



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:50 PM UTC

PDB ID : 1KOG / pdb_00001kog
Title : Crystal structure of E. coli threonyl-tRNA synthetase interacting with the essential domain of its mRNA operator
Authors : Torres-Larrios, A.; Dock-Bregeon, A.C.; Romby, P.; Rees, B.; Sankaranarayanan, R.; Caillet, J.; Springer, M.; Ehresmann, C.; Ehresmann, B.; Moras, D.
Deposited on : 2001-12-20
Resolution : 3.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

The following versions of software and data (see [references](#) ①) were used in the production of this report:

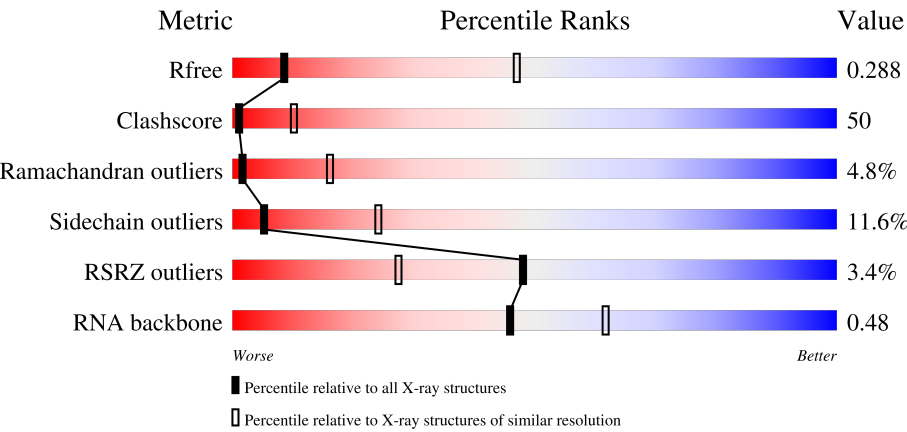
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtrriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)
RNA backbone	3983	1010 (3.98-3.02)

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

Mol	Chain	Length	Quality of chain
1	I	37	11% 32% 38% 22% 8%
1	J	37	5% 16% 54% 16% 14%
1	K	37	35% 14% 54% 16% 16%
1	L	37	3% 27% 43% 19% 11%

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Mol	Chain	Length	Quality of chain
1	M	37	
1	N	37	
1	O	37	
1	P	37	
2	A	401	
2	B	401	
2	C	401	
2	D	401	
2	E	401	
2	F	401	
2	G	401	
2	H	401	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 32922 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 Threonyl-tRNA synthetase mRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	I	37	Total	C	N	O	P	0	0	0
			785	349	132	267	37			
1	J	37	Total	C	N	O	P	0	0	0
			785	349	132	267	37			
1	K	37	Total	C	N	O	P	0	0	0
			785	349	132	267	37			
1	L	37	Total	C	N	O	P	0	0	0
			785	349	132	267	37			
1	M	37	Total	C	N	O	P	0	0	0
			785	349	132	267	37			
1	N	37	Total	C	N	O	P	0	0	0
			785	349	132	267	37			
1	O	37	Total	C	N	O	P	0	0	0
			785	349	132	267	37			
1	P	37	Total	C	N	O	P	0	0	0
			785	349	132	267	37			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	401	Total	C	N	O	S	0	0	0
			3278	2069	576	610	23			
2	B	401	Total	C	N	O	S	0	0	0
			3278	2069	576	610	23			
2	C	401	Total	C	N	O	S	0	0	0
			3278	2069	576	610	23			
2	D	401	Total	C	N	O	S	0	0	0
			3278	2069	576	610	23			
2	E	401	Total	C	N	O	S	0	0	0
			3278	2069	576	610	23			
2	F	401	Total	C	N	O	S	0	0	0
			3278	2069	576	610	23			

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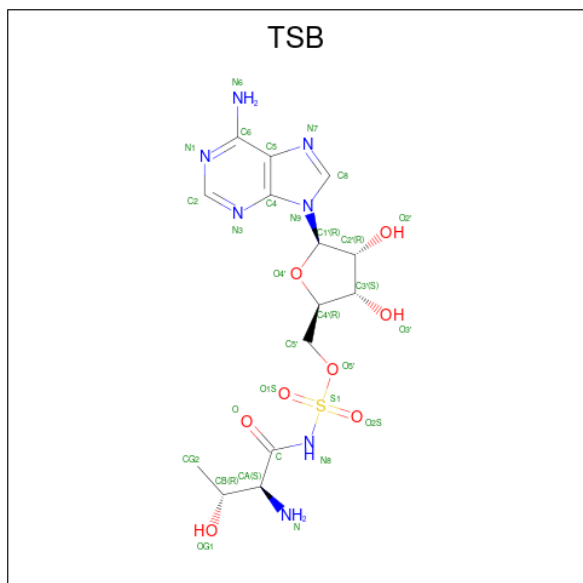
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	401	Total 3278	C 2069	N 576	O 610	S 23	0	0	0
2	H	401	Total 3278	C 2069	N 576	O 610	S 23	0	0	0

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Zn 1 1	0	0
3	B	1	Total Zn 1 1	0	0
3	C	1	Total Zn 1 1	0	0
3	D	1	Total Zn 1 1	0	0
3	E	1	Total Zn 1 1	0	0
3	F	1	Total Zn 1 1	0	0
3	G	1	Total Zn 1 1	0	0
3	H	1	Total Zn 1 1	0	0

- Molecule 4 is 5'-O-(N-(L-THREONYL)-SULFAMOYL)ADENOSINE (CCD ID: TSB) (formula: $C_{14}H_{21}N_7O_8S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			30	14	7	8	1		
4	B	1	Total	C	N	O	S	0	0
			30	14	7	8	1		
4	C	1	Total	C	N	O	S	0	0
			30	14	7	8	1		
4	D	1	Total	C	N	O	S	0	0
			30	14	7	8	1		
4	E	1	Total	C	N	O	S	0	0
			30	14	7	8	1		
4	F	1	Total	C	N	O	S	0	0
			30	14	7	8	1		
4	G	1	Total	C	N	O	S	0	0
			30	14	7	8	1		
4	H	1	Total	C	N	O	S	0	0
			30	14	7	8	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	I	11	Total	O	0	0
			11	11		
5	J	12	Total	O	0	0
			12	12		
5	K	1	Total	O	0	0
			1	1		
5	L	8	Total	O	0	0
			8	8		
5	M	6	Total	O	0	0
			6	6		
5	N	9	Total	O	0	0
			9	9		
5	O	11	Total	O	0	0
			11	11		
5	P	9	Total	O	0	0
			9	9		
5	A	10	Total	O	0	0
			10	10		
5	B	15	Total	O	0	0
			15	15		
5	C	11	Total	O	0	0
			11	11		
5	D	14	Total	O	0	0
			14	14		

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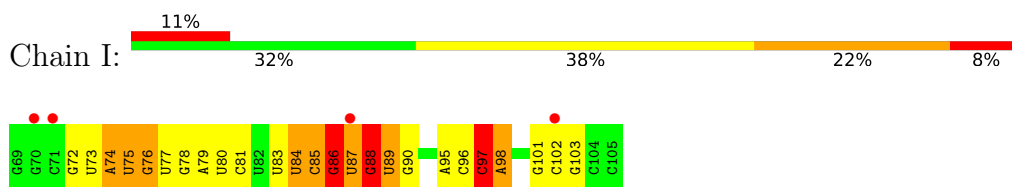
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	E	11	Total 11	O 11	0	0
5	F	13	Total 13	O 13	0	0
5	G	13	Total 13	O 13	0	0
5	H	16	Total 16	O 16	0	0

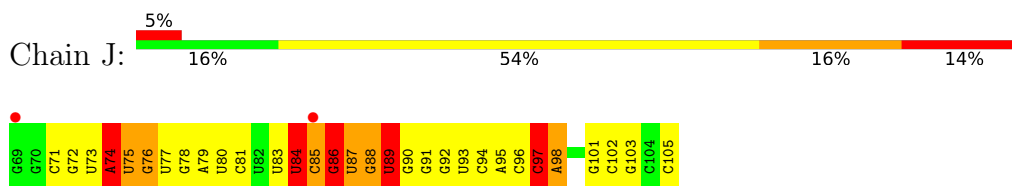
3 Residue-property plots [i](#)

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

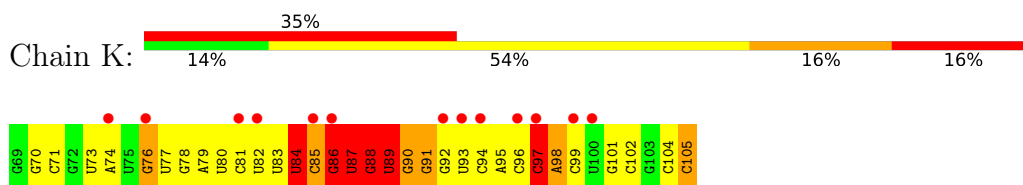
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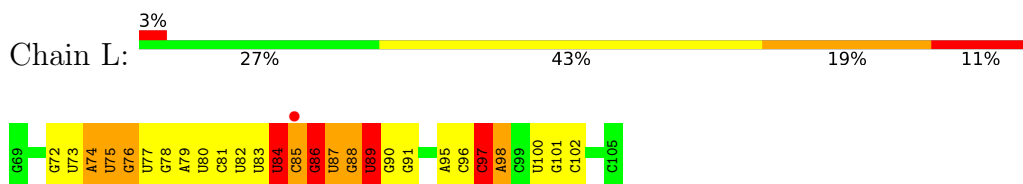
- Molecule 1: Threonyl-tRNA synthetase mRNA



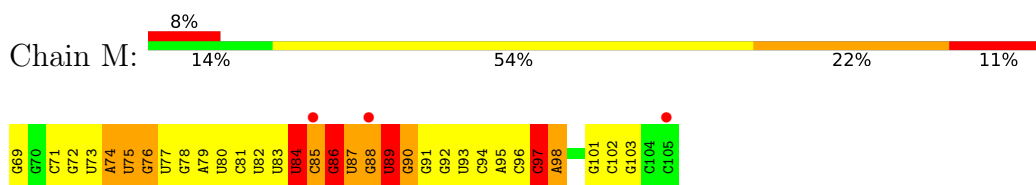
- Molecule 1: Threonyl-tRNA synthetase mRNA



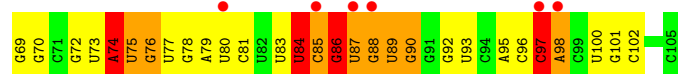
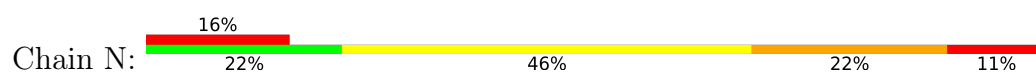
- Molecule 1: Threonyl-tRNA synthetase mRNA



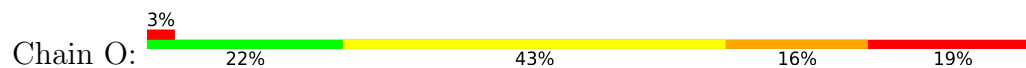
- Molecule 1: Threonyl-tRNA synthetase mRNA



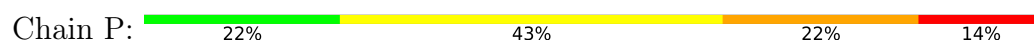
- Molecule 1: Threonyl-tRNA synthetase mRNA



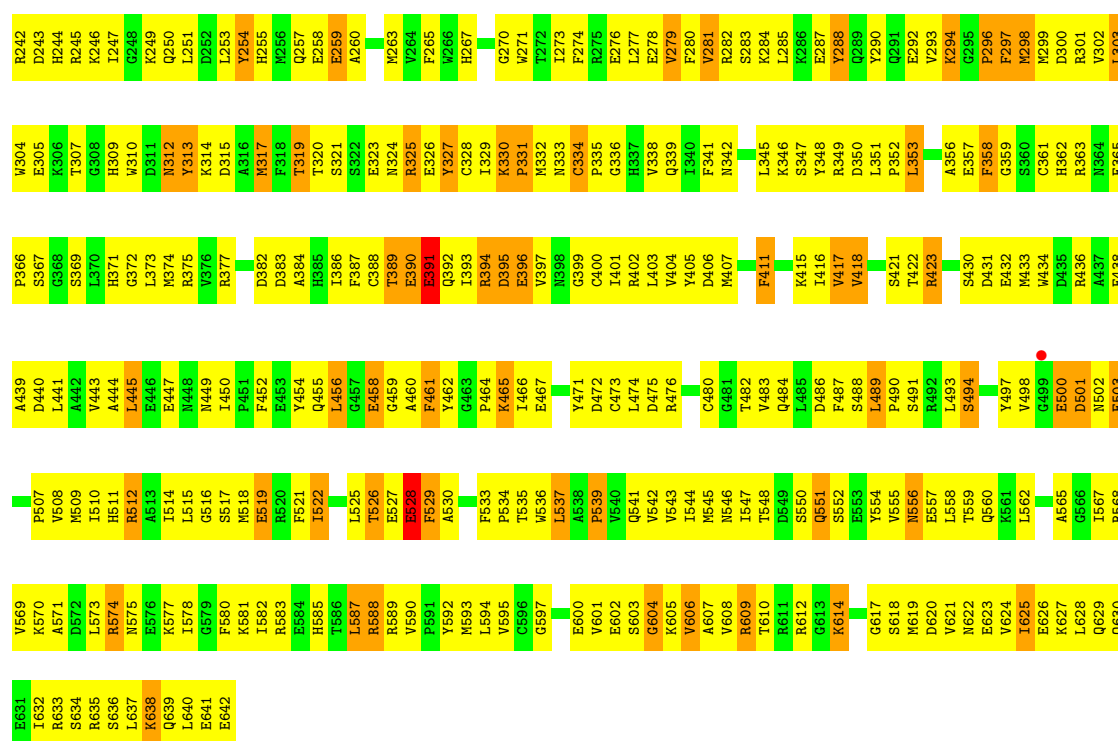
• Molecule 1: Threonyl-tRNA synthetase mRNA



• Molecule 1: Threonyl-tRNA synthetase mRNA

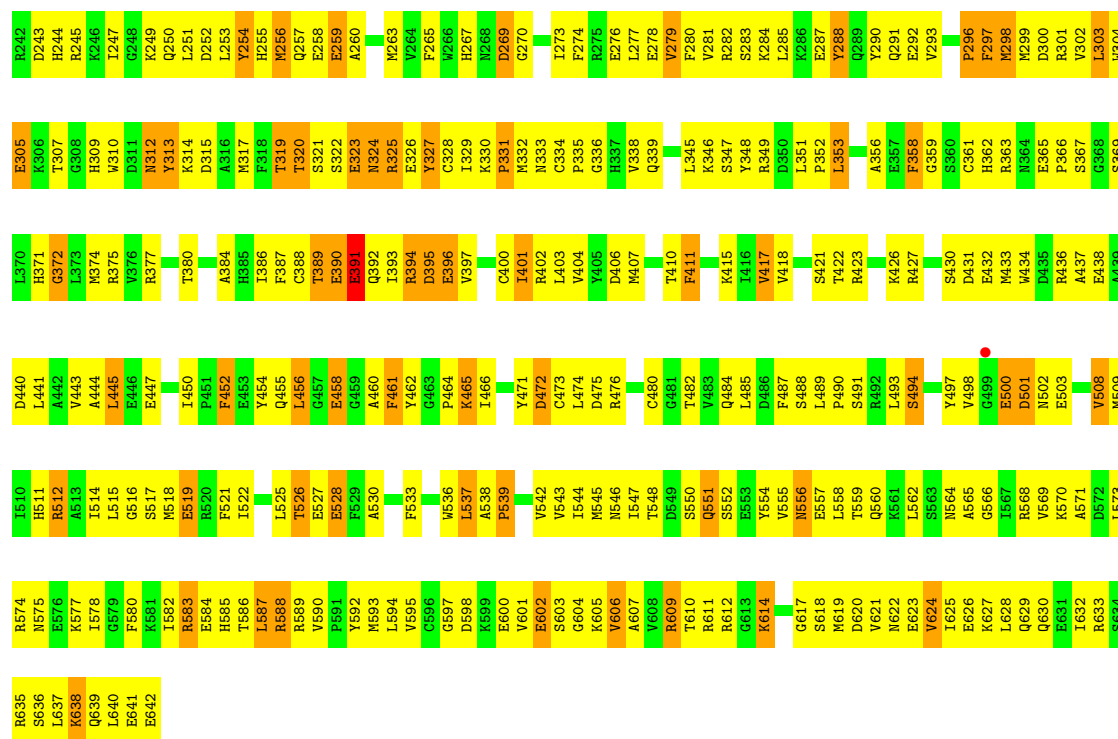


• Molecule 2: Threonyl-tRNA synthetase

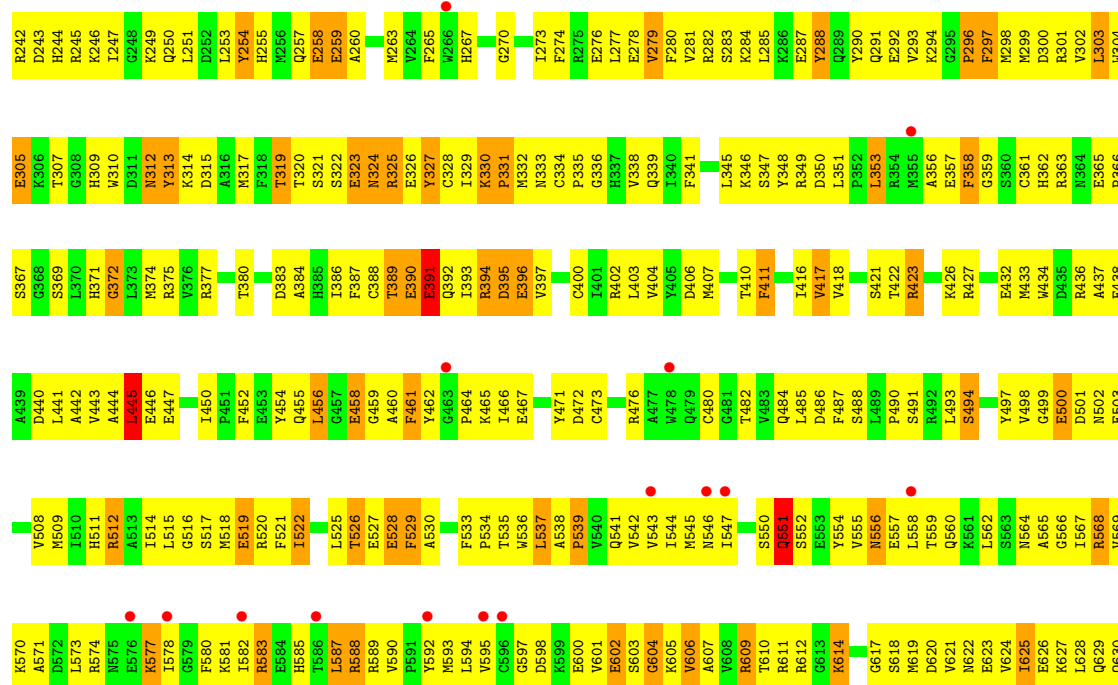


• Molecule 2: Threonyl-tRNA synthetase

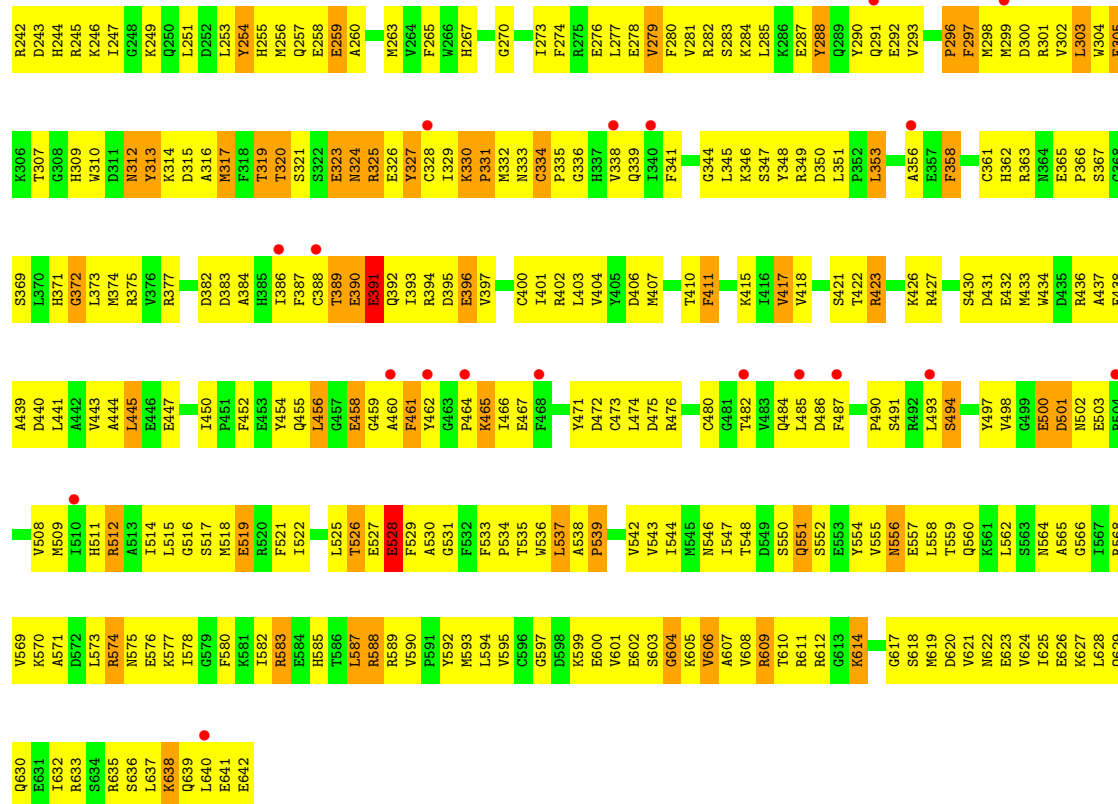




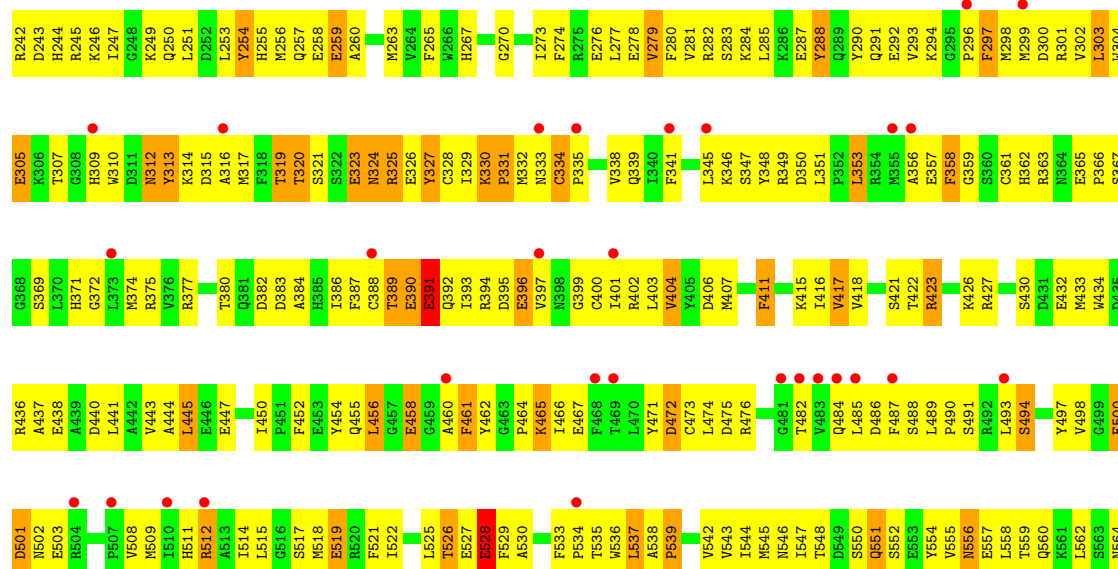
• Molecule 2: Threonyl-tRNA synthetase

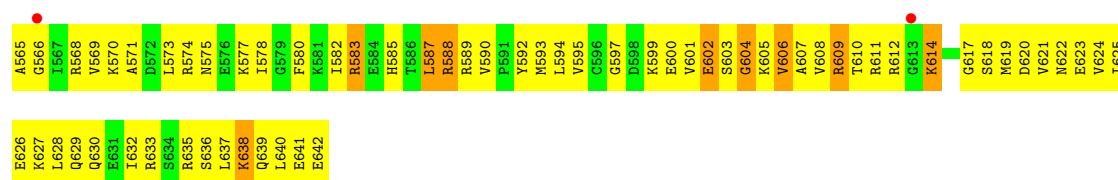


• Molecule 2: Threonyl-tRNA synthetase

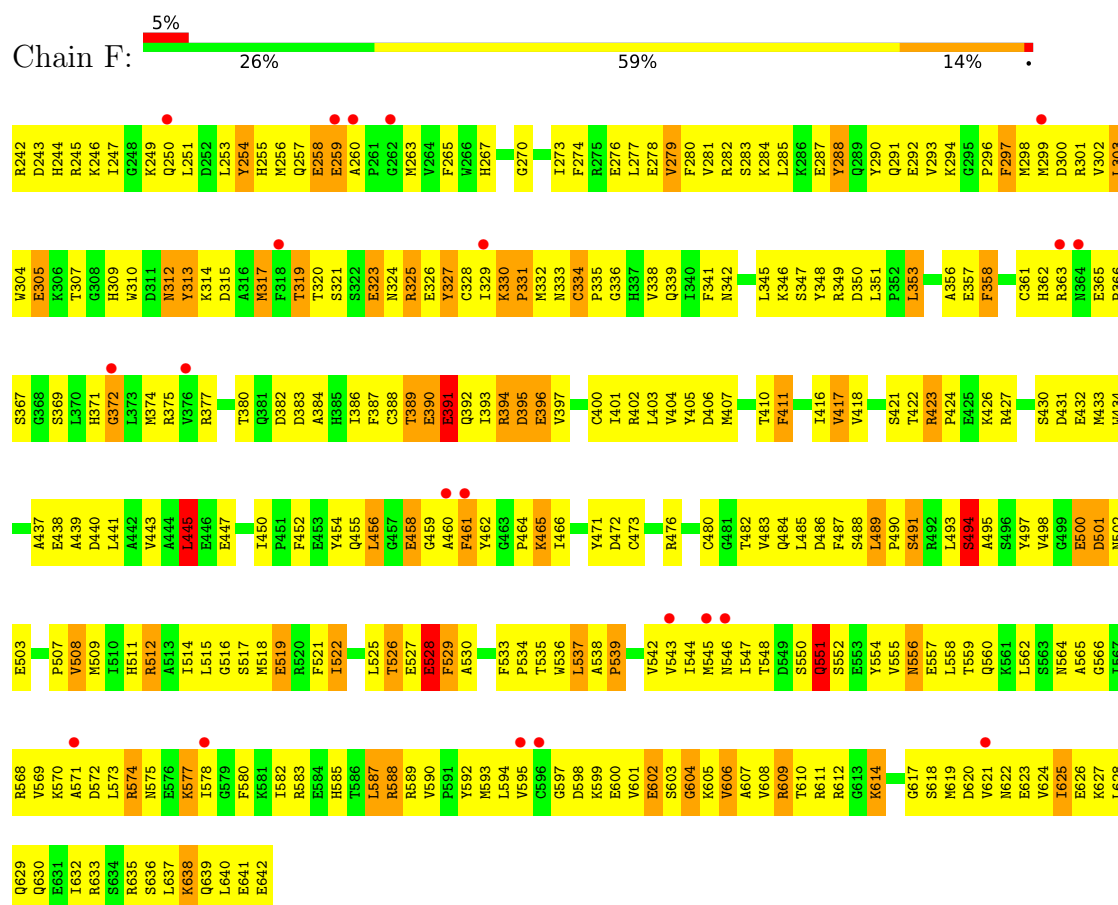


• Molecule 2: Threonyl-tRNA synthetase

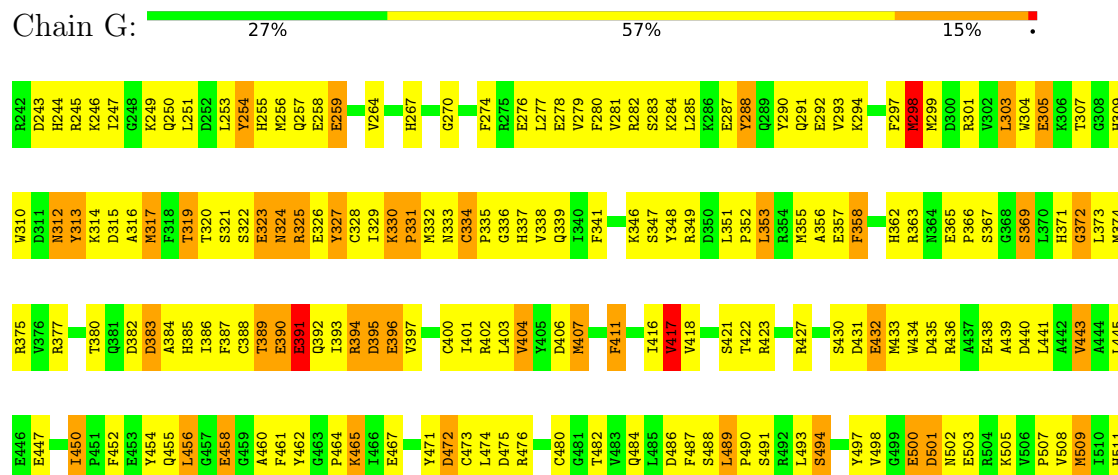


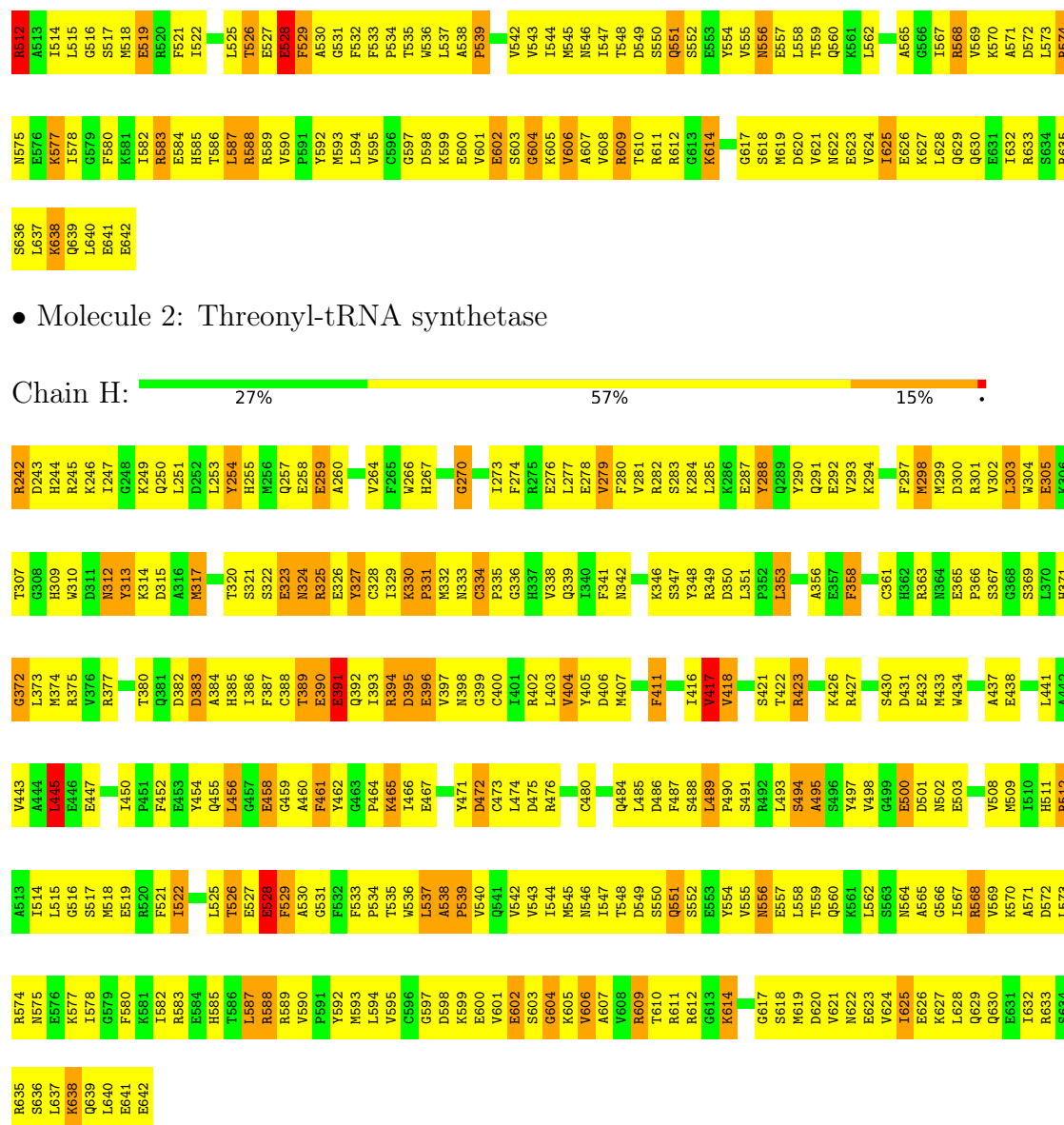


• Molecule 2: Threonyl-tRNA synthetase



• Molecule 2: Threonyl-tRNA synthetase





• Molecule 2: Threonyl-tRNA synthetase

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	188.45Å 101.74Å 199.34Å 90.00° 114.40° 90.00°	Depositor
Resolution (Å)	29.80 – 3.50 29.80 – 3.50	Depositor EDS
% Data completeness (in resolution range)	90.0 (29.80-3.50) 94.2 (29.80-3.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.63 (at 3.47Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.251 , 0.287 0.256 , 0.288	Depositor DCC
R_{free} test set	8671 reflections (9.81%)	wwPDB-VP
Wilson B-factor (Å ²)	96.2	Xtriage
Anisotropy	0.477	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 114.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	32922	wwPDB-VP
Average B, all atoms (Å ²)	94.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.58% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	I	0.49	0/874	0.99	7/1358 (0.5%)
1	J	0.56	0/874	1.00	7/1358 (0.5%)
1	K	0.66	0/874	0.98	7/1358 (0.5%)
1	L	0.46	0/874	0.95	6/1358 (0.4%)
1	M	0.53	0/874	1.00	7/1358 (0.5%)
1	N	0.47	0/874	0.88	6/1358 (0.4%)
1	O	0.67	0/874	1.08	9/1358 (0.7%)
1	P	0.51	0/874	1.01	8/1358 (0.6%)
2	A	0.78	2/3349 (0.1%)	1.13	23/4508 (0.5%)
2	B	0.85	3/3349 (0.1%)	1.12	24/4508 (0.5%)
2	C	0.59	0/3349	1.06	21/4508 (0.5%)
2	D	0.55	0/3349	1.06	21/4508 (0.5%)
2	E	0.55	0/3349	1.06	22/4508 (0.5%)
2	F	0.61	0/3349	1.09	23/4508 (0.5%)
2	G	0.97	3/3349 (0.1%)	1.19	24/4508 (0.5%)
2	H	0.78	1/3349 (0.0%)	1.14	29/4508 (0.6%)
All	All	0.69	9/33784 (0.0%)	1.08	244/46928 (0.5%)

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	I	0	2
1	J	0	4
1	K	0	1
1	L	0	3
1	M	0	2
1	N	0	2
1	O	0	3

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Mol	Chain	#Chirality outliers	#Planarity outliers
1	P	0	3
All	All	0	20

