB2	2.02	0.41
1:A:625:VAL:HA	1:A:629:VAL:HG21	2.03	0.41
1:A:610:LYS:O	1:A:611:LEU:O	2.39	0.41
1:A:20:PRO:O	1:A:24:LEU:HB2	2.19	0.41
2:N:113:DC:H4'	2:N:114:DT:OP2	2.21	0.41
1:A:567:VAL:HG22	1:A:880:PHE:CG	2.55	0.41
1:A:611:LEU:HD12	1:A:669:GLN:HB2	2.03	0.41
1:A:573:ILE:HD11	1:A:577:LYS:HE2	2.02	0.41
1:A:698:TRP:HH2	1:A:846:TYR:CD2	2.39	0.41
1:A:486:HIS:CD2	1:A:518:TYR:OH	2.66	0.41
1:A:100:PRO:HG2	1:A:103:PHE:HB2	2.02	0.41
1:A:570:ILE:HA	1:A:573:ILE:HG23	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:842:LEU:O	1:A:845:PHE:HB3	2.20	0.41
4:T:10:DG:H2'	4:T:11:DT:H72	2.03	0.41
1:A:110:LYS:HA	1:A:111:PRO:HD2	1.93	0.41
1:A:428:ALA:H	1:A:435:GLN:NE2	2.19	0.41
1:A:16:LEU:HD23	1:A:16:LEU:HA	1.82	0.41
1:A:825:PHE:CZ	1:A:829:ARG:NH2	2.88	0.41
1:A:137:VAL:HG12	1:A:217:ILE:HD11	2.03	0.41
1:A:274:PRO:HA	1:A:275:PRO:HD3	1.72	0.41
1:A:173:ARG:C	1:A:175:GLY:H	2.24	0.40
1:A:862:PRO:O	1:A:863:ALA:CB	2.69	0.40
1:A:421:ASP:C	1:A:421:ASP:OD1	2.60	0.40
1:A:126:LEU:C	1:A:128:SER:N	2.74	0.40
1:A:298:ARG:HD2	4:T:6:DT:C5'	2.26	0.40
1:A:729:THR:CG2	1:A:733:PHE:H	2.35	0.40
1:A:425:ARG:HD3	1:A:811:HIS:CD2	2.56	0.40
1:A:350:GLU:HG3	5:A:2177:HOH:O	2.20	0.40
1:A:206:LYS:HE2	4:T:9:DA:C2	2.55	0.40
1:A:13:ASP:CG	1:A:153:ARG:HH22	2.25	0.40
1:A:544:GLN:HG2	1:A:561:LEU:CD1	2.52	0.40
4:T:5:DC:O2	4:T:5:DC:H2'	2.22	0.40
1:A:729:THR:CG2	1:A:733:PHE:CB	2.95	0.40
1:A:303:LYS:HD3	5:A:2149:HOH:O	2.20	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	858/883 (97%)	734 (86%)	88 (10%)	36 (4%)	<b>3</b> <b>2</b>

All (36) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	168	GLU
1	A	169	GLN
1	A	175	GLY
1	A	178	TYR
1	A	237	VAL
1	A	539	SER
1	A	594	VAL
1	A	604	GLU
1	A	606	SER
1	A	609	VAL
1	A	611	LEU
1	A	712	ASP
1	A	715	THR
1	A	857	GLN
1	A	862	PRO
1	A	10	ASP
1	A	177	VAL
1	A	180	LYS
1	A	597	VAL
1	A	716	GLY
1	A	721	LYS
1	A	559	VAL
1	A	711	LYS
1	A	714	LYS
1	A	370	ASN
1	A	863	ALA
1	A	7	ALA
1	A	8	LYS
1	A	632	ARG
1	A	856	SER
1	A	861	MET
1	A	625	VAL
1	A	508	PRO
1	A	470	VAL
1	A	710	VAL

### 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	680/729 (93%)	629 (92%)	51 (8%)	17	26

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	THR
1	A	27	HIS
1	A	46	MET
1	A	99	ARG
1	A	155	ARG
1	A	162	PHE
1	A	206	LYS
1	A	215	ARG
1	A	256	THR
1	A	261	LEU
1	A	289	ASN
1	A	291	ARG
1	A	292	ARG
1	A	305	LEU
1	A	313	MET
1	A	335	LEU
1	A	404	GLN
1	A	412	LYS
1	A	419	ASN
1	A	422	TRP
1	A	423	ARG
1	A	437	ASN
1	A	454	LYS
1	A	459	TRP
1	A	473	VAL
1	A	506	ASP
1	A	514	PHE
1	A	517	GLU
1	A	539	SER
1	A	540	CYS
1	A	553	GLU
1	A	554	VAL
1	A	564	SER
1	A	573	ILE
1	A	580	GLU
1	A	616	LEU

