



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:27 AM GMT

PDB ID : 3EPH
Title : Crystallographic snapshots of eukaryotic dimethylallyltransferase acting on tRNA: Insight into tRNA recognition and reaction mechanism
Authors : Huang, R.H.; Zhou, C.
Deposited on : 2008-09-29
Resolution : 2.95 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

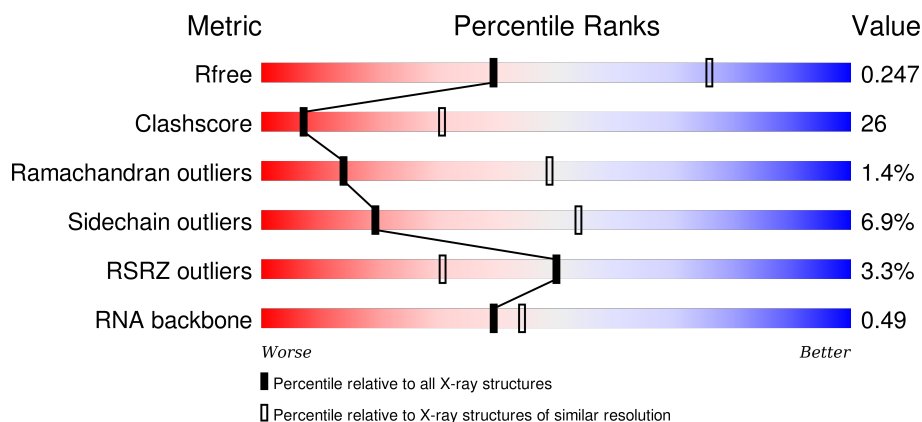
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	2184 (3.00-2.92)
Clashscore	102246	2552 (3.00-2.92)
Ramachandran outliers	100387	2468 (3.00-2.92)
Sidechain outliers	100360	2471 (3.00-2.92)
RSRZ outliers	91569	2201 (3.00-2.92)
RNA backbone	2183	1010 (3.36-2.56)

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	A	409	<div> <div>2%</div> <div>49%</div> <div>45%</div> <div>• •</div> </div>
1	B	409	<div> <div>3%</div> <div>46%</div> <div>48%</div> <div>• •</div> </div>
2	E	69	<div> <div>3%</div> <div>38%</div> <div>45%</div> <div>12%</div> <div>6%</div> </div>
2	F	69	<div> <div>9%</div> <div>35%</div> <div>43%</div> <div>16%</div> <div>6%</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	MG	A	4	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9934 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 tRNA isopentenyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	402	Total	C	N	O	S	0	0	0
			3339	2119	589	617	14			
1	B	402	Total	C	N	O	S	0	0	0
			3339	2119	589	617	14			

- Molecule 2 is a RNA chain called tRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	69	Total	C	N	O	P	0	0	0
			1472	656	258	489	69			
2	F	69	Total	C	N	O	P	0	0	0
			1472	656	258	489	69			

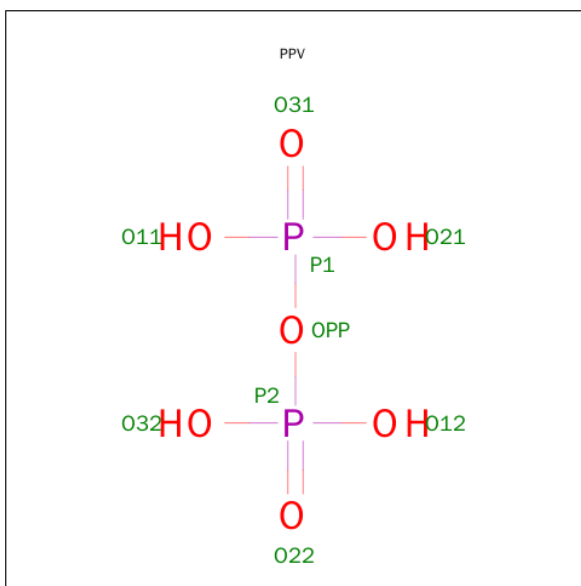
- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	3	Total	Mg	0	0
			3	3		
4	A	4	Total	Mg	0	0
			4	4		
4	F	1	Total	Mg	0	0
			1	1		
4	E	2	Total	Mg	0	0
			2	2		

- Molecule 5 is PYROPHOSPHATE (three-letter code: PPV) (formula: $\text{H}_4\text{O}_7\text{P}_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	P	0	0
			9	7	2		
5	B	1	Total	O	P	0	0
			9	7	2		

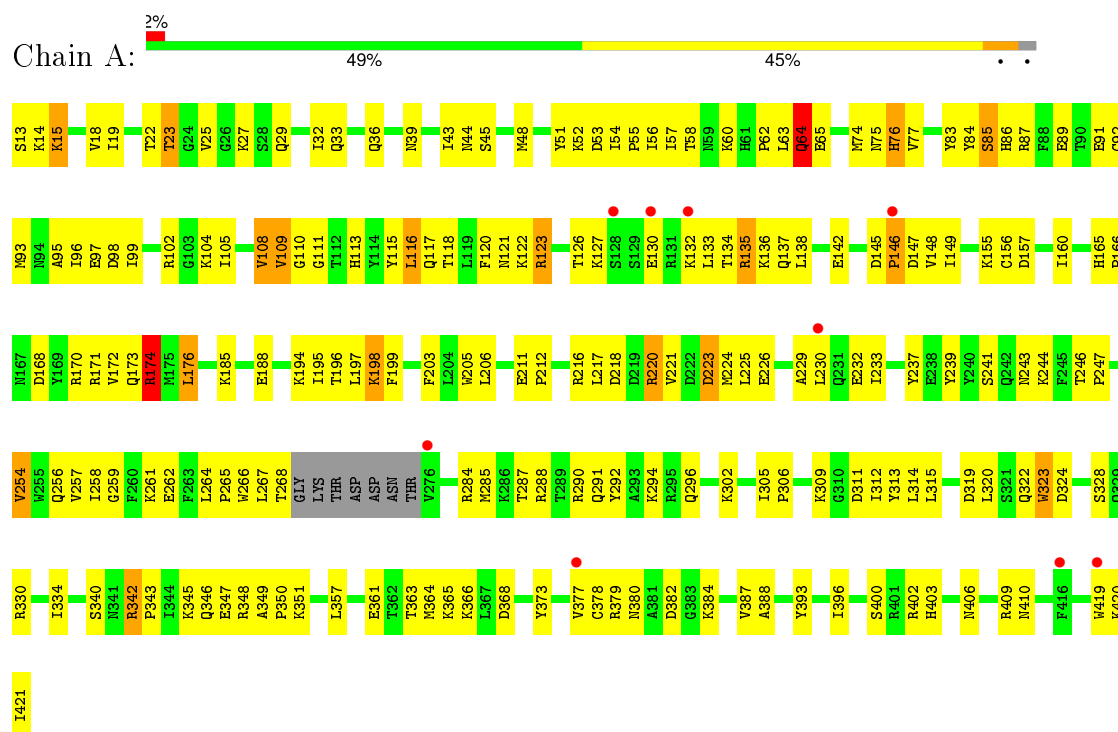
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	85	Total	O	0	0
			85	85		
6	E	58	Total	O	0	0
			58	58		
6	B	104	Total	O	0	0
			104	104		
6	F	35	Total	O	0	0
			35	35		