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	407	MET	SD-CE	10.25	2.05	1.79
2	G	509	MET	SD-CE	6.77	1.96	1.79
2	B	256	MET	SD-CE	6.49	1.95	1.79
2	G	355	MET	SD-CE	6.45	1.95	1.79
2	B	269	ASP	CG-OD1	-6.06	1.13	1.25
2	B	269	ASP	CG-OD2	-5.84	1.14	1.25
2	H	538	ALA	CA-C	-5.47	1.47	1.52
2	A	281	VAL	CA-CB	-5.33	1.48	1.54
2	A	298	MET	SD-CE	5.14	1.92	1.79

All (244) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	O	74	A	N9-C1'-C2'	10.28	129.42	114.00
2	B	391	GLU	N-CA-C	-9.50	101.11	112.89
1	L	74	A	N9-C1'-C2'	9.36	126.04	112.00
1	N	74	A	N9-C1'-C2'	9.18	125.77	112.00
1	I	74	A	N9-C1'-C2'	9.18	125.76	112.00
2	E	391	GLU	N-CA-C	-9.03	101.69	112.89
2	F	391	GLU	N-CA-C	-8.96	101.61	112.54
1	M	74	A	N9-C1'-C2'	8.93	127.39	114.00
2	G	391	GLU	N-CA-C	-8.86	101.73	112.54
2	C	391	GLU	N-CA-C	-8.80	101.98	112.89
2	D	391	GLU	N-CA-C	-8.80	101.98	112.89
2	H	391	GLU	N-CA-C	-8.61	102.03	112.54
1	P	74	A	N9-C1'-C2'	8.56	126.84	114.00
2	A	391	GLU	N-CA-C	-8.42	102.27	112.54
1	K	86	G	N9-C1'-C2'	8.31	126.46	114.00
2	E	389	THR	N-CA-C	-8.29	100.22	110.41
2	D	389	THR	N-CA-C	-8.15	100.38	110.41
2	G	389	THR	N-CA-C	-8.02	100.19	110.53
2	B	270	GLY	N-CA-C	-7.99	102.83	112.49
1	P	86	G	N9-C1'-C2'	7.76	123.64	112.00
1	J	74	A	N9-C1'-C2'	7.71	125.57	114.00
1	P	105	C	C2'-C3'-O3'	7.66	125.18	113.70
2	H	389	THR	N-CA-C	-7.63	101.02	110.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	382	ASP	N-CA-C	-7.62	102.19	112.26
1	K	97	C	C2'-C3'-O3'	-7.53	102.41	113.70
1	K	81	C	C3'-C2'-O2'	7.51	121.97	110.70
1	O	87	U	N1-C1'-C2'	7.51	123.27	112.00
2	A	389	THR	N-CA-C	-7.39	101.32	110.41
2	H	604	GLY	N-CA-C	-7.37	105.82	115.47
2	B	389	THR	N-CA-C	-7.33	101.08	110.53
2	F	389	THR	N-CA-C	-7.31	101.42	110.41
1	I	86	G	N9-C1'-C2'	7.28	122.91	112.00
2	C	604	GLY	N-CA-C	-7.21	106.03	115.47
2	A	359	GLY	N-CA-C	7.19	122.18	110.97
1	I	87	U	N1-C1'-C2'	7.10	122.66	112.00
2	C	500	GLU	N-CA-C	-7.09	103.60	111.82
2	H	330	LYS	N-CA-C	7.08	118.28	109.57
1	J	86	G	N9-C1'-C2'	7.06	122.59	112.00
2	F	461	PHE	N-CA-C	-7.06	104.40	112.87
2	F	604	GLY	N-CA-C	-7.06	106.22	115.47
2	A	604	GLY	N-CA-C	-7.03	106.26	115.47
2	G	604	GLY	N-CA-C	-7.02	106.27	115.47
2	D	588	ARG	N-CA-C	-6.99	104.57	113.23
2	G	588	ARG	N-CA-C	-6.98	104.57	113.23
2	H	382	ASP	N-CA-C	-6.97	103.06	112.26
2	B	604	GLY	N-CA-C	-6.91	106.42	115.47
1	N	86	G	N9-C1'-C2'	6.89	122.33	112.00
2	C	389	THR	N-CA-C	-6.88	101.95	110.41
2	E	604	GLY	N-CA-C	-6.78	106.59	115.47
1	J	87	U	N1-C1'-C2'	6.77	122.15	112.00
2	A	383	ASP	N-CA-C	6.67	119.80	109.60
2	E	588	ARG	N-CA-C	-6.64	104.99	113.23
2	D	500	GLU	N-CA-C	-6.62	104.14	111.82
1	J	97	C	C2'-C3'-O3'	-6.62	99.58	109.50
2	G	330	LYS	N-CA-C	6.61	117.70	109.57
2	E	500	GLU	N-CA-C	-6.60	104.16	111.82
2	C	588	ARG	N-CA-C	-6.60	105.05	113.23
2	E	330	LYS	N-CA-C	6.60	117.68	109.57
1	M	87	U	N1-C1'-C2'	6.59	121.89	112.00
1	O	86	G	N9-C1'-C2'	6.59	121.89	112.00
2	F	330	LYS	N-CA-C	6.58	117.67	109.57
1	M	86	G	N9-C1'-C2'	6.58	121.87	112.00
1	O	97	C	N1-C1'-C2'	6.58	123.87	114.00
2	A	330	LYS	CA-C-N	6.58	128.06	119.84
2	A	330	LYS	C-N-CA	6.58	128.06	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	588	ARG	N-CA-C	-6.57	105.08	113.23
2	A	500	GLU	N-CA-C	-6.53	104.24	111.82
2	F	588	ARG	N-CA-C	-6.53	105.13	113.23
1	O	74	A	O4'-C1'-C2'	6.53	112.33	105.80
1	L	86	G	N9-C1'-C2'	6.49	121.73	112.00
2	D	604	GLY	N-CA-C	-6.49	106.97	115.47
2	H	330	LYS	CA-C-N	6.49	127.95	119.84
2	H	330	LYS	C-N-CA	6.49	127.95	119.84
2	G	404	VAL	CB-CA-C	-6.40	103.50	112.14
2	D	358	PHE	N-CA-C	-6.39	100.04	109.62
2	G	472	ASP	N-CA-C	6.37	120.63	109.80
1	N	87	U	N1-C1'-C2'	6.34	121.51	112.00
2	G	383	ASP	N-CA-C	6.30	119.62	109.79
2	F	500	GLU	N-CA-C	-6.29	104.52	111.82
1	K	84	U	C2'-C3'-O3'	6.24	118.86	109.50
1	P	89	U	C2'-C3'-O3'	-6.24	104.35	113.70
2	G	500	GLU	N-CA-C	-6.22	104.60	111.82
2	B	260	ALA	CA-C-N	6.22	127.88	120.23
2	B	260	ALA	C-N-CA	6.22	127.88	120.23
2	D	461	PHE	N-CA-C	-6.22	105.41	112.87
2	B	330	LYS	N-CA-C	6.17	118.36	109.48
1	L	87	U	N1-C1'-C2'	6.15	121.22	112.00
2	B	556	ASN	N-CA-C	-6.14	104.64	111.71
1	P	84	U	C2'-C3'-O3'	6.14	118.71	109.50
1	P	87	U	N1-C1'-C2'	6.13	121.19	112.00
1	I	97	C	C2'-C3'-O3'	-6.13	100.31	109.50
2	B	500	GLU	N-CA-C	-6.11	104.73	111.82
2	G	556	ASN	N-CA-C	-6.11	104.62	111.28
2	F	383	ASP	N-CA-C	6.10	118.93	109.60
2	D	556	ASN	N-CA-C	-6.08	104.65	111.28
2	B	465	LYS	N-CA-C	6.08	118.19	107.61
2	A	588	ARG	N-CA-C	-6.06	105.72	113.23
1	K	87	U	C2'-C3'-O3'	-6.05	104.63	113.70
2	F	382	ASP	N-CA-C	-6.04	104.28	112.26
2	A	461	PHE	N-CA-C	-6.04	105.62	112.87
2	B	358	PHE	N-CA-C	-6.04	100.57	109.62
2	A	330	LYS	N-CA-C	6.00	116.96	109.57
2	D	330	LYS	N-CA-C	6.00	118.11	109.48
2	E	461	PHE	N-CA-C	-5.98	105.69	112.87
1	O	84	U	C2'-C3'-O3'	5.98	118.47	109.50
2	E	556	ASN	N-CA-C	-5.97	104.77	111.28
2	C	330	LYS	CA-C-N	5.96	127.28	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	330	LYS	C-N-CA	5.96	127.28	119.84
1	L	84	U	C2'-C3'-O3'	5.95	118.42	109.50
2	G	264	VAL	N-CA-C	5.93	117.28	109.21
2	E	472	ASP	N-CA-C	5.92	118.45	109.63
1	M	97	C	N1-C1'-C2'	5.91	122.87	114.00
2	B	330	LYS	CA-C-N	5.91	127.23	119.84
2	B	330	LYS	C-N-CA	5.91	127.23	119.84
2	F	350	ASP	N-CA-C	-5.91	105.84	113.16
2	H	305	GLU	N-CA-C	-5.89	104.86	111.28
2	H	500	GLU	N-CA-C	-5.89	104.98	111.82
2	H	588	ARG	N-CA-C	-5.87	105.95	113.23
2	A	565	ALA	N-CA-C	-5.86	105.40	112.54
1	M	97	C	C2'-C3'-O3'	-5.85	100.73	109.50
2	E	382	ASP	N-CA-C	-5.85	104.54	112.26
2	C	260	ALA	CA-C-N	5.84	126.76	119.98
2	C	260	ALA	C-N-CA	5.84	126.76	119.98
1	N	84	U	C2'-C3'-O3'	5.83	118.25	109.50
2	C	461	PHE	N-CA-C	-5.81	105.90	112.87
2	H	540	VAL	N-CA-C	-5.81	100.11	108.48
1	M	74	A	O4'-C1'-C2'	5.81	111.61	105.80
2	E	383	ASP	N-CA-C	5.80	118.48	109.60
2	A	465	LYS	N-CA-C	5.80	117.70	107.61
1	J	84	U	C2'-C3'-O3'	5.79	118.18	109.50
2	H	472	ASP	N-CA-C	5.78	119.63	109.80
2	H	404	VAL	CB-CA-C	-5.76	104.60	111.97
2	D	565	ALA	N-CA-C	-5.74	105.98	112.87
2	F	565	ALA	N-CA-C	-5.74	105.38	112.38
2	F	319	THR	N-CA-C	5.72	118.79	109.06
2	C	330	LYS	N-CA-C	5.70	117.69	109.48
2	E	565	ALA	N-CA-C	-5.70	106.03	112.87
2	G	465	LYS	N-CA-C	5.70	117.53	107.61
2	B	472	ASP	N-CA-C	5.69	118.11	109.63
2	A	260	ALA	CA-C-N	5.68	126.57	119.98
2	A	260	ALA	C-N-CA	5.68	126.57	119.98
2	H	383	ASP	N-CA-C	5.67	118.64	109.79
2	E	358	PHE	N-CA-C	-5.66	101.13	109.62
2	C	556	ASN	N-CA-C	-5.66	105.11	111.28
2	C	319	THR	N-CA-C	5.62	118.36	108.75
2	D	383	ASP	N-CA-C	5.60	118.17	109.60
2	B	565	ALA	N-CA-C	-5.59	106.16	112.87
2	H	538	ALA	N-CA-C	5.59	117.70	109.42
2	H	465	LYS	N-CA-C	5.58	117.32	107.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	319	THR	N-CA-C	5.57	118.27	108.75
2	F	330	LYS	CA-C-N	5.57	126.80	119.84
2	F	330	LYS	C-N-CA	5.57	126.80	119.84
2	B	359	GLY	N-CA-C	5.56	119.64	110.97
2	D	305	GLU	N-CA-C	-5.55	105.13	111.07
1	P	97	C	N1-C1'-C2'	5.54	122.31	114.00
2	D	260	ALA	CA-C-N	5.54	126.41	119.98
2	D	260	ALA	C-N-CA	5.54	126.41	119.98
2	C	358	PHE	N-CA-C	-5.53	101.32	109.62
2	E	465	LYS	N-CA-C	5.52	117.21	107.61
1	N	74	A	C4'-C3'-O3'	-5.51	104.73	113.00
2	B	452	PHE	N-CA-C	5.51	115.83	108.38
2	D	382	ASP	N-CA-C	-5.51	104.98	112.26
2	E	305	GLU	N-CA-C	-5.51	105.28	111.28
2	G	512	ARG	NE-CZ-NH1	-5.50	116.00	121.50
2	F	305	GLU	N-CA-C	-5.49	105.29	111.28
2	A	319	THR	N-CA-C	5.49	118.13	108.75
2	D	330	LYS	CA-C-N	5.47	126.68	119.84
2	D	330	LYS	C-N-CA	5.47	126.68	119.84
2	H	565	ALA	N-CA-C	-5.47	106.30	112.87
2	D	350	ASP	N-CA-C	-5.46	106.61	113.28
2	C	305	GLU	N-CA-C	-5.46	105.33	111.28
2	C	383	ASP	N-CA-C	5.46	116.96	110.19
2	A	556	ASN	N-CA-C	-5.45	105.34	111.28
2	F	260	ALA	CA-C-N	5.45	126.93	120.23
2	F	260	ALA	C-N-CA	5.45	126.93	120.23
2	F	556	ASN	N-CA-C	-5.45	105.34	111.28
1	I	88	G	N9-C1'-C2'	5.44	122.17	114.00
2	B	404	VAL	CB-CA-C	-5.44	105.01	111.97
2	F	465	LYS	N-CA-C	5.43	117.05	107.61
2	G	417	VAL	N-CA-C	5.43	116.12	108.36
1	K	89	U	N1-C1'-C2'	5.42	120.13	112.00
2	C	359	GLY	N-CA-C	5.42	119.42	110.97
2	E	325	ARG	CB-CA-C	-5.42	107.63	114.40
2	F	404	VAL	CB-CA-C	-5.41	105.04	111.97
2	H	350	ASP	N-CA-C	-5.41	106.45	113.16
2	B	319	THR	N-CA-C	5.41	117.99	108.75
2	E	359	GLY	N-CA-C	5.40	119.40	110.97
1	L	74	A	C2'-C3'-O3'	-5.40	105.60	113.70
2	A	404	VAL	CB-CA-C	-5.40	105.06	111.97
1	J	97	C	N1-C1'-C2'	5.39	122.09	114.00
2	A	382	ASP	N-CA-C	-5.39	105.14	112.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	264	VAL	N-CA-C	5.39	116.54	109.21
2	G	325	ARG	CB-CA-C	-5.39	107.67	114.40
2	F	325	ARG	CB-CA-C	-5.38	107.68	114.40
2	B	305	GLU	N-CA-C	-5.36	105.43	111.28
2	H	358	PHE	N-CA-C	-5.36	101.58	109.62
1	J	74	A	O4'-C1'-C2'	5.34	111.14	105.80
2	B	461	PHE	N-CA-C	-5.34	106.46	112.87
2	B	325	ARG	CB-CA-C	-5.34	107.73	114.40
2	A	358	PHE	N-CA-C	-5.33	101.62	109.62
2	H	260	ALA	CA-C-N	5.33	126.79	120.23
2	H	260	ALA	C-N-CA	5.33	126.79	120.23
2	H	556	ASN	N-CA-C	-5.33	105.47	111.28
2	G	319	THR	N-CA-C	5.32	117.34	108.99
2	A	350	ASP	N-CA-C	-5.31	106.58	113.16
2	D	325	ARG	CB-CA-C	-5.31	107.77	114.40
1	O	89	U	C2'-C3'-O3'	-5.30	105.75	113.70
2	F	508	VAL	N-CA-C	-5.28	100.78	109.12
2	C	325	ARG	CB-CA-C	-5.28	107.80	114.40
2	E	319	THR	N-CA-C	5.28	117.78	108.75
1	K	88	G	N9-C1'-C2'	5.27	121.90	114.00
2	H	495	ALA	N-CA-C	-5.25	101.56	109.85
1	M	84	U	C2'-C3'-O3'	5.23	117.35	109.50
1	N	97	C	N1-C1'-C2'	5.23	121.85	114.00
2	G	330	LYS	CA-C-N	5.23	126.38	119.84
2	G	330	LYS	C-N-CA	5.23	126.38	119.84
2	D	465	LYS	N-CA-C	5.22	116.70	107.61
2	H	417	VAL	N-CA-C	5.22	115.47	108.17
2	B	508	VAL	N-CA-C	-5.20	100.91	109.12
2	G	358	PHE	N-CA-C	-5.19	101.84	109.62
2	C	350	ASP	N-CA-C	-5.18	106.96	113.28
2	G	443	VAL	CB-CA-C	-5.18	102.79	111.29
2	G	305	GLU	N-CA-C	-5.17	105.54	111.07
2	C	404	VAL	CB-CA-C	-5.16	105.37	111.97
1	I	74	A	C2'-C3'-O3'	-5.13	106.00	113.70
2	H	461	PHE	N-CA-C	-5.10	106.75	112.87
2	E	404	VAL	CB-CA-C	-5.10	105.45	111.97
2	F	358	PHE	N-CA-C	-5.10	101.97	109.62
2	C	565	ALA	N-CA-C	-5.09	106.17	112.38
2	H	418	VAL	N-CA-C	5.08	116.59	108.87
2	H	325	ARG	CB-CA-C	-5.07	108.06	114.40
2	A	418	VAL	N-CA-C	5.07	116.57	108.87
1	L	97	C	N1-C1'-C2'	5.06	121.59	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	369	SER	N-CA-C	5.06	119.21	113.19
2	H	270	GLY	N-CA-C	-5.06	106.66	112.73
1	I	97	C	N1-C1'-C2'	5.06	121.59	114.00
1	O	87	U	C4'-C3'-O3'	-5.06	105.42	113.00
2	A	503	GLU	N-CA-C	5.04	117.79	109.72
2	G	450	ILE	CA-C-O	5.04	122.47	119.19
1	O	75	U	C4'-C3'-O3'	-5.03	105.46	113.00
2	E	260	ALA	CA-C-N	5.03	126.41	120.23
2	E	260	ALA	C-N-CA	5.03	126.41	120.23
2	E	350	ASP	N-CA-C	-5.02	106.93	113.16
1	P	74	A	O4'-C1'-C2'	5.00	110.80	105.80

There are no chirality outliers.

All (20) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	I	74	A	Sidechain
1	I	86	G	Sidechain
1	J	74	A	Sidechain
1	J	86	G	Sidechain
1	J	89	U	Sidechain
1	J	97	C	Sidechain
1	K	86	G	Sidechain
1	L	74	A	Sidechain
1	L	86	G	Sidechain
1	L	89	U	Sidechain
1	M	86	G	Sidechain
1	M	89	U	Sidechain
1	N	74	A	Sidechain
1	N	86	G	Sidechain
1	O	74	A	Sidechain
1	O	86	G	Sidechain
1	O	88	G	Sidechain
1	P	74	A	Sidechain
1	P	86	G	Sidechain
1	P	89	U	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	I	785	0	397	45	0
1	J	785	0	397	53	0
1	K	785	0	397	78	0
1	L	785	0	397	42	0
1	M	785	0	397	73	0
1	N	785	0	397	43	0
1	O	785	0	397	55	0
1	P	785	0	397	65	0
2	A	3278	0	3208	364	0
2	B	3278	0	3208	341	0
2	C	3278	0	3208	356	0
2	D	3278	0	3208	359	0
2	E	3278	0	3208	355	0
2	F	3278	0	3208	367	0
2	G	3278	0	3208	343	0
2	H	3278	0	3208	335	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
4	A	30	0	21	3	0
4	B	30	0	20	4	0
4	C	30	0	21	3	0
4	D	30	0	21	4	0
4	E	30	0	21	4	0
4	F	30	0	21	5	0
4	G	30	0	21	4	0
4	H	30	0	21	4	0
5	A	10	0	0	0	0
5	B	15	0	0	8	0
5	C	11	0	0	3	0
5	D	14	0	0	4	0
5	E	11	0	0	5	0
5	F	13	0	0	2	0
5	G	13	0	0	4	0
5	H	16	0	0	3	0
5	I	11	0	0	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	J	12	0	0	8	0
5	K	1	0	0	0	0
5	L	8	0	0	3	0
5	M	6	0	0	10	0
5	N	9	0	0	2	0
5	O	11	0	0	9	0
5	P	9	0	0	14	0
All	All	32922	0	29007	3085	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 50.