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Mol	Chain	Res	Type
1	A	633	SER
1	A	647	ARG
1	A	698	TRP
1	A	706	LEU
1	A	730	PRO
1	A	735	VAL
1	A	763	THR
1	A	768	GLU
1	A	783	VAL
1	A	816	THR
1	A	828	VAL
1	A	831	THR
1	A	832	MET
1	A	842	LEU
1	A	846	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	ASN
1	A	27	HIS
1	A	36	GLN
1	A	86	ASN
1	A	107	GLN
1	A	161	HIS
1	A	165	ASN
1	A	171	ASN
1	A	176	HIS
1	A	184	GLN
1	A	233	ASN
1	A	239	GLN
1	A	269	GLN
1	A	289	ASN
1	A	300	HIS
1	A	321	ASN
1	A	324	GLN
1	A	339	ASN
1	A	370	ASN
1	A	406	ASN
1	A	410	ASN
1	A	411	HIS
1	A	435	GLN

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Mol	Chain	Res	Type
1	A	437	ASN
1	A	463	HIS
1	A	485	ASN
1	A	486	HIS
1	A	522	GLN
1	A	523	HIS
1	A	583	GLN
1	A	669	GLN
1	A	671	ASN
1	A	697	ASN
1	A	726	HIS
1	A	737	GLN
1	A	744	GLN
1	A	764	ASN
1	A	781	ASN
1	A	786	GLN
1	A	811	HIS
1	A	823	ASN
1	A	869	ASN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	R	1/3 (33%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

## 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
3	GTP	R	1	3,4	25,34,34	1.30	4 (16%)	34,54,54	4.43	9 (26%)

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	GTP	R	1	3,4	-	0/18/38/38	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	R	1	GTP	O4'-C1'	2.16	1.43	1.41
3	R	1	GTP	PG-O3G	2.26	1.62	1.54
3	R	1	GTP	PG-O2G	2.63	1.64	1.54
3	R	1	GTP	C6-N1	3.64	1.39	1.33

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	R	1	GTP	C5-C6-N1	-8.61	111.82	123.59
3	R	1	GTP	O2B-PB-O3B	-6.14	77.22	105.09
3	R	1	GTP	O2G-PG-O3B	-2.91	91.88	105.09
3	R	1	GTP	N3-C2-N1	-2.46	123.70	127.44
3	R	1	GTP	PB-O3B-PG	-2.21	125.25	132.67
3	R	1	GTP	PA-O3A-PB	5.70	148.75	132.73
3	R	1	GTP	C6-N1-C2	6.57	125.06	115.94
3	R	1	GTP	O3G-PG-O3B	10.64	153.38	105.09
3	R	1	GTP	O3A-PA-O5'	17.91	150.46	102.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	R	1	GTP	2	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	862/883 (97%)	0.39	65 (7%) 17 17	18, 41, 85, 105	0
2	N	14/17 (82%)	0.23	1 (7%) 19 19	31, 49, 71, 108	0
3	R	2/3 (66%)	0.37	0 100 100	59, 59, 59, 61	0
4	T	22/22 (100%)	0.24	1 (4%) 37 38	39, 51, 95, 101	0
All	All	900/925 (97%)	0.38	67 (7%) 17 17	18, 42, 86, 108	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	362	MET	10.6
1	A	175	GLY	8.6
1	A	599	ASP	8.3
1	A	601	ASN	6.7
1	A	600	GLU	6.6
1	A	602	THR	6.6
1	A	598	THR	6.1
1	A	603	GLY	6.1
2	N	114	DT	6.0
1	A	608	LYS	5.5
1	A	363	LYS	5.5
1	A	364	PRO	5.3
1	A	361	PRO	5.3
1	A	606	SER	5.2
1	A	372	GLU	5.0
1	A	858	LEU	4.8
1	A	166	VAL	4.8
4	T	6	DT	4.7
1	A	374	LEU	4.6
1	A	180	LYS	4.6
1	A	596	THR	4.5

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Mol	Chain	Res	Type	RSRZ
1	A	597	VAL	4.3
1	A	713	LYS	4.3
1	A	851	ASP	4.3
1	A	178	TYR	4.1
1	A	719	LEU	4.1
1	A	174	VAL	4.0
1	A	370	ASN	3.9
1	A	167	GLU	3.8
1	A	367	ILE	3.7
1	A	376	ALA	3.7
1	A	385	TYR	3.6
1	A	52	ARG	3.6
1	A	368	ASP	3.5
1	A	607	GLU	3.3
1	A	358	GLU	3.2
1	A	369	MET	3.1
1	A	8	LYS	3.0
1	A	883	ALA	2.9
1	A	6	ILE	2.8
1	A	595	VAL	2.8
1	A	380	ALA	2.8
1	A	196	LEU	2.8
1	A	609	VAL	2.7
1	A	714	LYS	2.7
1	A	715	THR	2.7
1	A	381	ALA	2.6
1	A	850	ALA	2.5
1	A	168	GLU	2.5
1	A	861	MET	2.4
1	A	860	LYS	2.4
1	A	9	ASN	2.4
1	A	21	PHE	2.4
1	A	377	TRP	2.3
1	A	541	SER	2.3
1	A	365	GLU	2.3
1	A	300	HIS	2.3
1	A	360	LEU	2.3
1	A	356	GLU	2.2
1	A	882	PHE	2.2
1	A	375	THR	2.2
1	A	723	CYS	2.2
1	A	50	ARG	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	176	HIS	2.2
1	A	849	PHE	2.1
1	A	382	ALA	2.1
1	A	720	ARG	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. 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	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	GTP	R	1	32/32	0.56	0.32	-	81,92,115,115	0

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.