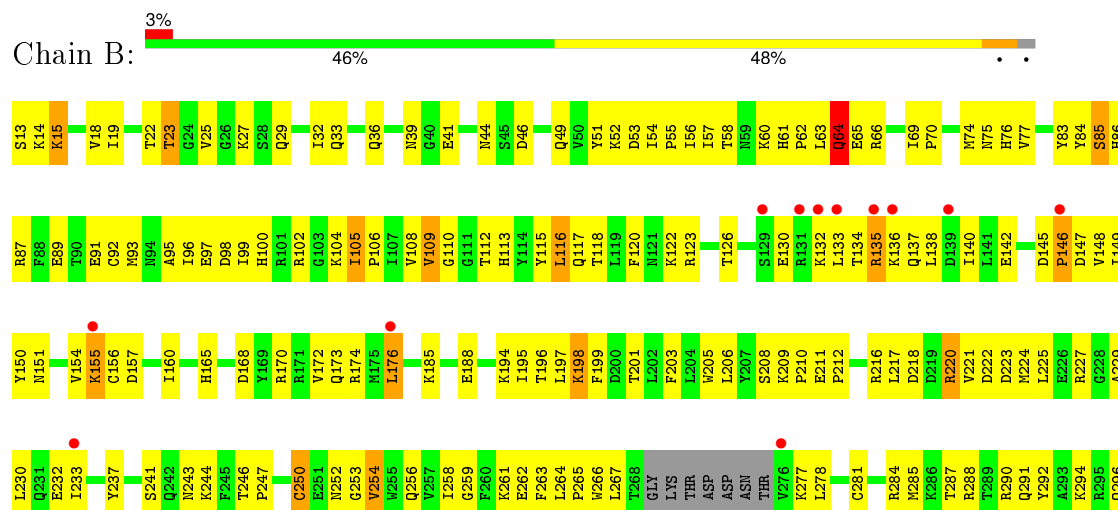
3 Residue-property plots

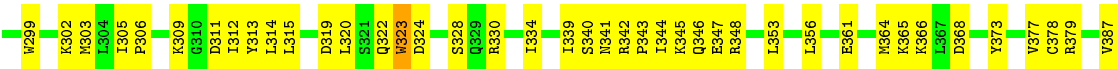
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ($\text{RSRZ} > 2$). 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 isopentenyltransferase

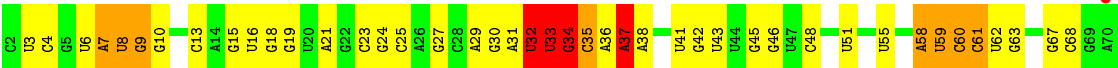


• Molecule 1: tRNA isopentenyltransferase

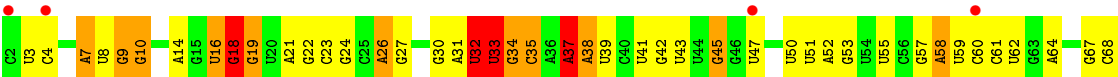




• Molecule 2: tRNA



• Molecule 2: tRNA



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	162.46Å 211.60Å 125.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.95 48.75 – 2.94	Depositor EDS
% Data completeness (in resolution range)	81.2 (50.00-2.95) 87.7 (48.75-2.94)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.43 (at 2.96Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.197 , 0.236 0.210 , 0.247	Depositor DCC
R_{free} test set	3288 reflections (8.12%)	DCC
Wilson B-factor (Å ²)	62.0	Xtriage
Anisotropy	0.848	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 60.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 42940 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9934	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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: ZN, MG, PPV

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.41	0/3411	0.59	4/4598 (0.1%)
1	B	0.41	0/3411	0.59	2/4598 (0.0%)
2	E	0.47	0/1644	0.88	11/2561 (0.4%)
2	F	0.43	0/1644	0.86	8/2561 (0.3%)
All	All	0.43	0/10110	0.71	25/14318 (0.2%)

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	E	1	0
2	F	1	0
All	All	2	0

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	32	U	C2'-C3'-O3'	8.97	129.23	109.50
2	F	37	A	C2'-C3'-O3'	8.79	128.83	109.50
2	E	37	A	C2'-C3'-O3'	8.73	128.71	109.50
2	E	34	G	C2'-C3'-O3'	8.11	127.33	109.50
2	F	34	G	C2'-C3'-O3'	8.05	127.22	109.50
2	F	18	G	C2'-C3'-O3'	7.83	126.71	109.50
2	E	33	U	C2'-C3'-O3'	7.45	125.88	109.50
2	F	32	U	C2'-C3'-O3'	7.05	125.01	109.50
2	E	60	C	N1-C1'-C2'	6.90	122.97	114.00
1	A	174	ARG	NE-CZ-NH1	6.62	123.61	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	174	ARG	NE-CZ-NH2	-6.54	117.03	120.30
2	F	33	U	C2'-C3'-O3'	6.27	123.73	113.70
2	F	34	G	C4'-C3'-O3'	6.03	125.05	113.00
1	A	342	ARG	NE-CZ-NH2	-5.90	117.35	120.30
1	B	174	ARG	NE-CZ-NH2	5.77	123.19	120.30
2	E	33	U	C4'-C3'-C2'	5.59	108.19	102.60
2	F	18	G	C4'-C3'-C2'	5.59	108.19	102.60
1	B	174	ARG	NE-CZ-NH1	-5.57	117.51	120.30
2	E	34	G	C4'-C3'-C2'	5.43	108.03	102.60
2	E	37	A	C4'-C3'-C2'	5.42	108.02	102.60
2	E	34	G	C4'-C3'-O3'	5.41	123.82	113.00
2	E	32	U	C4'-C3'-C2'	5.30	107.90	102.60
2	E	9	G	O4'-C1'-N9	5.22	112.38	108.20
1	A	342	ARG	NE-CZ-NH1	5.16	122.88	120.30
2	F	32	U	C4'-C3'-C2'	5.09	107.69	102.60

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	E	34	G	C3'
2	F	34	G	C3'

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3339	0	3346	185	0
1	B	3339	0	3345	201	0
2	E	1472	0	741	48	0
2	F	1472	0	741	52	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	4	0	0	0	0
4	B	3	0	0	0	0
4	E	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	F	1	0	0	0	0
5	A	9	0	0	2	0
5	B	9	0	0	1	0
6	A	85	0	0	38	0
6	B	104	0	0	39	0
6	E	58	0	0	15	0
6	F	35	0	0	16	0
All	All	9934	0	8173	467	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 26.