All (3085) 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)
2:G:407:MET:CE	2:G:407:MET:SD	2.05	1.44
2:D:559:THR:HG21	2:D:571:ALA:HB2	1.29	1.15
2:G:559:THR:HG21	2:G:571:ALA:HB2	1.27	1.12
2:E:559:THR:HG21	2:E:571:ALA:HB2	1.30	1.12
2:A:559:THR:HG21	2:A:571:ALA:HB2	1.27	1.11
2:C:559:THR:HG21	2:C:571:ALA:HB2	1.30	1.10
2:F:559:THR:HG21	2:F:571:ALA:HB2	1.26	1.09
2:H:559:THR:HG21	2:H:571:ALA:HB2	1.31	1.09
2:B:559:THR:HG21	2:B:571:ALA:HB2	1.32	1.07
2:E:411:PHE:HD2	2:E:525:LEU:HD21	1.21	1.06
2:D:411:PHE:HE2	2:D:521:PHE:CE1	1.74	1.05
2:B:473:CYS:HA	5:B:221:HOH:O	1.57	1.05
1:O:87:U:H2'	5:O:1519:HOH:O	1.58	1.03
2:B:411:PHE:HD2	2:B:525:LEU:HD21	1.23	1.02
2:A:411:PHE:HD2	2:A:525:LEU:HD21	1.25	1.02
2:G:421:SER:HB3	2:G:458:GLU:HB3	1.40	1.01
2:H:585:HIS:HB2	2:H:593:MET:HE3	1.46	0.98
1:K:86:G:H5'	1:K:87:U:OP1	1.63	0.97
2:E:421:SER:HB3	2:E:458:GLU:HB3	1.47	0.97
2:H:411:PHE:HD2	2:H:525:LEU:HD21	1.28	0.96
2:B:421:SER:HB3	2:B:458:GLU:HB3	1.45	0.95
2:A:298:MET:HG2	2:B:263:MET:HE2	1.50	0.94
2:G:585:HIS:HB2	2:G:593:MET:HE3	1.48	0.94
2:E:585:HIS:HB2	2:E:593:MET:HE3	1.48	0.94
2:A:421:SER:HB3	2:A:458:GLU:HB3	1.45	0.93
1:I:103:G:H5''	5:I:113:HOH:O	1.68	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:421:SER:HB3	2:D:458:GLU:HB3	1.49	0.92
2:F:585:HIS:HB2	2:F:593:MET:HE3	1.48	0.92
2:D:411:PHE:CE2	2:D:521:PHE:CE1	2.57	0.92
2:E:411:PHE:CD2	2:E:525:LEU:HD21	2.04	0.92
2:C:421:SER:HB3	2:C:458:GLU:HB3	1.50	0.91
2:D:543:VAL:HG23	2:D:590:VAL:HG11	1.52	0.91
2:H:421:SER:HB3	2:H:458:GLU:HB3	1.50	0.91
1:K:84:U:H4'	1:K:85:C:H5'	1.53	0.91
2:C:585:HIS:HB2	2:C:593:MET:HE3	1.53	0.91
1:O:77:U:O3'	5:O:1539:HOH:O	1.89	0.91
2:C:389:THR:OG1	2:C:392:GLN:HG3	1.71	0.91
2:C:411:PHE:HE2	2:C:521:PHE:CE1	1.87	0.90
2:A:309:HIS:HB3	2:A:317:MET:HE1	1.54	0.90
2:C:411:PHE:HE2	2:C:521:PHE:HE1	1.18	0.90
2:F:309:HIS:HB3	2:F:317:MET:HE1	1.53	0.90
2:F:411:PHE:HE2	2:F:521:PHE:CE1	1.88	0.90
1:K:88:G:H4'	1:K:89:U:OP1	1.71	0.90
2:B:411:PHE:CD2	2:B:525:LEU:HD21	2.06	0.89
2:D:585:HIS:HB2	2:D:593:MET:HE3	1.53	0.89
2:G:309:HIS:HB3	2:G:317:MET:HE1	1.54	0.89
1:O:85:C:H5	2:G:547:ILE:O	1.55	0.89
2:G:543:VAL:HG23	2:G:590:VAL:HG11	1.55	0.89
2:A:389:THR:OG1	2:A:392:GLN:HG3	1.71	0.89
2:B:389:THR:OG1	2:B:392:GLN:HG3	1.72	0.89
2:F:543:VAL:HG23	2:F:590:VAL:HG11	1.54	0.88
2:A:411:PHE:CD2	2:A:525:LEU:HD21	2.08	0.88
2:C:411:PHE:HD2	2:C:525:LEU:HD21	1.39	0.88
1:K:101:G:O2'	1:K:102:C:H5'	1.73	0.88
2:H:543:VAL:HG23	2:H:590:VAL:HG11	1.53	0.88
1:J:98:A:H4'	5:J:1027:HOH:O	1.74	0.88
1:K:87:U:H2'	1:K:87:U:O2	1.73	0.87
2:A:587:LEU:HD23	2:A:587:LEU:H	1.38	0.87
2:B:585:HIS:HB2	2:B:593:MET:HE3	1.56	0.87
2:G:452:PHE:HE2	2:G:454:TYR:HE1	1.22	0.87
2:H:309:HIS:HB3	2:H:317:MET:HE1	1.55	0.87
2:G:551:GLN:HE21	2:G:597:GLY:HA2	1.39	0.87
2:F:421:SER:HB3	2:F:458:GLU:HB3	1.55	0.87
2:A:307:THR:HG22	2:A:493:LEU:HD21	1.54	0.87
2:A:543:VAL:HG23	2:A:590:VAL:HG11	1.56	0.86
2:E:307:THR:HG22	2:E:493:LEU:HD21	1.57	0.86
2:H:389:THR:OG1	2:H:392:GLN:HG3	1.75	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:84:U:H4'	1:K:85:C:C5'	2.05	0.86
2:C:543:VAL:HG23	2:C:590:VAL:HG11	1.55	0.86
2:D:411:PHE:CE2	2:D:521:PHE:CZ	2.62	0.86
2:A:263:MET:HE2	2:B:298:MET:HG2	1.57	0.86
2:A:585:HIS:HB2	2:A:593:MET:HE3	1.57	0.86
2:E:389:THR:OG1	2:E:392:GLN:HG3	1.75	0.86
2:H:411:PHE:CD2	2:H:525:LEU:HD21	2.09	0.86
2:A:276:GLU:O	2:A:279:VAL:HG22	1.75	0.86
2:G:389:THR:OG1	2:G:392:GLN:HG3	1.75	0.86
2:G:587:LEU:HD23	2:G:587:LEU:H	1.40	0.86
2:C:307:THR:HG22	2:C:493:LEU:HD21	1.57	0.86
1:I:86:G:H2'	5:I:109:HOH:O	1.74	0.86
2:B:309:HIS:HB3	2:B:317:MET:HE1	1.56	0.86
2:F:411:PHE:CD2	2:F:525:LEU:HD21	2.10	0.86
2:C:255:HIS:ND1	2:C:267:HIS:HE1	1.75	0.85
2:E:309:HIS:HB3	2:E:317:MET:HE1	1.57	0.85
2:H:473:CYS:HA	5:H:821:HOH:O	1.76	0.85
2:G:276:GLU:O	2:G:279:VAL:HG22	1.77	0.84
2:H:551:GLN:HE21	2:H:597:GLY:HA2	1.41	0.84
2:E:543:VAL:HG23	2:E:590:VAL:HG11	1.59	0.84
2:C:411:PHE:CD2	2:C:525:LEU:HD21	2.12	0.84
1:O:89:U:O2	1:O:89:U:H2'	1.78	0.84
2:D:411:PHE:HE2	2:D:521:PHE:HE1	1.25	0.84
2:F:389:THR:OG1	2:F:392:GLN:HG3	1.77	0.84
2:F:411:PHE:HD2	2:F:525:LEU:HD21	1.41	0.84
2:G:411:PHE:HD2	2:G:525:LEU:HD21	1.43	0.84
1:M:89:U:O2	1:M:89:U:H2'	1.75	0.83
2:A:633:ARG:NH1	2:H:638:LYS:HZ1	1.75	0.83
2:H:407:MET:SD	2:H:514:ILE:HG21	2.18	0.83
2:C:309:HIS:HB3	2:C:317:MET:HE1	1.59	0.83
2:D:307:THR:HG22	2:D:493:LEU:HD21	1.58	0.83
1:I:88:G:H8	5:I:112:HOH:O	1.60	0.82
2:C:587:LEU:HD23	2:C:587:LEU:H	1.43	0.82
2:D:309:HIS:HB3	2:D:317:MET:HE1	1.61	0.82
2:H:587:LEU:HD23	2:H:587:LEU:H	1.43	0.82
2:B:587:LEU:HD23	2:B:587:LEU:H	1.42	0.82
2:G:551:GLN:NE2	2:G:597:GLY:HA2	1.93	0.82
2:C:411:PHE:CE2	2:C:521:PHE:CE1	2.66	0.82
1:N:79:A:H1'	2:E:502:ASN:OD1	1.80	0.82
2:D:389:THR:OG1	2:D:392:GLN:HG3	1.80	0.82
1:N:100:U:H4'	5:F:7014:HOH:O	1.77	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:255:HIS:ND1	2:F:267:HIS:HE1	1.78	0.81
2:F:411:PHE:HE2	2:F:521:PHE:HE1	1.26	0.81
2:C:276:GLU:O	2:C:279:VAL:HG22	1.80	0.81
2:E:411:PHE:CE2	2:E:521:PHE:CE1	2.69	0.81
2:C:551:GLN:HE21	2:C:597:GLY:HA2	1.43	0.81
2:A:551:GLN:HE21	2:A:597:GLY:HA2	1.44	0.81
2:F:551:GLN:HE21	2:F:597:GLY:HA2	1.44	0.81
1:O:86:G:O6	2:G:599:LYS:CB	2.29	0.81
2:H:307:THR:HG22	2:H:493:LEU:HD21	1.63	0.81
1:N:86:G:H2'	5:N:1413:HOH:O	1.81	0.81
2:H:255:HIS:ND1	2:H:267:HIS:HE1	1.79	0.81
2:B:501:ASP:OD2	2:B:503:GLU:HB2	1.81	0.80
2:B:543:VAL:HG23	2:B:590:VAL:HG11	1.63	0.80
2:E:587:LEU:HD23	2:E:587:LEU:H	1.46	0.80
2:H:348:TYR:CD1	2:H:349:ARG:N	2.49	0.80
1:O:97:C:H5''	1:O:98:A:OP1	1.81	0.80
2:C:320:THR:HG22	2:C:321:SER:H	1.45	0.80
1:P:83:U:H3	2:H:575:ASN:HD21	1.30	0.80
2:B:307:THR:HG22	2:B:493:LEU:HD21	1.61	0.80
2:H:551:GLN:NE2	2:H:597:GLY:HA2	1.97	0.80
1:N:88:G:H4'	1:N:89:U:OP1	1.80	0.80
2:G:621:VAL:O	2:G:624:VAL:HG22	1.80	0.80
2:B:255:HIS:ND1	2:B:267:HIS:HE1	1.80	0.80
2:E:319:THR:O	2:F:320:THR:HG23	1.80	0.80
2:F:411:PHE:CE2	2:F:521:PHE:CE1	2.69	0.80
2:G:254:TYR:CZ	2:G:373:LEU:HD21	2.16	0.80
2:B:411:PHE:N	2:B:411:PHE:HD1	1.80	0.80
2:F:402:ARG:HH11	2:F:402:ARG:HB2	1.47	0.80
2:G:411:PHE:CD2	2:G:525:LEU:HD21	2.17	0.80
2:C:452:PHE:HE2	2:C:454:TYR:HE1	1.30	0.80
1:P:97:C:H5''	1:P:98:A:OP1	1.81	0.79
2:A:321:SER:HB3	2:A:326:GLU:HA	1.64	0.79
2:F:307:THR:HG22	2:F:493:LEU:HD21	1.62	0.79
2:H:251:LEU:HA	2:H:588:ARG:HH12	1.47	0.79
2:G:251:LEU:HA	2:G:588:ARG:HH12	1.47	0.79
1:M:98:A:O3'	5:M:1327:HOH:O	1.99	0.79
2:A:501:ASP:OD2	2:A:503:GLU:HB2	1.81	0.79
2:A:633:ARG:CZ	2:H:638:LYS:HZ3	1.96	0.79
2:D:255:HIS:ND1	2:D:267:HIS:HE1	1.80	0.79
2:D:501:ASP:OD2	2:D:503:GLU:HB2	1.82	0.79
2:A:320:THR:HG23	2:B:319:THR:O	1.83	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:320:THR:HG22	2:B:321:SER:H	1.48	0.79
2:E:501:ASP:OD2	2:E:503:GLU:HB2	1.83	0.79
1:P:89:U:O2	1:P:89:U:H2'	1.81	0.79
2:D:551:GLN:HE21	2:D:597:GLY:HA2	1.48	0.79
2:F:320:THR:HG22	2:F:321:SER:H	1.48	0.79
2:F:551:GLN:NE2	2:F:597:GLY:HA2	1.98	0.79
2:G:578:ILE:O	2:G:582:ILE:HG12	1.83	0.79
2:A:255:HIS:ND1	2:A:267:HIS:HE1	1.80	0.78
2:A:319:THR:O	2:B:320:THR:HG23	1.83	0.78
2:E:251:LEU:HA	2:E:588:ARG:HH12	1.48	0.78
2:E:411:PHE:N	2:E:411:PHE:HD1	1.80	0.78
2:G:403:LEU:HG	2:G:407:MET:CE	2.14	0.78
2:H:276:GLU:O	2:H:279:VAL:HG22	1.83	0.78
2:D:276:GLU:O	2:D:279:VAL:HG22	1.83	0.78
2:E:411:PHE:HE2	2:E:521:PHE:HE1	1.29	0.78
2:E:255:HIS:ND1	2:E:267:HIS:HE1	1.80	0.78
2:A:633:ARG:CZ	2:H:638:LYS:NZ	2.47	0.78
2:G:348:TYR:CD1	2:G:349:ARG:N	2.52	0.78
1:M:89:U:C5	2:E:583:ARG:CZ	2.66	0.78
2:C:501:ASP:OD2	2:C:503:GLU:HB2	1.83	0.78
2:F:587:LEU:H	2:F:587:LEU:HD23	1.48	0.78
2:A:249:LYS:HB3	2:A:249:LYS:HZ3	1.47	0.78
2:G:251:LEU:HA	2:G:588:ARG:NH1	1.99	0.78
2:G:402:ARG:HB2	2:G:402:ARG:HH11	1.49	0.78
2:B:251:LEU:HA	2:B:588:ARG:HH12	1.48	0.78
2:C:298:MET:HG2	2:D:263:MET:HE2	1.66	0.78
2:D:621:VAL:O	2:D:624:VAL:HG22	1.83	0.78
2:D:411:PHE:CD2	2:D:525:LEU:HD21	2.18	0.77
2:G:320:THR:HG22	2:G:321:SER:H	1.49	0.77
2:A:407:MET:SD	2:A:514:ILE:HG21	2.24	0.77
2:E:298:MET:HG2	2:F:263:MET:HE2	1.66	0.77
2:F:585:HIS:HB2	2:F:593:MET:CE	2.15	0.77
2:C:407:MET:SD	2:C:514:ILE:HG21	2.24	0.77
2:E:320:THR:HG22	2:E:321:SER:H	1.49	0.77
1:J:89:U:O2	1:J:89:U:H2'	1.82	0.77
2:A:303:LEU:HD11	2:A:339:GLN:HE21	1.48	0.77
2:E:551:GLN:HE21	2:E:597:GLY:HA2	1.50	0.77
2:H:578:ILE:O	2:H:582:ILE:HG12	1.84	0.77
2:H:501:ASP:OD2	2:H:503:GLU:HB2	1.84	0.77
2:A:390:GLU:HA	2:A:393:ILE:HG13	1.66	0.77
1:O:85:C:C5	2:G:547:ILE:O	2.37	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:578:ILE:O	2:A:582:ILE:HG12	1.83	0.77
2:H:402:ARG:HB2	2:H:402:ARG:HH11	1.50	0.77
1:N:89:U:H2'	1:N:89:U:O2	1.84	0.77
2:G:307:THR:HG22	2:G:493:LEU:HD21	1.67	0.77
2:H:585:HIS:HB2	2:H:593:MET:CE	2.15	0.77
2:E:411:PHE:CE2	2:E:521:PHE:HE1	2.02	0.76
2:F:276:GLU:O	2:F:279:VAL:HG22	1.84	0.76
2:B:452:PHE:HE2	2:B:454:TYR:HE1	1.30	0.76
2:C:551:GLN:NE2	2:C:597:GLY:HA2	1.98	0.76
2:F:251:LEU:HA	2:F:588:ARG:NH1	2.00	0.76
2:D:587:LEU:HD23	2:D:587:LEU:H	1.50	0.76
2:G:365:GLU:H	2:G:377:ARG:HD3	1.50	0.76
1:O:86:G:O6	2:G:599:LYS:HB2	1.86	0.76
2:D:452:PHE:HE2	2:D:454:TYR:HE1	1.31	0.76
2:F:251:LEU:HA	2:F:588:ARG:HH12	1.49	0.76
2:F:585:HIS:CB	2:F:593:MET:HE3	2.15	0.76
2:A:411:PHE:N	2:A:411:PHE:HD1	1.83	0.76
2:D:320:THR:HG22	2:D:321:SER:H	1.50	0.76
2:F:621:VAL:O	2:F:624:VAL:HG22	1.86	0.76
2:G:255:HIS:ND1	2:G:267:HIS:HE1	1.82	0.76
1:J:75:U:O2'	1:J:76:G:OP1	2.02	0.76
1:M:97:C:H5''	1:M:98:A:OP1	1.84	0.76
2:A:551:GLN:NE2	2:A:597:GLY:HA2	2.01	0.76
1:I:88:G:H4'	1:I:89:U:OP1	1.85	0.76
2:A:411:PHE:CE2	2:A:521:PHE:CE1	2.74	0.76
2:C:637:LEU:H	2:C:637:LEU:HD12	1.50	0.76
2:D:407:MET:SD	2:D:514:ILE:HG21	2.25	0.76
2:E:320:THR:HG23	2:F:319:THR:O	1.86	0.76
2:E:390:GLU:HA	2:E:393:ILE:HG13	1.68	0.76
2:C:263:MET:HE2	2:D:298:MET:HG2	1.65	0.76
2:D:251:LEU:HA	2:D:588:ARG:HH12	1.51	0.76
2:F:460:ALA:HB1	2:F:462:TYR:CD1	2.21	0.76
1:L:88:G:H4'	1:L:89:U:OP1	1.85	0.75
2:A:325:ARG:HH11	2:A:325:ARG:HG3	1.51	0.75
2:E:452:PHE:HE2	2:E:454:TYR:HE1	1.34	0.75
2:A:517:SER:HB2	4:A:2002:TSB:H1'	1.68	0.75
2:E:621:VAL:O	2:E:624:VAL:HG22	1.85	0.75
1:J:97:C:H5''	1:J:98:A:OP1	1.86	0.75
2:A:402:ARG:HH11	2:A:402:ARG:HB2	1.50	0.75
2:H:299:MET:O	2:H:327:TYR:HB3	1.86	0.75
2:B:421:SER:CB	2:B:458:GLU:HB3	2.17	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:84:U:H4'	1:I:85:C:C5'	2.16	0.75
2:C:251:LEU:HA	2:C:588:ARG:NH1	2.02	0.75
1:K:76:G:N2	1:K:97:C:C2	2.55	0.75
2:A:298:MET:HG2	2:B:263:MET:CE	2.16	0.74
2:H:251:LEU:HA	2:H:588:ARG:NH1	2.03	0.74
2:E:473:CYS:HA	5:E:6010:HOH:O	1.86	0.74
2:H:411:PHE:HD2	2:H:525:LEU:CD2	1.99	0.74
2:A:411:PHE:HE2	2:A:521:PHE:HE1	1.33	0.74
1:L:88:G:H8	5:D:5016:HOH:O	1.69	0.74
2:A:452:PHE:HE2	2:A:454:TYR:HE1	1.33	0.74
2:E:411:PHE:N	2:E:411:PHE:CD1	2.54	0.74
2:F:321:SER:HB3	2:F:326:GLU:HA	1.70	0.74
2:H:465:LYS:HG3	2:H:484:GLN:HG2	1.68	0.74
1:P:86:G:H3'	5:P:1613:HOH:O	1.86	0.74
2:F:390:GLU:HA	2:F:393:ILE:HG13	1.69	0.74
1:J:103:G:OP2	5:J:1025:HOH:O	2.05	0.74
2:A:411:PHE:CE2	2:A:521:PHE:HE1	2.06	0.74
2:A:465:LYS:HG3	2:A:484:GLN:HG2	1.69	0.74
2:H:403:LEU:HG	2:H:407:MET:CE	2.18	0.74
2:C:251:LEU:HA	2:C:588:ARG:HH12	1.51	0.74
2:H:585:HIS:CB	2:H:593:MET:HE3	2.17	0.74
1:J:86:G:H3'	5:J:1013:HOH:O	1.88	0.74
2:C:621:VAL:O	2:C:624:VAL:HG22	1.86	0.74
1:K:77:U:H1'	1:K:98:A:N1	2.03	0.74
2:B:251:LEU:HA	2:B:588:ARG:NH1	2.02	0.74
2:B:585:HIS:HB2	2:B:593:MET:CE	2.17	0.74
2:C:411:PHE:N	2:C:411:PHE:HD1	1.85	0.74
2:H:452:PHE:HE2	2:H:454:TYR:HE1	1.35	0.74
2:D:390:GLU:HA	2:D:393:ILE:HG13	1.69	0.73
2:F:578:ILE:O	2:F:582:ILE:HG12	1.88	0.73
2:G:390:GLU:HA	2:G:393:ILE:HG13	1.68	0.73
1:O:88:G:H4'	1:O:89:U:OP1	1.86	0.73
2:G:362:HIS:HD2	5:G:732:HOH:O	1.71	0.73
1:K:82:U:H2'	1:K:83:U:H5'	1.70	0.73
2:E:276:GLU:O	2:E:279:VAL:HG22	1.88	0.73
2:H:637:LEU:H	2:H:637:LEU:HD12	1.53	0.73
1:L:98:A:H4'	5:L:1227:HOH:O	1.87	0.73
2:E:585:HIS:CB	2:E:593:MET:HE3	2.17	0.73
2:G:249:LYS:HB3	2:G:249:LYS:NZ	2.02	0.73
2:C:411:PHE:CE2	2:C:521:PHE:HE1	2.05	0.73
2:F:348:TYR:CD1	2:F:349:ARG:N	2.57	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:402:ARG:HB2	2:B:402:ARG:HH11	1.52	0.73
2:C:402:ARG:HH11	2:C:402:ARG:HB2	1.54	0.73
2:E:321:SER:HB3	2:E:326:GLU:HA	1.70	0.73
2:A:633:ARG:NH1	2:H:638:LYS:NZ	2.37	0.73
2:D:637:LEU:H	2:D:637:LEU:HD12	1.53	0.73
2:A:583:ARG:O	2:A:587:LEU:HD23	1.88	0.73
2:D:402:ARG:HB2	2:D:402:ARG:HH11	1.53	0.73
2:D:551:GLN:NE2	2:D:597:GLY:HA2	2.03	0.73
2:D:625:ILE:O	2:D:629:GLN:HG3	1.89	0.73
2:E:251:LEU:HA	2:E:588:ARG:NH1	2.04	0.73
2:E:411:PHE:HD2	2:E:525:LEU:CD2	1.99	0.73
2:F:452:PHE:HE2	2:F:454:TYR:HE1	1.35	0.73
2:A:251:LEU:HA	2:A:588:ARG:HH12	1.53	0.72
2:C:411:PHE:N	2:C:411:PHE:CD1	2.57	0.72
2:H:621:VAL:O	2:H:624:VAL:HG22	1.88	0.72
2:D:251:LEU:HA	2:D:588:ARG:NH1	2.04	0.72
2:B:411:PHE:HD2	2:B:525:LEU:CD2	2.02	0.72
1:M:69:G:OP3	1:P:105:C:O3'	2.08	0.72
2:F:501:ASP:OD2	2:F:503:GLU:HB2	1.89	0.72
1:L:78:G:O4'	2:C:345:LEU:HD23	1.89	0.72
1:M:88:G:H4'	1:M:89:U:OP1	1.87	0.72
2:B:411:PHE:N	2:B:411:PHE:CD1	2.51	0.72
2:D:578:ILE:O	2:D:582:ILE:HG12	1.89	0.72
2:E:585:HIS:HB2	2:E:593:MET:CE	2.18	0.72
2:F:637:LEU:HD12	2:F:637:LEU:H	1.54	0.72
2:G:547:ILE:HD12	2:G:597:GLY:HA3	1.71	0.72
2:H:390:GLU:HA	2:H:393:ILE:HG13	1.72	0.72
1:K:78:G:H2'	1:K:79:A:H8	1.54	0.72
2:C:403:LEU:HG	2:C:407:MET:CE	2.19	0.72
2:E:403:LEU:HG	2:E:407:MET:CE	2.20	0.72
2:F:517:SER:HB2	4:F:7002:TSB:H1'	1.72	0.72
2:A:277:LEU:O	2:A:281:VAL:HG23	1.89	0.72
2:A:637:LEU:HD12	2:A:637:LEU:H	1.54	0.72
2:F:403:LEU:HG	2:F:407:MET:CE	2.20	0.72
1:O:86:G:H2'	5:O:1513:HOH:O	1.88	0.72
2:F:411:PHE:CE2	2:F:521:PHE:CZ	2.77	0.72
2:C:578:ILE:O	2:C:582:ILE:HG12	1.90	0.72
2:G:417:VAL:CG1	2:G:471:TYR:HE1	2.03	0.72
1:K:90:G:H8	2:C:577:LYS:CD	2.02	0.71
1:L:89:U:O2	1:L:89:U:H2'	1.88	0.71
1:P:86:G:C4	2:H:547:ILE:HD13	2.25	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:407:MET:SD	2:E:514:ILE:HG21	2.30	0.71
1:J:88:G:H4'	1:J:89:U:OP1	1.90	0.71
1:L:79:A:H1'	2:C:502:ASN:OD1	1.90	0.71
2:A:320:THR:HG22	2:A:321:SER:H	1.54	0.71
2:A:411:PHE:N	2:A:411:PHE:CD1	2.56	0.71
2:A:462:TYR:CE2	2:A:511:HIS:CE1	2.78	0.71
2:E:625:ILE:O	2:E:629:GLN:HG3	1.90	0.71
2:F:547:ILE:HD12	2:F:597:GLY:HA3	1.72	0.71
2:G:392:GLN:O	2:G:396:GLU:HB2	1.89	0.71
2:G:637:LEU:H	2:G:637:LEU:HD12	1.54	0.71
2:H:411:PHE:N	2:H:411:PHE:HD1	1.88	0.71
1:O:91:G:O3'	5:O:1518:HOH:O	2.08	0.71
2:E:299:MET:O	2:E:327:TYR:HB3	1.90	0.71
2:A:280:PHE:O	2:A:283:SER:HB3	1.91	0.71
2:B:625:ILE:O	2:B:629:GLN:HG3	1.91	0.71
2:C:321:SER:HB3	2:C:326:GLU:HA	1.73	0.71
2:A:251:LEU:HA	2:A:588:ARG:NH1	2.05	0.71
2:C:243:ASP:OD2	2:C:245:ARG:HB2	1.91	0.71
2:C:319:THR:O	2:D:320:THR:HG23	1.91	0.71
2:C:390:GLU:HA	2:C:393:ILE:HG13	1.72	0.71
2:C:473:CYS:HA	5:C:4013:HOH:O	1.90	0.71
2:G:514:ILE:HG22	2:G:515:LEU:N	2.06	0.71
2:H:244:HIS:CE1	2:H:374:MET:HE2	2.25	0.71
1:K:82:U:C2'	1:K:83:U:H5'	2.21	0.71
2:E:578:ILE:O	2:E:582:ILE:HG12	1.91	0.71
1:K:70:G:H22	1:K:105:C:H1'	1.56	0.71
2:G:551:GLN:HE21	2:G:597:GLY:CA	2.03	0.71
2:H:277:LEU:O	2:H:281:VAL:HG23	1.91	0.71
2:B:637:LEU:H	2:B:637:LEU:HD12	1.56	0.71
2:E:402:ARG:HB2	2:E:402:ARG:HH11	1.54	0.71
1:K:97:C:H4'	1:K:98:A:OP1	1.91	0.70
1:O:89:U:C5	2:G:583:ARG:CZ	2.73	0.70
2:B:585:HIS:CB	2:B:593:MET:HE3	2.21	0.70
2:H:547:ILE:HD12	2:H:597:GLY:HA3	1.73	0.70
1:L:77:U:H1'	1:L:98:A:N1	2.06	0.70
2:A:585:HIS:HB2	2:A:593:MET:CE	2.21	0.70
2:E:637:LEU:H	2:E:637:LEU:HD12	1.56	0.70
2:G:321:SER:HB3	2:G:326:GLU:HA	1.72	0.70
2:C:517:SER:HB2	4:C:4002:TSB:H1'	1.72	0.70
1:M:84:U:H4'	1:M:85:C:C5'	2.22	0.70
2:C:411:PHE:CE2	2:C:521:PHE:CZ	2.79	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:320:THR:HG22	2:H:321:SER:H	1.53	0.70
1:L:97:C:H5'	1:L:98:A:OP1	1.91	0.70
2:C:585:HIS:HB2	2:C:593:MET:CE	2.20	0.70
2:A:621:VAL:O	2:A:624:VAL:HG22	1.92	0.70
2:B:276:GLU:O	2:B:279:VAL:HG22	1.91	0.70
2:B:578:ILE:O	2:B:582:ILE:HG12	1.92	0.70
2:F:411:PHE:N	2:F:411:PHE:HD1	1.89	0.70
1:J:77:U:H1'	1:J:98:A:N1	2.06	0.70
2:A:462:TYR:HD2	2:A:486:ASP:OD2	1.73	0.70
2:C:544:ILE:HD11	2:C:562:LEU:HD22	1.72	0.70
2:C:585:HIS:CB	2:C:593:MET:HE3	2.21	0.70
2:G:433:MET:HE3	2:G:487:PHE:HB3	1.73	0.70
2:G:501:ASP:OD2	2:G:503:GLU:HB2	1.92	0.70
1:N:84:U:H4'	1:N:85:C:C5'	2.22	0.70
2:C:460:ALA:HB1	2:C:462:TYR:CD1	2.27	0.70
1:J:91:G:H4'	5:J:1018:HOH:O	1.92	0.70
2:B:321:SER:HB3	2:B:326:GLU:HA	1.72	0.70
2:D:299:MET:O	2:D:327:TYR:HB3	1.91	0.70
2:H:321:SER:HB3	2:H:326:GLU:HA	1.74	0.70
2:A:290:TYR:CE2	2:A:512:ARG:NH1	2.60	0.70
2:A:585:HIS:CB	2:A:593:MET:HE3	2.22	0.70
2:D:249:LYS:NZ	2:D:249:LYS:HB3	2.07	0.70
2:F:460:ALA:HB1	2:F:462:TYR:CE1	2.26	0.70
1:K:92:G:H2'	1:K:93:U:O4'	1.92	0.69
2:D:411:PHE:N	2:D:411:PHE:CD1	2.58	0.69
2:F:592:TYR:OH	2:F:639:GLN:HB3	1.92	0.69
2:A:421:SER:HA	2:A:455:GLN:HB3	1.74	0.69
2:C:465:LYS:HG3	2:C:484:GLN:HG2	1.72	0.69
2:H:249:LYS:NZ	2:H:249:LYS:HB3	2.06	0.69
2:H:411:PHE:N	2:H:411:PHE:CD1	2.59	0.69
2:B:551:GLN:HE21	2:B:597:GLY:HA2	1.56	0.69
2:B:583:ARG:NH2	5:B:216:HOH:O	2.26	0.69
2:F:583:ARG:O	2:F:587:LEU:HD23	1.92	0.69
1:I:75:U:O2'	1:I:76:G:OP1	2.09	0.69
2:A:547:ILE:HD12	2:A:597:GLY:HA3	1.72	0.69
2:B:402:ARG:HB2	2:B:402:ARG:NH1	2.07	0.69
1:I:89:U:H2'	1:I:89:U:O2	1.90	0.69
2:G:303:LEU:HD11	2:G:339:GLN:HE21	1.57	0.69
1:I:87:U:H2'	5:I:110:HOH:O	1.92	0.69
1:P:88:G:H4'	1:P:89:U:OP1	1.92	0.69
2:A:460:ALA:HB1	2:A:462:TYR:CD1	2.28	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:70:G:H2'	1:K:71:C:H6	1.56	0.69
2:D:585:HIS:HB2	2:D:593:MET:CE	2.23	0.69
2:F:402:ARG:HB2	2:F:402:ARG:NH1	2.06	0.69
2:D:547:ILE:HD12	2:D:597:GLY:HA3	1.74	0.69
2:E:551:GLN:NE2	2:E:597:GLY:HA2	2.06	0.69
2:F:346:LYS:O	2:F:497:TYR:HA	1.93	0.69
2:G:411:PHE:HD2	2:G:525:LEU:CD2	2.04	0.69
2:H:554:TYR:O	2:H:558:LEU:HD13	1.92	0.69
1:P:77:U:H1'	1:P:98:A:N1	2.08	0.69
2:C:592:TYR:OH	2:C:639:GLN:HB3	1.92	0.69
2:E:421:SER:CB	2:E:458:GLU:HB3	2.21	0.69
2:E:583:ARG:O	2:E:587:LEU:HD23	1.92	0.69
1:K:70:G:N2	1:K:105:C:H1'	2.08	0.69
2:B:388:CYS:HB2	2:B:392:GLN:OE1	1.93	0.69
2:C:583:ARG:O	2:C:587:LEU:HD23	1.92	0.69
2:D:592:TYR:OH	2:D:639:GLN:HB3	1.93	0.69
2:F:421:SER:HA	2:F:455:GLN:HB3	1.75	0.69
2:B:551:GLN:NE2	2:B:597:GLY:HA2	2.08	0.68
1:J:78:G:O4'	2:A:345:LEU:HD23	1.93	0.68
2:B:411:PHE:CE2	2:B:521:PHE:CE1	2.81	0.68
1:J:88:G:H1'	2:B:583:ARG:HG2	1.76	0.68
1:K:105:C:H2'	1:K:105:C:O2	1.91	0.68
2:F:462:TYR:CE2	2:F:511:HIS:CE1	2.81	0.68
2:F:544:ILE:HD11	2:F:562:LEU:HD22	1.75	0.68
1:M:77:U:H1'	1:M:98:A:N1	2.08	0.68
1:P:100:U:H4'	5:P:1630:HOH:O	1.93	0.68
2:D:321:SER:HB3	2:D:326:GLU:HA	1.75	0.68
2:E:421:SER:HA	2:E:455:GLN:HB3	1.74	0.68
1:L:91:G:H4'	5:L:1218:HOH:O	1.92	0.68
2:A:460:ALA:HB1	2:A:462:TYR:CE1	2.28	0.68
2:D:421:SER:CB	2:D:458:GLU:HB3	2.22	0.68
2:H:551:GLN:HE21	2:H:597:GLY:CA	2.07	0.68
1:L:84:U:H4'	1:L:85:C:C5'	2.23	0.68
2:C:460:ALA:HB1	2:C:462:TYR:CE1	2.28	0.68
2:G:417:VAL:HG13	2:G:471:TYR:HE1	1.59	0.68
2:A:517:SER:CB	4:A:2002:TSB:H1'	2.24	0.68
2:D:585:HIS:CB	2:D:593:MET:HE3	2.24	0.68
2:F:500:GLU:C	2:F:502:ASN:H	2.02	0.68
1:I:98:A:H4'	5:I:114:HOH:O	1.93	0.68
1:K:91:G:O2'	1:K:92:G:H5'	1.94	0.68
1:N:77:U:H1'	1:N:98:A:N1	2.09	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:465:LYS:HG3	2:F:484:GLN:HG2	1.75	0.67
2:H:595:VAL:HB	2:H:607:ALA:HB3	1.76	0.67
1:M:91:G:O3'	5:M:1318:HOH:O	2.12	0.67
1:O:89:U:H5	2:G:583:ARG:CZ	2.07	0.67
2:A:348:TYR:CD1	2:A:349:ARG:N	2.61	0.67
2:B:403:LEU:HG	2:B:407:MET:CE	2.22	0.67
2:B:622:ASN:ND2	5:B:226:HOH:O	2.26	0.67
2:G:411:PHE:N	2:G:411:PHE:CD1	2.62	0.67
2:C:421:SER:HA	2:C:455:GLN:HB3	1.75	0.67
2:C:535:THR:HG22	2:C:632:ILE:HG12	1.75	0.67
2:D:595:VAL:HB	2:D:607:ALA:HB3	1.77	0.67
2:G:544:ILE:HD11	2:G:562:LEU:HD22	1.76	0.67
2:H:304:TRP:CZ3	2:H:328:CYS:HB2	2.28	0.67
2:H:346:LYS:O	2:H:497:TYR:HA	1.93	0.67
2:A:462:TYR:HE2	2:A:511:HIS:CE1	2.12	0.67
2:D:402:ARG:HB2	2:D:402:ARG:NH1	2.10	0.67
2:F:551:GLN:HE21	2:F:597:GLY:CA	2.08	0.67
2:E:460:ALA:HB1	2:E:462:TYR:CE1	2.30	0.67
2:E:554:TYR:O	2:E:557:GLU:HB3	1.94	0.67
2:F:290:TYR:CE2	2:F:512:ARG:NH1	2.62	0.67
2:G:407:MET:SD	2:G:514:ILE:HG21	2.34	0.67
1:J:89:U:C5	2:B:583:ARG:NH1	2.62	0.67
2:C:277:LEU:O	2:C:281:VAL:HG23	1.95	0.67
2:C:348:TYR:CD1	2:C:349:ARG:N	2.63	0.67
1:N:97:C:H5''	1:N:98:A:OP1	1.94	0.67
2:B:339:GLN:OE1	2:B:339:GLN:HA	1.93	0.67
2:D:403:LEU:HG	2:D:407:MET:CE	2.25	0.67
2:D:411:PHE:HD2	2:D:525:LEU:HD21	1.59	0.67
2:D:433:MET:HE3	2:D:487:PHE:HB3	1.76	0.67
2:E:460:ALA:HB1	2:E:462:TYR:CD1	2.30	0.67
2:F:411:PHE:N	2:F:411:PHE:CD1	2.59	0.67
2:A:346:LYS:O	2:A:497:TYR:HA	1.95	0.67
2:G:299:MET:O	2:G:327:TYR:HB3	1.95	0.67
2:G:402:ARG:HB2	2:G:402:ARG:NH1	2.09	0.67
2:A:402:ARG:HB2	2:A:402:ARG:NH1	2.10	0.67
2:A:403:LEU:HG	2:A:407:MET:CE	2.25	0.67
2:A:411:PHE:HD2	2:A:525:LEU:CD2	2.04	0.67
2:H:421:SER:HA	2:H:455:GLN:HB3	1.77	0.67
2:A:592:TYR:OH	2:A:639:GLN:HB3	1.94	0.67
2:B:390:GLU:HA	2:B:393:ILE:HG13	1.77	0.67
2:D:421:SER:HA	2:D:455:GLN:HB3	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:392:GLN:O	2:H:396:GLU:HB2	1.95	0.67
1:J:89:U:C5	2:B:583:ARG:CZ	2.79	0.66
2:E:402:ARG:HB2	2:E:402:ARG:NH1	2.10	0.66
2:G:329:ILE:N	2:G:329:ILE:HD12	2.10	0.66
1:M:103:G:H5''	5:M:1325:HOH:O	1.95	0.66
1:O:84:U:H4'	1:O:85:C:C5'	2.25	0.66
1:P:84:U:H4'	1:P:85:C:C5'	2.25	0.66
2:C:320:THR:HG23	2:D:319:THR:O	1.95	0.66
2:F:299:MET:O	2:F:327:TYR:HB3	1.95	0.66
2:G:585:HIS:HB2	2:G:593:MET:CE	2.25	0.66
1:L:75:U:O2'	1:L:76:G:OP1	2.13	0.66
2:B:441:LEU:HD12	2:B:464:PRO:HB2	1.77	0.66
2:B:621:VAL:O	2:B:624:VAL:HG22	1.94	0.66
2:G:304:TRP:CZ3	2:G:328:CYS:HB2	2.30	0.66
2:G:421:SER:CB	2:G:458:GLU:HB3	2.21	0.66
2:H:417:VAL:CG1	2:H:471:TYR:HE1	2.09	0.66
2:C:547:ILE:HD12	2:C:597:GLY:HA3	1.76	0.66
2:A:299:MET:O	2:A:327:TYR:HB3	1.96	0.66
2:B:517:SER:HB2	4:B:3002:TSB:H1'	1.78	0.66
2:C:551:GLN:HE21	2:C:597:GLY:CA	2.09	0.66
2:D:249:LYS:HB3	2:D:249:LYS:HZ3	1.61	0.66
2:E:544:ILE:HD11	2:E:562:LEU:HD22	1.78	0.66
2:G:625:ILE:O	2:G:629:GLN:HG3	1.96	0.66
1:O:89:U:O2	1:O:89:U:C2'	2.44	0.66
2:B:544:ILE:HD11	2:B:562:LEU:HD22	1.77	0.66
2:E:411:PHE:HE2	2:E:521:PHE:CE1	2.10	0.66
2:F:517:SER:CB	4:F:7002:TSB:H1'	2.25	0.66
2:G:421:SER:HA	2:G:455:GLN:HB3	1.78	0.66
2:H:402:ARG:HB2	2:H:402:ARG:NH1	2.10	0.66
2:B:433:MET:HE3	2:B:487:PHE:HB3	1.77	0.66
2:D:517:SER:HB2	4:D:5002:TSB:H1'	1.78	0.66
2:E:348:TYR:CD1	2:E:349:ARG:N	2.64	0.66
2:F:514:ILE:HG22	2:F:515:LEU:N	2.11	0.66
2:G:500:GLU:C	2:G:502:ASN:H	2.03	0.66
1:I:87:U:H1'	2:A:582:ILE:HG13	1.77	0.66
2:D:583:ARG:O	2:D:587:LEU:HD23	1.95	0.66
2:E:547:ILE:HD12	2:E:597:GLY:HA3	1.77	0.66
2:E:592:TYR:OH	2:E:639:GLN:HB3	1.95	0.66
2:F:243:ASP:OD2	2:F:245:ARG:HB2	1.95	0.66
1:K:101:G:H2'	1:K:102:C:H6	1.61	0.66
2:C:462:TYR:CE2	2:C:511:HIS:CE1	2.84	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:580:PHE:O	2:G:580:PHE:HD1	1.79	0.66
2:D:544:ILE:HD11	2:D:562:LEU:HD22	1.77	0.65
2:G:332:MET:HG2	2:G:363:ARG:HH21	1.61	0.65
2:G:417:VAL:CG1	2:G:471:TYR:CE1	2.79	0.65
2:H:325:ARG:HH11	2:H:325:ARG:HG3	1.61	0.65
2:A:544:ILE:HD11	2:A:562:LEU:HD22	1.76	0.65
2:C:421:SER:CB	2:C:458:GLU:HB3	2.24	0.65
2:C:433:MET:HE3	2:C:487:PHE:HB3	1.77	0.65
2:D:411:PHE:N	2:D:411:PHE:HD1	1.94	0.65
2:D:465:LYS:HG3	2:D:484:GLN:HG2	1.77	0.65
2:F:625:ILE:O	2:F:629:GLN:HG3	1.96	0.65
2:H:514:ILE:HG22	2:H:515:LEU:HD23	1.78	0.65
1:J:84:U:H4'	1:J:85:C:C5'	2.26	0.65
1:L:89:U:C5	2:D:583:ARG:CZ	2.79	0.65
2:A:500:GLU:C	2:A:502:ASN:H	2.03	0.65
2:B:243:ASP:OD2	2:B:245:ARG:HB2	1.96	0.65
2:B:325:ARG:HG3	2:B:325:ARG:HH11	1.61	0.65
2:B:547:ILE:HD12	2:B:597:GLY:HA3	1.78	0.65
2:E:434:TRP:O	2:E:438:GLU:HG3	1.97	0.65
1:O:77:U:H1'	1:O:98:A:N1	2.12	0.65
2:B:249:LYS:NZ	2:B:249:LYS:HB3	2.12	0.65
2:B:554:TYR:O	2:B:558:LEU:HD13	1.96	0.65
2:E:595:VAL:HB	2:E:607:ALA:HB3	1.79	0.65
2:F:580:PHE:O	2:F:580:PHE:HD1	1.79	0.65
2:B:299:MET:O	2:B:327:TYR:HB3	1.96	0.65
2:C:554:TYR:O	2:C:557:GLU:HB3	1.97	0.65
2:H:544:ILE:HD11	2:H:562:LEU:HD22	1.78	0.65
2:H:612:ARG:HH11	2:H:612:ARG:HG3	1.62	0.65
1:K:104:C:H2'	1:K:105:C:O4'	1.97	0.65
1:M:85:C:H5	2:E:547:ILE:O	1.79	0.65
2:F:407:MET:SD	2:F:514:ILE:HG21	2.35	0.65
2:G:554:TYR:O	2:G:558:LEU:HD13	1.97	0.65
1:K:90:G:OP2	2:C:577:LYS:HB2	1.96	0.65
1:N:83:U:H3	2:F:575:ASN:HD21	1.45	0.65
2:D:441:LEU:HD12	2:D:464:PRO:HB2	1.78	0.65
1:K:70:G:H2'	1:K:71:C:C6	2.31	0.65
1:N:86:G:O6	2:F:599:LYS:HB2	1.97	0.65
2:A:253:LEU:O	2:A:254:TYR:HB3	1.97	0.65
2:B:411:PHE:CE2	2:B:521:PHE:HE1	2.14	0.65
2:E:441:LEU:HD12	2:E:464:PRO:HB2	1.78	0.65
2:E:554:TYR:O	2:E:558:LEU:HD13	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:580:PHE:HD1	2:E:580:PHE:O	1.78	0.65
2:H:582:ILE:O	2:H:593:MET:HE1	1.97	0.65
2:F:421:SER:CB	2:F:458:GLU:HB3	2.27	0.65
2:G:277:LEU:O	2:G:281:VAL:HG23	1.96	0.65
2:G:280:PHE:CE2	2:G:407:MET:HG2	2.32	0.65
2:H:433:MET:HE3	2:H:487:PHE:HB3	1.78	0.65
2:E:517:SER:HB2	4:E:6002:TSB:H1'	1.79	0.65
2:H:500:GLU:C	2:H:502:ASN:H	2.05	0.65
1:M:89:U:H5	2:E:583:ARG:CZ	2.10	0.64
2:A:257:GLN:OE1	2:B:339:GLN:HB3	1.96	0.64
2:H:514:ILE:HG22	2:H:515:LEU:N	2.12	0.64
2:H:583:ARG:O	2:H:587:LEU:HD23	1.97	0.64
2:C:595:VAL:HB	2:C:607:ALA:HB3	1.78	0.64
2:E:433:MET:HE3	2:E:487:PHE:HB3	1.79	0.64
2:F:392:GLN:O	2:F:396:GLU:HB2	1.97	0.64
1:J:88:G:C8	5:J:1023:HOH:O	2.50	0.64
1:K:87:U:O2	1:K:87:U:C2'	2.45	0.64
1:K:88:G:C4'	1:K:89:U:OP1	2.45	0.64
2:A:249:LYS:HB3	2:A:249:LYS:NZ	2.11	0.64
2:B:334:CYS:N	2:B:335:PRO:HD2	2.12	0.64
2:B:500:GLU:C	2:B:502:ASN:H	2.04	0.64
1:K:90:G:H8	2:C:577:LYS:HD3	1.61	0.64
1:M:89:U:O2	1:M:89:U:C2'	2.43	0.64
2:C:299:MET:O	2:C:327:TYR:HB3	1.97	0.64
2:E:462:TYR:CE2	2:E:511:HIS:CE1	2.85	0.64
2:E:500:GLU:C	2:E:502:ASN:H	2.04	0.64
2:F:303:LEU:HD11	2:F:339:GLN:HE21	1.62	0.64
2:F:433:MET:HE3	2:F:487:PHE:HB3	1.78	0.64
1:K:86:G:C4	2:C:547:ILE:HD13	2.33	0.64
2:A:421:SER:CB	2:A:458:GLU:HB3	2.23	0.64
2:A:514:ILE:HG22	2:A:515:LEU:N	2.11	0.64
2:D:500:GLU:C	2:D:502:ASN:H	2.04	0.64
2:F:535:THR:HG22	2:F:632:ILE:HG12	1.79	0.64
2:H:391:GLU:OE1	2:H:391:GLU:N	2.30	0.64
2:H:417:VAL:HG13	2:H:471:TYR:HE1	1.62	0.64
2:B:392:GLN:O	2:B:396:GLU:HB2	1.98	0.64
2:D:339:GLN:OE1	2:D:339:GLN:HA	1.98	0.64
2:B:348:TYR:CD1	2:B:349:ARG:N	2.65	0.64
2:C:625:ILE:O	2:C:629:GLN:HG3	1.98	0.64
2:D:473:CYS:HA	5:D:5013:HOH:O	1.97	0.64
2:G:325:ARG:HG3	2:G:325:ARG:HH11	1.63	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:472:ASP:OD1	2:C:476:ARG:HB2	1.98	0.64
2:E:392:GLN:O	2:E:396:GLU:HB2	1.97	0.64
2:G:585:HIS:CB	2:G:593:MET:HE3	2.24	0.64
1:P:89:U:O2	1:P:89:U:C2'	2.46	0.64
2:A:244:HIS:CE1	2:A:374:MET:HE2	2.33	0.64
2:A:551:GLN:HE21	2:A:597:GLY:CA	2.11	0.64
2:C:339:GLN:HB3	2:D:257:GLN:OE1	1.98	0.64
2:D:422:THR:OG1	2:D:456:LEU:HA	1.98	0.64
2:E:465:LYS:HG3	2:E:484:GLN:HG2	1.80	0.64
2:F:365:GLU:H	2:F:377:ARG:HD3	1.63	0.64
2:H:625:ILE:O	2:H:629:GLN:HG3	1.98	0.64
1:N:86:G:C4	2:F:547:ILE:HD13	2.33	0.64
2:B:253:LEU:O	2:B:254:TYR:HB3	1.97	0.64
2:G:411:PHE:N	2:G:411:PHE:HD1	1.94	0.64
1:O:83:U:H3	2:G:575:ASN:HD21	1.45	0.63
2:D:348:TYR:CD1	2:D:349:ARG:N	2.67	0.63
2:F:253:LEU:O	2:F:254:TYR:HB3	1.98	0.63
2:F:411:PHE:CE2	2:F:521:PHE:HE1	2.11	0.63
2:D:460:ALA:HB1	2:D:462:TYR:CE1	2.33	0.63
2:F:462:TYR:HE2	2:F:511:HIS:CE1	2.16	0.63
1:O:97:C:C5'	1:O:98:A:OP1	2.45	0.63
2:A:369:SER:O	2:A:375:ARG:HD2	1.98	0.63
2:C:346:LYS:O	2:C:497:TYR:HA	1.98	0.63
2:D:243:ASP:OD2	2:D:245:ARG:HB2	1.98	0.63
2:C:402:ARG:HB2	2:C:402:ARG:NH1	2.14	0.63
2:B:411:PHE:HE2	2:B:521:PHE:HE1	1.47	0.63
2:B:421:SER:HA	2:B:455:GLN:HB3	1.80	0.63
2:D:417:VAL:HG13	2:D:471:TYR:HE1	1.63	0.63
2:G:334:CYS:N	2:G:335:PRO:HD2	2.13	0.63
1:J:87:U:H1'	2:B:582:ILE:HG13	1.81	0.63
1:P:84:U:H4'	1:P:85:C:O5'	1.99	0.63
2:B:417:VAL:HG13	2:B:471:TYR:HE1	1.64	0.63
2:C:257:GLN:HB3	2:C:259:GLU:OE2	1.98	0.63
2:D:554:TYR:O	2:D:558:LEU:HD13	1.98	0.63
2:C:612:ARG:HH11	2:C:612:ARG:HG3	1.63	0.63
1:N:78:G:O4'	2:E:345:LEU:HD23	1.98	0.63
1:O:89:U:O5'	5:O:1528:HOH:O	2.15	0.63
2:B:554:TYR:O	2:B:557:GLU:HB3	1.99	0.63
2:B:628:LEU:O	2:B:632:ILE:HG13	1.99	0.63
1:I:78:G:O2'	1:I:79:A:H5'	1.99	0.62
2:F:417:VAL:HG13	2:F:471:TYR:HE1	1.62	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:595:VAL:HB	2:G:607:ALA:HB3	1.80	0.62
1:I:77:U:H1'	1:I:98:A:N1	2.14	0.62
1:K:79:A:C2	1:K:94:C:C2	2.87	0.62
2:C:526:THR:HG23	2:C:533:PHE:HZ	1.63	0.62
2:D:460:ALA:HB1	2:D:462:TYR:CD1	2.35	0.62
2:A:417:VAL:CG1	2:A:471:TYR:HE1	2.12	0.62
2:A:417:VAL:HG13	2:A:471:TYR:HE1	1.62	0.62
2:A:630:GLN:HA	2:A:633:ARG:NH1	2.14	0.62
2:B:595:VAL:HB	2:B:607:ALA:HB3	1.79	0.62
2:C:249:LYS:NZ	2:C:249:LYS:HB3	2.15	0.62
2:G:391:GLU:OE1	2:G:391:GLU:N	2.32	0.62
2:B:582:ILE:O	2:B:593:MET:HE1	1.98	0.62
2:B:583:ARG:O	2:B:587:LEU:HD23	1.99	0.62
2:C:580:PHE:O	2:C:580:PHE:HD1	1.82	0.62
1:K:98:A:C5	2:D:345:LEU:HB2	2.34	0.62
2:A:332:MET:HE2	4:A:2002:TSB:HN12	1.63	0.62
2:C:517:SER:CB	4:C:4002:TSB:H1'	2.28	0.62
2:H:580:PHE:O	2:H:580:PHE:HD1	1.81	0.62
1:M:84:U:H4'	1:M:85:C:O5'	2.00	0.62
2:D:392:GLN:O	2:D:396:GLU:HB2	1.98	0.62
2:E:612:ARG:HH11	2:E:612:ARG:HG3	1.63	0.62
2:A:392:GLN:O	2:A:396:GLU:HB2	1.98	0.62
2:D:610:THR:CG2	2:D:614:LYS:HB3	2.29	0.62
2:E:249:LYS:NZ	2:E:249:LYS:HB3	2.15	0.62
2:E:277:LEU:O	2:E:281:VAL:HG23	2.00	0.62
2:E:582:ILE:O	2:E:593:MET:HE1	1.99	0.62
2:F:559:THR:HG21	2:F:571:ALA:CB	2.17	0.62
2:F:627:LYS:HB3	2:F:640:LEU:HD21	1.82	0.62
2:C:630:GLN:HA	2:C:633:ARG:NH1	2.14	0.62
2:F:277:LEU:O	2:F:281:VAL:HG23	2.00	0.62
2:H:388:CYS:HB2	2:H:392:GLN:OE1	2.00	0.62
2:H:418:VAL:HG21	2:H:450:ILE:HG21	1.82	0.62
2:H:421:SER:CB	2:H:458:GLU:HB3	2.26	0.62
1:I:84:U:H4'	1:I:85:C:O5'	2.00	0.62
2:A:433:MET:HE3	2:A:487:PHE:HB3	1.81	0.62
2:E:417:VAL:HG13	2:E:471:TYR:HE1	1.64	0.62
2:G:554:TYR:O	2:G:557:GLU:HB3	1.99	0.62
2:F:341:PHE:O	2:F:346:LYS:NZ	2.33	0.62
2:H:253:LEU:O	2:H:254:TYR:HB3	2.00	0.62
1:K:73:U:C5	1:K:97:C:N4	2.67	0.61
1:K:95:A:C1'	2:D:502:ASN:HD21	2.13	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:84:U:H4'	1:O:85:C:O5'	2.00	0.61
2:A:284:LYS:HG3	2:A:407:MET:HG3	1.82	0.61
2:B:542:VAL:HG12	2:B:543:VAL:N	2.15	0.61
2:E:627:LYS:HB3	2:E:640:LEU:HD21	1.81	0.61
2:F:249:LYS:NZ	2:F:249:LYS:HB3	2.13	0.61
2:H:411:PHE:CE2	2:H:521:PHE:CE1	2.88	0.61
2:H:411:PHE:CE2	2:H:521:PHE:HE1	2.17	0.61
2:A:580:PHE:O	2:A:580:PHE:HD1	1.83	0.61
2:B:465:LYS:HG3	2:B:484:GLN:HG2	1.82	0.61
2:E:253:LEU:O	2:E:254:TYR:HB3	1.99	0.61
2:H:592:TYR:OH	2:H:639:GLN:HB3	2.00	0.61
2:A:303:LEU:HD11	2:A:339:GLN:NE2	2.16	0.61
2:A:595:VAL:HB	2:A:607:ALA:HB3	1.80	0.61
2:D:303:LEU:HD11	2:D:339:GLN:HE21	1.65	0.61
2:E:374:MET:HE1	2:E:527:GLU:HG3	1.82	0.61
1:K:86:G:C5'	1:K:87:U:OP1	2.45	0.61
1:M:69:G:O5'	1:P:105:C:O2'	2.18	0.61
2:A:388:CYS:HB2	2:A:392:GLN:OE1	2.00	0.61
2:D:277:LEU:O	2:D:281:VAL:HG23	2.00	0.61
2:G:517:SER:CB	4:G:8002:TSB:H1'	2.30	0.61
2:H:310:TRP:O	2:H:314:LYS:HB3	2.01	0.61
2:D:374:MET:HE1	2:D:527:GLU:HG3	1.82	0.61
2:D:582:ILE:O	2:D:593:MET:HE1	2.00	0.61
2:E:263:MET:HE2	2:F:298:MET:HG2	1.82	0.61
1:N:75:U:O2'	1:N:76:G:OP1	2.19	0.61
2:C:543:VAL:HG11	2:C:585:HIS:CE1	2.36	0.61
2:D:580:PHE:O	2:D:580:PHE:HD1	1.84	0.61
2:G:582:ILE:O	2:G:593:MET:HE1	2.01	0.61
2:B:257:GLN:HB3	2:B:259:GLU:OE2	2.01	0.61
2:C:374:MET:HE1	2:C:527:GLU:HG3	1.82	0.61
2:B:580:PHE:HD1	2:B:580:PHE:O	1.83	0.61
2:C:417:VAL:CG1	2:C:471:TYR:HE1	2.14	0.61
2:C:417:VAL:HG13	2:C:471:TYR:HE1	1.66	0.61
2:D:257:GLN:HB3	2:D:259:GLU:OE2	2.01	0.61
2:E:422:THR:OG1	2:E:456:LEU:HA	2.00	0.61
2:F:375:ARG:NH2	5:F:7007:HOH:O	2.27	0.61
2:G:310:TRP:O	2:G:314:LYS:HB3	2.00	0.61
2:G:346:LYS:O	2:G:497:TYR:HA	2.01	0.61
2:G:517:SER:HB2	4:G:8002:TSB:H1'	1.81	0.61
2:G:544:ILE:HG22	2:G:555:VAL:HG13	1.81	0.61
1:O:86:G:O6	2:G:599:LYS:HB3	1.99	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:86:G:C2'	5:O:1513:HOH:O	2.45	0.60
1:O:101:G:O2'	1:O:102:C:H5'	2.01	0.60
2:B:259:GLU:H	2:B:259:GLU:CD	2.07	0.60
2:B:280:PHE:CE2	2:B:407:MET:HG2	2.36	0.60
2:B:517:SER:CB	4:B:3002:TSB:H1'	2.31	0.60
2:C:392:GLN:O	2:C:396:GLU:HB2	2.01	0.60
2:C:500:GLU:C	2:C:502:ASN:H	2.06	0.60
2:G:583:ARG:O	2:G:587:LEU:HD23	2.01	0.60
2:C:296:PRO:HG2	2:D:265:PHE:CE1	2.35	0.60
2:E:339:GLN:OE1	2:E:339:GLN:HA	2.00	0.60
2:E:346:LYS:O	2:E:497:TYR:HA	2.01	0.60
1:M:87:U:H3'	5:M:1319:HOH:O	2.00	0.60
2:C:244:HIS:CE1	2:C:374:MET:HE2	2.36	0.60
2:D:411:PHE:CE2	2:D:521:PHE:HZ	2.14	0.60
2:E:310:TRP:O	2:E:314:LYS:HB3	2.01	0.60
2:F:595:VAL:HB	2:F:607:ALA:HB3	1.82	0.60
2:G:630:GLN:HA	2:G:633:ARG:NH1	2.15	0.60
2:H:609:ARG:O	2:H:609:ARG:HG3	2.01	0.60
2:D:612:ARG:HG3	2:D:612:ARG:HH11	1.66	0.60
2:F:284:LYS:HG3	2:F:407:MET:HG3	1.83	0.60
2:F:422:THR:OG1	2:F:456:LEU:HA	2.00	0.60
2:F:554:TYR:O	2:F:558:LEU:HD13	2.01	0.60
1:M:83:U:H3	2:E:575:ASN:HD21	1.49	0.60
2:A:434:TRP:O	2:A:438:GLU:HG3	2.01	0.60
2:D:551:GLN:HE21	2:D:597:GLY:CA	2.14	0.60
2:H:369:SER:O	2:H:375:ARG:HD2	2.02	0.60
2:H:441:LEU:HD12	2:H:464:PRO:HB2	1.83	0.60
2:A:257:GLN:HB3	2:A:259:GLU:OE2	2.02	0.60
2:A:462:TYR:CD2	2:A:486:ASP:OD2	2.54	0.60
2:D:369:SER:O	2:D:375:ARG:HD2	2.01	0.60
2:F:606:VAL:O	2:F:618:SER:HA	2.01	0.60
2:G:281:VAL:HG12	2:G:285:LEU:HD12	1.82	0.60
1:M:86:G:O6	2:E:599:LYS:HB2	2.00	0.60
1:M:87:U:H2'	5:M:1319:HOH:O	2.00	0.60
1:P:87:U:H2'	5:P:1619:HOH:O	2.01	0.60
2:D:310:TRP:O	2:D:314:LYS:HB3	2.02	0.60
2:D:627:LYS:HB3	2:D:640:LEU:HD21	1.82	0.60
2:E:280:PHE:CE2	2:E:407:MET:HG2	2.37	0.60
2:F:462:TYR:HD2	2:F:486:ASP:OD2	1.85	0.60
2:F:543:VAL:HG11	2:F:585:HIS:CE1	2.36	0.60
2:E:417:VAL:CG1	2:E:471:TYR:HE1	2.14	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:403:LEU:HG	2:G:407:MET:HE3	1.83	0.60
2:G:452:PHE:CE2	2:G:454:TYR:HE1	2.13	0.60
2:H:280:PHE:O	2:H:283:SER:HB3	2.01	0.60
2:H:339:GLN:OE1	2:H:339:GLN:HA	2.02	0.60
2:H:517:SER:CB	4:H:9002:TSB:H1'	2.31	0.60
1:J:80:U:H2'	1:J:81:C:C6	2.37	0.60
2:B:630:GLN:HA	2:B:633:ARG:NH1	2.17	0.60
2:H:280:PHE:CE2	2:H:407:MET:HG2	2.37	0.60
1:N:84:U:H4'	1:N:85:C:O5'	2.02	0.60
2:C:627:LYS:HB3	2:C:640:LEU:HD21	1.82	0.60
2:D:393:ILE:HD11	2:D:508:VAL:HG11	1.82	0.60
2:E:303:LEU:HD11	2:E:339:GLN:HE21	1.67	0.60
2:G:254:TYR:CE1	2:G:373:LEU:HD21	2.37	0.60
2:G:620:ASP:HB3	2:G:623:GLU:HB2	1.84	0.60
1:I:101:G:O2'	1:I:102:C:H5'	2.02	0.59
1:M:69:G:P	1:P:105:C:O3'	2.60	0.59
2:A:325:ARG:HG3	2:A:325:ARG:NH1	2.10	0.59
2:A:423:ARG:HA	2:A:434:TRP:CZ3	2.37	0.59
2:A:484:GLN:HB2	2:A:511:HIS:HB2	1.84	0.59
2:D:346:LYS:O	2:D:497:TYR:HA	2.02	0.59
2:E:620:ASP:HB3	2:E:623:GLU:HB2	1.84	0.59
2:F:334:CYS:N	2:F:335:PRO:HD2	2.16	0.59
2:F:388:CYS:HB2	2:F:392:GLN:OE1	2.02	0.59
2:H:259:GLU:H	2:H:259:GLU:CD	2.07	0.59
1:I:89:U:C5	2:A:583:ARG:NH1	2.70	0.59
1:P:101:G:O2'	1:P:102:C:H5'	2.02	0.59
2:E:259:GLU:H	2:E:259:GLU:CD	2.08	0.59
2:H:517:SER:HB2	4:H:9002:TSB:H1'	1.83	0.59
2:A:472:ASP:OD1	2:A:476:ARG:HB2	2.01	0.59
2:A:606:VAL:O	2:A:618:SER:HA	2.02	0.59
2:B:421:SER:HB3	2:B:458:GLU:C	2.27	0.59
2:B:592:TYR:OH	2:B:639:GLN:HB3	2.03	0.59
2:B:609:ARG:HG3	2:B:609:ARG:O	2.02	0.59
2:C:339:GLN:OE1	2:C:339:GLN:HA	2.02	0.59
2:C:554:TYR:O	2:C:558:LEU:HD13	2.03	0.59
2:E:243:ASP:OD2	2:E:245:ARG:HB2	2.00	0.59
2:E:369:SER:O	2:E:375:ARG:HD2	2.00	0.59
2:E:546:ASN:OD1	2:E:573:LEU:HA	2.02	0.59
2:F:280:PHE:O	2:F:283:SER:HB3	2.02	0.59
2:H:257:GLN:HB3	2:H:259:GLU:OE2	2.00	0.59
2:A:332:MET:HG2	2:A:363:ARG:HH21	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:535:THR:HG22	2:E:632:ILE:HG12	1.85	0.59
2:F:417:VAL:CG1	2:F:471:TYR:HE1	2.15	0.59
2:H:535:THR:HG22	2:H:632:ILE:HG12	1.84	0.59
1:J:84:U:H4'	1:J:85:C:O5'	2.02	0.59
1:J:89:U:C6	2:B:583:ARG:CZ	2.85	0.59
1:N:101:G:O2'	1:N:102:C:H5'	2.02	0.59
2:C:325:ARG:HH11	2:C:325:ARG:HG3	1.67	0.59
2:E:388:CYS:HB2	2:E:392:GLN:OE1	2.02	0.59
2:G:247:ILE:O	2:G:250:GLN:HB2	2.02	0.59
2:G:259:GLU:CD	2:G:259:GLU:H	2.11	0.59
2:G:369:SER:O	2:G:375:ARG:HD2	2.01	0.59
2:A:514:ILE:HG22	2:A:515:LEU:HD23	1.84	0.59
2:B:514:ILE:HG22	2:B:515:LEU:N	2.18	0.59
2:E:411:PHE:CE2	2:E:521:PHE:CZ	2.90	0.59
2:F:612:ARG:HH11	2:F:612:ARG:HG3	1.67	0.59
2:G:281:VAL:HG12	2:G:285:LEU:CD1	2.32	0.59
2:G:434:TRP:O	2:G:438:GLU:HG3	2.02	0.59
2:G:612:ARG:HH11	2:G:612:ARG:HG3	1.66	0.59
2:A:259:GLU:CD	2:A:259:GLU:H	2.09	0.59
2:A:544:ILE:HG22	2:A:555:VAL:HG13	1.85	0.59
2:B:278:GLU:CB	2:B:358:PHE:HE2	2.15	0.59
2:D:273:ILE:HG23	2:D:537:LEU:HD22	1.85	0.59
2:D:325:ARG:HG3	2:D:325:ARG:HH11	1.66	0.59
2:H:304:TRP:CE3	2:H:328:CYS:HB2	2.38	0.59
1:P:87:U:C2'	5:P:1619:HOH:O	2.51	0.59
2:A:422:THR:OG1	2:A:456:LEU:HA	2.01	0.59
2:B:346:LYS:O	2:B:497:TYR:HA	2.03	0.59
2:B:407:MET:SD	2:B:514:ILE:HG21	2.43	0.59
2:D:411:PHE:CE2	2:D:521:PHE:HE1	2.05	0.59
2:D:434:TRP:O	2:D:438:GLU:HG3	2.02	0.59
2:E:393:ILE:HD11	2:E:508:VAL:HG11	1.85	0.59
2:E:517:SER:CB	4:E:6002:TSB:H1'	2.33	0.59
2:F:257:GLN:HB3	2:F:259:GLU:OE2	2.02	0.59
2:F:610:THR:CG2	2:F:614:LYS:HB3	2.33	0.59
2:G:592:TYR:OH	2:G:639:GLN:HB3	2.03	0.59
2:H:281:VAL:HG12	2:H:285:LEU:HD12	1.84	0.59
2:H:417:VAL:CG1	2:H:471:TYR:CE1	2.86	0.59
1:J:78:G:O2'	1:J:79:A:H5'	2.03	0.59
1:P:86:G:C4	2:H:547:ILE:CD1	2.85	0.59
2:A:329:ILE:HD12	2:A:329:ILE:N	2.17	0.59
2:B:310:TRP:O	2:B:314:LYS:HB3	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:544:ILE:HG22	2:C:555:VAL:HG13	1.85	0.59
2:C:620:ASP:HB3	2:C:623:GLU:HB2	1.84	0.59
2:D:417:VAL:CG1	2:D:471:TYR:HE1	2.16	0.59
2:D:554:TYR:O	2:D:557:GLU:HB3	2.02	0.59
1:M:75:U:O2'	1:M:76:G:OP1	2.21	0.59
2:A:365:GLU:H	2:A:377:ARG:HD3	1.67	0.59
2:A:554:TYR:O	2:A:558:LEU:HD13	2.03	0.59
2:C:278:GLU:HB3	2:C:358:PHE:HE2	1.68	0.59
2:D:244:HIS:CE1	2:D:374:MET:HE2	2.38	0.59
2:D:253:LEU:O	2:D:254:TYR:HB3	2.03	0.59
2:G:253:LEU:O	2:G:254:TYR:HB3	2.01	0.59
2:B:546:ASN:OD1	2:B:573:LEU:HA	2.03	0.58
2:C:303:LEU:HD11	2:C:339:GLN:HE21	1.68	0.58
1:K:98:A:N6	2:D:345:LEU:HD22	2.18	0.58
2:A:374:MET:HE1	2:A:527:GLU:HG3	1.85	0.58
2:B:627:LYS:HB3	2:B:640:LEU:HD21	1.85	0.58
2:C:556:ASN:ND2	2:C:573:LEU:HD11	2.18	0.58
2:D:421:SER:HB3	2:D:458:GLU:C	2.28	0.58
2:E:610:THR:CG2	2:E:614:LYS:HB3	2.33	0.58
2:F:351:LEU:HD22	2:F:387:PHE:O	2.03	0.58
2:F:374:MET:HE1	2:F:527:GLU:HG3	1.85	0.58
2:G:244:HIS:CE1	2:G:374:MET:HE2	2.38	0.58
2:G:290:TYR:CE2	2:G:512:ARG:NH1	2.71	0.58
2:G:375:ARG:NH2	5:G:710:HOH:O	2.26	0.58
1:O:77:U:O5'	1:O:77:U:H6	1.85	0.58
2:F:434:TRP:O	2:F:438:GLU:HG3	2.02	0.58
2:G:546:ASN:OD1	2:G:573:LEU:HA	2.04	0.58
1:M:78:G:H5'	5:M:1339:HOH:O	2.03	0.58
2:A:535:THR:HG22	2:A:632:ILE:HG12	1.84	0.58
2:A:587:LEU:HD23	2:A:587:LEU:N	2.14	0.58
2:A:625:ILE:O	2:A:629:GLN:HG3	2.03	0.58
2:C:369:SER:O	2:C:375:ARG:HD2	2.03	0.58
2:D:517:SER:CB	4:D:5002:TSB:H1'	2.33	0.58
2:E:551:GLN:HE21	2:E:597:GLY:CA	2.16	0.58
2:F:410:THR:C	2:F:411:PHE:HD1	2.12	0.58
2:F:526:THR:HG23	2:F:533:PHE:HZ	1.68	0.58
2:A:417:VAL:CG1	2:A:471:TYR:CE1	2.87	0.58
2:B:418:VAL:HG21	2:B:450:ILE:HG21	1.86	0.58
2:D:411:PHE:CD2	2:D:521:PHE:CZ	2.91	0.58
2:D:418:VAL:HG21	2:D:450:ILE:HG21	1.85	0.58
2:E:257:GLN:HB3	2:E:259:GLU:OE2	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:580:PHE:CD1	2:E:580:PHE:C	2.82	0.58
2:H:461:PHE:C	2:H:461:PHE:CD1	2.81	0.58
1:L:89:U:C5	2:D:583:ARG:NH1	2.72	0.58
2:A:612:ARG:HG3	2:A:612:ARG:HH11	1.67	0.58
2:D:535:THR:HG22	2:D:632:ILE:HG12	1.83	0.58
2:F:325:ARG:HH11	2:F:325:ARG:HG3	1.67	0.58
2:F:369:SER:O	2:F:375:ARG:HD2	2.03	0.58
2:H:325:ARG:HG3	2:H:325:ARG:NH1	2.18	0.58
1:I:89:U:O2	1:I:89:U:C2'	2.51	0.58
1:N:75:U:O2'	1:N:89:U:O4	2.21	0.58
2:C:365:GLU:H	2:C:377:ARG:HD3	1.68	0.58
2:C:422:THR:OG1	2:C:456:LEU:HA	2.03	0.58
2:E:273:ILE:HG23	2:E:537:LEU:HD22	1.85	0.58
2:E:622:ASN:O	2:E:626:GLU:HG3	2.04	0.58
2:A:536:TRP:CZ3	2:A:635:ARG:HD3	2.38	0.58
2:B:612:ARG:HH11	2:B:612:ARG:HG3	1.69	0.58
2:C:280:PHE:CE2	2:C:407:MET:HG2	2.38	0.58
2:C:514:ILE:HG22	2:C:515:LEU:N	2.19	0.58
2:F:472:ASP:OD1	2:F:476:ARG:HB2	2.03	0.58
2:F:630:GLN:HA	2:F:633:ARG:NH1	2.19	0.58
2:G:303:LEU:HD11	2:G:339:GLN:NE2	2.19	0.58
2:G:620:ASP:O	2:G:623:GLU:HB3	2.02	0.58
2:H:278:GLU:HB3	2:H:358:PHE:HE2	1.68	0.58
1:M:69:G:OP2	1:P:105:C:C3'	2.52	0.58
1:O:87:U:O4	2:G:607:ALA:HB1	2.04	0.58
2:A:441:LEU:HD12	2:A:464:PRO:HB2	1.85	0.58
2:E:542:VAL:HG12	2:E:543:VAL:N	2.18	0.58
2:H:290:TYR:CE2	2:H:512:ARG:NH1	2.71	0.58
2:H:348:TYR:CD1	2:H:348:TYR:C	2.82	0.58
2:A:304:TRP:CZ3	2:A:328:CYS:HB2	2.39	0.58
2:D:422:THR:O	2:D:458:GLU:O	2.21	0.58
2:F:281:VAL:HG12	2:F:285:LEU:HD12	1.86	0.58
2:F:484:GLN:HB2	2:F:511:HIS:HB2	1.85	0.58
2:G:527:GLU:O	2:G:528:GLU:O	2.21	0.58
2:H:620:ASP:O	2:H:623:GLU:HB3	2.04	0.58
1:N:89:U:O2	1:N:89:U:C2'	2.50	0.57
2:A:362:HIS:HE1	2:B:297:PHE:CD1	2.22	0.57
2:A:391:GLU:OE1	2:A:391:GLU:N	2.36	0.57
2:B:278:GLU:HB3	2:B:358:PHE:HE2	1.69	0.57
2:F:554:TYR:O	2:F:557:GLU:HB3	2.02	0.57
2:G:421:SER:HB3	2:G:458:GLU:CB	2.26	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:536:TRP:O	2:H:568:ARG:NH2	2.36	0.57
2:H:610:THR:CG2	2:H:614:LYS:HB3	2.34	0.57
1:M:89:U:H5	2:E:583:ARG:NH2	2.01	0.57
1:P:97:C:C5'	1:P:98:A:OP1	2.51	0.57
2:A:554:TYR:O	2:A:557:GLU:HB3	2.04	0.57