All (467) 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:B:342:ARG:HB3	6:B:473:HOH:O	1.56	1.02
1:A:60:LYS:HE3	1:A:75:ASN:HD22	1.24	1.01
2:E:9:G:H3'	6:E:89:HOH:O	1.62	1.00
1:A:195:ILE:HG12	6:A:479:HOH:O	1.65	0.95
2:E:6:U:H3	2:E:67:G:H1	1.15	0.95
1:A:198:LYS:HA	6:A:499:HOH:O	1.69	0.93
1:B:60:LYS:HE3	1:B:75:ASN:HD22	1.32	0.92
2:F:50:U:H3	2:F:64:A:H61	0.95	0.91
1:A:147:ASP:HA	6:A:477:HOH:O	1.70	0.89
2:F:19:G:H5'	2:F:60:C:H42	1.38	0.89
6:A:494:HOH:O	1:B:195:ILE:HD12	1.72	0.88
1:A:292:TYR:CE1	2:E:37:A:H5'	2.10	0.86
1:A:342:ARG:HB3	6:A:466:HOH:O	1.76	0.85
6:A:467:HOH:O	1:B:364:MET:HE3	1.77	0.85
1:A:86:HIS:HE1	6:E:113:HOH:O	1.61	0.84
2:F:9:G:H3'	6:F:78:HOH:O	1.79	0.83
1:B:406:ASN:HB3	6:B:476:HOH:O	1.78	0.83
1:B:211:GLU:HB2	1:B:212:PRO:HD3	1.61	0.83
1:B:292:TYR:CE1	2:F:37:A:H5'	2.13	0.82
1:B:365:LYS:HG2	6:B:450:HOH:O	1.80	0.81
1:B:334:ILE:HG12	1:B:346:GLN:HG3	1.62	0.81
1:A:322:GLN:HB3	6:A:465:HOH:O	1.80	0.79
1:A:221:VAL:HG13	1:A:285:MET:HE2	1.62	0.79
1:A:211:GLU:HB2	1:A:212:PRO:HD3	1.64	0.78
2:E:51:U:H3	2:E:63:G:H1	1.27	0.78
1:A:52:LYS:HG2	6:A:429:HOH:O	1.84	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:406:ASN:HB3	6:A:442:HOH:O	1.84	0.77
2:E:13:C:H5''	6:E:126:HOH:O	1.84	0.77
1:A:134:THR:HG22	1:A:136:LYS:H	1.50	0.77
1:A:322:GLN:HG3	6:A:464:HOH:O	1.87	0.75
1:A:267:LEU:HD11	6:A:500:HOH:O	1.87	0.75
2:E:45:G:H3'	6:E:75:HOH:O	1.87	0.74
1:B:221:VAL:HG13	1:B:285:MET:HE2	1.67	0.74
1:A:302:LYS:O	1:A:366:LYS:HE2	1.88	0.73
2:E:34:G:P	6:E:113:HOH:O	2.48	0.72
2:F:4:C:H42	2:F:69:G:H1	1.36	0.72
1:A:334:ILE:HG12	1:A:346:GLN:HG3	1.70	0.72
2:E:19:G:H5'	2:E:60:C:H42	1.54	0.72
1:B:134:THR:HB	1:B:137:GLN:HG3	1.72	0.72
1:B:23:THR:HG22	5:B:422:PPV:O22	1.90	0.72
1:A:117:GLN:HB3	1:A:122:LYS:HB3	1.72	0.72
1:B:117:GLN:HB3	1:B:122:LYS:HB3	1.72	0.71
1:B:112:THR:N	6:B:483:HOH:O	2.24	0.71
1:A:134:THR:HB	1:A:137:GLN:HG3	1.73	0.71
2:E:31:A:C2'	2:E:32:U:H5'	2.20	0.71
1:A:361:GLU:HA	1:A:365:LYS:HE3	1.72	0.70
1:B:225:LEU:HD23	1:B:229:ALA:HB3	1.74	0.70
1:B:123:ARG:C	1:B:123:ARG:HD2	2.13	0.69
1:A:39:ASN:OD1	1:A:104:LYS:HE2	1.92	0.69
1:B:302:LYS:O	1:B:366:LYS:HE2	1.91	0.69
2:E:55:U:OP2	6:E:116:HOH:O	2.10	0.69
1:B:205:TRP:HB2	1:B:312:ILE:HD11	1.75	0.69
1:B:134:THR:HG22	1:B:136:LYS:H	1.56	0.68
1:B:60:LYS:HB3	6:B:521:HOH:O	1.92	0.68
1:B:361:GLU:HA	1:B:365:LYS:HE3	1.74	0.68
1:B:148:VAL:HG13	6:B:460:HOH:O	1.92	0.68
1:B:256:GLN:O	1:B:261:LYS:HE3	1.94	0.68
1:A:194:LYS:HD2	1:A:196:THR:OG1	1.94	0.68
2:F:16:U:H2'	2:F:18:G:H4'	1.76	0.68
1:A:89:GLU:O	1:A:93:MET:HB2	1.94	0.67
1:B:51:TYR:OH	6:B:485:HOH:O	2.10	0.67
1:B:194:LYS:HB3	6:B:461:HOH:O	1.93	0.67
1:B:39:ASN:OD1	1:B:104:LYS:HE2	1.94	0.67
1:B:52:LYS:HG3	1:B:77:VAL:O	1.93	0.67
2:E:45:G:H5'	2:E:46:G:OP2	1.95	0.67
2:E:31:A:O2'	2:E:32:U:H5'	1.95	0.67
1:A:104:LYS:HD3	6:A:496:HOH:O	1.94	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:267:LEU:HD21	6:A:500:HOH:O	1.95	0.66
1:B:104:LYS:HD3	6:B:505:HOH:O	1.96	0.66
1:B:322:GLN:HB3	6:B:447:HOH:O	1.95	0.66
1:A:220:ARG:O	1:A:224:MET:HB2	1.95	0.66
1:B:199:PHE:N	6:B:464:HOH:O	2.05	0.66
1:B:89:GLU:O	1:B:93:MET:HB2	1.96	0.65
1:A:256:GLN:O	1:A:261:LYS:HE3	1.96	0.65
2:E:58:A:H3'	6:E:99:HOH:O	1.96	0.65
1:A:225:LEU:HD23	1:A:229:ALA:HB3	1.76	0.65
1:A:57:ILE:HA	1:A:224:MET:HE3	1.77	0.65
2:F:50:U:H3	2:F:64:A:N6	1.80	0.65
1:A:195:ILE:HD13	6:A:482:HOH:O	1.97	0.65
1:A:56:ILE:HD13	1:A:232:GLU:OE2	1.97	0.65
1:A:33:GLN:HE22	1:A:323:TRP:HE1	1.44	0.65
1:A:93:MET:O	1:A:97:GLU:HG2	1.96	0.65
1:B:303:MET:HE1	6:F:98:HOH:O	1.97	0.64
2:F:52:A:H61	2:F:62:U:H3	1.45	0.64
1:B:57:ILE:HA	1:B:224:MET:HE3	1.80	0.64
2:F:45:G:H3'	6:F:92:HOH:O	1.98	0.64
2:F:67:G:H2'	2:F:68:C:C6	2.33	0.64
1:A:258:ILE:HG13	2:E:37:A:H1'	1.80	0.64
1:B:198:LYS:HG3	6:B:464:HOH:O	1.95	0.64
2:E:27:G:N3	6:E:79:HOH:O	2.30	0.64
1:B:220:ARG:O	1:B:224:MET:HB2	1.97	0.64
2:F:9:G:C8	6:F:78:HOH:O	2.50	0.63
1:B:377:VAL:HG12	6:B:476:HOH:O	1.98	0.63