2:C:388:CYS:HB2	2:C:392:GLN:OE1	2.04	0.57
2:F:280:PHE:CE2	2:F:407:MET:HG2	2.38	0.57
2:G:441:LEU:HD12	2:G:464:PRO:HB2	1.84	0.57
2:G:528:GLU:O	2:G:530:ALA:N	2.37	0.57
2:H:620:ASP:HB3	2:H:623:GLU:HB2	1.84	0.57
1:L:89:U:O2	1:L:89:U:C2'	2.51	0.57
2:B:421:SER:HB3	2:B:458:GLU:CB	2.27	0.57
2:C:556:ASN:HD21	2:C:573:LEU:HD11	1.69	0.57
2:E:421:SER:HB3	2:E:458:GLU:C	2.28	0.57
2:E:514:ILE:HG22	2:E:515:LEU:N	2.18	0.57
2:F:423:ARG:HA	2:F:434:TRP:CZ3	2.38	0.57
2:H:249:LYS:HB3	2:H:249:LYS:HZ3	1.68	0.57
2:C:259:GLU:H	2:C:259:GLU:CD	2.12	0.57
2:C:338:VAL:HG11	2:C:493:LEU:HB2	1.85	0.57
2:D:472:ASP:OD1	2:D:476:ARG:HB2	2.04	0.57
2:E:265:PHE:CE1	2:F:296:PRO:HG2	2.39	0.57
2:F:391:GLU:N	2:F:391:GLU:OE1	2.37	0.57
2:G:341:PHE:O	2:G:346:LYS:NZ	2.36	0.57
2:A:449:ASN:ND2	2:H:452:PHE:CD2	2.72	0.57
2:B:391:GLU:OE1	2:B:391:GLU:N	2.38	0.57
2:B:434:TRP:O	2:B:438:GLU:HG3	2.04	0.57
2:C:620:ASP:O	2:C:623:GLU:HB3	2.04	0.57
2:D:630:GLN:HA	2:D:633:ARG:NH1	2.19	0.57
2:F:421:SER:HB3	2:F:458:GLU:C	2.30	0.57
2:F:582:ILE:O	2:F:593:MET:HE1	2.05	0.57
1:I:79:A:O2'	2:B:500:GLU:O	2.23	0.57
1:L:84:U:H4'	1:L:85:C:O5'	2.04	0.57
1:P:85:C:H5	2:H:547:ILE:O	1.86	0.57
2:E:462:TYR:HE2	2:E:511:HIS:CE1	2.23	0.57
2:H:332:MET:HG2	2:H:363:ARG:HH21	1.69	0.57
1:J:77:U:O5'	1:J:77:U:H6	1.87	0.57
1:L:89:U:H5	2:D:583:ARG:CZ	2.16	0.57
2:A:243:ASP:OD2	2:A:245:ARG:HB2	2.04	0.57
2:A:636:SER:OG	2:A:638:LYS:HD3	2.04	0.57
2:D:259:GLU:CD	2:D:259:GLU:H	2.10	0.57
2:D:606:VAL:O	2:D:618:SER:HA	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:403:LEU:O	2:E:407:MET:HB2	2.05	0.57
2:F:514:ILE:HG22	2:F:515:LEU:HD23	1.86	0.57
1:J:101:G:O2'	1:J:102:C:H5'	2.05	0.57
2:A:536:TRP:CE3	2:A:635:ARG:HD3	2.39	0.57
2:C:375:ARG:NH2	5:C:4008:HOH:O	2.36	0.57
2:D:411:PHE:CD2	2:D:521:PHE:HZ	2.23	0.57
2:D:620:ASP:O	2:D:624:VAL:HG13	2.04	0.57
2:G:348:TYR:CD1	2:G:348:TYR:C	2.82	0.57
1:I:84:U:H4'	1:I:85:C:H5'	1.86	0.57
1:K:79:A:O2'	2:D:500:GLU:O	2.22	0.57
1:L:101:G:O2'	1:L:102:C:H5'	2.05	0.57
2:B:325:ARG:HG3	2:B:325:ARG:NH1	2.19	0.57
2:B:550:SER:C	2:B:552:SER:H	2.12	0.57
2:C:249:LYS:HB3	2:C:249:LYS:HZ3	1.68	0.57
2:G:403:LEU:O	2:G:407:MET:HB2	2.04	0.57
2:B:620:ASP:HB3	2:B:623:GLU:HB2	1.87	0.57
2:E:620:ASP:O	2:E:624:VAL:HG13	2.05	0.57
2:H:243:ASP:OD2	2:H:245:ARG:HB2	2.04	0.57
2:H:422:THR:OG1	2:H:456:LEU:HA	2.05	0.57
2:H:547:ILE:HD12	2:H:597:GLY:CA	2.35	0.57
2:A:580:PHE:CD1	2:A:580:PHE:C	2.83	0.56
2:B:332:MET:HG2	2:B:363:ARG:HH21	1.70	0.56
2:B:417:VAL:CG1	2:B:471:TYR:HE1	2.18	0.56
2:B:500:GLU:OE1	2:B:500:GLU:N	2.37	0.56
2:B:610:THR:CG2	2:B:614:LYS:HB3	2.34	0.56
2:G:325:ARG:HG3	2:G:325:ARG:NH1	2.20	0.56
2:H:341:PHE:O	2:H:346:LYS:NZ	2.35	0.56
2:H:587:LEU:HD23	2:H:587:LEU:N	2.19	0.56
1:M:80:U:H2'	1:M:81:C:C6	2.40	0.56
1:P:75:U:O2'	1:P:76:G:OP1	2.23	0.56
2:A:633:ARG:NH2	2:H:638:LYS:HD2	2.18	0.56
2:B:638:LYS:HB2	2:B:642:GLU:HG3	1.87	0.56
2:C:536:TRP:O	2:C:568:ARG:NH2	2.38	0.56
2:D:484:GLN:HB2	2:D:511:HIS:HB2	1.87	0.56
2:E:418:VAL:HG21	2:E:450:ILE:HG21	1.86	0.56
2:G:285:LEU:HD22	2:G:290:TYR:CG	2.40	0.56
2:G:460:ALA:C	2:G:462:TYR:H	2.13	0.56
2:G:536:TRP:O	2:G:568:ARG:NH2	2.38	0.56
2:H:365:GLU:H	2:H:377:ARG:HD3	1.69	0.56
2:H:421:SER:HB3	2:H:458:GLU:C	2.31	0.56
1:J:80:U:H2'	1:J:81:C:H6	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:89:U:O2	1:J:89:U:C2'	2.48	0.56
1:O:93:U:N3	1:O:94:C:C5	2.74	0.56
2:A:297:PHE:CD1	2:B:362:HIS:HE1	2.22	0.56
2:B:277:LEU:O	2:B:281:VAL:HG23	2.06	0.56
2:C:253:LEU:O	2:C:254:TYR:HB3	2.05	0.56
2:C:397:VAL:O	2:C:400:CYS:HB2	2.04	0.56
2:C:462:TYR:HE2	2:C:511:HIS:CE1	2.21	0.56
2:C:546:ASN:OD1	2:C:573:LEU:HA	2.05	0.56
2:D:620:ASP:HB3	2:D:623:GLU:HB2	1.87	0.56
2:G:332:MET:HG2	2:G:363:ARG:NH2	2.21	0.56
2:G:411:PHE:CE2	2:G:521:PHE:CE1	2.94	0.56
2:G:542:VAL:HG12	2:G:543:VAL:N	2.20	0.56
2:H:472:ASP:OD1	2:H:476:ARG:HB2	2.04	0.56
2:H:542:VAL:HG12	2:H:543:VAL:N	2.20	0.56
2:A:304:TRP:CE3	2:A:328:CYS:HB2	2.40	0.56
2:A:620:ASP:HB3	2:A:623:GLU:HB2	1.87	0.56
2:C:421:SER:HB3	2:C:458:GLU:C	2.31	0.56
2:E:312:ASN:O	2:E:313:TYR:HB2	2.06	0.56
2:F:259:GLU:CD	2:F:259:GLU:H	2.12	0.56
2:F:278:GLU:HB3	2:F:358:PHE:HE2	1.70	0.56
2:F:332:MET:HG2	2:F:363:ARG:HH21	1.70	0.56
2:G:624:VAL:HG23	2:G:625:ILE:N	2.20	0.56
1:M:89:U:C5	2:E:583:ARG:NH1	2.73	0.56
2:D:391:GLU:OE1	2:D:391:GLU:N	2.38	0.56
2:D:490:PRO:HD3	2:D:509:MET:HE2	1.86	0.56
2:D:514:ILE:HG22	2:D:515:LEU:N	2.20	0.56
2:E:304:TRP:CZ3	2:E:328:CYS:HB2	2.41	0.56
2:F:338:VAL:HG11	2:F:493:LEU:HB2	1.88	0.56
2:F:609:ARG:HG3	2:F:609:ARG:O	2.05	0.56
2:G:580:PHE:C	2:G:580:PHE:CD1	2.81	0.56
1:K:95:A:H1'	2:D:502:ASN:HD21	1.71	0.56
1:P:87:U:C2'	1:P:88:G:OP2	2.53	0.56
2:B:544:ILE:HG12	2:B:594:LEU:HD12	1.88	0.56
2:D:462:TYR:CE2	2:D:511:HIS:CE1	2.93	0.56
2:D:544:ILE:HG22	2:D:555:VAL:HG13	1.87	0.56
2:G:374:MET:HE1	2:G:527:GLU:HG3	1.87	0.56
2:G:514:ILE:HG22	2:G:515:LEU:HD23	1.87	0.56
2:G:587:LEU:HD23	2:G:587:LEU:N	2.16	0.56
2:G:624:VAL:CG2	2:G:625:ILE:N	2.67	0.56
2:H:580:PHE:CD1	2:H:580:PHE:C	2.82	0.56
1:M:75:U:O2'	1:M:89:U:O4	2.22	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:620:ASP:O	2:C:624:VAL:HG13	2.06	0.56
2:E:338:VAL:HG11	2:E:493:LEU:HB2	1.88	0.56
2:F:312:ASN:O	2:F:313:TYR:HB2	2.06	0.56
2:H:273:ILE:HG23	2:H:537:LEU:CD2	2.36	0.56
2:H:281:VAL:HG12	2:H:285:LEU:CD1	2.35	0.56
2:H:400:CYS:SG	2:H:512:ARG:HG3	2.46	0.56
1:M:87:U:C2'	5:M:1319:HOH:O	2.54	0.56
2:B:369:SER:O	2:B:375:ARG:HD2	2.06	0.56
2:C:555:VAL:HB	2:C:573:LEU:HD21	1.88	0.56
2:D:312:ASN:O	2:D:313:TYR:HB2	2.05	0.56
2:F:403:LEU:HG	2:F:407:MET:HE3	1.88	0.56
2:F:418:VAL:HG21	2:F:450:ILE:HG21	1.88	0.56
1:I:78:G:O4'	2:B:345:LEU:HD23	2.06	0.56
1:K:98:A:HO2'	2:D:344:GLY:H	1.50	0.56
1:M:77:U:H6	1:M:77:U:O5'	1.89	0.56
2:A:547:ILE:HD12	2:A:597:GLY:CA	2.36	0.56
2:C:312:ASN:O	2:C:313:TYR:HB2	2.06	0.56
2:C:580:PHE:CD1	2:C:580:PHE:C	2.84	0.56
2:D:280:PHE:CE2	2:D:407:MET:HG2	2.41	0.56
2:D:334:CYS:N	2:D:335:PRO:HD2	2.21	0.56
2:G:461:PHE:CD1	2:G:461:PHE:C	2.84	0.56
2:H:403:LEU:O	2:H:407:MET:HB2	2.06	0.56
1:I:97:C:H5''	1:I:98:A:OP1	2.05	0.56
2:B:526:THR:O	2:B:530:ALA:N	2.37	0.56
2:D:332:MET:HG2	2:D:363:ARG:HH21	1.70	0.56
2:F:304:TRP:CZ3	2:F:328:CYS:HB2	2.40	0.56
2:G:609:ARG:HG3	2:G:609:ARG:O	2.06	0.56
2:A:301:ARG:O	2:A:305:GLU:HG3	2.06	0.55
2:C:410:THR:C	2:C:411:PHE:HD1	2.13	0.55
2:E:391:GLU:OE1	2:E:391:GLU:N	2.39	0.55
2:H:418:VAL:HG21	2:H:450:ILE:CG2	2.35	0.55
2:H:630:GLN:HA	2:H:633:ARG:NH1	2.21	0.55
1:P:86:G:O6	2:H:599:LYS:CB	2.54	0.55
2:F:422:THR:O	2:F:458:GLU:O	2.22	0.55
2:G:304:TRP:CE3	2:G:328:CYS:HB2	2.41	0.55
1:M:78:G:C5'	5:M:1339:HOH:O	2.53	0.55
2:B:556:ASN:ND2	2:B:573:LEU:HD11	2.21	0.55
2:E:490:PRO:HD3	2:E:509:MET:HE2	1.87	0.55
2:E:628:LEU:O	2:E:632:ILE:HG13	2.07	0.55
2:E:630:GLN:HA	2:E:633:ARG:NH1	2.21	0.55
2:G:580:PHE:HD1	2:G:580:PHE:C	2.14	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:73:U:O4	1:K:97:C:N3	2.39	0.55
2:C:310:TRP:O	2:C:314:LYS:HB3	2.06	0.55
2:C:547:ILE:HD12	2:C:597:GLY:CA	2.37	0.55
2:D:580:PHE:C	2:D:580:PHE:CD1	2.84	0.55
2:E:386:ILE:HG21	2:E:396:GLU:HG3	1.89	0.55
2:E:472:ASP:OD1	2:E:476:ARG:HB2	2.06	0.55
2:G:339:GLN:OE1	2:G:339:GLN:HA	2.05	0.55
2:G:411:PHE:CE2	2:G:521:PHE:HE1	2.24	0.55
2:G:421:SER:HB3	2:G:458:GLU:C	2.31	0.55
2:A:310:TRP:O	2:A:314:LYS:HB3	2.06	0.55
2:A:578:ILE:HD11	2:A:595:VAL:CG2	2.36	0.55
2:B:556:ASN:HD21	2:B:573:LEU:HD11	1.72	0.55
2:D:386:ILE:HG21	2:D:396:GLU:CG	2.37	0.55
2:E:362:HIS:HE1	2:F:297:PHE:CD1	2.24	0.55
2:E:461:PHE:C	2:E:461:PHE:CD1	2.85	0.55
2:F:304:TRP:CE3	2:F:328:CYS:HB2	2.42	0.55
2:G:422:THR:HB	2:G:438:GLU:OE2	2.06	0.55
2:H:397:VAL:O	2:H:400:CYS:HB2	2.07	0.55
2:H:443:VAL:O	2:H:447:GLU:HB2	2.05	0.55
1:K:97:C:C4'	1:K:98:A:OP1	2.55	0.55
1:M:93:U:H2'	1:M:94:C:H6	1.71	0.55
2:A:620:ASP:O	2:A:623:GLU:HB3	2.07	0.55
2:B:418:VAL:HG21	2:B:450:ILE:CG2	2.36	0.55
2:B:606:VAL:O	2:B:618:SER:HA	2.06	0.55
2:F:461:PHE:CD1	2:F:461:PHE:C	2.83	0.55
2:F:546:ASN:OD1	2:F:573:LEU:HA	2.06	0.55
2:G:278:GLU:HG2	2:G:518:MET:HE2	1.88	0.55
2:G:416:ILE:O	2:G:416:ILE:HG22	2.06	0.55
2:H:312:ASN:O	2:H:313:TYR:HB2	2.06	0.55
1:J:97:C:C5'	1:J:98:A:OP1	2.54	0.55
1:K:83:U:C2'	1:K:84:U:H5'	2.36	0.55
2:C:301:ARG:O	2:C:305:GLU:HG3	2.07	0.55
2:D:542:VAL:HG12	2:D:543:VAL:N	2.21	0.55
2:E:278:GLU:HG2	2:E:518:MET:HE2	1.89	0.55
2:E:580:PHE:HD1	2:E:580:PHE:C	2.15	0.55
2:F:536:TRP:O	2:F:568:ARG:NH2	2.40	0.55
2:F:620:ASP:O	2:F:624:VAL:HG13	2.06	0.55
2:G:254:TYR:CD2	2:G:255:HIS:N	2.74	0.55
2:B:374:MET:HE1	2:B:527:GLU:HG3	1.89	0.55
2:B:422:THR:O	2:B:423:ARG:C	2.50	0.55
2:C:265:PHE:CE1	2:D:296:PRO:HG2	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:417:VAL:CG1	2:E:471:TYR:CE1	2.89	0.55
2:F:244:HIS:CE1	2:F:374:MET:HE2	2.40	0.55
2:F:417:VAL:CG1	2:F:471:TYR:CE1	2.89	0.55
2:F:551:GLN:HG2	2:F:597:GLY:C	2.32	0.55
2:F:556:ASN:ND2	2:F:573:LEU:HD11	2.22	0.55
2:G:329:ILE:N	2:G:329:ILE:CD1	2.70	0.55
2:G:569:VAL:HG22	2:G:570:LYS:N	2.21	0.55
1:K:90:G:C8	2:C:577:LYS:CD	2.87	0.55
2:B:365:GLU:H	2:B:377:ARG:HD3	1.72	0.55
2:C:423:ARG:HA	2:C:434:TRP:CZ3	2.41	0.55
2:C:609:ARG:HG3	2:C:609:ARG:O	2.07	0.55
2:G:388:CYS:HB2	2:G:392:GLN:OE1	2.05	0.55
2:G:610:THR:CG2	2:G:614:LYS:HB3	2.37	0.55
2:H:544:ILE:HG22	2:H:555:VAL:HG13	1.88	0.55
1:J:72:G:H2'	1:J:73:U:H6	1.72	0.55
1:K:83:U:O2'	1:K:84:U:H5'	2.07	0.55
1:O:75:U:O2'	1:O:76:G:OP1	2.24	0.55
2:A:462:TYR:HE2	2:A:511:HIS:NE2	2.04	0.55
2:B:298:MET:HE2	2:B:327:TYR:CD1	2.42	0.55
2:D:351:LEU:HD22	2:D:387:PHE:O	2.07	0.55
2:E:339:GLN:HB3	2:F:257:GLN:OE1	2.06	0.55
2:F:580:PHE:CD1	2:F:580:PHE:C	2.83	0.55
2:G:452:PHE:HE2	2:G:454:TYR:CE1	2.13	0.55
1:L:78:G:O2'	1:L:79:A:H5'	2.06	0.54
2:B:484:GLN:HB2	2:B:511:HIS:HB2	1.89	0.54
2:D:278:GLU:HG2	2:D:518:MET:HE2	1.89	0.54
2:D:304:TRP:CZ3	2:D:328:CYS:HB2	2.43	0.54
2:D:388:CYS:HB2	2:D:392:GLN:OE1	2.06	0.54
2:G:627:LYS:HB3	2:G:640:LEU:HD21	1.89	0.54
1:M:97:C:C5'	1:M:98:A:OP1	2.52	0.54
2:C:334:CYS:N	2:C:335:PRO:HD2	2.22	0.54
2:D:284:LYS:HG3	2:D:407:MET:HG3	1.90	0.54
2:F:547:ILE:HD12	2:F:597:GLY:CA	2.36	0.54
2:G:407:MET:CE	2:G:407:MET:CG	2.85	0.54
2:H:303:LEU:HD11	2:H:339:GLN:HE21	1.72	0.54
2:H:338:VAL:HG11	2:H:493:LEU:HB2	1.89	0.54
2:H:393:ILE:HD11	2:H:508:VAL:HG11	1.89	0.54
2:H:554:TYR:O	2:H:557:GLU:HB3	2.06	0.54
1:K:90:G:P	2:C:577:LYS:HZ2	2.30	0.54
2:A:546:ASN:OD1	2:A:573:LEU:HA	2.08	0.54
2:B:441:LEU:O	2:B:444:ALA:HB3	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:461:PHE:CD1	2:C:461:PHE:C	2.86	0.54
2:D:403:LEU:O	2:D:407:MET:HB2	2.07	0.54
2:D:417:VAL:CG1	2:D:471:TYR:CE1	2.91	0.54
2:E:638:LYS:HB2	2:E:642:GLU:HG3	1.89	0.54
2:F:597:GLY:O	2:F:601:VAL:HG23	2.06	0.54
2:F:638:LYS:HB2	2:F:642:GLU:HG3	1.90	0.54
2:G:257:GLN:HB3	2:G:259:GLU:OE2	2.07	0.54
2:G:312:ASN:O	2:G:313:TYR:HB2	2.05	0.54
2:G:484:GLN:HB2	2:G:511:HIS:HB2	1.88	0.54
2:H:334:CYS:N	2:H:335:PRO:HD2	2.23	0.54
2:A:341:PHE:O	2:A:346:LYS:NZ	2.38	0.54
2:A:526:THR:HG23	2:A:533:PHE:HZ	1.72	0.54
2:A:578:ILE:HD11	2:A:595:VAL:HG22	1.88	0.54
2:C:348:TYR:CD1	2:C:348:TYR:C	2.84	0.54
2:C:417:VAL:CG1	2:C:471:TYR:CE1	2.89	0.54
2:D:461:PHE:CD1	2:D:461:PHE:C	2.86	0.54
2:E:365:GLU:H	2:E:377:ARG:HD3	1.71	0.54
2:E:386:ILE:HG21	2:E:396:GLU:CG	2.38	0.54
2:F:536:TRP:CE3	2:F:635:ARG:HD3	2.43	0.54
2:G:335:PRO:HG2	2:G:336:GLY:H	1.72	0.54
2:H:460:ALA:HB1	2:H:462:TYR:CD1	2.42	0.54
2:H:580:PHE:HD1	2:H:580:PHE:C	2.15	0.54
1:M:78:G:O4'	2:F:345:LEU:HD23	2.08	0.54
2:A:536:TRP:O	2:A:568:ARG:NH2	2.40	0.54
2:B:348:TYR:CD1	2:B:348:TYR:C	2.86	0.54
2:D:325:ARG:HG3	2:D:325:ARG:NH1	2.22	0.54
2:G:383:ASP:OD2	2:G:385:HIS:CE1	2.60	0.54
1:K:101:G:H2'	1:K:102:C:C6	2.41	0.54
2:A:312:ASN:O	2:A:313:TYR:HB2	2.06	0.54
2:A:351:LEU:HD22	2:A:387:PHE:O	2.07	0.54
2:B:284:LYS:HG3	2:B:407:MET:HG3	1.90	0.54
2:C:331:PRO:O	2:C:361:CYS:HB3	2.07	0.54
2:D:610:THR:HG21	2:D:614:LYS:HB3	1.90	0.54
2:F:460:ALA:C	2:F:462:TYR:N	2.64	0.54
2:F:556:ASN:HD21	2:F:573:LEU:HD11	1.73	0.54
2:G:278:GLU:HB3	2:G:358:PHE:HE2	1.72	0.54
2:G:422:THR:O	2:G:458:GLU:O	2.25	0.54
2:H:284:LYS:HG3	2:H:407:MET:HG3	1.90	0.54
2:H:434:TRP:O	2:H:438:GLU:HG3	2.08	0.54
2:A:288:TYR:N	2:A:288:TYR:CD1	2.76	0.54
2:C:332:MET:HG2	2:C:363:ARG:HH21	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:391:GLU:N	2:C:391:GLU:OE1	2.40	0.54
2:C:610:THR:CG2	2:C:614:LYS:HB3	2.38	0.54
2:D:290:TYR:CE2	2:D:512:ARG:NH1	2.75	0.54
2:D:546:ASN:OD1	2:D:573:LEU:HA	2.08	0.54
2:E:422:THR:O	2:E:458:GLU:O	2.25	0.54
2:E:606:VAL:O	2:E:618:SER:HA	2.08	0.54
2:F:539:PRO:O	2:F:568:ARG:HD2	2.07	0.54
2:F:557:GLU:O	2:F:560:GLN:HB3	2.08	0.54
2:A:278:GLU:HG2	2:A:518:MET:HE2	1.89	0.54
2:A:461:PHE:CD1	2:A:461:PHE:C	2.85	0.54
2:A:490:PRO:HD3	2:A:509:MET:HE2	1.90	0.54
2:B:273:ILE:HG23	2:B:537:LEU:HD22	1.90	0.54
2:E:330:LYS:O	5:E:6011:HOH:O	2.18	0.54
2:E:423:ARG:HA	2:E:434:TRP:CZ3	2.42	0.54
2:F:441:LEU:HD12	2:F:464:PRO:HB2	1.89	0.54
2:H:462:TYR:CE2	2:H:511:HIS:CE1	2.96	0.54
1:K:83:U:H1'	1:K:90:G:C6	2.43	0.54
2:A:281:VAL:HG12	2:A:285:LEU:HD12	1.90	0.54
2:A:421:SER:HB3	2:A:458:GLU:C	2.32	0.54
2:A:544:ILE:HG12	2:A:594:LEU:HD12	1.89	0.54
2:A:627:LYS:HB3	2:A:640:LEU:HD21	1.90	0.54
2:B:423:ARG:HD2	2:B:438:GLU:OE2	2.07	0.54
2:B:460:ALA:C	2:B:462:TYR:H	2.16	0.54
2:D:411:PHE:HE2	2:D:521:PHE:CZ	2.10	0.54
2:D:536:TRP:CE3	2:D:635:ARG:HD3	2.43	0.54
2:E:362:HIS:HD2	5:E:6011:HOH:O	1.91	0.54
2:F:335:PRO:HG2	2:F:336:GLY:H	1.72	0.54
2:G:365:GLU:N	2:G:377:ARG:HD3	2.21	0.54
2:G:465:LYS:HG3	2:G:484:GLN:HG2	1.90	0.54
2:G:578:ILE:HD11	2:G:595:VAL:HG22	1.90	0.54
2:H:274:PHE:CE1	2:H:518:MET:HG3	2.43	0.54
1:J:75:U:O2'	1:J:89:U:O4	2.24	0.54
2:A:298:MET:CG	2:B:263:MET:HE2	2.30	0.54
2:A:411:PHE:CE2	2:A:521:PHE:CZ	2.96	0.54
2:A:542:VAL:HG12	2:A:543:VAL:N	2.23	0.54
2:B:620:ASP:O	2:B:623:GLU:HB3	2.08	0.54
2:D:273:ILE:HG23	2:D:537:LEU:CD2	2.38	0.54
2:E:325:ARG:HH11	2:E:325:ARG:HG3	1.73	0.54
2:E:484:GLN:HB2	2:E:511:HIS:HB2	1.90	0.54
2:F:544:ILE:HG22	2:F:555:VAL:HG13	1.89	0.54
2:F:580:PHE:HD1	2:F:580:PHE:C	2.16	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:620:ASP:HB3	2:F:623:GLU:HB2	1.89	0.54
2:G:380:THR:N	2:G:519:GLU:OE2	2.41	0.54
2:H:578:ILE:HD11	2:H:595:VAL:HG22	1.90	0.54
2:D:278:GLU:HB3	2:D:358:PHE:HE2	1.73	0.53
2:D:410:THR:C	2:D:411:PHE:HD1	2.15	0.53
2:E:284:LYS:HG3	2:E:407:MET:HG3	1.88	0.53
2:E:332:MET:HG2	2:E:363:ARG:HH21	1.72	0.53
2:E:604:GLY:O	2:E:621:VAL:HG23	2.09	0.53
2:H:335:PRO:HG2	2:H:336:GLY:H	1.73	0.53
2:A:622:ASN:O	2:A:626:GLU:HG3	2.07	0.53
2:D:338:VAL:HG11	2:D:493:LEU:HB2	1.89	0.53
2:E:281:VAL:HG12	2:E:285:LEU:HD12	1.91	0.53
2:F:310:TRP:O	2:F:314:LYS:HB3	2.08	0.53
2:G:400:CYS:SG	2:G:512:ARG:HG3	2.48	0.53
2:H:423:ARG:HA	2:H:434:TRP:CZ3	2.44	0.53
2:H:460:ALA:HB1	2:H:462:TYR:CE1	2.43	0.53
2:H:606:VAL:O	2:H:618:SER:HA	2.08	0.53
1:I:86:G:C2'	5:I:109:HOH:O	2.45	0.53
1:O:75:U:O2'	1:O:89:U:O4	2.25	0.53
2:G:422:THR:OG1	2:G:456:LEU:HA	2.08	0.53
1:M:89:U:C6	2:E:583:ARG:CZ	2.92	0.53
1:P:103:G:C5'	5:P:1625:HOH:O	2.56	0.53
2:A:278:GLU:HB3	2:A:358:PHE:HE2	1.74	0.53
2:A:583:ARG:O	2:A:587:LEU:CD2	2.55	0.53
2:B:338:VAL:HG11	2:B:493:LEU:HB2	1.91	0.53
2:B:422:THR:OG1	2:B:456:LEU:HA	2.09	0.53
2:B:637:LEU:HD12	2:B:637:LEU:N	2.22	0.53
2:C:325:ARG:HG3	2:C:325:ARG:NH1	2.23	0.53
2:D:473:CYS:HB2	2:D:527:GLU:OE1	2.09	0.53
2:E:624:VAL:CG2	2:E:625:ILE:N	2.72	0.53
2:F:572:ASP:OD2	2:F:574:ARG:NH2	2.31	0.53
2:G:278:GLU:CB	2:G:358:PHE:HE2	2.22	0.53
2:H:374:MET:HE1	2:H:527:GLU:HG3	1.90	0.53
1:O:80:U:H2'	1:O:81:C:C6	2.44	0.53
1:P:86:G:O6	2:H:599:LYS:HB2	2.09	0.53
1:P:103:G:H5''	5:P:1625:HOH:O	2.08	0.53
2:B:375:ARG:NH1	5:B:210:HOH:O	2.35	0.53
2:B:386:ILE:HG21	2:B:396:GLU:HG3	1.90	0.53
2:B:460:ALA:C	2:B:462:TYR:N	2.66	0.53
2:C:559:THR:HG21	2:C:571:ALA:CB	2.20	0.53
2:C:600:GLU:HB3	2:C:605:LYS:O	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:638:LYS:HB2	2:C:642:GLU:HG3	1.90	0.53
2:E:274:PHE:CD1	2:E:274:PHE:C	2.86	0.53
2:E:298:MET:HG2	2:F:263:MET:CE	2.36	0.53
2:F:325:ARG:HG3	2:F:325:ARG:NH1	2.23	0.53
2:F:403:LEU:O	2:F:407:MET:HB2	2.08	0.53
2:G:547:ILE:HD12	2:G:597:GLY:CA	2.37	0.53
2:H:422:THR:O	2:H:458:GLU:O	2.26	0.53
2:H:622:ASN:ND2	5:H:826:HOH:O	2.36	0.53
1:I:88:G:HI'	2:A:583:ARG:HG2	1.90	0.53
1:K:84:U:H4'	1:K:85:C:O5'	2.09	0.53
1:N:79:A:O2'	2:E:502:ASN:HB2	2.07	0.53
2:A:320:THR:CG2	2:B:319:THR:H	2.22	0.53
2:B:620:ASP:O	2:B:624:VAL:HG13	2.08	0.53
2:D:460:ALA:C	2:D:462:TYR:H	2.16	0.53
2:D:556:ASN:ND2	2:D:573:LEU:HD11	2.23	0.53
2:E:620:ASP:O	2:E:623:GLU:HB3	2.08	0.53
2:F:339:GLN:HA	2:F:339:GLN:OE1	2.08	0.53
2:F:555:VAL:HB	2:F:573:LEU:HD21	1.89	0.53
2:G:301:ARG:O	2:G:305:GLU:HG3	2.08	0.53
2:H:624:VAL:CG2	2:H:625:ILE:N	2.71	0.53
2:H:637:LEU:HD12	2:H:637:LEU:N	2.22	0.53
1:K:98:A:C6	2:D:345:LEU:HB2	2.42	0.53
1:P:77:U:H6	1:P:77:U:O5'	1.92	0.53
2:A:620:ASP:O	2:A:624:VAL:HG13	2.08	0.53
2:B:288:TYR:N	2:B:288:TYR:HD1	2.07	0.53
2:C:539:PRO:O	2:C:568:ARG:HD2	2.08	0.53
2:F:288:TYR:N	2:F:288:TYR:CD1	2.77	0.53
2:F:348:TYR:CD1	2:F:348:TYR:C	2.86	0.53
2:A:356:ALA:HA	2:A:384:ALA:HA	1.91	0.53
2:A:580:PHE:HD1	2:A:580:PHE:C	2.16	0.53
2:B:422:THR:HB	2:B:438:GLU:OE2	2.09	0.53
2:B:461:PHE:C	2:B:461:PHE:CD1	2.86	0.53
2:B:519:GLU:CD	2:B:519:GLU:H	2.17	0.53
2:D:288:TYR:N	2:D:288:TYR:CD1	2.77	0.53
2:D:288:TYR:N	2:D:288:TYR:HD1	2.06	0.53
2:E:600:GLU:HB3	2:E:605:LYS:O	2.09	0.53
2:H:278:GLU:CB	2:H:358:PHE:HE2	2.22	0.53
2:A:339:GLN:OE1	2:A:339:GLN:HA	2.09	0.53
2:B:472:ASP:OD1	2:B:476:ARG:HB2	2.09	0.53
2:C:403:LEU:HG	2:C:407:MET:HE3	1.90	0.53
2:D:329:ILE:HD12	2:D:329:ILE:N	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:334:CYS:N	2:E:335:PRO:HD2	2.24	0.53
2:F:460:ALA:C	2:F:462:TYR:H	2.15	0.53
2:H:329:ILE:HD12	2:H:329:ILE:N	2.24	0.53
1:M:69:G:OP2	1:P:105:C:H2'	2.09	0.53
2:A:254:TYR:CZ	2:A:373:LEU:HD21	2.44	0.53
2:A:338:VAL:HG11	2:A:493:LEU:HB2	1.91	0.53
2:A:348:TYR:CD1	2:A:348:TYR:C	2.86	0.53
2:A:555:VAL:HB	2:A:573:LEU:HD21	1.90	0.53
2:A:638:LYS:HB2	2:A:642:GLU:HG3	1.90	0.53
2:B:281:VAL:HG12	2:B:285:LEU:HD12	1.90	0.53
2:B:288:TYR:N	2:B:288:TYR:CD1	2.76	0.53
2:C:284:LYS:HG3	2:C:407:MET:HG3	1.90	0.53
2:C:403:LEU:O	2:C:407:MET:HB2	2.08	0.53
2:G:243:ASP:OD2	2:G:245:ARG:HB2	2.08	0.53
2:H:309:HIS:HB3	2:H:317:MET:CE	2.36	0.53
2:H:627:LYS:HB3	2:H:640:LEU:HD21	1.89	0.53
1:P:80:U:H2'	1:P:81:C:C6	2.44	0.52
2:A:334:CYS:N	2:A:335:PRO:HD2	2.24	0.52
2:C:423:ARG:HD2	2:C:438:GLU:OE2	2.08	0.52
2:D:281:VAL:HG12	2:D:285:LEU:HD12	1.91	0.52
2:E:288:TYR:CD1	2:E:288:TYR:N	2.77	0.52
2:G:462:TYR:CE2	2:G:511:HIS:CE1	2.96	0.52
2:H:247:ILE:O	2:H:250:GLN:HB2	2.09	0.52
2:H:375:ARG:NH2	5:H:810:HOH:O	2.17	0.52
1:O:93:U:C2	1:O:94:C:C5	2.97	0.52
1:P:75:U:O2'	1:P:89:U:O4	2.26	0.52
2:A:288:TYR:N	2:A:288:TYR:HD1	2.07	0.52
2:A:362:HIS:CE1	2:B:297:PHE:CG	2.97	0.52
2:A:416:ILE:O	2:A:416:ILE:HG22	2.09	0.52
2:B:551:GLN:HE21	2:B:597:GLY:CA	2.20	0.52
2:B:580:PHE:CD1	2:B:580:PHE:C	2.85	0.52
2:C:281:VAL:HG12	2:C:285:LEU:HD12	1.91	0.52
2:D:301:ARG:O	2:D:305:GLU:HG3	2.09	0.52
2:D:536:TRP:O	2:D:568:ARG:NH2	2.42	0.52
2:D:556:ASN:HD21	2:D:573:LEU:HD11	1.74	0.52
2:E:278:GLU:HB3	2:E:358:PHE:HE2	1.74	0.52
2:E:297:PHE:CD1	2:F:362:HIS:HE1	2.27	0.52
2:F:423:ARG:O	2:F:423:ARG:HG2	2.10	0.52
2:G:418:VAL:HG21	2:G:450:ILE:HG21	1.90	0.52
1:P:87:U:H2'	1:P:88:G:OP2	2.09	0.52
2:A:462:TYR:CE2	2:A:511:HIS:NE2	2.77	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:288:TYR:N	2:C:288:TYR:CD1	2.78	0.52
2:C:422:THR:O	2:C:458:GLU:O	2.27	0.52
2:D:622:ASN:O	2:D:626:GLU:HG3	2.09	0.52
2:E:288:TYR:N	2:E:288:TYR:HD1	2.07	0.52
2:E:543:VAL:HG11	2:E:585:HIS:CE1	2.44	0.52
2:F:331:PRO:O	2:F:361:CYS:HB3	2.09	0.52
2:F:473:CYS:HB2	2:F:527:GLU:OE1	2.08	0.52
2:G:535:THR:HG22	2:G:632:ILE:HG12	1.92	0.52
2:H:578:ILE:HD11	2:H:595:VAL:CG2	2.38	0.52
1:N:86:G:H1'	2:F:547:ILE:HG21	1.92	0.52
2:A:529:PHE:CD2	2:A:534:PRO:HD3	2.45	0.52
2:B:490:PRO:HD3	2:B:509:MET:HE2	1.92	0.52
2:B:587:LEU:HD23	2:B:587:LEU:N	2.19	0.52
2:E:403:LEU:HG	2:E:407:MET:HE3	1.91	0.52
2:F:329:ILE:HD12	2:F:329:ILE:N	2.25	0.52
1:L:100:U:H4'	5:L:1230:HOH:O	2.08	0.52
1:M:93:U:C2	1:M:94:C:C6	2.97	0.52
2:A:254:TYR:CD2	2:A:255:HIS:N	2.77	0.52
2:A:500:GLU:OE1	2:A:500:GLU:N	2.41	0.52
2:C:542:VAL:HG12	2:C:543:VAL:N	2.25	0.52
2:C:578:ILE:HD11	2:C:595:VAL:HG22	1.92	0.52
2:C:622:ASN:O	2:C:626:GLU:HG3	2.10	0.52
2:E:557:GLU:O	2:E:560:GLN:HB3	2.09	0.52
2:G:284:LYS:HG3	2:G:407:MET:HG3	1.91	0.52
2:G:386:ILE:HG21	2:G:396:GLU:CG	2.39	0.52
2:H:278:GLU:HG2	2:H:518:MET:HE2	1.91	0.52
1:J:87:U:H3'	5:J:1019:HOH:O	2.10	0.52
1:P:88:G:P	5:P:1623:HOH:O	2.66	0.52
2:C:462:TYR:HD2	2:C:486:ASP:OD2	1.93	0.52
2:D:386:ILE:HG21	2:D:396:GLU:HG3	1.90	0.52
2:D:460:ALA:C	2:D:462:TYR:N	2.65	0.52
2:D:536:TRP:CZ3	2:D:635:ARG:HD3	2.44	0.52
2:H:282:ARG:HD3	2:H:292:GLU:OE1	2.10	0.52
2:H:557:GLU:O	2:H:560:GLN:HB3	2.09	0.52
2:C:258:GLU:O	2:D:325:ARG:NH2	2.42	0.52
2:E:569:VAL:HG22	2:E:570:LYS:N	2.25	0.52
2:G:366:PRO:O	2:G:367:SER:C	2.53	0.52
2:H:365:GLU:HA	2:H:365:GLU:OE1	2.10	0.52
2:A:569:VAL:HG22	2:A:570:LYS:N	2.25	0.52
2:C:536:TRP:CE3	2:C:635:ARG:HD3	2.44	0.52
2:C:544:ILE:HG12	2:C:594:LEU:HD12	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:597:GLY:O	2:C:601:VAL:HG23	2.08	0.52
2:D:365:GLU:H	2:D:377:ARG:HD3	1.73	0.52
2:D:543:VAL:HG11	2:D:585:HIS:CE1	2.44	0.52
2:E:244:HIS:CE1	2:E:374:MET:HE2	2.45	0.52
2:F:490:PRO:HD3	2:F:509:MET:HE2	1.91	0.52
2:F:536:TRP:CZ3	2:F:635:ARG:HD3	2.44	0.52
2:F:578:ILE:HD11	2:F:595:VAL:CG2	2.40	0.52
2:G:637:LEU:HD12	2:G:637:LEU:N	2.24	0.52
2:H:403:LEU:HG	2:H:407:MET:HE1	1.89	0.52
2:H:544:ILE:HG12	2:H:594:LEU:HD12	1.91	0.52
1:O:93:U:C2	1:O:94:C:C6	2.98	0.52
1:P:84:U:O2	2:H:549:ASP:OD1	2.28	0.52
2:A:543:VAL:HG11	2:A:585:HIS:CE1	2.45	0.52
2:B:636:SER:OG	2:B:638:LYS:HD3	2.10	0.52
2:C:606:VAL:O	2:C:618:SER:HA	2.10	0.52
2:C:620:ASP:HB3	2:C:623:GLU:CB	2.40	0.52
2:D:638:LYS:HB2	2:D:642:GLU:HG3	1.91	0.52
2:E:329:ILE:N	2:E:329:ILE:HD12	2.25	0.52
2:E:332:MET:HE2	4:E:6002:TSB:HN12	1.75	0.52
2:E:556:ASN:ND2	2:E:573:LEU:HD11	2.25	0.52
2:G:386:ILE:HG21	2:G:396:GLU:HG3	1.92	0.52
2:B:332:MET:HE2	4:B:3002:TSB:HN12	1.75	0.52
2:D:620:ASP:O	2:D:623:GLU:HB3	2.10	0.52
2:E:544:ILE:HG22	2:E:555:VAL:HG13	1.91	0.52
2:F:288:TYR:N	2:F:288:TYR:HD1	2.07	0.52
2:F:624:VAL:CG2	2:F:625:ILE:N	2.73	0.52
2:G:375:ARG:NH1	5:G:710:HOH:O	2.32	0.52
1:K:78:G:H2'	1:K:79:A:C8	2.42	0.51
2:D:544:ILE:HG12	2:D:594:LEU:HD12	1.91	0.51
2:D:636:SER:OG	2:D:638:LYS:HD3	2.10	0.51
2:E:610:THR:HG21	2:E:614:LYS:HB3	1.92	0.51
2:F:278:GLU:HG2	2:F:518:MET:HE2	1.92	0.51
2:F:628:LEU:O	2:F:632:ILE:HG13	2.10	0.51
2:G:443:VAL:O	2:G:447:GLU:HB2	2.09	0.51
2:H:488:SER:C	2:H:490:PRO:HD2	2.35	0.51
1:O:89:U:H2'	1:O:90:G:OP1	2.10	0.51
2:A:362:HIS:HE1	2:B:297:PHE:CG	2.28	0.51
2:A:403:LEU:HG	2:A:407:MET:HE3	1.92	0.51
2:B:304:TRP:CZ3	2:B:328:CYS:HB2	2.45	0.51
2:D:418:VAL:HG21	2:D:450:ILE:CG2	2.39	0.51
2:D:423:ARG:HD2	2:D:438:GLU:OE2	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:624:VAL:CG2	2:D:625:ILE:N	2.72	0.51
2:E:290:TYR:CE2	2:E:512:ARG:NH1	2.78	0.51
2:E:462:TYR:HD2	2:E:486:ASP:OD2	1.92	0.51
2:F:622:ASN:O	2:F:626:GLU:HG3	2.09	0.51
2:G:578:ILE:HD11	2:G:595:VAL:CG2	2.39	0.51
2:H:273:ILE:HG23	2:H:537:LEU:HD22	1.91	0.51
2:A:280:PHE:CE2	2:A:407:MET:HG2	2.44	0.51
2:A:559:THR:HG21	2:A:571:ALA:CB	2.19	0.51
2:C:278:GLU:CB	2:C:358:PHE:HE2	2.22	0.51
2:C:582:ILE:O	2:C:593:MET:HE1	2.10	0.51
2:D:285:LEU:HD22	2:D:290:TYR:CG	2.45	0.51
2:D:555:VAL:HB	2:D:573:LEU:HD21	1.91	0.51
2:E:304:TRP:CE3	2:E:328:CYS:HB2	2.45	0.51
2:G:620:ASP:O	2:G:624:VAL:HG13	2.10	0.51
2:B:557:GLU:O	2:B:560:GLN:HB3	2.10	0.51
2:D:514:ILE:HG22	2:D:515:LEU:HD23	1.92	0.51
2:E:296:PRO:HG2	2:F:265:PHE:CE1	2.45	0.51
2:E:418:VAL:HG21	2:E:450:ILE:CG2	2.40	0.51
2:E:482:THR:HG21	4:E:6002:TSB:H5'2	1.92	0.51
2:G:515:LEU:HD22	2:G:521:PHE:CE2	2.46	0.51
2:H:636:SER:OG	2:H:638:LYS:HD3	2.11	0.51
1:I:89:U:C6	2:A:583:ARG:CZ	2.93	0.51
2:A:331:PRO:O	2:A:361:CYS:HB3	2.09	0.51
2:A:339:GLN:HB3	2:B:257:GLN:OE1	2.11	0.51
2:A:422:THR:O	2:A:423:ARG:C	2.53	0.51
2:A:624:VAL:CG2	2:A:625:ILE:N	2.72	0.51
2:B:269:ASP:HB3	2:B:539:PRO:HB3	1.92	0.51
2:B:277:LEU:HB2	2:B:518:MET:HE1	1.92	0.51
2:B:423:ARG:HE	2:B:438:GLU:CD	2.19	0.51
2:C:442:ALA:O	2:C:446:GLU:HG2	2.11	0.51
2:C:580:PHE:HD1	2:C:580:PHE:C	2.18	0.51
2:E:547:ILE:HD12	2:E:597:GLY:CA	2.41	0.51
2:F:287:GLU:HB2	2:F:288:TYR:HD1	1.75	0.51
2:G:338:VAL:HG11	2:G:493:LEU:HB2	1.92	0.51
2:G:606:VAL:O	2:G:618:SER:HA	2.11	0.51
2:H:301:ARG:O	2:H:305:GLU:HG3	2.09	0.51
1:I:84:U:O2'	1:I:85:C:OP2	2.28	0.51
1:J:95:A:H2'	1:J:96:C:C6	2.46	0.51
1:K:85:C:C1'	2:C:577:LYS:HG2	2.40	0.51
1:P:89:U:O5'	5:P:1628:HOH:O	2.19	0.51
2:B:247:ILE:O	2:B:250:GLN:HB2	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:443:VAL:O	2:D:447:GLU:HB2	2.10	0.51
2:E:620:ASP:HB3	2:E:623:GLU:CB	2.41	0.51
2:F:620:ASP:O	2:F:623:GLU:HB3	2.11	0.51
1:M:80:U:H2'	1:M:81:C:H6	1.75	0.51
2:B:543:VAL:HG11	2:B:585:HIS:CE1	2.46	0.51
2:B:562:LEU:HD12	2:B:625:ILE:HD11	1.93	0.51
2:C:280:PHE:O	2:C:283:SER:HB3	2.09	0.51
2:C:297:PHE:CD1	2:D:362:HIS:HE1	2.28	0.51
2:C:462:TYR:HE2	2:C:511:HIS:NE2	2.09	0.51
2:C:578:ILE:HD11	2:C:595:VAL:CG2	2.41	0.51
2:E:348:TYR:CD1	2:E:348:TYR:C	2.89	0.51
2:G:622:ASN:O	2:G:626:GLU:HG3	2.10	0.51
2:H:251:LEU:O	2:H:253:LEU:HD23	2.11	0.51
2:A:421:SER:CA	2:A:455:GLN:HB3	2.40	0.51
2:B:452:PHE:CE2	2:B:454:TYR:HE1	2.20	0.51
2:C:304:TRP:CZ3	2:C:328:CYS:HB2	2.46	0.51
2:D:462:TYR:HD2	2:D:486:ASP:OD2	1.93	0.51
2:F:610:THR:HG21	2:F:614:LYS:HB3	1.93	0.51
2:A:423:ARG:O	2:A:423:ARG:HG2	2.10	0.51
2:C:290:TYR:CE2	2:C:512:ARG:NH1	2.79	0.51
2:D:526:THR:HG23	2:D:533:PHE:HZ	1.75	0.51
2:D:604:GLY:O	2:D:621:VAL:HG23	2.10	0.51
2:G:638:LYS:HB2	2:G:642:GLU:HG3	1.93	0.51
2:H:555:VAL:HB	2:H:573:LEU:HD21	1.92	0.51
2:A:637:LEU:HD12	2:A:637:LEU:N	2.23	0.51
2:B:274:PHE:CD1	2:B:274:PHE:C	2.89	0.51
2:B:580:PHE:HD1	2:B:580:PHE:C	2.19	0.51
2:C:624:VAL:CG2	2:C:625:ILE:N	2.74	0.51
2:C:636:SER:OG	2:C:638:LYS:HD3	2.11	0.51
2:D:452:PHE:HE2	2:D:454:TYR:CE1	2.22	0.51
2:D:547:ILE:HD12	2:D:597:GLY:CA	2.39	0.51
2:F:462:TYR:HE2	2:F:511:HIS:NE2	2.09	0.51
2:H:638:LYS:HB2	2:H:642:GLU:HG3	1.92	0.51
1:M:86:G:O6	2:E:599:LYS:CB	2.59	0.50
2:C:257:GLN:OE1	2:D:339:GLN:HB3	2.10	0.50
2:D:278:GLU:CB	2:D:358:PHE:HE2	2.24	0.50
2:D:348:TYR:CD1	2:D:348:TYR:C	2.88	0.50
2:E:351:LEU:HD22	2:E:387:PHE:O	2.10	0.50
2:E:578:ILE:HD11	2:E:595:VAL:HG22	1.93	0.50
2:F:301:ARG:O	2:F:305:GLU:HG3	2.10	0.50
2:H:620:ASP:O	2:H:624:VAL:HG13	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:78:G:N3	1:K:79:A:C8	2.78	0.50
2:A:488:SER:C	2:A:490:PRO:HD2	2.36	0.50
2:B:244:HIS:CE1	2:B:374:MET:HE2	2.45	0.50
2:D:580:PHE:HD1	2:D:580:PHE:C	2.17	0.50
2:E:249:LYS:HB3	2:E:249:LYS:HZ3	1.75	0.50
2:E:422:THR:O	2:E:423:ARG:C	2.54	0.50
2:E:443:VAL:O	2:E:447:GLU:HB2	2.11	0.50
2:H:411:PHE:HE2	2:H:521:PHE:HE1	1.56	0.50
2:H:422:THR:HB	2:H:438:GLU:OE2	2.12	0.50
2:H:603:SER:C	2:H:605:LYS:H	2.18	0.50
2:C:329:ILE:HD12	2:C:329:ILE:N	2.26	0.50
2:C:529:PHE:CD2	2:C:534:PRO:HD3	2.47	0.50
2:C:547:ILE:HG12	2:C:578:ILE:HD12	1.92	0.50
2:D:341:PHE:O	2:D:346:LYS:NZ	2.38	0.50
2:E:305:GLU:HA	2:E:310:TRP:HB3	1.94	0.50
2:E:460:ALA:C	2:E:462:TYR:H	2.19	0.50
2:E:550:SER:C	2:E:552:SER:H	2.18	0.50
2:G:636:SER:OG	2:G:638:LYS:HD3	2.11	0.50
1:J:105:C:H2'	1:O:69:G:O4'	2.12	0.50
1:K:104:C:H2'	1:K:105:C:C4'	2.42	0.50
2:B:335:PRO:HG2	2:B:336:GLY:H	1.76	0.50
2:C:557:GLU:O	2:C:560:GLN:HB3	2.11	0.50
2:D:578:ILE:HD11	2:D:595:VAL:HG22	1.93	0.50
2:E:556:ASN:HD21	2:E:573:LEU:HD11	1.77	0.50
2:F:281:VAL:HG12	2:F:285:LEU:CD1	2.42	0.50
2:F:462:TYR:CD2	2:F:486:ASP:OD2	2.64	0.50
2:G:309:HIS:HB3	2:G:317:MET:CE	2.36	0.50
2:G:500:GLU:C	2:G:502:ASN:N	2.69	0.50
2:H:331:PRO:O	2:H:361:CYS:HB3	2.12	0.50
1:P:78:G:O2'	1:P:79:A:H5'	2.12	0.50
2:A:411:PHE:HE2	2:A:521:PHE:CE1	2.17	0.50
2:A:418:VAL:HG21	2:A:450:ILE:HG21	1.93	0.50
2:A:500:GLU:C	2:A:502:ASN:N	2.70	0.50
2:B:455:GLN:O	2:B:456:LEU:C	2.53	0.50
2:C:441:LEU:HD12	2:C:464:PRO:HB2	1.93	0.50
2:E:280:PHE:O	2:E:283:SER:HB3	2.11	0.50
2:F:529:PHE:H	2:F:529:PHE:HD1	1.58	0.50
2:G:274:PHE:CD1	2:G:274:PHE:C	2.90	0.50
2:H:342:ASN:HD21	2:H:495:ALA:HA	1.75	0.50
2:H:548:THR:HG23	2:H:551:GLN:OE1	2.12	0.50
2:B:417:VAL:CG1	2:B:471:TYR:CE1	2.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:288:TYR:N	2:C:288:TYR:HD1	2.08	0.50
2:C:473:CYS:HB2	2:C:527:GLU:OE1	2.12	0.50
2:D:637:LEU:HD12	2:D:637:LEU:N	2.23	0.50
2:E:325:ARG:NH2	2:F:258:GLU:O	2.45	0.50
2:E:331:PRO:O	2:E:361:CYS:HB3	2.12	0.50
2:F:411:PHE:CE2	2:F:521:PHE:HZ	2.27	0.50
2:F:488:SER:C	2:F:490:PRO:HD2	2.36	0.50
2:F:544:ILE:HG12	2:F:594:LEU:HD12	1.92	0.50
2:G:556:ASN:ND2	2:G:573:LEU:HD11	2.27	0.50
2:H:452:PHE:HE2	2:H:454:TYR:CE1	2.23	0.50
1:K:85:C:C2	2:C:577:LYS:HG3	2.46	0.50
2:A:297:PHE:CG	2:B:362:HIS:CE1	3.00	0.50
2:A:329:ILE:N	2:A:329:ILE:CD1	2.75	0.50
2:B:243:ASP:O	2:B:247:ILE:HG13	2.12	0.50
2:B:277:LEU:CB	2:B:518:MET:HE1	2.42	0.50
2:B:362:HIS:HD2	5:B:232:HOH:O	1.94	0.50
2:B:600:GLU:HB3	2:B:605:LYS:O	2.11	0.50
2:C:434:TRP:O	2:C:438:GLU:HG3	2.11	0.50
2:C:588:ARG:C	2:C:589:ARG:HG2	2.36	0.50
2:D:397:VAL:O	2:D:401:ILE:HG12	2.12	0.50
2:G:550:SER:C	2:G:552:SER:H	2.19	0.50
1:I:89:U:C5	2:A:583:ARG:CZ	2.95	0.50
1:O:92:G:H2'	1:O:93:U:H6	1.76	0.50
2:B:285:LEU:HD22	2:B:290:TYR:CG	2.47	0.50
2:B:536:TRP:CE3	2:B:635:ARG:HD3	2.47	0.50
2:B:569:VAL:HG22	2:B:570:LYS:N	2.26	0.50
2:C:247:ILE:O	2:C:250:GLN:HB2	2.12	0.50
2:D:304:TRP:CE3	2:D:328:CYS:HB2	2.46	0.50
2:D:500:GLU:OE1	2:D:500:GLU:N	2.44	0.50
2:E:273:ILE:HG23	2:E:537:LEU:CD2	2.41	0.50
2:H:550:SER:C	2:H:552:SER:H	2.20	0.50
2:H:620:ASP:HB3	2:H:623:GLU:CB	2.42	0.50
1:L:77:U:O5'	1:L:77:U:H6	1.95	0.50
1:M:92:G:H2'	1:M:93:U:H6	1.77	0.50
2:B:578:ILE:HD11	2:B:595:VAL:HG22	1.94	0.50
2:C:445:LEU:HD21	2:C:466:ILE:HG12	1.94	0.50
2:C:551:GLN:HG2	2:C:597:GLY:C	2.37	0.50
2:E:403:LEU:HG	2:E:407:MET:HE1	1.94	0.50
2:E:500:GLU:C	2:E:502:ASN:N	2.70	0.50
2:E:536:TRP:O	2:E:568:ARG:NH2	2.45	0.50
2:F:624:VAL:HG23	2:F:625:ILE:N	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:278:GLU:HG2	2:B:518:MET:HE2	1.93	0.49
2:B:334:CYS:SG	2:B:511:HIS:HE1	2.34	0.49
2:C:298:MET:HE2	2:C:327:TYR:CD1	2.47	0.49
2:C:484:GLN:HB2	2:C:511:HIS:HB2	1.94	0.49
2:D:578:ILE:HD11	2:D:595:VAL:CG2	2.42	0.49
2:E:514:ILE:HG22	2:E:515:LEU:HD23	1.93	0.49
2:E:536:TRP:CE3	2:E:635:ARG:HD3	2.47	0.49
2:E:637:LEU:HD12	2:E:637:LEU:N	2.26	0.49
2:F:247:ILE:O	2:F:250:GLN:HB2	2.11	0.49
1:L:80:U:H2'	1:L:81:C:C6	2.47	0.49
2:B:393:ILE:HD11	2:B:508:VAL:HG11	1.94	0.49
2:B:403:LEU:HG	2:B:407:MET:HE3	1.93	0.49
2:F:483:VAL:HG22	2:F:512:ARG:HB2	1.94	0.49
2:G:280:PHE:CZ	2:G:407:MET:HG2	2.47	0.49
2:G:472:ASP:OD1	2:G:476:ARG:HB2	2.12	0.49
2:G:536:TRP:CZ3	2:G:635:ARG:HD3	2.47	0.49
2:G:620:ASP:HB3	2:G:623:GLU:CB	2.42	0.49
2:H:356:ALA:HA	2:H:384:ALA:HA	1.93	0.49
1:J:72:G:H2'	1:J:73:U:C6	2.47	0.49
1:M:101:G:O2'	1:M:102:C:H5'	2.12	0.49
1:N:79:A:C1'	2:E:502:ASN:OD1	2.57	0.49
2:C:278:GLU:HG2	2:C:518:MET:HE2	1.92	0.49
2:C:403:LEU:HG	2:C:407:MET:HE1	1.92	0.49
2:E:309:HIS:HB3	2:E:317:MET:CE	2.38	0.49
2:F:542:VAL:HG12	2:F:543:VAL:N	2.27	0.49
2:G:559:THR:HG21	2:G:571:ALA:CB	2.20	0.49
2:G:582:ILE:HA	2:G:593:MET:CE	2.42	0.49
2:H:621:VAL:HG12	2:H:622:ASN:N	2.27	0.49
1:K:93:U:H2'	1:K:94:C:C6	2.46	0.49
2:A:455:GLN:O	2:A:456:LEU:C	2.55	0.49
2:A:473:CYS:HB2	2:A:527:GLU:OE1	2.12	0.49
2:C:393:ILE:HD11	2:C:508:VAL:HG11	1.95	0.49
2:C:480:CYS:O	2:C:516:GLY:HA3	2.12	0.49
2:D:557:GLU:O	2:D:560:GLN:HB3	2.12	0.49
2:E:624:VAL:HG23	2:E:625:ILE:N	2.27	0.49
2:F:402:ARG:NH1	2:F:402:ARG:CB	2.75	0.49
2:G:291:GLN:NE2	2:G:353:LEU:HD21	2.26	0.49
2:G:551:GLN:HG2	2:G:597:GLY:C	2.36	0.49
2:H:254:TYR:CZ	2:H:373:LEU:HD21	2.48	0.49
2:A:609:ARG:HG3	2:A:609:ARG:O	2.11	0.49
2:C:587:LEU:HD23	2:C:587:LEU:N	2.20	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:637:LEU:HD12	2:C:637:LEU:N	2.21	0.49
2:D:280:PHE:O	2:D:283:SER:HB3	2.12	0.49
2:D:619:MET:HG3	2:D:624:VAL:HG12	1.94	0.49
2:F:249:LYS:HB3	2:F:249:LYS:HZ3	1.76	0.49
2:G:526:THR:HG23	2:G:533:PHE:HZ	1.77	0.49
2:G:539:PRO:O	2:G:568:ARG:HD2	2.13	0.49
2:H:280:PHE:CZ	2:H:407:MET:HG2	2.48	0.49
2:H:291:GLN:NE2	2:H:353:LEU:HD21	2.28	0.49
1:N:84:U:O2'	1:N:85:C:OP2	2.30	0.49
2:A:305:GLU:HA	2:A:310:TRP:HB3	1.94	0.49
2:A:443:VAL:O	2:A:447:GLU:HB2	2.13	0.49
2:C:332:MET:HE2	4:C:4002:TSB:HN12	1.77	0.49
2:C:536:TRP:CZ3	2:C:635:ARG:HD3	2.47	0.49
2:D:423:ARG:HA	2:D:434:TRP:CZ3	2.47	0.49
2:D:620:ASP:HB3	2:D:623:GLU:CB	2.42	0.49
2:E:590:VAL:O	2:E:611:ARG:HB3	2.13	0.49
2:G:418:VAL:HG21	2:G:450:ILE:CG2	2.43	0.49
2:G:427:ARG:HB2	2:G:434:TRP:CZ2	2.47	0.49
2:G:474:LEU:O	2:G:475:ASP:C	2.55	0.49
2:H:386:ILE:HG21	2:H:396:GLU:CG	2.43	0.49
2:H:400:CYS:O	2:H:404:VAL:HG23	2.13	0.49
2:H:551:GLN:HG2	2:H:597:GLY:C	2.37	0.49
1:J:84:U:O2'	1:J:85:C:OP2	2.31	0.49
2:B:415:LYS:HG2	2:B:471:TYR:CD1	2.48	0.49
2:B:603:SER:C	2:B:605:LYS:H	2.20	0.49
2:C:243:ASP:O	2:C:247:ILE:HG13	2.13	0.49
2:D:550:SER:C	2:D:552:SER:H	2.20	0.49
2:D:600:GLU:HB3	2:D:605:LYS:O	2.12	0.49
2:E:254:TYR:HE2	2:E:256:MET:HE2	1.77	0.49
2:F:416:ILE:O	2:F:416:ILE:HG22	2.10	0.49
2:F:578:ILE:HD11	2:F:595:VAL:HG22	1.93	0.49
2:H:500:GLU:C	2:H:502:ASN:N	2.71	0.49
2:B:329:ILE:N	2:B:329:ILE:HD12	2.27	0.49
2:B:590:VAL:O	2:B:611:ARG:HB3	2.12	0.49
2:C:298:MET:HG2	2:D:263:MET:CE	2.41	0.49
2:F:455:GLN:O	2:F:456:LEU:C	2.56	0.49
2:F:480:CYS:O	2:F:516:GLY:HA3	2.13	0.49
2:F:588:ARG:C	2:F:589:ARG:HG2	2.37	0.49
2:F:620:ASP:HB3	2:F:623:GLU:CB	2.43	0.49
2:H:323:GLU:H	2:H:323:GLU:CD	2.19	0.49
2:H:624:VAL:HG23	2:H:625:ILE:N	2.26	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:105:C:O2	1:K:105:C:C2'	2.60	0.49
1:L:97:C:C5'	1:L:98:A:OP1	2.60	0.49
1:M:69:G:OP2	1:P:105:C:O3'	2.31	0.49
1:M:69:G:OP2	1:P:105:C:C2'	2.61	0.49
2:A:297:PHE:CG	2:B:362:HIS:HE1	2.31	0.49
2:B:443:VAL:O	2:B:447:GLU:HB2	2.11	0.49
2:C:443:VAL:O	2:C:447:GLU:HB2	2.13	0.49
2:C:619:MET:HG3	2:C:624:VAL:HG12	1.93	0.49
2:D:624:VAL:HG23	2:D:625:ILE:N	2.27	0.49
2:E:300:ASP:OD2	2:E:302:VAL:HB	2.13	0.49
2:E:609:ARG:O	2:E:609:ARG:HG3	2.13	0.49
2:F:309:HIS:HB3	2:F:317:MET:CE	2.35	0.49
2:G:621:VAL:HG12	2:G:622:ASN:N	2.28	0.49
2:H:274:PHE:C	2:H:274:PHE:CD1	2.90	0.49
2:H:610:THR:HG21	2:H:614:LYS:HB3	1.94	0.49
1:M:88:G:H8	5:E:6013:HOH:O	1.96	0.49
2:B:588:ARG:C	2:B:589:ARG:HG2	2.38	0.49
2:B:624:VAL:CG2	2:B:625:ILE:N	2.75	0.49
2:C:304:TRP:CE3	2:C:328:CYS:HB2	2.48	0.49
2:C:624:VAL:O	2:C:628:LEU:HG	2.12	0.49
2:E:460:ALA:C	2:E:462:TYR:N	2.68	0.49
2:F:282:ARG:HD3	2:F:292:GLU:OE1	2.13	0.49
2:F:332:MET:HE2	4:F:7002:TSB:HN12	1.78	0.49
1:I:86:G:P	1:I:86:G:O4'	2.71	0.48
2:A:277:LEU:O	2:A:280:PHE:HB3	2.12	0.48
2:A:421:SER:HB3	2:A:458:GLU:CB	2.30	0.48
2:A:551:GLN:HG2	2:A:597:GLY:C	2.38	0.48
2:C:514:ILE:HG22	2:C:515:LEU:HD23	1.94	0.48
2:C:529:PHE:H	2:C:529:PHE:HD1	1.58	0.48
2:D:597:GLY:O	2:D:601:VAL:HG23	2.12	0.48
2:G:460:ALA:C	2:G:462:TYR:N	2.65	0.48
2:G:555:VAL:HB	2:G:573:LEU:HD21	1.94	0.48
2:H:330:LYS:HA	2:H:331:PRO:HD3	1.60	0.48
2:A:403:LEU:O	2:A:407:MET:HB2	2.12	0.48
2:B:547:ILE:HD12	2:B:597:GLY:CA	2.43	0.48
2:C:583:ARG:O	2:C:587:LEU:CD2	2.60	0.48
2:D:274:PHE:CD1	2:D:274:PHE:C	2.91	0.48
2:D:603:SER:C	2:D:605:LYS:H	2.21	0.48
2:D:609:ARG:HG3	2:D:609:ARG:O	2.13	0.48
2:E:500:GLU:OE1	2:E:500:GLU:N	2.43	0.48
2:F:500:GLU:OE1	2:F:500:GLU:N	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:512:ARG:NE	2:F:512:ARG:O	2.45	0.48
1:M:69:G:P	1:P:105:C:HO3'	2.31	0.48
1:O:87:U:C2'	1:O:88:G:OP2	2.60	0.48
2:A:263:MET:CE	2:B:298:MET:HG2	2.38	0.48
2:A:624:VAL:HG23	2:A:625:ILE:N	2.26	0.48
2:C:462:TYR:CE2	2:C:511:HIS:NE2	2.81	0.48
2:C:545:MET:SD	2:C:582:ILE:HD13	2.52	0.48
2:E:285:LEU:HD22	2:E:290:TYR:CG	2.49	0.48
2:E:325:ARG:HG3	2:E:325:ARG:NH1	2.27	0.48
2:E:603:SER:C	2:E:605:LYS:H	2.19	0.48
2:G:282:ARG:HD3	2:G:292:GLU:OE1	2.13	0.48
2:G:351:LEU:HD22	2:G:387:PHE:O	2.13	0.48
2:G:422:THR:O	2:G:423:ARG:C	2.56	0.48
1:K:78:G:O2'	1:K:79:A:H5'	2.12	0.48
2:B:303:LEU:HD11	2:B:339:GLN:HE21	1.78	0.48
2:B:422:THR:O	2:B:458:GLU:O	2.31	0.48
2:B:610:THR:HG22	2:B:614:LYS:O	2.13	0.48
2:D:462:TYR:HE2	2:D:511:HIS:CE1	2.31	0.48
2:E:278:GLU:CB	2:E:358:PHE:HE2	2.26	0.48
2:E:427:ARG:HB2	2:E:434:TRP:CZ2	2.49	0.48
2:E:597:GLY:O	2:E:601:VAL:HG23	2.14	0.48
2:F:491:SER:O	2:F:494:SER:N	2.41	0.48
2:G:251:LEU:O	2:G:253:LEU:HD23	2.14	0.48
2:G:462:TYR:HD2	2:G:486:ASP:OD2	1.97	0.48
2:G:500:GLU:OE1	2:G:500:GLU:N	2.44	0.48
2:H:253:LEU:HD23	2:H:253:LEU:N	2.29	0.48
2:H:351:LEU:HD22	2:H:387:PHE:O	2.13	0.48
2:H:386:ILE:HG21	2:H:396:GLU:HG3	1.95	0.48
2:H:423:ARG:O	2:H:423:ARG:HG2	2.13	0.48
1:N:89:U:O2'	2:F:580:PHE:HB2	2.13	0.48
1:P:86:G:C2'	5:P:1613:HOH:O	2.61	0.48
2:A:288:TYR:CZ	2:A:403:LEU:HD13	2.48	0.48
2:A:620:ASP:HB3	2:A:623:GLU:CB	2.43	0.48
2:B:365:GLU:OE1	2:B:365:GLU:HA	2.13	0.48
2:B:610:THR:HG21	2:B:614:LYS:HB3	1.96	0.48
2:C:320:THR:HG22	2:C:321:SER:N	2.22	0.48
2:C:628:LEU:O	2:C:632:ILE:HG13	2.13	0.48
2:D:331:PRO:O	2:D:361:CYS:HB3	2.14	0.48
2:D:519:GLU:H	2:D:519:GLU:CD	2.20	0.48
2:D:551:GLN:HG2	2:D:597:GLY:C	2.38	0.48
2:E:636:SER:OG	2:E:638:LYS:HD3	2.11	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:480:CYS:O	2:H:516:GLY:HA3	2.13	0.48
1:J:93:U:H2'	1:J:94:C:H6	1.79	0.48
1:L:88:G:H1'	2:D:583:ARG:HG2	1.96	0.48
1:N:77:U:O5'	1:N:77:U:H6	1.96	0.48
1:N:84:U:H4'	1:N:85:C:H5'	1.96	0.48
2:A:319:THR:H	2:B:320:THR:CG2	2.26	0.48
2:A:597:GLY:O	2:A:601:VAL:HG23	2.13	0.48
2:A:610:THR:CG2	2:A:614:LYS:HB3	2.42	0.48
2:B:320:THR:HG22	2:B:321:SER:N	2.25	0.48
2:B:403:LEU:O	2:B:407:MET:HB2	2.14	0.48
2:B:500:GLU:C	2:B:502:ASN:N	2.70	0.48
2:B:536:TRP:CZ3	2:B:635:ARG:HD3	2.48	0.48
2:C:263:MET:CE	2:D:298:MET:HG2	2.39	0.48
2:C:512:ARG:O	2:C:512:ARG:NE	2.44	0.48
2:E:265:PHE:CZ	2:F:296:PRO:HG2	2.48	0.48
2:E:362:HIS:CE1	2:F:297:PHE:CG	3.02	0.48
2:F:305:GLU:HA	2:F:310:TRP:HB3	1.94	0.48
2:F:411:PHE:CD2	2:F:521:PHE:CZ	3.01	0.48
2:F:418:VAL:HG21	2:F:450:ILE:CG2	2.44	0.48
2:G:280:PHE:O	2:G:283:SER:HB3	2.12	0.48
2:H:244:HIS:ND1	2:H:374:MET:HE2	2.29	0.48
1:M:80:U:OP2	2:F:349:ARG:NH2	2.47	0.48
2:A:441:LEU:O	2:A:444:ALA:HB3	2.13	0.48
2:A:539:PRO:O	2:A:568:ARG:HD2	2.12	0.48
2:B:339:GLN:OE1	2:B:339:GLN:CA	2.60	0.48
2:C:400:CYS:SG	2:C:512:ARG:HG3	2.54	0.48
2:C:452:PHE:HE2	2:C:454:TYR:CE1	2.20	0.48
2:C:465:LYS:HG3	2:C:484:GLN:CG	2.42	0.48
2:C:604:GLY:O	2:C:621:VAL:HG23	2.13	0.48
2:D:254:TYR:HA	2:D:267:HIS:ND1	2.27	0.48
2:D:303:LEU:HD11	2:D:339:GLN:NE2	2.28	0.48
2:D:582:ILE:HA	2:D:593:MET:CE	2.44	0.48
2:E:555:VAL:HB	2:E:573:LEU:HD21	1.95	0.48
2:F:545:MET:SD	2:F:582:ILE:HD13	2.54	0.48
2:G:362:HIS:CD2	5:G:732:HOH:O	2.56	0.48
2:H:546:ASN:OD1	2:H:573:LEU:HA	2.13	0.48
2:H:569:VAL:HG22	2:H:570:LYS:N	2.29	0.48
2:H:603:SER:C	2:H:605:LYS:N	2.71	0.48
1:I:88:G:C8	5:I:112:HOH:O	2.48	0.48
1:I:95:A:H2'	1:I:96:C:O4'	2.14	0.48
1:K:98:A:O2'	2:D:344:GLY:N	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:582:ILE:O	2:A:593:MET:HE1	2.14	0.48
2:B:389:THR:HG1	2:B:392:GLN:HG3	1.74	0.48
2:C:254:TYR:HA	2:C:267:HIS:ND1	2.29	0.48
2:C:273:ILE:HG23	2:C:537:LEU:HD22	1.96	0.48
2:E:554:TYR:CE1	2:E:558:LEU:HD11	2.49	0.48
2:E:578:ILE:HD11	2:E:595:VAL:CG2	2.43	0.48
2:F:386:ILE:HG21	2:F:396:GLU:HG3	1.95	0.48
2:G:556:ASN:HD21	2:G:573:LEU:HD11	1.79	0.48
2:G:572:ASP:OD2	2:G:574:ARG:NH2	2.37	0.48
1:I:72:G:H2'	1:I:73:U:H6	1.78	0.48
1:K:79:A:N1	1:K:94:C:N3	2.62	0.48
1:M:69:G:C8	1:P:105:C:O2	2.67	0.48
1:O:95:A:H2'	1:O:96:C:O4'	2.14	0.48
2:A:335:PRO:HG2	2:A:336:GLY:H	1.79	0.48
2:A:386:ILE:O	2:A:509:MET:HA	2.13	0.48
2:C:253:LEU:O	2:C:270:GLY:HA3	2.14	0.48
2:C:421:SER:CA	2:C:455:GLN:HB3	2.43	0.48
2:E:515:LEU:HD22	2:E:521:PHE:CE2	2.48	0.48
2:E:536:TRP:CD1	2:E:635:ARG:HA	2.49	0.48
2:G:397:VAL:O	2:G:400:CYS:HB2	2.13	0.48
2:G:423:ARG:HD2	2:G:438:GLU:OE2	2.14	0.48
2:H:254:TYR:CD2	2:H:255:HIS:N	2.82	0.48
2:H:332:MET:HE1	2:H:462:TYR:OH	2.12	0.48
1:L:75:U:O2'	1:L:89:U:O4	2.28	0.48
1:L:87:U:H1'	2:D:582:ILE:HG13	1.96	0.48
2:C:437:ALA:HA	2:C:485:LEU:HD23	1.96	0.48
2:E:411:PHE:CD2	2:E:525:LEU:CD2	2.85	0.48
2:E:423:ARG:HD2	2:E:438:GLU:OE2	2.14	0.48
2:G:332:MET:HE1	2:G:462:TYR:OH	2.13	0.48
2:H:526:THR:O	2:H:530:ALA:HA	2.14	0.48
2:H:545:MET:HE3	2:H:593:MET:HB3	1.96	0.48
1:I:79:A:H1'	2:B:502:ASN:OD1	2.13	0.47
1:J:103:G:P	5:J:1025:HOH:O	2.71	0.47
2:A:634:SER:O	2:A:635:ARG:C	2.57	0.47
2:C:309:HIS:HB3	2:C:317:MET:CE	2.39	0.47
2:D:323:GLU:CD	2:D:323:GLU:H	2.22	0.47
2:D:403:LEU:HG	2:D:407:MET:HE3	1.96	0.47
2:D:423:ARG:HE	2:D:438:GLU:CD	2.22	0.47
2:G:544:ILE:HG12	2:G:594:LEU:HD12	1.96	0.47
2:H:430:SER:O	2:H:431:ASP:C	2.55	0.47
2:H:538:ALA:O	2:H:539:PRO:C	2.57	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:92:G:H2'	1:J:93:U:H6	1.79	0.47