1:A:115:TYR:O	1:A:118:THR:HG22	1.98	0.63
1:A:420:LYS:O	1:A:421:ILE:HD13	1.99	0.63
1:A:54:ILE:HD12	1:A:232:GLU:HG2	1.81	0.63
1:B:54:ILE:HD12	1:B:232:GLU:HG2	1.80	0.63
1:B:62:PRO:HB2	1:B:64:GLN:HE21	1.62	0.63
2:F:19:G:H5'	2:F:60:C:N4	2.12	0.62
1:A:205:TRP:HB2	1:A:312:ILE:HD11	1.80	0.62
1:B:334:ILE:HG12	1:B:346:GLN:CG	2.29	0.62
2:E:27:G:H1	2:E:43:U:H3	1.48	0.62
1:B:168:ASP:O	1:B:172:VAL:HG23	1.99	0.61
1:A:313:TYR:CD2	1:A:348:ARG:HA	2.35	0.61
1:B:284:ARG:CZ	1:B:288:ARG:HD3	2.30	0.61
1:A:62:PRO:HB2	1:A:64:GLN:HE21	1.65	0.61
1:A:334:ILE:HG12	1:A:346:GLN:CG	2.30	0.60
2:F:18:G:H21	2:F:57:G:H2'	1.65	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:PHE:N	6:A:499:HOH:O	2.17	0.60
6:B:453:HOH:O	2:F:42:G:H5'	2.00	0.60
1:A:322:GLN:O	1:A:324:ASP:N	2.34	0.60
1:B:313:TYR:CD2	1:B:348:ARG:HA	2.36	0.60
1:B:302:LYS:HE3	6:F:76:HOH:O	2.01	0.60
1:B:387:VAL:O	1:B:387:VAL:HG23	2.01	0.60
1:B:314:LEU:O	1:B:330:ARG:NH1	2.35	0.59
1:B:58:THR:CG2	1:B:60:LYS:HG3	2.32	0.59
1:A:345:LYS:HE3	1:B:368:ASP:OD2	2.01	0.59
1:A:198:LYS:HG3	6:A:499:HOH:O	2.02	0.59
1:B:401:ARG:HD3	6:F:100:HOH:O	2.02	0.59
1:A:244:LYS:N	6:A:462:HOH:O	2.36	0.59
1:A:120:PHE:HB3	1:A:197:LEU:HA	1.83	0.59
2:E:58:A:H2'	2:E:60:C:OP2	2.02	0.59
1:B:33:GLN:HE22	1:B:323:TRP:HE1	1.48	0.58
1:A:229:ALA:O	1:A:233:ILE:HG13	2.04	0.58
1:B:157:ASP:O	1:B:160:ILE:HG22	2.04	0.58
1:B:44:ASN:HB3	1:B:74:MET:HB2	1.84	0.58
1:B:198:LYS:HA	6:B:464:HOH:O	2.03	0.58
1:B:168:ASP:HB2	6:B:438:HOH:O	2.04	0.58
1:A:138:LEU:O	1:A:142:GLU:HG2	2.04	0.58
1:A:22:THR:O	1:A:25:VAL:HG23	2.03	0.58
1:B:115:TYR:O	1:B:118:THR:HG22	2.04	0.58
1:A:314:LEU:O	1:A:330:ARG:NH1	2.36	0.58
1:B:292:TYR:O	1:B:296:GLN:HG3	2.04	0.58
1:B:264:LEU:N	1:B:265:PRO:HD2	2.18	0.58
2:F:4:C:N4	2:F:69:G:H1	2.00	0.58
1:A:44:ASN:HB3	1:A:74:MET:HB2	1.84	0.58
1:B:347:GLU:O	1:B:348:ARG:HB3	2.04	0.58
1:B:305:ILE:HB	1:B:306:PRO:HD3	1.86	0.58
2:F:31:A:O2'	2:F:32:U:H5'	2.04	0.57
1:B:149:ILE:HD11	1:B:173:GLN:HA	1.86	0.57
2:F:52:A:H2'	2:F:53:G:O4'	2.04	0.57
1:B:322:GLN:O	1:B:324:ASP:N	2.38	0.57
1:A:241:SER:HA	6:A:462:HOH:O	2.02	0.57
1:B:138:LEU:O	1:B:142:GLU:HG2	2.05	0.57
1:A:23:THR:HG22	5:A:422:PPV:O22	2.04	0.57
1:A:292:TYR:O	1:A:296:GLN:HG3	2.04	0.57
1:A:18:VAL:HG13	1:A:203:PHE:HA	1.87	0.57
1:B:61:HIS:N	6:B:521:HOH:O	2.37	0.56
2:F:27:G:H1	2:F:43:U:H3	1.53	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:170:ARG:HD3	6:F:91:HOH:O	2.05	0.56
1:B:63:LEU:C	1:B:65:GLU:H	2.09	0.56
2:E:9:G:H8	6:E:89:HOH:O	1.88	0.56
1:A:170:ARG:HD3	6:E:86:HOH:O	2.05	0.56
1:A:185:LYS:HD2	1:A:188:GLU:OE2	2.06	0.56
1:A:86:HIS:CE1	6:E:113:HOH:O	2.44	0.56
1:B:287:THR:O	1:B:291:GLN:HG3	2.06	0.56
1:B:120:PHE:HB3	1:B:197:LEU:HA	1.87	0.56
1:A:63:LEU:C	1:A:65:GLU:H	2.09	0.56
2:F:18:G:N2	2:F:57:G:H2'	2.20	0.56
1:B:92:CYS:O	1:B:96:ILE:HG12	2.05	0.55
1:A:198:LYS:CA	6:A:499:HOH:O	2.39	0.55
2:F:60:C:H5''	2:F:61:C:OP2	2.06	0.55
1:B:93:MET:O	1:B:97:GLU:HG2	2.07	0.55
1:A:54:ILE:O	1:A:54:ILE:HG23	2.07	0.55
1:A:149:ILE:HD11	1:A:173:GLN:HA	1.87	0.55
1:B:227:ARG:NH1	6:B:520:HOH:O	2.39	0.55
1:B:229:ALA:O	1:B:233:ILE:HG13	2.05	0.55
1:B:194:LYS:HD2	1:B:196:THR:OG1	2.08	0.54
1:A:145:ASP:CG	1:A:146:PRO:HD2	2.27	0.54
1:A:366:LYS:HD3	1:B:343:PRO:HG3	1.89	0.54
1:A:347:GLU:O	1:A:348:ARG:HB3	2.06	0.54
1:A:388:ALA:HA	6:A:444:HOH:O	2.07	0.54
1:B:66:ARG:NH2	6:B:521:HOH:O	2.39	0.54
2:E:15:G:H2'	2:E:16:U:C6	2.43	0.54
1:A:206:LEU:HD12	1:A:315:LEU:HB2	1.90	0.54
1:B:345:LYS:N	6:B:474:HOH:O	2.40	0.54
2:E:9:G:O2'	2:E:45:G:H2'	2.08	0.54
1:A:368:ASP:OD2	1:B:345:LYS:HE3	2.08	0.54
1:A:264:LEU:N	1:A:265:PRO:HD2	2.24	0.53
1:A:123:ARG:HD2	1:A:123:ARG:C	2.29	0.53
1:A:268:THR:HA	6:A:432:HOH:O	2.09	0.53
1:B:55:PRO:HA	1:B:58:THR:HG22	1.89	0.53
2:E:27:G:N2	6:E:79:HOH:O	2.33	0.53
1:A:122:LYS:HE3	6:A:482:HOH:O	2.08	0.53
2:E:29:A:H2'	2:E:30:G:H8	1.72	0.53
1:A:284:ARG:CZ	1:A:288:ARG:HD3	2.39	0.53
1:B:148:VAL:HG22	6:B:460:HOH:O	2.08	0.53
1:A:387:VAL:HG23	1:A:387:VAL:O	2.09	0.53
1:B:259:GLY:HA2	1:B:285:MET:SD	2.50	0.52
1:B:56:ILE:HD13	1:B:232:GLU:OE2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:284:ARG:NH2	1:B:288:ARG:HH11	2.07	0.52
1:B:117:GLN:HB2	6:B:523:HOH:O	2.09	0.52
2:F:52:A:N6	2:F:62:U:H3	2.06	0.52
1:B:95:ALA:O	1:B:99:ILE:HG13	2.10	0.52
2:F:31:A:C2'	2:F:32:U:H5'	2.40	0.52
1:A:55:PRO:HA	1:A:58:THR:HG22	1.91	0.52
1:A:13:SER:O	1:A:15:LYS:HD2	2.10	0.52
1:A:262:GLU:HG3	1:A:284:ARG:HG3	1.92	0.52
1:B:62:PRO:HB2	1:B:64:GLN:NE2	2.25	0.52
1:A:52:LYS:HG3	1:A:77:VAL:O	2.10	0.52
2:F:35:C:H5	6:F:99:HOH:O	1.92	0.52
1:A:58:THR:CG2	1:A:60:LYS:HG3	2.39	0.51
1:B:54:ILE:O	1:B:54:ILE:HG23	2.10	0.51
1:A:83:TYR:N	6:A:448:HOH:O	2.42	0.51
2:F:7:A:H2'	6:F:96:HOH:O	2.10	0.51
1:A:323:TRP:N	6:A:465:HOH:O	2.44	0.51
2:F:9:G:H5'	2:F:10:G:OP2	2.10	0.51
1:B:53:ASP:O	1:B:54:ILE:HG22	2.10	0.51
1:A:157:ASP:O	1:A:160:ILE:HG22	2.10	0.51
1:A:92:CYS:O	1:A:96:ILE:HG12	2.11	0.51
1:A:60:LYS:HE3	1:A:75:ASN:ND2	2.09	0.51
2:F:38:A:H2'	2:F:38:A:N3	2.26	0.51
1:A:259:GLY:HA2	1:A:285:MET:SD	2.50	0.51
1:B:322:GLN:C	6:B:447:HOH:O	2.50	0.51
1:A:32:ILE:O	1:A:36:GLN:HG3	2.10	0.51
2:F:50:U:H2'	2:F:51:U:O4'	2.12	0.50
2:F:58:A:H3'	6:F:88:HOH:O	2.11	0.50
1:B:258:ILE:HG13	2:F:37:A:H1'	1.92	0.50
1:A:53:ASP:O	1:A:54:ILE:HG22	2.11	0.50
2:E:7:A:H4'	2:E:8:U:OP2	2.12	0.50
1:A:55:PRO:HG2	1:A:56:ILE:HD12	1.94	0.50
1:B:23:THR:HG23	6:B:503:HOH:O	2.11	0.50
1:A:400:SER:HB3	2:E:43:U:H5"	1.93	0.50
1:A:85:SER:N	6:A:431:HOH:O	2.41	0.50
1:A:378:CYS:SG	1:A:403:HIS:HA	2.51	0.50
1:A:51:TYR:OH	6:A:448:HOH:O	2.19	0.50
1:A:95:ALA:O	1:A:99:ILE:HG13	2.11	0.50
1:B:150:TYR:O	1:B:154:VAL:HG23	2.12	0.50
2:F:41:U:H2'	2:F:42:G:H8	1.77	0.50
1:B:237:TYR:CG	1:B:267:LEU:HD13	2.46	0.50
1:A:168:ASP:O	1:A:172:VAL:HG23	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:TYR:CG	1:A:267:LEU:HD13	2.45	0.50
1:A:305:ILE:HB	1:A:306:PRO:HD3	1.94	0.50
1:A:409:ARG:HG3	1:A:410:ASN:N	2.25	0.50
1:A:54:ILE:HG13	1:A:57:ILE:HG12	1.93	0.50
1:A:284:ARG:NH2	1:A:288:ARG:HH11	2.09	0.49
1:B:323:TRP:N	6:B:447:HOH:O	2.44	0.49
1:B:262:GLU:HG3	1:B:284:ARG:HG3	1.94	0.49
2:F:41:U:H2'	2:F:42:G:C8	2.47	0.49
2:E:29:A:H2'	2:E:30:G:C8	2.47	0.49
1:B:393:TYR:HB2	6:B:499:HOH:O	2.12	0.49
2:E:23:C:H2'	2:E:24:G:H8	1.77	0.49
2:E:9:G:C3'	6:E:89:HOH:O	2.38	0.49
1:A:377:VAL:HG12	6:A:442:HOH:O	2.11	0.49
1:B:241:SER:HA	6:B:466:HOH:O	2.12	0.49
1:A:254:VAL:O	1:A:257:VAL:HG23	2.11	0.49
1:A:340:SER:CB	1:A:342:ARG:HD3	2.42	0.49
1:B:156:CYS:O	1:B:157:ASP:HB3	2.13	0.49
1:B:206:LEU:HD12	1:B:315:LEU:HB2	1.95	0.49
1:B:54:ILE:HD12	1:B:232:GLU:CG	2.43	0.49
1:B:409:ARG:HG3	1:B:410:ASN:N	2.28	0.49
1:B:87:ARG:NE	1:B:91:GLU:OE2	2.46	0.49
1:A:237:TYR:CE2	1:A:267:LEU:HB3	2.48	0.48
2:E:19:G:H5'	2:E:60:C:N4	2.23	0.48
2:F:14:A:N1	2:F:22:G:H1'	2.28	0.48
1:A:165:HIS:CD2	2:E:41:U:H4'	2.48	0.48
1:A:216:ARG:NH1	1:A:320:LEU:HG	2.29	0.48
1:B:54:ILE:HG13	1:B:57:ILE:HG12	1.95	0.48
1:A:27:LYS:N	5:A:422:PPV:O11	2.38	0.48
1:B:290:ARG:O	1:B:294:LYS:HG3	2.12	0.48
1:B:18:VAL:HG13	1:B:203:PHE:HA	1.95	0.48
1:A:44:ASN:CB	1:A:74:MET:HB2	2.42	0.48
1:B:185:LYS:HD2	1:B:188:GLU:OE2	2.13	0.48
1:B:32:ILE:O	1:B:36:GLN:HG3	2.14	0.48
2:E:35:C:OP2	2:E:36:A:N7	2.46	0.48
1:B:44:ASN:HD21	1:B:46:ASP:CB	2.27	0.48
1:B:135:ARG:HA	1:B:135:ARG:CZ	2.44	0.48
1:A:63:LEU:HD12	1:A:63:LEU:H	1.79	0.48
1:B:98:ASP:O	1:B:102:ARG:HG3	2.13	0.48
1:B:343:PRO:HD2	6:B:473:HOH:O	2.12	0.48
1:B:225:LEU:HA	1:B:229:ALA:HB3	1.95	0.48
1:B:145:ASP:HB3	1:B:148:VAL:CG1	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:58:A:N6	2:E:61:C:C2	2.82	0.48
1:B:197:LEU:HD21	1:B:201:THR:HB	1.96	0.48
6:B:510:HOH:O	2:F:35:C:H1'	2.13	0.48
1:B:299:TRP:CE3	1:B:303:MET:HG3	2.49	0.47
2:F:35:C:H5''	6:F:74:HOH:O	2.13	0.47
1:A:19:ILE:HB	1:A:109:VAL:HG13	1.96	0.47
1:A:225:LEU:HA	1:A:229:ALA:HB3	1.96	0.47
1:B:237:TYR:CE2	1:B:267:LEU:HB3	2.49	0.47
1:B:324:ASP:HA	1:B:328:SER:HB2	1.97	0.47
1:B:392:LYS:O	1:B:396:ILE:HG13	2.14	0.47
1:B:113:HIS:O	1:B:116:LEU:HB2	2.14	0.47