2:B:282:ARG:HD3	2:B:292:GLU:OE1	2.14	0.47
2:B:331:PRO:O	2:B:361:CYS:HB3	2.14	0.47
2:B:473:CYS:HB2	2:B:527:GLU:OE1	2.14	0.47
2:B:555:VAL:HB	2:B:573:LEU:HD21	1.95	0.47
2:C:291:GLN:NE2	2:C:353:LEU:HD21	2.29	0.47
2:C:386:ILE:HG21	2:C:396:GLU:CG	2.44	0.47
2:E:287:GLU:HB2	2:E:288:TYR:HD1	1.79	0.47
2:F:278:GLU:CB	2:F:358:PHE:HE2	2.26	0.47
2:F:371:HIS:O	2:F:372:GLY:C	2.56	0.47
2:F:547:ILE:HG12	2:F:578:ILE:HD12	1.96	0.47
2:F:637:LEU:HD12	2:F:637:LEU:N	2.26	0.47
2:G:298:MET:HE2	2:G:327:TYR:CD1	2.49	0.47
2:H:484:GLN:HB2	2:H:511:HIS:HB2	1.96	0.47
1:M:78:G:O2'	1:M:79:A:H5'	2.14	0.47
2:A:356:ALA:HA	2:A:384:ALA:CB	2.45	0.47
2:A:356:ALA:CB	2:A:384:ALA:HB2	2.44	0.47
2:C:500:GLU:C	2:C:502:ASN:N	2.72	0.47
2:C:609:ARG:HD3	2:C:610:THR:O	2.14	0.47
2:F:394:ARG:O	2:F:395:ASP:C	2.57	0.47
2:F:636:SER:OG	2:F:638:LYS:HD3	2.14	0.47
2:H:288:TYR:N	2:H:288:TYR:CD1	2.82	0.47
2:H:403:LEU:HG	2:H:407:MET:HE3	1.93	0.47
2:H:536:TRP:CZ3	2:H:635:ARG:HD3	2.49	0.47
1:K:85:C:O2	1:K:90:G:N7	2.47	0.47
1:K:98:A:N3	1:K:98:A:H2'	2.30	0.47
2:A:556:ASN:ND2	2:A:573:LEU:HD11	2.28	0.47
2:B:621:VAL:HG12	2:B:622:ASN:N	2.29	0.47
2:C:624:VAL:HG23	2:C:625:ILE:N	2.29	0.47
2:E:301:ARG:O	2:E:305:GLU:HG3	2.14	0.47
2:G:253:LEU:O	2:G:270:GLY:HA3	2.14	0.47
2:G:557:GLU:O	2:G:560:GLN:HB3	2.14	0.47
2:G:569:VAL:O	2:G:570:LYS:HG2	2.14	0.47
2:H:423:ARG:HD2	2:H:438:GLU:OE2	2.15	0.47
2:H:500:GLU:OE1	2:H:500:GLU:N	2.45	0.47
1:J:88:G:H1'	2:B:583:ARG:CG	2.43	0.47
2:A:296:PRO:HG2	2:B:265:PHE:CE1	2.50	0.47
2:B:423:ARG:HA	2:B:434:TRP:CZ3	2.49	0.47
2:B:578:ILE:HD11	2:B:595:VAL:CG2	2.45	0.47
2:C:411:PHE:CD2	2:C:521:PHE:CZ	3.03	0.47
2:C:423:ARG:HE	2:C:438:GLU:CD	2.22	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:544:ILE:HG12	2:E:594:LEU:HD12	1.96	0.47
2:E:621:VAL:HG12	2:E:622:ASN:N	2.29	0.47
2:F:274:PHE:CD1	2:F:274:PHE:C	2.91	0.47
2:F:281:VAL:O	2:F:282:ARG:C	2.56	0.47
2:F:393:ILE:HD11	2:F:508:VAL:HG11	1.97	0.47
2:G:536:TRP:CE3	2:G:635:ARG:HD3	2.49	0.47
2:H:556:ASN:ND2	2:H:573:LEU:HD11	2.29	0.47
1:I:95:A:H2'	1:I:96:C:C6	2.48	0.47
1:K:78:G:H21	2:D:502:ASN:CG	2.20	0.47
1:N:78:G:O2'	1:N:79:A:H5'	2.15	0.47
2:A:273:ILE:HG22	2:A:522:ILE:HD11	1.96	0.47
2:A:287:GLU:HB2	2:A:288:TYR:HD1	1.80	0.47
2:C:281:VAL:HG12	2:C:285:LEU:CD1	2.45	0.47
2:C:287:GLU:HB2	2:C:288:TYR:HD1	1.79	0.47
2:C:526:THR:HG23	2:C:533:PHE:CZ	2.46	0.47
2:D:281:VAL:HG12	2:D:285:LEU:CD1	2.45	0.47
2:E:569:VAL:O	2:E:570:LYS:HG2	2.14	0.47
2:H:421:SER:HB3	2:H:458:GLU:CB	2.35	0.47
2:H:539:PRO:O	2:H:568:ARG:HD2	2.15	0.47
1:K:88:G:H5''	1:K:89:U:P	2.55	0.47
2:A:247:ILE:O	2:A:250:GLN:HB2	2.14	0.47
2:A:562:LEU:HD12	2:A:625:ILE:HD11	1.96	0.47
2:A:619:MET:HG3	2:A:624:VAL:HG12	1.96	0.47
2:B:278:GLU:HB3	2:B:358:PHE:CE2	2.49	0.47
2:B:280:PHE:O	2:B:283:SER:HB3	2.15	0.47
2:B:323:GLU:H	2:B:323:GLU:CD	2.22	0.47
2:B:375:ARG:NH2	5:B:208:HOH:O	2.26	0.47
2:B:474:LEU:O	2:B:475:ASP:C	2.57	0.47
2:B:550:SER:O	2:B:552:SER:N	2.47	0.47
2:B:554:TYR:CE1	2:B:558:LEU:HD11	2.50	0.47
2:B:609:ARG:HD3	2:B:610:THR:O	2.15	0.47
2:C:418:VAL:HG21	2:C:450:ILE:HG21	1.97	0.47
2:C:460:ALA:C	2:C:462:TYR:N	2.72	0.47
2:C:488:SER:C	2:C:490:PRO:HD2	2.40	0.47
2:C:490:PRO:HD3	2:C:509:MET:HE2	1.95	0.47
2:C:603:SER:C	2:C:605:LYS:H	2.21	0.47
2:D:397:VAL:O	2:D:400:CYS:HB2	2.15	0.47
2:D:445:LEU:HD21	2:D:466:ILE:HG12	1.97	0.47
2:D:580:PHE:HE1	5:D:5014:HOH:O	1.96	0.47
2:E:473:CYS:HB2	2:E:527:GLU:OE1	2.15	0.47
2:E:588:ARG:C	2:E:589:ARG:HG2	2.40	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:243:ASP:O	2:F:247:ILE:HG13	2.15	0.47
2:F:298:MET:CE	2:F:327:TYR:HB2	2.45	0.47
2:F:443:VAL:O	2:F:447:GLU:HB2	2.15	0.47
2:F:624:VAL:O	2:F:628:LEU:HG	2.14	0.47
2:G:246:LYS:O	2:G:250:GLN:HG2	2.15	0.47
2:G:423:ARG:HA	2:G:434:TRP:CZ3	2.50	0.47
2:G:543:VAL:HG11	2:G:585:HIS:CE1	2.50	0.47
1:K:79:A:C2	1:K:94:C:O2	2.68	0.47
2:A:386:ILE:HG21	2:A:396:GLU:HG3	1.95	0.47
2:A:548:THR:HG23	2:A:551:GLN:OE1	2.15	0.47
2:B:301:ARG:O	2:B:305:GLU:HG3	2.14	0.47
2:B:411:PHE:CE2	2:B:521:PHE:CZ	3.03	0.47
2:C:422:THR:O	2:C:423:ARG:C	2.57	0.47
2:D:422:THR:O	2:D:423:ARG:C	2.58	0.47
2:E:274:PHE:C	2:E:274:PHE:HD1	2.22	0.47
2:E:421:SER:CA	2:E:455:GLN:HB3	2.43	0.47
2:E:436:ARG:HG2	2:E:440:ASP:OD2	2.15	0.47
2:G:278:GLU:OE2	2:G:358:PHE:HD2	1.98	0.47
2:G:403:LEU:HG	2:G:407:MET:HE1	1.93	0.47
2:H:288:TYR:N	2:H:288:TYR:HD1	2.12	0.47
1:N:80:U:H2'	1:N:81:C:C6	2.50	0.47
1:N:86:G:C4	2:F:547:ILE:CD1	2.97	0.47
2:A:265:PHE:CE1	2:B:296:PRO:HG2	2.50	0.47
2:A:281:VAL:HG12	2:A:285:LEU:CD1	2.45	0.47
2:B:423:ARG:NE	2:B:438:GLU:OE1	2.48	0.47
2:C:562:LEU:HD12	2:C:625:ILE:HD11	1.96	0.47
2:E:365:GLU:HA	2:E:365:GLU:OE1	2.15	0.47
2:E:519:GLU:CD	2:E:519:GLU:H	2.22	0.47
2:F:303:LEU:HD11	2:F:339:GLN:NE2	2.27	0.47
2:F:543:VAL:HG11	2:F:585:HIS:ND1	2.30	0.47
2:G:421:SER:OG	2:G:467:GLU:OE2	2.33	0.47
2:G:600:GLU:HB3	2:G:605:LYS:O	2.14	0.47
2:H:266:TRP:CH2	2:H:522:ILE:HD12	2.50	0.47
2:H:416:ILE:O	2:H:416:ILE:HG22	2.14	0.47
2:H:589:ARG:O	2:H:611:ARG:HD3	2.15	0.47
2:H:624:VAL:O	2:H:628:LEU:HG	2.15	0.47
1:L:86:G:C4	2:D:547:ILE:HD13	2.49	0.47
1:P:84:U:O2'	1:P:85:C:OP2	2.29	0.47
2:B:400:CYS:O	2:B:401:ILE:C	2.58	0.47
2:C:501:ASP:O	2:C:503:GLU:HG2	2.15	0.47
2:E:323:GLU:H	2:E:323:GLU:CD	2.23	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:332:MET:HE1	2:E:462:TYR:OH	2.15	0.47
2:E:526:THR:HG23	2:E:533:PHE:HZ	1.79	0.47
2:E:583:ARG:O	2:E:587:LEU:CD2	2.63	0.47
2:F:452:PHE:HE2	2:F:454:TYR:CE1	2.24	0.47
2:G:332:MET:HE2	4:G:8002:TSB:HN12	1.80	0.47
2:G:460:ALA:HB1	2:G:462:TYR:CE1	2.50	0.47
2:G:562:LEU:HB3	2:G:569:VAL:HG11	1.97	0.47
2:H:484:GLN:HE22	4:H:9002:TSB:HN8	1.62	0.47
2:H:526:THR:HG23	2:H:533:PHE:HZ	1.80	0.47
2:H:621:VAL:O	2:H:622:ASN:C	2.57	0.47
2:H:628:LEU:O	2:H:632:ILE:HG13	2.14	0.47
1:K:89:U:O2	1:K:89:U:H2'	2.15	0.46
1:M:93:U:N3	1:M:94:C:C5	2.82	0.46
1:O:84:U:H3	2:G:549:ASP:CG	2.23	0.46
2:B:536:TRP:O	2:B:568:ARG:NH2	2.48	0.46
2:B:624:VAL:O	2:B:628:LEU:HG	2.14	0.46
2:C:335:PRO:HG2	2:C:336:GLY:H	1.79	0.46
2:C:529:PHE:CD1	2:C:529:PHE:N	2.78	0.46
2:D:624:VAL:O	2:D:628:LEU:HG	2.15	0.46
2:E:277:LEU:HB2	2:E:518:MET:HE1	1.97	0.46
2:E:281:VAL:HG12	2:E:285:LEU:CD1	2.43	0.46
2:E:534:PRO:O	2:E:535:THR:C	2.58	0.46
2:E:536:TRP:CZ3	2:E:635:ARG:HD3	2.50	0.46
2:E:547:ILE:HG12	2:E:578:ILE:HD12	1.97	0.46
2:E:619:MET:HG3	2:E:624:VAL:HG12	1.97	0.46
2:F:528:GLU:O	2:F:530:ALA:N	2.48	0.46
2:F:583:ARG:O	2:F:587:LEU:CD2	2.61	0.46
2:G:356:ALA:HA	2:G:384:ALA:HA	1.97	0.46
2:G:421:SER:CA	2:G:455:GLN:HB3	2.44	0.46
2:G:531:GLY:O	2:G:589:ARG:HG3	2.16	0.46
2:G:597:GLY:O	2:G:601:VAL:HG23	2.15	0.46
1:N:95:A:H2'	1:N:96:C:O4'	2.15	0.46
2:A:397:VAL:O	2:A:400:CYS:HB2	2.14	0.46
2:A:600:GLU:HB3	2:A:605:LYS:O	2.15	0.46
2:A:621:VAL:O	2:A:622:ASN:C	2.57	0.46
2:B:281:VAL:HG12	2:B:285:LEU:CD1	2.44	0.46
2:B:603:SER:C	2:B:605:LYS:N	2.72	0.46
2:D:559:THR:HG21	2:D:571:ALA:CB	2.21	0.46
2:F:280:PHE:CZ	2:F:407:MET:HG2	2.50	0.46
2:F:298:MET:HE2	2:F:327:TYR:CD1	2.50	0.46
2:F:386:ILE:HG21	2:F:396:GLU:CG	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:603:SER:C	2:F:605:LYS:H	2.23	0.46
2:H:545:MET:CE	2:H:593:MET:HB3	2.45	0.46
1:M:71:C:N3	1:M:72:G:N7	2.64	0.46
1:O:93:U:H2'	1:O:94:C:H6	1.80	0.46
2:A:608:VAL:HG21	2:A:619:MET:HE3	1.97	0.46
2:B:386:ILE:HG21	2:B:396:GLU:CG	2.45	0.46
2:E:400:CYS:SG	2:E:512:ARG:HG3	2.55	0.46
2:F:403:LEU:HG	2:F:407:MET:HE1	1.97	0.46
2:A:449:ASN:ND2	2:H:452:PHE:CE2	2.83	0.46
2:B:291:GLN:NE2	2:B:353:LEU:HD21	2.30	0.46
2:B:482:THR:O	2:B:512:ARG:HA	2.16	0.46
2:B:515:LEU:HD22	2:B:521:PHE:CE2	2.50	0.46
2:C:501:ASP:CG	2:C:503:GLU:HB2	2.40	0.46
2:C:519:GLU:H	2:C:519:GLU:CD	2.22	0.46
2:C:610:THR:HG21	2:C:614:LYS:HB3	1.97	0.46
2:D:305:GLU:HA	2:D:310:TRP:HB3	1.97	0.46
2:D:365:GLU:OE1	2:D:365:GLU:HA	2.15	0.46
2:F:285:LEU:HD21	2:F:512:ARG:HH12	1.80	0.46
2:H:287:GLU:HB2	2:H:288:TYR:HD1	1.79	0.46
1:J:89:U:H5	2:B:583:ARG:CZ	2.27	0.46
1:L:95:A:H2'	1:L:96:C:C6	2.50	0.46
2:B:501:ASP:O	2:B:503:GLU:HG2	2.16	0.46
2:B:536:TRP:CD1	2:B:635:ARG:HA	2.50	0.46
2:B:547:ILE:HG12	2:B:578:ILE:HD12	1.97	0.46
2:C:305:GLU:HA	2:C:310:TRP:HB3	1.98	0.46
2:C:386:ILE:HG21	2:C:396:GLU:HG3	1.96	0.46
2:D:621:VAL:HG12	2:D:622:ASN:N	2.31	0.46
2:E:423:ARG:HE	2:E:438:GLU:CD	2.23	0.46
2:F:600:GLU:HB3	2:F:605:LYS:O	2.15	0.46
2:H:394:ARG:O	2:H:395:ASP:C	2.59	0.46
1:I:95:A:H1'	2:B:502:ASN:HD21	1.81	0.46
1:M:93:U:C2	1:M:94:C:C5	3.03	0.46
2:A:371:HIS:O	2:A:372:GLY:C	2.57	0.46
2:A:545:MET:SD	2:A:582:ILE:HD13	2.56	0.46
2:B:542:VAL:CG1	2:B:543:VAL:N	2.77	0.46
2:C:274:PHE:C	2:C:274:PHE:CD1	2.92	0.46
2:C:356:ALA:HA	2:C:384:ALA:HA	1.98	0.46
2:C:550:SER:C	2:C:552:SER:H	2.23	0.46
2:D:500:GLU:C	2:D:502:ASN:N	2.70	0.46
2:F:445:LEU:HD21	2:F:466:ILE:HG12	1.98	0.46
2:F:619:MET:HG3	2:F:624:VAL:HG12	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:460:ALA:HB1	2:G:462:TYR:CD1	2.50	0.46
2:H:274:PHE:CZ	2:H:518:MET:HG3	2.51	0.46
2:H:421:SER:OG	2:H:467:GLU:OE2	2.33	0.46
2:H:460:ALA:C	2:H:462:TYR:N	2.72	0.46
1:K:93:U:C4	1:K:94:C:N4	2.84	0.46
1:M:77:U:O2'	1:M:98:A:N1	2.48	0.46
1:M:86:G:C4	2:E:547:ILE:HD13	2.51	0.46
1:O:86:G:P	1:O:86:G:O4'	2.73	0.46
2:A:527:GLU:O	2:A:528:GLU:O	2.34	0.46
2:C:612:ARG:HH11	2:C:612:ARG:CG	2.29	0.46
2:D:366:PRO:O	2:D:367:SER:C	2.59	0.46
2:E:255:HIS:HE2	2:E:257:GLN:CG	2.29	0.46
2:E:341:PHE:O	2:E:346:LYS:NZ	2.40	0.46
2:E:548:THR:HG23	2:E:551:GLN:OE1	2.15	0.46
2:F:423:ARG:HD2	2:F:438:GLU:OE2	2.16	0.46
2:G:347:SER:HA	2:G:498:VAL:H	1.81	0.46
2:G:582:ILE:HA	2:G:593:MET:HE2	1.98	0.46
2:H:285:LEU:HD21	2:H:512:ARG:HH12	1.81	0.46
2:H:300:ASP:OD2	2:H:302:VAL:HB	2.16	0.46
1:I:84:U:O2'	1:I:85:C:P	2.73	0.46
1:K:95:A:C2	1:K:96:C:C2	3.03	0.46
2:A:288:TYR:CE2	2:A:403:LEU:HD13	2.51	0.46
2:A:423:ARG:NE	2:A:438:GLU:OE1	2.48	0.46
2:B:254:TYR:HE2	2:B:256:MET:HE2	1.81	0.46
2:B:452:PHE:HE2	2:B:454:TYR:CE1	2.21	0.46
2:C:422:THR:HB	2:C:438:GLU:OE2	2.15	0.46
2:C:621:VAL:HG12	2:C:622:ASN:N	2.31	0.46
2:D:628:LEU:O	2:D:632:ILE:HG13	2.14	0.46
2:E:247:ILE:O	2:E:250:GLN:HB2	2.16	0.46
2:F:356:ALA:HA	2:F:384:ALA:CB	2.46	0.46
2:F:500:GLU:C	2:F:502:ASN:N	2.69	0.46
2:G:322:SER:O	2:G:323:GLU:C	2.57	0.46
2:G:603:SER:C	2:G:605:LYS:H	2.22	0.46
2:H:366:PRO:O	2:H:367:SER:C	2.59	0.46
2:H:422:THR:O	2:H:423:ARG:C	2.58	0.46
2:H:455:GLN:O	2:H:456:LEU:C	2.59	0.46
2:H:462:TYR:HD2	2:H:486:ASP:OD2	1.98	0.46
2:H:562:LEU:HD12	2:H:625:ILE:HD11	1.98	0.46
2:H:602:GLU:OE1	2:H:602:GLU:N	2.49	0.46
1:L:86:G:P	1:L:86:G:O4'	2.73	0.46
1:P:86:G:C3'	5:P:1613:HOH:O	2.54	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:274:PHE:CD1	2:A:274:PHE:C	2.94	0.46
2:A:512:ARG:NE	2:A:512:ARG:O	2.48	0.46
2:C:273:ILE:HG22	2:C:522:ILE:HD11	1.98	0.46
2:C:297:PHE:CG	2:D:362:HIS:HE1	2.34	0.46
2:D:389:THR:HG1	2:D:392:GLN:HG3	1.79	0.46
2:E:329:ILE:N	2:E:329:ILE:CD1	2.79	0.46
2:F:397:VAL:O	2:F:401:ILE:HG12	2.15	0.46
2:G:402:ARG:NH1	2:G:402:ARG:CB	2.77	0.46
2:G:562:LEU:HD12	2:G:625:ILE:HD11	1.98	0.46
2:H:329:ILE:N	2:H:329:ILE:CD1	2.79	0.46
2:H:588:ARG:C	2:H:589:ARG:HG2	2.40	0.46
1:J:86:G:P	1:J:86:G:O4'	2.73	0.46
2:A:278:GLU:CB	2:A:358:PHE:HE2	2.29	0.46
2:A:462:TYR:CD1	2:A:462:TYR:N	2.84	0.46
2:A:621:VAL:HG12	2:A:622:ASN:N	2.29	0.46
2:B:280:PHE:CZ	2:B:407:MET:HG2	2.51	0.46
2:B:526:THR:HG23	2:B:533:PHE:HZ	1.81	0.46
2:B:602:GLU:OE1	2:B:602:GLU:N	2.49	0.46
2:D:347:SER:HA	2:D:498:VAL:H	1.80	0.46
2:D:548:THR:HG23	2:D:551:GLN:OE1	2.16	0.46
2:E:277:LEU:O	2:E:280:PHE:HB3	2.16	0.46
2:E:299:MET:HE1	2:E:335:PRO:HB2	1.97	0.46
2:E:386:ILE:O	2:E:509:MET:HA	2.16	0.46
2:E:602:GLU:OE1	2:E:602:GLU:N	2.49	0.46
2:F:550:SER:C	2:F:552:SER:H	2.23	0.46
2:H:274:PHE:C	2:H:274:PHE:HD1	2.24	0.46
2:H:619:MET:HG3	2:H:624:VAL:HG12	1.98	0.46
1:I:80:U:H2'	1:I:81:C:C6	2.51	0.45
1:I:87:U:C2'	1:I:88:G:OP2	2.64	0.45
1:O:87:U:C2'	5:O:1519:HOH:O	2.39	0.45
1:P:87:U:H3'	5:P:1619:HOH:O	2.16	0.45
2:A:528:GLU:O	2:A:530:ALA:N	2.49	0.45
2:B:555:VAL:O	2:B:559:THR:HG23	2.16	0.45
2:B:620:ASP:HB3	2:B:623:GLU:CB	2.46	0.45
2:C:500:GLU:OE1	2:C:500:GLU:N	2.48	0.45
2:D:339:GLN:OE1	2:D:339:GLN:CA	2.64	0.45
2:D:490:PRO:HD3	2:D:509:MET:CE	2.46	0.45
2:F:330:LYS:HA	2:F:331:PRO:HD3	1.64	0.45
2:H:411:PHE:HE2	2:H:525:LEU:HD11	1.81	0.45
2:A:402:ARG:NH1	2:A:402:ARG:CB	2.78	0.45
2:A:501:ASP:CG	2:A:503:GLU:HB2	2.42	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:550:SER:C	2:A:552:SER:H	2.24	0.45
2:B:375:ARG:NH2	5:B:210:HOH:O	2.36	0.45
2:C:389:THR:HG1	2:C:392:GLN:HG3	1.77	0.45
2:C:538:ALA:O	2:C:539:PRO:C	2.59	0.45
2:D:356:ALA:HA	2:D:384:ALA:HA	1.98	0.45
2:D:427:ARG:HB2	2:D:434:TRP:CZ2	2.51	0.45
2:E:423:ARG:O	2:E:423:ARG:HG2	2.15	0.45
2:E:455:GLN:O	2:E:456:LEU:C	2.59	0.45
2:F:273:ILE:HG22	2:F:522:ILE:HD11	1.97	0.45
2:F:290:TYR:CZ	2:F:512:ARG:NH1	2.85	0.45
2:G:569:VAL:CG2	2:G:570:LYS:N	2.79	0.45
1:M:87:U:C3'	5:M:1319:HOH:O	2.62	0.45
1:O:97:C:C4'	1:O:98:A:OP1	2.64	0.45
2:A:294:LYS:HA	2:A:357:GLU:HG3	1.98	0.45
2:B:278:GLU:CA	2:B:358:PHE:HE2	2.29	0.45
2:B:329:ILE:N	2:B:329:ILE:CD1	2.80	0.45
2:B:545:MET:SD	2:B:582:ILE:HD13	2.56	0.45
2:C:589:ARG:O	2:C:611:ARG:HD3	2.16	0.45
2:D:300:ASP:OD2	2:D:302:VAL:HB	2.16	0.45
2:D:526:THR:O	2:D:530:ALA:N	2.50	0.45
2:D:526:THR:O	2:D:530:ALA:HA	2.16	0.45
2:D:538:ALA:O	2:D:539:PRO:C	2.58	0.45
2:F:253:LEU:O	2:F:270:GLY:HA3	2.16	0.45
2:F:422:THR:HB	2:F:438:GLU:OE2	2.17	0.45
2:H:526:THR:O	2:H:530:ALA:N	2.49	0.45
2:A:462:TYR:HE2	2:A:511:HIS:CD2	2.35	0.45
2:A:541:GLN:HE22	2:A:632:ILE:HD13	1.82	0.45
2:D:536:TRP:CD1	2:D:635:ARG:HA	2.52	0.45
2:D:569:VAL:HG22	2:D:570:LYS:N	2.30	0.45
2:F:298:MET:HE2	2:F:327:TYR:HB2	1.97	0.45
2:F:323:GLU:CD	2:F:323:GLU:H	2.22	0.45
2:F:462:TYR:CE2	2:F:511:HIS:NE2	2.83	0.45
2:G:279:VAL:CG2	2:G:280:PHE:N	2.80	0.45
2:G:280:PHE:HE2	2:G:407:MET:HG2	1.78	0.45
2:G:320:THR:HG22	2:G:321:SER:N	2.25	0.45
1:I:88:G:H1'	2:A:583:ARG:CG	2.46	0.45
1:K:85:C:C2'	2:C:577:LYS:HG2	2.46	0.45
1:K:93:U:H2'	1:K:94:C:H6	1.82	0.45
1:M:71:C:C2	1:M:72:G:C8	3.04	0.45
1:M:81:C:O2'	1:M:82:U:H5'	2.16	0.45
1:O:77:U:O2'	1:O:98:A:N1	2.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:471:TYR:CD1	2:A:471:TYR:N	2.85	0.45
2:B:304:TRP:CE3	2:B:328:CYS:HB2	2.52	0.45
2:B:550:SER:C	2:B:552:SER:N	2.72	0.45
2:C:281:VAL:O	2:C:282:ARG:C	2.60	0.45
2:C:282:ARG:HD3	2:C:292:GLU:OE1	2.16	0.45
2:D:437:ALA:HA	2:D:485:LEU:HD23	1.97	0.45
2:E:285:LEU:HD21	2:E:512:ARG:HH12	1.82	0.45
2:E:462:TYR:CE2	2:E:511:HIS:NE2	2.85	0.45
2:E:624:VAL:O	2:E:628:LEU:HG	2.17	0.45
2:F:380:THR:N	2:F:519:GLU:OE2	2.50	0.45
2:F:386:ILE:O	2:F:509:MET:HA	2.17	0.45
2:G:490:PRO:HD3	2:G:509:MET:CE	2.46	0.45
2:G:627:LYS:O	2:G:630:GLN:HB3	2.16	0.45
1:K:98:A:N7	2:D:345:LEU:HD13	2.31	0.45
2:A:445:LEU:HD21	2:A:466:ILE:HG12	1.98	0.45
2:B:277:LEU:O	2:B:280:PHE:HB3	2.17	0.45
2:C:411:PHE:CE2	2:C:521:PHE:HZ	2.31	0.45
2:C:569:VAL:HG22	2:C:570:LYS:N	2.31	0.45
2:D:421:SER:HB3	2:D:458:GLU:CB	2.35	0.45
2:E:362:HIS:HE1	2:F:297:PHE:CG	2.34	0.45
2:E:397:VAL:O	2:E:400:CYS:HB2	2.16	0.45
2:G:330:LYS:HA	2:G:331:PRO:HD3	1.63	0.45
2:G:538:ALA:O	2:G:539:PRO:C	2.59	0.45
2:H:371:HIS:O	2:H:372:GLY:C	2.59	0.45
2:H:380:THR:O	2:H:380:THR:HG22	2.17	0.45
1:J:96:C:H2'	1:J:97:C:C6	2.51	0.45
1:N:89:U:O3'	2:F:577:LYS:NZ	2.44	0.45
2:A:273:ILE:HG23	2:A:537:LEU:CD2	2.46	0.45
2:A:405:TYR:HE1	2:A:416:ILE:HG21	1.82	0.45
2:A:423:ARG:HE	2:A:438:GLU:CD	2.23	0.45
2:A:534:PRO:O	2:A:535:THR:C	2.59	0.45
2:A:603:SER:C	2:A:605:LYS:H	2.24	0.45
2:B:251:LEU:HD22	2:B:588:ARG:O	2.17	0.45
2:B:300:ASP:OD2	2:B:302:VAL:HB	2.16	0.45
2:B:312:ASN:O	2:B:313:TYR:HB2	2.16	0.45
2:C:621:VAL:O	2:C:622:ASN:C	2.59	0.45
2:F:294:LYS:HA	2:F:357:GLU:HG3	1.98	0.45
2:F:397:VAL:O	2:F:400:CYS:HB2	2.17	0.45
2:G:323:GLU:CD	2:G:323:GLU:H	2.24	0.45
2:H:460:ALA:C	2:H:462:TYR:H	2.25	0.45
2:H:554:TYR:CE1	2:H:558:LEU:HD11	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:422:THR:HB	2:A:438:GLU:OE2	2.17	0.45
2:A:482:THR:O	2:A:512:ARG:HA	2.16	0.45
2:B:501:ASP:O	2:B:502:ASN:C	2.58	0.45
2:B:551:GLN:HE21	2:B:598:ASP:N	2.14	0.45
2:C:323:GLU:CD	2:C:323:GLU:H	2.23	0.45
2:C:441:LEU:O	2:C:444:ALA:HB3	2.16	0.45
2:D:609:ARG:HD3	2:D:610:THR:O	2.17	0.45
2:E:445:LEU:HD21	2:E:466:ILE:HG12	1.98	0.45
2:E:610:THR:HG22	2:E:614:LYS:O	2.17	0.45
2:F:555:VAL:O	2:F:559:THR:HG23	2.17	0.45
1:J:93:U:C2	1:J:94:C:C6	3.05	0.45
1:N:86:G:O6	2:F:599:LYS:CB	2.65	0.45
1:O:89:U:P	5:O:1528:HOH:O	2.75	0.45
1:P:80:U:H2'	1:P:81:C:H6	1.80	0.45
2:A:417:VAL:HG13	2:A:471:TYR:CE1	2.47	0.45
2:A:422:THR:O	2:A:458:GLU:O	2.34	0.45
2:A:438:GLU:O	2:A:439:ALA:C	2.60	0.45
2:B:427:ARG:HB2	2:B:434:TRP:CZ2	2.52	0.45
2:B:490:PRO:HD3	2:B:509:MET:CE	2.47	0.45
2:C:294:LYS:HA	2:C:294:LYS:HD2	1.84	0.45
2:C:585:HIS:HA	2:C:588:ARG:HB2	1.99	0.45
2:D:253:LEU:O	2:D:270:GLY:HA3	2.17	0.45
2:D:610:THR:HG22	2:D:614:LYS:O	2.17	0.45
2:E:280:PHE:CZ	2:E:407:MET:HG2	2.51	0.45
2:F:526:THR:HG23	2:F:533:PHE:CZ	2.50	0.45
2:F:562:LEU:HB3	2:F:569:VAL:HG11	1.97	0.45
2:G:383:ASP:OD2	2:G:385:HIS:HE1	2.00	0.45
2:G:567:ILE:O	2:G:569:VAL:N	2.50	0.45
1:K:90:G:C8	2:C:577:LYS:HD2	2.51	0.45
1:M:87:U:C2'	1:M:88:G:OP2	2.65	0.45
2:A:282:ARG:HD3	2:A:292:GLU:OE1	2.16	0.45
2:A:460:ALA:C	2:A:462:TYR:N	2.72	0.45
2:A:480:CYS:O	2:A:516:GLY:HA3	2.17	0.45
2:A:488:SER:O	2:A:489:LEU:C	2.60	0.45
2:A:588:ARG:C	2:A:589:ARG:HG2	2.43	0.45
2:A:628:LEU:O	2:A:632:ILE:HG13	2.17	0.45
2:B:279:VAL:CG2	2:B:280:PHE:N	2.80	0.45
2:B:423:ARG:HG2	2:B:423:ARG:O	2.17	0.45
2:C:294:LYS:HA	2:C:357:GLU:HG3	1.99	0.45
2:C:297:PHE:CG	2:D:362:HIS:CE1	3.05	0.45
2:C:371:HIS:O	2:C:372:GLY:C	2.60	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:436:ARG:HG2	2:C:440:ASP:OD2	2.17	0.45
2:C:455:GLN:O	2:C:456:LEU:C	2.60	0.45
2:D:277:LEU:HB2	2:D:518:MET:HE1	1.98	0.45
2:E:251:LEU:HD22	2:E:588:ARG:O	2.17	0.45
2:F:421:SER:CA	2:F:455:GLN:HB3	2.45	0.45
1:N:87:U:H2'	5:N:1419:HOH:O	2.16	0.44
2:B:551:GLN:CD	2:B:551:GLN:H	2.22	0.44
2:C:298:MET:HE2	2:C:327:TYR:HB2	1.99	0.44
2:C:303:LEU:HD11	2:C:339:GLN:NE2	2.31	0.44
2:C:543:VAL:HG11	2:C:585:HIS:ND1	2.31	0.44
2:E:243:ASP:O	2:E:247:ILE:HG13	2.17	0.44
2:E:298:MET:HE2	2:E:327:TYR:CD1	2.52	0.44
2:E:320:THR:HG22	2:E:321:SER:N	2.25	0.44
2:F:255:HIS:HE2	2:F:257:GLN:CG	2.30	0.44
2:F:422:THR:O	2:F:423:ARG:C	2.60	0.44
2:F:465:LYS:HG3	2:F:484:GLN:CG	2.44	0.44
2:G:397:VAL:O	2:G:401:ILE:HG12	2.17	0.44
2:H:569:VAL:O	2:H:570:LYS:HG2	2.17	0.44
1:I:83:U:H3	2:A:575:ASN:HD21	1.65	0.44
1:L:84:U:H4'	1:L:85:C:H5'	1.97	0.44
1:O:89:U:H5	2:G:583:ARG:NH2	2.14	0.44
2:A:423:ARG:HD2	2:A:438:GLU:OE2	2.18	0.44
2:B:253:LEU:N	2:B:253:LEU:HD23	2.32	0.44
2:B:624:VAL:HG23	2:B:625:ILE:N	2.32	0.44
2:C:242:ARG:HE	2:C:246:LYS:HD3	1.82	0.44
2:C:323:GLU:O	2:C:324:ASN:O	2.35	0.44
2:C:421:SER:HB3	2:C:458:GLU:CB	2.35	0.44
2:D:329:ILE:N	2:D:329:ILE:CD1	2.80	0.44
2:F:285:LEU:HD22	2:F:290:TYR:CG	2.52	0.44
2:F:366:PRO:O	2:F:367:SER:C	2.59	0.44
2:G:249:LYS:HB3	2:G:249:LYS:HZ3	1.78	0.44
2:G:254:TYR:HE2	2:G:256:MET:HE2	1.82	0.44
2:H:551:GLN:HE21	2:H:598:ASP:N	2.16	0.44
1:I:86:G:H3'	5:I:109:HOH:O	2.17	0.44
1:M:84:U:H4'	1:M:85:C:H5'	1.96	0.44
2:A:556:ASN:HD21	2:A:573:LEU:HD11	1.82	0.44
2:B:309:HIS:HB3	2:B:317:MET:CE	2.37	0.44
2:B:421:SER:CA	2:B:455:GLN:HB3	2.47	0.44
2:C:423:ARG:NE	2:C:438:GLU:OE1	2.49	0.44
2:C:437:ALA:HB2	2:C:487:PHE:CD2	2.52	0.44
2:D:353:LEU:HB3	2:D:387:PHE:HB2	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:426:LYS:HE2	2:D:426:LYS:HB3	1.84	0.44
2:D:585:HIS:HA	2:D:588:ARG:HB2	1.98	0.44
2:F:291:GLN:NE2	2:F:353:LEU:HD21	2.32	0.44
2:F:365:GLU:N	2:F:377:ARG:HD3	2.30	0.44
2:G:288:TYR:N	2:G:288:TYR:CD1	2.85	0.44
2:G:400:CYS:O	2:G:404:VAL:HG23	2.17	0.44
2:G:488:SER:O	2:G:489:LEU:C	2.60	0.44
2:G:604:GLY:O	2:G:621:VAL:HG23	2.18	0.44
2:H:501:ASP:O	2:H:503:GLU:HG2	2.18	0.44
1:N:72:G:H2'	1:N:73:U:H6	1.82	0.44
1:P:85:C:C5	2:H:547:ILE:O	2.68	0.44
2:A:285:LEU:HD21	2:A:512:ARG:HH12	1.83	0.44
2:B:278:GLU:HA	2:B:358:PHE:HE2	1.83	0.44
2:C:254:TYR:CD2	2:C:255:HIS:N	2.85	0.44
2:C:298:MET:HG2	2:C:298:MET:O	2.16	0.44
2:C:351:LEU:HD22	2:C:387:PHE:O	2.17	0.44
2:C:362:HIS:HE1	2:D:297:PHE:CD1	2.35	0.44
2:C:460:ALA:C	2:C:462:TYR:H	2.24	0.44
2:C:590:VAL:O	2:C:611:ARG:HB3	2.17	0.44
2:D:281:VAL:O	2:D:282:ARG:C	2.61	0.44
2:D:501:ASP:O	2:D:503:GLU:HG2	2.18	0.44
2:F:411:PHE:CD2	2:F:521:PHE:HZ	2.35	0.44
2:F:423:ARG:HE	2:F:438:GLU:CD	2.24	0.44
2:G:251:LEU:CA	2:G:588:ARG:HH12	2.25	0.44