1:A:48:MET:SD	1:A:257:VAL:HG22	2.55	0.47
1:A:174:ARG:NH1	2:E:34:G:OP1	2.45	0.47
1:A:246:THR:HB	1:A:247:PRO:HD2	1.96	0.47
1:A:54:ILE:HD12	1:A:232:GLU:CG	2.43	0.47
1:B:230:LEU:HD23	1:B:233:ILE:HD12	1.96	0.47
1:A:340:SER:HB2	1:A:342:ARG:HD3	1.97	0.47
1:B:261:LYS:HG2	6:B:442:HOH:O	2.15	0.47
1:B:39:ASN:HD21	1:B:104:LYS:HD2	1.79	0.47
1:B:198:LYS:CA	6:B:464:HOH:O	2.61	0.47
1:B:393:TYR:HD2	1:B:396:ILE:HD12	1.80	0.47
1:B:246:THR:HB	1:B:247:PRO:HD2	1.96	0.47
1:A:343:PRO:HG3	1:B:366:LYS:HD3	1.96	0.47
2:F:42:G:H4'	6:F:77:HOH:O	2.14	0.47
1:B:63:LEU:HD12	1:B:63:LEU:H	1.79	0.47
2:E:62:U:O2'	2:E:63:G:H5'	2.15	0.47
1:A:25:VAL:HA	6:A:472:HOH:O	2.14	0.46
1:A:135:ARG:CZ	1:A:135:ARG:HA	2.45	0.46
1:B:195:ILE:HG22	1:B:195:ILE:O	2.15	0.46
1:B:252:ASN:HA	6:B:510:HOH:O	2.15	0.46
1:B:13:SER:O	1:B:15:LYS:HD2	2.15	0.46
1:B:145:ASP:CG	1:B:146:PRO:HD2	2.35	0.46
2:E:67:G:H2'	2:E:68:C:C6	2.50	0.46
2:E:31:A:H2'	2:E:32:U:H5'	1.94	0.46
1:A:342:ARG:NH1	6:A:437:HOH:O	2.48	0.46
1:A:170:ARG:HG2	2:E:35:C:N4	2.30	0.46
1:A:364:MET:CE	1:B:345:LYS:HB3	2.45	0.46
1:A:306:PRO:HB3	1:A:363:THR:HG22	1.97	0.46
1:B:393:TYR:CD2	1:B:396:ILE:HD12	2.50	0.46
1:B:244:LYS:N	6:B:466:HOH:O	2.48	0.46
2:F:18:G:O2'	2:F:57:G:N2	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:26:A:H3'	2:F:27:G:H8	1.80	0.45
1:B:99:ILE:CG2	1:B:104:LYS:HB3	2.47	0.45
2:F:51:U:H2'	6:F:83:HOH:O	2.15	0.45
1:A:22:THR:HA	1:A:296:GLN:OE1	2.17	0.45
1:B:19:ILE:HB	1:B:109:VAL:HG13	1.99	0.45
1:B:319:ASP:OD2	1:B:322:GLN:HG2	2.17	0.45
2:F:30:G:O2'	2:F:31:A:H5'	2.17	0.45
1:A:111:GLY:N	6:A:427:HOH:O	2.50	0.45
2:F:23:C:O2'	2:F:24:G:H5'	2.17	0.45
1:B:258:ILE:CG1	2:F:38:A:OP1	2.64	0.45
2:E:41:U:H2'	2:E:42:G:H8	1.82	0.45
1:B:46:ASP:HB3	1:B:49:GLN:HB2	1.97	0.45
2:F:38:A:H5'	2:F:39:U:OP2	2.17	0.45
1:A:121:ASN:HD22	1:A:123:ARG:HH21	1.65	0.45
1:B:266:TRP:O	1:B:266:TRP:CG	2.69	0.45
1:B:99:ILE:HG23	1:B:104:LYS:HB3	1.97	0.45
1:B:216:ARG:NH1	6:B:457:HOH:O	2.50	0.45
1:A:113:HIS:O	1:A:116:LEU:HB2	2.17	0.45
1:B:250:CYS:SG	1:B:261:LYS:HB3	2.56	0.45
1:B:165:HIS:CD2	2:F:41:U:H4'	2.52	0.45
1:A:156:CYS:O	1:A:157:ASP:HB3	2.17	0.45
1:A:87:ARG:NE	1:A:91:GLU:OE2	2.50	0.45
1:A:98:ASP:O	1:A:102:ARG:HG3	2.17	0.45
1:B:133:LEU:HB3	1:B:138:LEU:HG	1.99	0.44
1:B:277:LYS:HB3	6:B:525:HOH:O	2.17	0.44
1:A:171:ARG:NE	2:E:33:U:OP1	2.51	0.44
1:B:25:VAL:HG12	1:B:25:VAL:O	2.16	0.44
1:A:216:ARG:O	1:A:220:ARG:HB2	2.17	0.44
1:B:211:GLU:HB2	1:B:212:PRO:CD	2.42	0.44
1:A:43:ILE:O	1:A:108:VAL:HA	2.17	0.44
1:A:62:PRO:HB2	1:A:64:GLN:NE2	2.32	0.44
1:A:364:MET:O	1:B:345:LYS:NZ	2.51	0.44
1:B:347:GLU:HB2	6:B:513:HOH:O	2.18	0.44
1:A:345:LYS:N	6:A:438:HOH:O	2.45	0.44
2:E:41:U:H2'	2:E:42:G:C8	2.52	0.44
1:A:45:SER:OG	1:A:110:GLY:HA3	2.18	0.44
1:A:51:TYR:CD1	1:A:254:VAL:HG13	2.53	0.44
1:B:284:ARG:NH2	1:B:288:ARG:HD3	2.31	0.44
1:A:195:ILE:N	6:A:455:HOH:O	2.51	0.44
1:B:230:LEU:HA	1:B:233:ILE:HD12	1.99	0.44
1:B:57:ILE:HG13	1:B:58:THR:N	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:16:U:H3	2:E:59:U:H3	1.65	0.44
1:A:230:LEU:HD23	1:A:233:ILE:HD12	1.99	0.43
1:B:63:LEU:O	1:B:65:GLU:N	2.51	0.43
1:A:224:MET:O	1:A:229:ALA:HB2	2.17	0.43
1:B:378:CYS:SG	1:B:403:HIS:HA	2.58	0.43
1:B:147:ASP:O	1:B:151:ASN:ND2	2.51	0.43
1:A:89:GLU:HG3	1:A:118:THR:OG1	2.18	0.43
1:B:51:TYR:CD1	1:B:254:VAL:HG13	2.53	0.43
1:A:420:LYS:HG2	1:A:420:LYS:O	2.19	0.43
2:E:24:G:C6	2:E:25:C:C4	3.06	0.43
1:B:339:ILE:C	1:B:341:ASN:H	2.21	0.43
1:A:84:TYR:CE1	1:A:86:HIS:HB2	2.53	0.43
1:B:77:VAL:HG21	1:B:83:TYR:CD1	2.53	0.43
1:A:319:ASP:OD2	1:A:322:GLN:HG2	2.19	0.43
1:A:39:ASN:HD21	1:A:104:LYS:HD2	1.84	0.43
1:A:311:ASP:OD1	1:A:348:ARG:NH2	2.51	0.43
1:A:320:LEU:C	1:A:322:GLN:N	2.70	0.43
1:A:57:ILE:HG13	1:A:58:THR:N	2.33	0.43
1:A:75:ASN:C	6:A:429:HOH:O	2.56	0.43
1:B:54:ILE:N	1:B:55:PRO:CD	2.82	0.43
1:A:287:THR:O	1:A:291:GLN:HG3	2.19	0.43
1:B:105:ILE:HD12	1:B:106:PRO:HD2	2.00	0.43
2:F:31:A:P	6:F:98:HOH:O	2.76	0.43
1:B:63:LEU:HD12	1:B:63:LEU:N	2.33	0.43
1:A:145:ASP:HB3	1:A:148:VAL:CG1	2.49	0.43
2:E:58:A:O2'	2:E:59:U:P	2.77	0.43
1:B:224:MET:O	1:B:229:ALA:HB2	2.18	0.43
1:B:216:ARG:NH1	1:B:320:LEU:HG	2.34	0.43
1:B:340:SER:O	1:B:342:ARG:HG3	2.19	0.42
1:B:58:THR:HG23	1:B:60:LYS:HG3	1.99	0.42
2:F:58:A:N6	2:F:61:C:C4	2.87	0.42
1:B:218:ASP:O	1:B:221:VAL:HB	2.19	0.42
2:F:45:G:N2	6:F:100:HOH:O	2.49	0.42
2:E:23:C:H2'	2:E:24:G:C8	2.54	0.42
1:B:84:TYR:CE1	1:B:86:HIS:HB2	2.54	0.42
1:A:58:THR:HG23	1:A:60:LYS:HG3	2.00	0.42
2:F:18:G:N2	2:F:57:G:C4	2.88	0.42
1:B:89:GLU:HG3	1:B:118:THR:OG1	2.20	0.42
1:A:133:LEU:HB3	1:A:138:LEU:HG	2.01	0.42
1:A:382:ASP:OD2	1:A:384:LYS:HD2	2.19	0.42
1:B:27:LYS:HE3	1:B:110:GLY:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:135:ARG:HG2	1:B:135:ARG:HH11	1.83	0.42
1:B:400:SER:HB3	2:F:43:U:H5'	2.02	0.42
6:A:456:HOH:O	2:E:35:C:H5	2.01	0.42
1:A:148:VAL:HG13	1:A:149:ILE:N	2.33	0.42
1:B:155:LYS:HE2	6:B:459:HOH:O	2.19	0.42
1:A:350:PRO:O	1:A:351:LYS:C	2.57	0.42
1:B:46:ASP:CG	2:F:37:A:H62	2.23	0.42
1:B:365:LYS:HE2	6:B:450:HOH:O	2.20	0.42
1:B:148:VAL:HG13	1:B:149:ILE:N	2.33	0.42
1:B:320:LEU:C	1:B:322:GLN:N	2.70	0.42
1:A:63:LEU:C	1:A:65:GLU:N	2.73	0.42
1:A:349:ALA:HB2	1:A:357:LEU:HD12	2.01	0.42
1:B:135:ARG:NE	1:B:135:ARG:HA	2.35	0.42
1:A:134:THR:HG22	1:A:135:ARG:N	2.35	0.42
1:A:116:LEU:HD12	1:A:116:LEU:HA	1.76	0.42
1:A:217:LEU:O	1:A:221:VAL:HG23	2.19	0.42
1:B:155:LYS:NZ	1:B:155:LYS:HB2	2.35	0.42
1:A:380:ASN:CG	1:A:402:ARG:HH22	2.23	0.42
1:A:366:LYS:NZ	6:A:440:HOH:O	2.52	0.42
1:B:149:ILE:HG13	1:B:176:LEU:HD12	2.02	0.42
1:B:353:LEU:HB3	1:B:356:LEU:HD12	2.01	0.42
1:B:217:LEU:O	1:B:221:VAL:HG23	2.20	0.42
1:B:324:ASP:O	1:B:328:SER:HB2	2.20	0.42
1:B:315:LEU:HD23	1:B:330:ARG:HD3	2.02	0.42
1:A:324:ASP:HA	1:A:328:SER:HB2	2.02	0.41
1:B:41:GLU:OE2	1:B:70:PRO:HG2	2.20	0.41
2:E:9:G:C8	6:E:89:HOH:O	2.57	0.41
2:F:33:U:H5'	2:F:33:U:H6	1.85	0.41
2:F:16:U:O5'	2:F:16:U:H6	2.03	0.41
1:A:322:GLN:C	6:A:465:HOH:O	2.57	0.41
1:B:220:ARG:O	1:B:220:ARG:HD2	2.21	0.41
1:A:262:GLU:HG2	1:A:285:MET:HA	2.03	0.41
1:A:99:ILE:HG23	1:A:104:LYS:HB3	2.02	0.41
1:B:22:THR:O	1:B:25:VAL:HG23	2.20	0.41
1:A:393:TYR:HD2	1:A:396:ILE:HD12	1.85	0.41
1:B:208:SER:O	1:B:209:LYS:C	2.59	0.41
1:A:77:VAL:HG21	1:A:83:TYR:CD1	2.55	0.41
1:B:66:ARG:O	1:B:69:ILE:HG12	2.20	0.41
1:A:218:ASP:O	1:A:221:VAL:HB	2.21	0.41
1:B:145:ASP:HB3	1:B:148:VAL:HG11	2.01	0.41
1:B:264:LEU:N	1:B:265:PRO:CD	2.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:LYS:HG3	6:E:121:HOH:O	2.20	0.41
1:B:263:PHE:HA	1:B:281:CYS:SG	2.61	0.41
1:A:330:ARG:HH22	1:A:348:ARG:N	2.19	0.41
1:B:63:LEU:C	1:B:65:GLU:N	2.74	0.41
1:A:373:TYR:O	1:A:387:VAL:HA	2.21	0.41
1:B:266:TRP:CE2	1:B:278:LEU:HD13	2.55	0.41
1:A:290:ARG:O	1:A:294:LYS:HG3	2.20	0.41
2:E:30:G:O2'	2:E:31:A:H5'	2.21	0.41
1:A:99:ILE:CG2	1:A:104:LYS:HB3	2.51	0.41
1:B:373:TYR:O	1:B:387:VAL:HA	2.21	0.41
1:A:165:HIS:CG	1:A:166:PRO:HD2	2.56	0.41
1:B:100:HIS:C	1:B:102:ARG:H	2.24	0.41
1:A:393:TYR:CD2	1:A:396:ILE:HD12	2.56	0.41
1:A:266:TRP:CG	1:A:266:TRP:O	2.73	0.41
1:A:76:HIS:N	6:A:429:HOH:O	2.54	0.41
1:B:55:PRO:HG2	1:B:56:ILE:HD12	2.02	0.41
1:B:334:ILE:HG23	1:B:344:ILE:HG21	2.02	0.41
2:E:62:U:H2'	2:E:63:G:H8	1.86	0.41
1:A:334:ILE:CG1	1:A:346:GLN:HG3	2.46	0.41
1:B:311:ASP:OD1	1:B:348:ARG:NH2	2.53	0.41
1:B:134:THR:HG22	1:B:135:ARG:N	2.35	0.41
1:A:149:ILE:HG13	1:A:176:LEU:HD12	2.03	0.41
1:A:156:CYS:HB2	6:A:498:HOH:O	2.21	0.41
1:B:222:ASP:C	1:B:224:MET:H	2.23	0.40
1:B:116:LEU:HA	1:B:116:LEU:HD12	1.87	0.40
1:A:223:ASP:HA	1:A:226:GLU:HB2	2.04	0.40
1:B:210:PRO:HG2	6:B:439:HOH:O	2.21	0.40
2:F:9:G:H8	6:F:78:HOH:O	1.93	0.40
1:B:136:LYS:O	1:B:140:ILE:HG13	2.21	0.40
2:E:3:U:H2'	2:E:4:C:O4'	2.21	0.40
1:A:63:LEU:O	1:A:65:GLU:N	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	398/409 (97%)	341 (86%)	53 (13%)	4 (1%)	19	58
1	B	398/409 (97%)	339 (85%)	52 (13%)	7 (2%)	11	42
All	All	796/818 (97%)	680 (85%)	105 (13%)	11 (1%)	14	49

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	323	TRP
1	A	64	GLN
1	B	64	GLN
1	B	323	TRP
1	B	76	HIS
1	A	76	HIS
1	A	146	PRO
1	B	146	PRO
1	B	250	CYS
1	B	85	SER
1	B	253	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	
1	A	370/376 (98%)	343 (93%)	27 (7%)	17	50
1	B	370/376 (98%)	346 (94%)	24 (6%)	21	56
All	All	740/752 (98%)	689 (93%)	51 (7%)	19	54

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

Mol	Chain	Res	Type
1	A	14	LYS
1	A	15	LYS

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Mol	Chain	Res	Type
1	A	23	THR
1	A	29	GLN
1	A	64	GLN
1	A	85	SER
1	A	105	ILE
1	A	108	VAL
1	A	109	VAL
1	A	116	LEU
1	A	123	ARG
1	A	126	THR
1	A	130	GLU
1	A	132	LYS
1	A	135	ARG
1	A	155	LYS
1	A	174	ARG
1	A	176	LEU
1	A	198	LYS
1	A	220	ARG
1	A	223	ASP
1	A	239	TYR
1	A	243	ASN
1	A	254	VAL
1	A	309	LYS
1	A	379	ARG
1	A	419	TRP
1	B	14	LYS
1	B	15	LYS
1	B	23	THR
1	B	29	GLN
1	B	64	GLN
1	B	85	SER
1	B	105	ILE
1	B	108	VAL
1	B	109	VAL
1	B	116	LEU
1	B	126	THR
1	B	130	GLU
1	B	132	LYS
1	B	135	ARG
1	B	155	LYS
1	B	176	LEU
1	B	198	LYS

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Mol	Chain	Res	Type
1	B	220	ARG
1	B	223	ASP
1	B	243	ASN
1	B	254	VAL
1	B	309	LYS
1	B	379	ARG
1	B	420	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	44	ASN
1	A	64	GLN
1	A	75	ASN
1	A	117	GLN
1	A	121	ASN
1	A	151	ASN
1	A	249	GLN
1	A	256	GLN
1	A	341	ASN
1	B	33	GLN
1	B	44	ASN
1	B	64	GLN
1	B	75	ASN
1	B	117	GLN
1	B	121	ASN
1	B	151	ASN
1	B	249	GLN
1	B	341	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	E	68/69 (98%)	11 (16%)	7 (10%)
2	F	68/69 (98%)	21 (30%)	6 (8%)
All	All	136/138 (98%)	32 (23%)	13 (9%)

All (32) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	E	8	U
2	E	10	G
2	E	18	G
2	E	21	A
2	E	33	U
2	E	34	G
2	E	35	C
2	E	38	A
2	E	48	C
2	E	59	U
2	E	61	C
2	F	3	U
2	F	7	A
2	F	8	U
2	F	9	G
2	F	10	G
2	F	16	U
2	F	18	G
2	F	19	G
2	F	21	A
2	F	26	A
2	F	32	U
2	F	33	U
2	F	34	G
2	F	35	C
2	F	37	A
2	F	38	A
2	F	45	G
2	F	47	U
2	F	55	U
2	F	58	A
2	F	59	U

All (13) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	E	7	A
2	E	32	U
2	E	33	U
2	E	34	G
2	E	37	A
2	E	48	C
2	E	58	A

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Mol	Chain	Res	Type
2	F	18	G
2	F	32	U
2	F	33	U
2	F	34	G
2	F	37	A
2	F	58	A

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 14 ligands modelled in this entry, 12 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	$\# Z > 2$
5	PPV	A	422	4	6,8,8	2.36	2 (33%)	11,13,13	1.47	1 (9%)
5	PPV	B	422	4	6,8,8	2.33	2 (33%)	11,13,13	2.08	4 (36%)

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
5	PPV	A	422	4	-	0/6/6/6	0/0/0/0
5	PPV	B	422	4	-	0/6/6/6	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	422	PPV	P2-O22	3.10	1.61	1.51
5	A	422	PPV	P2-O22	3.12	1.61	1.51
5	B	422	PPV	P1-O31	3.15	1.61	1.51
5	A	422	PPV	P1-O31	3.15	1.61	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	422	PPV	P2-OPP-P1	-3.31	121.58	132.67
5	B	422	PPV	P2-OPP-P1	-2.74	123.49	132.67
5	B	422	PPV	O21-P1-OPP	-2.11	95.52	105.09
5	B	422	PPV	O11-P1-O31	3.47	121.73	110.58
5	B	422	PPV	O21-P1-O11	3.95	122.43	107.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	422	PPV	2	0
5	B	422	PPV	1	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	A	402/409 (98%)	0.20	9 (2%) 65 44	48, 78, 113, 141	0
1	B	402/409 (98%)	0.26	14 (3%) 48 29	51, 80, 118, 141	0
2	E	69/69 (100%)	0.43	2 (2%) 55 35	58, 98, 166, 184	0
2	F	69/69 (100%)	0.61	6 (8%) 13 6	61, 135, 196, 199	0
All	All	942/956 (98%)	0.27	31 (3%) 50 31	48, 81, 138, 199	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	276	VAL	4.4
1	B	276	VAL	3.8
1	B	135	ARG	3.6
1	B	132	LYS	3.4
1	B	136	LYS	3.3
1	B	155	LYS	3.3
1	B	418	LYS	3.2
2	E	71	G	3.1
1	B	146	PRO	3.0
2	F	47	U	2.9
1	B	129	SER	2.8
1	A	130	GLU	2.7
1	B	419	TRP	2.6
2	F	4	C	2.6
1	B	131	ARG	2.6
2	F	2	C	2.4
1	B	176	LEU	2.4
1	A	146	PRO	2.4
2	E	70	A	2.3
1	A	416	PHE	2.3
2	F	71	G	2.3

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Mol	Chain	Res	Type	RSRZ
2	F	60	C	2.2
1	B	233	ILE	2.2
1	A	419	TRP	2.2
1	A	128	SER	2.2
1	A	132	LYS	2.2
1	A	230	LEU	2.2
1	B	133	LEU	2.1
1	B	139	ASP	2.0
2	F	70	A	2.0
1	A	377	VAL	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	MG	A	4	1/1	0.90	0.38	13.89	64,64,64,64	0
5	PPV	B	422	9/9	0.98	0.18	-0.37	91,92,93,93	0
5	PPV	A	422	9/9	0.98	0.14	-1.30	93,94,96,96	0
3	ZN	B	1	1/1	0.99	0.16	-	72,72,72,72	0
4	MG	B	4	1/1	0.81	0.24	-	68,68,68,68	0
4	MG	B	5	1/1	0.69	0.14	-	69,69,69,69	0
4	MG	B	2	1/1	0.98	0.10	-	50,50,50,50	0
4	MG	A	2	1/1	0.95	0.19	-	46,46,46,46	0
4	MG	A	7	1/1	0.52	0.30	-	94,94,94,94	0
4	MG	E	73	1/1	0.73	0.13	-	76,76,76,76	0
4	MG	A	6	1/1	0.80	0.23	-	74,74,74,74	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	ZN	A	1	1/1	0.99	0.18	-	79,79,79,79	0
4	MG	F	72	1/1	0.38	0.17	-	92,92,92,92	0
4	MG	E	72	1/1	0.88	0.19	-	71,71,71,71	0

6.5 Other polymers [i](#)

There are no such residues in this entry.