2:G:534:PRO:O	2:G:535:THR:C	2.59	0.44
2:H:305:GLU:HA	2:H:310:TRP:HB3	1.98	0.44
2:H:542:VAL:CG1	2:H:543:VAL:N	2.80	0.44
1:K:89:U:O2'	2:C:580:PHE:CG	2.69	0.44
1:L:96:C:H2'	1:L:97:C:C6	2.52	0.44
2:A:255:HIS:HE2	2:A:257:GLN:CG	2.30	0.44
2:B:287:GLU:HB2	2:B:288:TYR:HD1	1.83	0.44
2:D:287:GLU:HB2	2:D:288:TYR:HD1	1.82	0.44
2:E:303:LEU:HD11	2:E:339:GLN:NE2	2.31	0.44
2:E:399:GLY:HA2	2:E:402:ARG:HH11	1.82	0.44
2:E:421:SER:OG	2:E:467:GLU:OE2	2.36	0.44
2:E:551:GLN:HG2	2:E:597:GLY:C	2.42	0.44
2:E:609:ARG:HD3	2:E:610:THR:O	2.18	0.44
2:F:471:TYR:N	2:F:471:TYR:CD1	2.86	0.44
2:G:455:GLN:O	2:G:456:LEU:C	2.61	0.44
2:H:323:GLU:O	2:H:324:ASN:O	2.36	0.44
1:J:88:G:H8	5:J:1023:HOH:O	1.97	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:89:U:O3'	2:G:577:LYS:NZ	2.46	0.44
1:P:72:G:H2'	1:P:73:U:H6	1.83	0.44
1:P:103:G:H5'	5:P:1625:HOH:O	2.18	0.44
2:A:273:ILE:HG23	2:A:537:LEU:HD22	1.99	0.44
2:A:562:LEU:HB3	2:A:569:VAL:HG11	2.00	0.44
2:B:403:LEU:HG	2:B:407:MET:HE1	1.99	0.44
2:B:619:MET:HG3	2:B:624:VAL:HG12	2.00	0.44
2:D:437:ALA:HB2	2:D:487:PHE:CD2	2.53	0.44
2:E:526:THR:O	2:E:530:ALA:HA	2.18	0.44
2:E:603:SER:C	2:E:605:LYS:N	2.73	0.44
2:F:437:ALA:HA	2:F:485:LEU:HD23	1.99	0.44
2:F:538:ALA:O	2:F:539:PRO:C	2.60	0.44
2:F:548:THR:HG23	2:F:551:GLN:OE1	2.18	0.44
2:F:569:VAL:HG22	2:F:570:LYS:N	2.33	0.44
2:G:365:GLU:OE1	2:G:365:GLU:HA	2.18	0.44
2:G:394:ARG:O	2:G:395:ASP:C	2.60	0.44
2:G:603:SER:C	2:G:605:LYS:N	2.75	0.44
2:H:395:ASP:OD1	2:H:395:ASP:N	2.47	0.44
1:M:86:G:P	1:M:86:G:O4'	2.76	0.44
1:P:86:G:N9	2:H:547:ILE:HD13	2.31	0.44
2:A:285:LEU:HD22	2:A:290:TYR:CG	2.52	0.44
2:A:356:ALA:HB2	2:A:384:ALA:HB2	2.00	0.44
2:A:386:ILE:HG21	2:A:396:GLU:CG	2.48	0.44
2:D:255:HIS:HE2	2:D:257:GLN:CG	2.31	0.44
2:D:562:LEU:HB3	2:D:569:VAL:HG11	2.00	0.44
2:F:562:LEU:HD12	2:F:625:ILE:HD11	2.00	0.44
2:G:287:GLU:HB2	2:G:288:TYR:HD1	1.82	0.44
2:G:288:TYR:N	2:G:288:TYR:HD1	2.15	0.44
2:G:371:HIS:O	2:G:372:GLY:C	2.59	0.44
2:H:405:TYR:HE1	2:H:416:ILE:HG21	1.83	0.44
2:H:427:ARG:HB2	2:H:434:TRP:CZ2	2.52	0.44
2:H:534:PRO:O	2:H:535:THR:C	2.60	0.44
2:H:609:ARG:HD3	2:H:610:THR:O	2.18	0.44
1:M:98:A:H4'	5:M:1327:HOH:O	2.17	0.44
2:A:603:SER:C	2:A:605:LYS:N	2.76	0.44
2:C:255:HIS:ND1	2:C:267:HIS:CE1	2.67	0.44
2:C:427:ARG:HA	2:C:461:PHE:O	2.18	0.44
2:D:291:GLN:NE2	2:D:353:LEU:HD21	2.33	0.44
2:D:528:GLU:O	2:D:530:ALA:N	2.51	0.44
2:E:339:GLN:OE1	2:E:339:GLN:CA	2.66	0.44
2:E:347:SER:HA	2:E:498:VAL:H	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:621:VAL:O	2:E:622:ASN:C	2.61	0.44
2:F:277:LEU:O	2:F:280:PHE:HB3	2.17	0.44
2:F:459:GLY:HA2	2:F:465:LYS:HE2	1.99	0.44
2:F:551:GLN:HE21	2:F:598:ASP:N	2.16	0.44
2:H:322:SER:O	2:H:323:GLU:C	2.61	0.44
2:H:471:TYR:N	2:H:471:TYR:CD1	2.86	0.44
2:H:474:LEU:O	2:H:475:ASP:C	2.60	0.44
1:J:75:U:O2'	1:J:76:G:P	2.76	0.44
1:J:83:U:H3	2:B:575:ASN:HD21	1.66	0.44
1:J:94:C:C2	1:J:95:A:C8	3.06	0.44
1:N:86:G:N3	2:F:547:ILE:CD1	2.81	0.44
2:A:320:THR:HG23	2:B:319:THR:H	1.83	0.44
2:A:332:MET:HG2	2:A:363:ARG:NH2	2.31	0.44
2:B:290:TYR:CE2	2:B:512:ARG:NH1	2.86	0.44
2:B:351:LEU:HD22	2:B:387:PHE:O	2.17	0.44
2:C:482:THR:O	2:C:512:ARG:HA	2.17	0.44
2:C:501:ASP:O	2:C:502:ASN:C	2.60	0.44
2:D:282:ARG:HD3	2:D:292:GLU:OE1	2.17	0.44
2:D:547:ILE:HG12	2:D:578:ILE:HD12	2.00	0.44
2:E:462:TYR:HE2	2:E:511:HIS:NE2	2.14	0.44
2:E:587:LEU:HD23	2:E:587:LEU:N	2.24	0.44
2:F:527:GLU:O	2:F:528:GLU:O	2.36	0.44
2:G:395:ASP:OD1	2:G:395:ASP:N	2.48	0.44
2:G:462:TYR:CD2	2:G:486:ASP:OD2	2.70	0.44
2:H:600:GLU:HB3	2:H:605:LYS:O	2.17	0.44
2:H:612:ARG:HH11	2:H:612:ARG:CG	2.28	0.44
2:A:250:GLN:NE2	2:A:250:GLN:HA	2.33	0.43
2:A:557:GLU:O	2:A:560:GLN:HB3	2.18	0.43
2:B:569:VAL:O	2:B:570:LYS:HG2	2.17	0.43
2:C:278:GLU:HB3	2:C:358:PHE:CE2	2.52	0.43
2:D:317:MET:HG3	2:D:331:PRO:HG3	2.00	0.43
2:D:403:LEU:HG	2:D:407:MET:HE1	1.97	0.43
2:D:441:LEU:O	2:D:444:ALA:HB3	2.18	0.43
2:D:455:GLN:O	2:D:456:LEU:C	2.60	0.43
2:D:501:ASP:CG	2:D:503:GLU:HB2	2.41	0.43
2:F:300:ASP:OD2	2:F:302:VAL:HB	2.18	0.43
2:F:587:LEU:HD23	2:F:587:LEU:N	2.25	0.43
2:F:603:SER:C	2:F:605:LYS:N	2.76	0.43
2:H:556:ASN:HD21	2:H:573:LEU:HD11	1.83	0.43
2:A:366:PRO:O	2:A:367:SER:C	2.61	0.43
2:A:388:CYS:O	2:A:508:VAL:HG23	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:418:VAL:HG21	2:A:450:ILE:CG2	2.49	0.43
2:B:473:CYS:CA	5:B:221:HOH:O	2.35	0.43
2:B:480:CYS:O	2:B:516:GLY:HA3	2.18	0.43
2:C:298:MET:CE	2:C:327:TYR:HB2	2.47	0.43
2:C:347:SER:HA	2:C:498:VAL:H	1.83	0.43
2:C:366:PRO:O	2:C:367:SER:C	2.61	0.43
2:D:555:VAL:O	2:D:559:THR:HG23	2.18	0.43
2:E:542:VAL:CG1	2:E:543:VAL:N	2.80	0.43
2:F:332:MET:HG2	2:F:363:ARG:NH2	2.33	0.43
2:F:353:LEU:HB3	2:F:387:PHE:HB2	1.99	0.43
2:G:386:ILE:O	2:G:509:MET:HA	2.18	0.43
2:G:551:GLN:HE21	2:G:597:GLY:C	2.25	0.43
2:H:388:CYS:O	2:H:508:VAL:HG23	2.18	0.43
2:H:625:ILE:O	2:H:626:GLU:C	2.61	0.43
1:I:77:U:O5'	1:I:77:U:H6	2.01	0.43
1:L:72:G:H2'	1:L:73:U:H6	1.81	0.43
1:L:79:A:O2'	2:C:500:GLU:O	2.35	0.43
1:M:89:U:H2'	1:M:90:G:OP1	2.18	0.43
1:P:86:G:P	1:P:86:G:O4'	2.76	0.43
2:A:332:MET:HE1	2:A:462:TYR:OH	2.18	0.43
2:A:348:TYR:CE2	2:A:500:GLU:HG3	2.53	0.43
2:A:529:PHE:H	2:A:529:PHE:HD1	1.55	0.43
2:B:402:ARG:NH1	2:B:402:ARG:CB	2.79	0.43
2:B:548:THR:C	2:B:550:SER:H	2.25	0.43
2:C:330:LYS:HA	2:C:331:PRO:HD3	1.61	0.43
2:C:394:ARG:O	2:C:395:ASP:C	2.61	0.43
2:D:254:TYR:CD2	2:D:255:HIS:N	2.85	0.43
2:D:299:MET:HE1	2:D:335:PRO:HB2	2.00	0.43
2:E:353:LEU:HD22	2:E:353:LEU:HA	1.82	0.43
2:E:437:ALA:HA	2:E:485:LEU:HD23	2.00	0.43
2:E:639:GLN:O	2:E:642:GLU:HG2	2.18	0.43
2:F:342:ASN:HD21	2:F:495:ALA:HA	1.83	0.43
2:F:356:ALA:HA	2:F:384:ALA:HA	2.00	0.43
2:F:604:GLY:O	2:F:621:VAL:HG23	2.18	0.43
2:F:627:LYS:CB	2:F:640:LEU:HD21	2.47	0.43
2:G:423:ARG:HE	2:G:438:GLU:CD	2.26	0.43
2:H:528:GLU:O	2:H:530:ALA:N	2.51	0.43
1:I:72:G:H2'	1:I:73:U:C6	2.52	0.43
1:N:87:U:C2'	1:N:88:G:OP2	2.66	0.43
2:A:387:PHE:CE1	2:A:509:MET:SD	3.11	0.43
2:A:569:VAL:O	2:A:570:LYS:HG2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:335:PRO:HG2	2:D:336:GLY:H	1.81	0.43
2:D:436:ARG:HG2	2:D:440:ASP:OD2	2.19	0.43
2:D:482:THR:HG21	4:D:5002:TSB:H5'2	1.99	0.43
2:E:282:ARG:HD3	2:E:292:GLU:OE1	2.19	0.43
2:E:582:ILE:HA	2:E:593:MET:CE	2.48	0.43
2:F:254:TYR:HA	2:F:267:HIS:ND1	2.33	0.43
2:F:353:LEU:HD22	2:F:353:LEU:HA	1.78	0.43
2:F:417:VAL:HG13	2:F:471:TYR:CE1	2.49	0.43
2:F:482:THR:O	2:F:512:ARG:HA	2.18	0.43
2:G:319:THR:HG22	2:G:328:CYS:HA	1.99	0.43
2:G:348:TYR:CE2	2:G:500:GLU:HG3	2.53	0.43
2:G:436:ARG:HG2	2:G:440:ASP:OD2	2.18	0.43
2:H:423:ARG:HE	2:H:438:GLU:CD	2.27	0.43
2:H:501:ASP:O	2:H:502:ASN:C	2.60	0.43
2:H:610:THR:HG22	2:H:614:LYS:O	2.18	0.43
1:J:71:C:C2	1:J:72:G:C8	3.06	0.43
1:J:87:U:C2'	1:J:88:G:OP2	2.66	0.43
2:A:423:ARG:HB3	2:A:434:TRP:CE3	2.54	0.43
2:B:285:LEU:HD21	2:B:512:ARG:HH12	1.83	0.43
2:C:300:ASP:OD2	2:C:302:VAL:HB	2.19	0.43
2:C:471:TYR:N	2:C:471:TYR:CD1	2.86	0.43
2:D:298:MET:HE2	2:D:327:TYR:CD1	2.53	0.43
2:D:521:PHE:HE1	2:D:525:LEU:HD11	1.82	0.43
2:D:554:TYR:CE1	2:D:558:LEU:HD11	2.54	0.43
2:D:562:LEU:HD12	2:D:625:ILE:HD11	2.00	0.43
2:E:254:TYR:HA	2:E:267:HIS:ND1	2.34	0.43
2:E:291:GLN:NE2	2:E:353:LEU:HD21	2.32	0.43
2:E:366:PRO:O	2:E:367:SER:C	2.60	0.43
2:F:609:ARG:HD3	2:F:610:THR:O	2.18	0.43
2:F:612:ARG:HH11	2:F:612:ARG:CG	2.32	0.43
2:G:439:ALA:O	2:G:440:ASP:C	2.62	0.43
2:G:521:PHE:HE1	2:G:525:LEU:HD11	1.84	0.43
2:G:583:ARG:O	2:G:586:THR:N	2.50	0.43
2:H:445:LEU:HD21	2:H:466:ILE:HG12	2.01	0.43
1:L:86:G:O6	2:D:599:LYS:HB2	2.18	0.43
1:N:69:G:O2'	1:N:70:G:H5'	2.18	0.43
1:P:72:G:H2'	1:P:73:U:C6	2.54	0.43
2:B:501:ASP:CG	2:B:503:GLU:HB2	2.42	0.43
2:C:380:THR:N	2:C:519:GLU:OE2	2.50	0.43
2:C:603:SER:C	2:C:605:LYS:N	2.75	0.43
2:C:610:THR:HG22	2:C:614:LYS:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:243:ASP:O	2:D:247:ILE:HG13	2.19	0.43
2:D:371:HIS:O	2:D:372:GLY:C	2.60	0.43
2:D:421:SER:OG	2:D:467:GLU:OE2	2.37	0.43
2:D:459:GLY:HA2	2:D:465:LYS:HE2	2.01	0.43
2:E:402:ARG:NH1	2:E:402:ARG:CB	2.80	0.43
2:E:462:TYR:CD2	2:E:486:ASP:OD2	2.71	0.43
2:E:564:ASN:C	2:E:566:GLY:N	2.77	0.43
2:E:612:ARG:HH11	2:E:612:ARG:CG	2.30	0.43
2:F:320:THR:HG22	2:F:321:SER:N	2.25	0.43
2:F:423:ARG:NE	2:F:438:GLU:OE1	2.52	0.43
2:F:526:THR:O	2:F:530:ALA:HA	2.18	0.43
2:G:609:ARG:HD3	2:G:610:THR:O	2.18	0.43
2:H:242:ARG:HE	2:H:246:LYS:HD3	1.84	0.43
1:M:95:A:H2'	1:M:96:C:O4'	2.18	0.43
1:P:87:U:O2'	1:P:88:G:P	2.77	0.43
2:A:327:TYR:CD1	2:A:327:TYR:N	2.87	0.43
2:A:465:LYS:HG3	2:A:484:GLN:CG	2.44	0.43
2:B:323:GLU:O	2:B:324:ASN:O	2.36	0.43
2:B:347:SER:HA	2:B:498:VAL:H	1.84	0.43
2:B:445:LEU:HD21	2:B:466:ILE:HG12	2.00	0.43
2:B:583:ARG:O	2:B:584:GLU:C	2.62	0.43
2:B:622:ASN:O	2:B:626:GLU:HG3	2.18	0.43
2:D:254:TYR:CZ	2:D:373:LEU:HD21	2.54	0.43
2:D:474:LEU:O	2:D:475:ASP:C	2.62	0.43
2:E:254:TYR:CE2	2:E:256:MET:HE2	2.52	0.43
2:E:521:PHE:HE1	2:E:525:LEU:HD11	1.84	0.43
2:E:608:VAL:HB	2:E:619:MET:HE3	1.99	0.43
2:F:484:GLN:HE22	4:F:7002:TSB:HN8	1.67	0.43
2:H:320:THR:HG22	2:H:321:SER:N	2.29	0.43
2:H:402:ARG:NH1	2:H:402:ARG:CB	2.79	0.43
2:H:597:GLY:O	2:H:601:VAL:HG23	2.17	0.43
1:O:86:G:H1'	2:G:547:ILE:HD13	2.01	0.43
2:A:254:TYR:CE1	2:A:373:LEU:HD21	2.54	0.43
2:B:255:HIS:HE2	2:B:257:GLN:CG	2.32	0.43
2:B:544:ILE:HG22	2:B:555:VAL:HG13	2.00	0.43
2:C:520:ARG:NH1	5:C:4008:HOH:O	2.50	0.43
2:D:603:SER:C	2:D:605:LYS:N	2.76	0.43
2:E:371:HIS:O	2:E:372:GLY:C	2.61	0.43
2:E:415:LYS:HG2	2:E:471:TYR:CD1	2.54	0.43
2:E:594:LEU:CD2	2:E:608:VAL:HG22	2.48	0.43
2:F:426:LYS:HE2	2:F:426:LYS:HB3	1.88	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:298:MET:HE2	2:H:327:TYR:CD1	2.54	0.43
1:J:79:A:H1'	2:A:502:ASN:OD1	2.19	0.43
1:J:89:U:H5	2:B:583:ARG:NH2	2.16	0.43
1:K:89:U:O2	1:K:89:U:C2'	2.66	0.43
1:L:95:A:H2'	1:L:96:C:O4'	2.19	0.43
1:N:87:U:H2'	1:N:88:G:OP2	2.19	0.43
1:N:97:C:C5'	1:N:98:A:OP1	2.64	0.43
1:P:87:U:H1'	2:H:582:ILE:HG13	2.00	0.43
2:A:351:LEU:HD13	2:A:388:CYS:HA	2.00	0.43
2:A:452:PHE:CE2	2:A:454:TYR:HE1	2.24	0.43
2:A:573:LEU:O	2:A:574:ARG:C	2.62	0.43
2:A:608:VAL:CB	2:A:619:MET:HE3	2.48	0.43
2:B:358:PHE:N	2:B:358:PHE:CD1	2.86	0.43
2:B:366:PRO:O	2:B:367:SER:C	2.60	0.43
2:B:622:ASN:HB3	2:H:612:ARG:O	2.19	0.43
2:C:280:PHE:CZ	2:C:407:MET:HG2	2.53	0.43
2:D:280:PHE:CZ	2:D:407:MET:HG2	2.53	0.43
2:E:242:ARG:HE	2:E:246:LYS:HD3	1.83	0.43
2:E:538:ALA:O	2:E:539:PRO:C	2.61	0.43
2:F:242:ARG:HE	2:F:246:LYS:HD3	1.83	0.43
2:G:281:VAL:O	2:G:282:ARG:C	2.62	0.43
2:G:542:VAL:CG1	2:G:543:VAL:N	2.82	0.43
2:H:243:ASP:O	2:H:247:ILE:HG13	2.19	0.43
1:L:87:U:C2'	1:L:88:G:OP2	2.66	0.43
1:L:88:G:H1'	2:D:583:ARG:CG	2.49	0.43
2:A:394:ARG:O	2:A:395:ASP:C	2.62	0.43
2:A:403:LEU:HG	2:A:407:MET:HE1	2.01	0.43
2:A:567:ILE:O	2:A:569:VAL:N	2.52	0.43
2:A:630:GLN:HA	2:A:633:ARG:HH12	1.84	0.43
2:B:353:LEU:HB3	2:B:387:PHE:HB2	2.00	0.43
2:B:410:THR:HB	2:B:411:PHE:CD1	2.53	0.43
2:D:332:MET:HE2	4:D:5002:TSB:HN12	1.83	0.43
2:D:588:ARG:C	2:D:589:ARG:HG2	2.44	0.43
2:E:526:THR:O	2:E:530:ALA:N	2.51	0.43
2:F:405:TYR:HE1	2:F:416:ILE:HG21	1.84	0.43
2:F:430:SER:O	2:F:431:ASP:C	2.62	0.43
2:G:313:TYR:CE1	2:G:316:ALA:HB3	2.54	0.43
2:G:417:VAL:HG11	2:G:471:TYR:CE1	2.52	0.43
2:G:430:SER:O	2:G:431:ASP:C	2.61	0.43
2:G:480:CYS:O	2:G:516:GLY:HA3	2.19	0.43
2:G:551:GLN:HE21	2:G:598:ASP:N	2.16	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:621:VAL:O	2:G:622:ASN:C	2.60	0.43
2:G:628:LEU:O	2:G:632:ILE:HG13	2.19	0.43
2:H:279:VAL:CG2	2:H:280:PHE:N	2.82	0.43
2:H:383:ASP:OD2	2:H:385:HIS:CE1	2.72	0.43
1:M:84:U:O2'	1:M:85:C:OP2	2.27	0.42
2:A:299:MET:O	2:A:304:TRP:HZ3	2.02	0.42
2:A:330:LYS:HA	2:A:331:PRO:HD3	1.60	0.42
2:A:353:LEU:HB3	2:A:387:PHE:HB2	1.99	0.42
2:A:585:HIS:HA	2:A:588:ARG:HB2	2.01	0.42
2:B:397:VAL:O	2:B:400:CYS:HB2	2.19	0.42
2:C:462:TYR:CD2	2:C:486:ASP:OD2	2.71	0.42
2:D:421:SER:CA	2:D:455:GLN:HB3	2.45	0.42
2:D:531:GLY:O	2:D:589:ARG:HG3	2.19	0.42
2:D:627:LYS:O	2:D:630:GLN:HB3	2.19	0.42
2:F:251:LEU:CA	2:F:588:ARG:HH12	2.25	0.42
2:F:254:TYR:CD2	2:F:255:HIS:N	2.87	0.42
2:F:273:ILE:HG23	2:F:537:LEU:HD22	2.01	0.42
2:F:529:PHE:CD2	2:F:534:PRO:HD3	2.54	0.42
2:F:585:HIS:HA	2:F:588:ARG:HB2	2.01	0.42
2:F:590:VAL:O	2:F:611:ARG:HB3	2.19	0.42
2:F:621:VAL:HG12	2:F:622:ASN:N	2.33	0.42
2:G:250:GLN:NE2	2:G:250:GLN:HA	2.34	0.42
2:G:572:ASP:OD1	2:G:574:ARG:HB2	2.18	0.42
2:H:255:HIS:HE2	2:H:257:GLN:CG	2.31	0.42
2:H:609:ARG:O	2:H:609:ARG:CG	2.67	0.42
2:A:281:VAL:O	2:A:282:ARG:C	2.61	0.42
2:A:309:HIS:HB3	2:A:317:MET:CE	2.37	0.42
2:C:249:LYS:HE2	2:C:249:LYS:HA	2.01	0.42
2:C:416:ILE:O	2:C:416:ILE:HG22	2.18	0.42
2:C:541:GLN:HE22	2:C:632:ILE:HD13	1.84	0.42
2:E:356:ALA:HA	2:E:384:ALA:HA	2.00	0.42
2:E:399:GLY:HA2	2:E:402:ARG:NH1	2.35	0.42
2:E:545:MET:SD	2:E:582:ILE:HD13	2.58	0.42
2:F:439:ALA:O	2:F:440:ASP:C	2.62	0.42
2:G:337:HIS:NE2	2:G:357:GLU:OE1	2.47	0.42
2:G:527:GLU:O	2:G:528:GLU:C	2.61	0.42
2:H:567:ILE:O	2:H:569:VAL:N	2.52	0.42
1:M:72:G:H2'	1:M:73:U:H6	1.84	0.42
1:N:92:G:H2'	1:N:93:U:H6	1.84	0.42
2:A:242:ARG:HE	2:A:246:LYS:HD3	1.85	0.42
2:A:280:PHE:CZ	2:A:407:MET:HG2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:290:TYR:CZ	2:A:512:ARG:NH1	2.86	0.42
2:A:400:CYS:SG	2:A:512:ARG:HG3	2.60	0.42
2:B:386:ILE:O	2:B:509:MET:HA	2.19	0.42
2:B:551:GLN:HG2	2:B:597:GLY:C	2.45	0.42
2:C:627:LYS:CB	2:C:640:LEU:HD21	2.47	0.42
2:D:423:ARG:NE	2:D:438:GLU:OE1	2.52	0.42
2:D:564:ASN:C	2:D:566:GLY:N	2.77	0.42
2:F:347:SER:HA	2:F:498:VAL:H	1.84	0.42
2:F:365:GLU:OE1	2:F:365:GLU:HA	2.19	0.42
2:F:438:GLU:O	2:F:439:ALA:C	2.62	0.42
2:G:393:ILE:O	2:G:397:VAL:HG23	2.19	0.42
2:G:610:THR:HG21	2:G:614:LYS:HB3	2.00	0.42
2:H:247:ILE:HD12	2:H:526:THR:HG22	2.00	0.42
2:H:277:LEU:HB2	2:H:518:MET:HE1	2.02	0.42
2:H:551:GLN:HE21	2:H:597:GLY:C	2.26	0.42
1:P:94:C:C2	1:P:95:A:C8	3.07	0.42
2:A:299:MET:HE3	2:A:303:LEU:HD13	2.01	0.42
2:A:604:GLY:O	2:A:621:VAL:HG23	2.20	0.42
2:C:365:GLU:OE1	2:C:365:GLU:HA	2.19	0.42
2:D:422:THR:HB	2:D:438:GLU:OE2	2.19	0.42
2:D:610:THR:HG22	2:D:614:LYS:HB3	2.00	0.42
2:E:330:LYS:HA	2:E:331:PRO:HD3	1.61	0.42
2:E:417:VAL:HG13	2:E:471:TYR:CE1	2.50	0.42
2:E:627:LYS:CB	2:E:640:LEU:HD21	2.47	0.42
2:F:569:VAL:O	2:F:570:LYS:HG2	2.20	0.42
2:G:351:LEU:HA	2:G:352:PRO:C	2.44	0.42
2:G:489:LEU:O	2:G:490:PRO:C	2.61	0.42
2:G:497:TYR:OH	2:G:505:LYS:HE2	2.19	0.42
2:G:526:THR:HG23	2:G:533:PHE:CZ	2.55	0.42
2:G:612:ARG:HH11	2:G:612:ARG:CG	2.32	0.42
2:H:332:MET:HE2	4:H:9002:TSB:HN12	1.84	0.42
2:H:347:SER:HA	2:H:498:VAL:H	1.85	0.42
2:H:529:PHE:CD2	2:H:534:PRO:HD3	2.55	0.42
1:I:86:G:C3'	5:I:109:HOH:O	2.66	0.42
1:M:87:U:H2'	1:M:88:G:OP2	2.19	0.42
1:N:78:G:C1'	2:E:345:LEU:HD23	2.49	0.42
1:O:78:G:P	5:O:1539:HOH:O	2.73	0.42
1:O:89:U:C2'	1:O:90:G:OP1	2.68	0.42
1:P:93:U:N3	1:P:94:C:C5	2.88	0.42
2:A:411:PHE:CD2	2:A:525:LEU:CD2	2.90	0.42
2:A:436:ARG:HG2	2:A:440:ASP:OD2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:501:ASP:O	2:A:503:GLU:HG2	2.19	0.42
2:B:305:GLU:HA	2:B:310:TRP:HB3	2.01	0.42
2:B:322:SER:O	2:B:323:GLU:C	2.61	0.42
2:C:411:PHE:CD2	2:C:521:PHE:HZ	2.37	0.42
2:C:459:GLY:HA2	2:C:465:LYS:HE2	2.01	0.42
2:E:317:MET:HG3	2:E:331:PRO:HG3	2.02	0.42
2:E:416:ILE:HG22	2:E:416:ILE:O	2.18	0.42
2:G:393:ILE:HD11	2:G:508:VAL:HG11	2.02	0.42
2:H:278:GLU:OE2	2:H:358:PHE:HD2	2.02	0.42
2:H:323:GLU:CD	2:H:323:GLU:N	2.77	0.42
2:H:589:ARG:O	2:H:611:ARG:NH1	2.48	0.42
1:M:69:G:N7	1:P:105:C:C2	2.87	0.42
1:P:96:C:H2'	1:P:97:C:C6	2.54	0.42
2:A:244:HIS:ND1	2:A:374:MET:HE2	2.34	0.42
2:A:254:TYR:HA	2:A:267:HIS:ND1	2.34	0.42
2:A:386:ILE:HB	2:A:510:ILE:HB	2.01	0.42
2:A:397:VAL:O	2:A:401:ILE:HG12	2.20	0.42
2:C:273:ILE:HG23	2:C:537:LEU:CD2	2.50	0.42
2:C:353:LEU:HB3	2:C:387:PHE:HB2	2.02	0.42
2:D:242:ARG:HE	2:D:246:LYS:HD3	1.85	0.42
2:D:254:TYR:HE2	2:D:256:MET:HE2	1.85	0.42
2:D:332:MET:HG2	2:D:363:ARG:NH2	2.34	0.42
2:D:400:CYS:O	2:D:404:VAL:HG23	2.19	0.42
2:D:439:ALA:O	2:D:440:ASP:C	2.62	0.42
2:F:427:ARG:HB2	2:F:434:TRP:CZ2	2.55	0.42
2:F:488:SER:O	2:F:489:LEU:C	2.63	0.42
2:F:602:GLU:N	2:F:602:GLU:OE1	2.53	0.42
2:G:548:THR:HG23	2:G:551:GLN:OE1	2.20	0.42
2:H:459:GLY:HA2	2:H:465:LYS:HE2	2.02	0.42
1:J:87:U:H2'	1:J:88:G:OP2	2.19	0.42
1:K:70:G:C4	1:K:71:C:C5	3.07	0.42
1:M:69:G:N7	1:P:105:C:O2	2.52	0.42
2:A:548:THR:C	2:A:550:SER:H	2.28	0.42
2:B:489:LEU:O	2:B:490:PRO:C	2.63	0.42
2:C:341:PHE:O	2:C:346:LYS:NZ	2.45	0.42
2:C:365:GLU:N	2:C:377:ARG:HD3	2.33	0.42
2:C:386:ILE:O	2:C:509:MET:HA	2.20	0.42
2:C:387:PHE:CE1	2:C:509:MET:SD	3.13	0.42
2:D:402:ARG:NH1	2:D:402:ARG:CB	2.80	0.42
2:D:501:ASP:O	2:D:502:ASN:C	2.61	0.42
2:D:548:THR:C	2:D:550:SER:H	2.28	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:297:PHE:CG	2:F:362:HIS:CE1	3.07	0.42
2:E:400:CYS:O	2:E:404:VAL:HG23	2.19	0.42
2:E:426:LYS:HE2	2:E:426:LYS:HB3	1.85	0.42
2:E:441:LEU:O	2:E:444:ALA:HB3	2.19	0.42
2:F:329:ILE:N	2:F:329:ILE:CD1	2.83	0.42
2:F:610:THR:HG22	2:F:614:LYS:O	2.19	0.42
2:G:590:VAL:O	2:G:611:ARG:HB3	2.19	0.42
2:H:543:VAL:HG11	2:H:585:HIS:CE1	2.53	0.42
2:H:604:GLY:O	2:H:621:VAL:HG23	2.20	0.42
1:I:89:U:C6	2:A:583:ARG:NH1	2.88	0.42
1:N:86:G:P	1:N:86:G:O4'	2.78	0.42
1:N:90:G:O5'	2:F:577:LYS:HD3	2.20	0.42
2:A:474:LEU:O	2:A:475:ASP:C	2.62	0.42
2:B:597:GLY:O	2:B:598:ASP:C	2.63	0.42
2:C:332:MET:HE1	2:C:462:TYR:OH	2.20	0.42
2:C:427:ARG:HB2	2:C:434:TRP:CZ2	2.55	0.42
2:C:562:LEU:HB3	2:C:569:VAL:HG11	2.01	0.42
2:C:602:GLU:N	2:C:602:GLU:OE1	2.53	0.42
2:D:249:LYS:HE2	2:D:249:LYS:HA	2.02	0.42
2:D:415:LYS:HG2	2:D:471:TYR:CD1	2.55	0.42
2:D:452:PHE:CE2	2:D:454:TYR:HE1	2.22	0.42
2:D:542:VAL:CG1	2:D:543:VAL:N	2.82	0.42
2:E:349:ARG:HE	2:E:349:ARG:HB2	1.61	0.42
2:E:474:LEU:O	2:E:475:ASP:C	2.63	0.42
2:E:550:SER:C	2:E:552:SER:N	2.78	0.42
2:G:305:GLU:HA	2:G:310:TRP:HB3	2.01	0.42
2:H:285:LEU:HD22	2:H:290:TYR:CG	2.54	0.42
2:H:462:TYR:CD2	2:H:486:ASP:OD2	2.72	0.42
2:H:531:GLY:O	2:H:589:ARG:HG3	2.20	0.42
1:K:80:U:H3'	1:K:80:U:C6	2.55	0.42
1:L:81:C:C4	1:L:82:U:C4	3.07	0.42
1:M:81:C:H2'	1:M:82:U:O4'	2.20	0.42
2:A:483:VAL:HG22	2:A:512:ARG:HB2	2.01	0.42
2:A:624:VAL:O	2:A:628:LEU:HG	2.20	0.42
2:B:252:ASP:OD1	2:B:267:HIS:CD2	2.72	0.42
2:B:323:GLU:CD	2:B:323:GLU:N	2.78	0.42
2:B:371:HIS:O	2:B:372:GLY:C	2.63	0.42
2:B:411:PHE:HE2	2:B:521:PHE:CE1	2.27	0.42
2:B:426:LYS:HE2	2:B:426:LYS:HB3	1.87	0.42
2:B:621:VAL:O	2:B:622:ASN:C	2.60	0.42
2:C:317:MET:HG3	2:C:331:PRO:HG3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:421:SER:OG	2:C:467:GLU:OE2	2.37	0.42
2:C:541:GLN:NE2	2:C:632:ILE:HD13	2.35	0.42
2:C:545:MET:HE3	2:C:593:MET:HB3	2.01	0.42
2:D:462:TYR:CD2	2:D:486:ASP:OD2	2.72	0.42
2:D:590:VAL:O	2:D:611:ARG:HB3	2.19	0.42
2:E:441:LEU:HD23	2:E:441:LEU:HA	1.90	0.42
2:E:482:THR:O	2:E:512:ARG:HA	2.19	0.42
2:E:501:ASP:CG	2:E:503:GLU:HB2	2.42	0.42
2:F:317:MET:HG3	2:F:331:PRO:HG3	2.01	0.42
2:F:327:TYR:CD1	2:F:327:TYR:N	2.87	0.42
2:F:421:SER:HB3	2:F:459:GLY:N	2.35	0.42
2:F:517:SER:HB3	4:F:7002:TSB:H1'	2.01	0.42
2:F:551:GLN:HE21	2:F:597:GLY:C	2.27	0.42
2:G:353:LEU:HB3	2:G:387:PHE:HB2	2.01	0.42
2:G:588:ARG:C	2:G:589:ARG:HG2	2.43	0.42
2:G:619:MET:HG3	2:G:624:VAL:HG12	2.01	0.42
2:H:394:ARG:O	2:H:397:VAL:N	2.53	0.42
1:J:84:U:H4'	1:J:85:C:H5'	2.01	0.42
1:J:95:A:H2'	1:J:96:C:O4'	2.20	0.42
2:A:610:THR:HG21	2:A:614:LYS:HB3	2.02	0.42
2:B:437:ALA:HA	2:B:485:LEU:HD23	2.02	0.42
2:B:564:ASN:C	2:B:566:GLY:N	2.78	0.42
2:B:597:GLY:O	2:B:601:VAL:HG23	2.20	0.42
2:C:551:GLN:HE21	2:C:598:ASP:N	2.18	0.42
2:D:430:SER:O	2:D:431:ASP:C	2.63	0.42
2:D:471:TYR:CD1	2:D:471:TYR:N	2.87	0.42
2:E:358:PHE:N	2:E:358:PHE:CD1	2.88	0.42
2:E:397:VAL:O	2:E:401:ILE:HG12	2.19	0.42
2:E:625:ILE:O	2:E:626:GLU:C	2.61	0.42
2:F:277:LEU:HB2	2:F:518:MET:HE1	2.02	0.42
2:F:626:GLU:HG3	2:F:626:GLU:H	1.73	0.42
2:G:411:PHE:HE2	2:G:521:PHE:HE1	1.67	0.42
2:G:431:ASP:O	2:G:432:GLU:C	2.63	0.42
1:M:92:G:H2'	1:M:93:U:C6	2.53	0.41
2:B:430:SER:O	2:B:431:ASP:C	2.61	0.41
2:B:488:SER:O	2:B:489:LEU:C	2.62	0.41
2:C:244:HIS:O	2:C:245:ARG:C	2.63	0.41
2:D:323:GLU:CD	2:D:323:GLU:N	2.78	0.41
2:D:583:ARG:HD3	5:D:5014:HOH:O	2.19	0.41
2:D:587:LEU:HD23	2:D:587:LEU:N	2.27	0.41
2:E:380:THR:N	2:E:519:GLU:OE2	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:627:LYS:O	2:E:630:GLN:HB3	2.20	0.41
2:F:254:TYR:HE2	2:F:256:MET:HE2	1.84	0.41
2:F:287:GLU:HB2	2:F:288:TYR:CD1	2.54	0.41
2:F:462:TYR:CD1	2:F:462:TYR:N	2.88	0.41
2:G:247:ILE:C	2:G:250:GLN:HB2	2.45	0.41
2:G:490:PRO:HD3	2:G:509:MET:HE2	2.02	0.41
2:H:501:ASP:CG	2:H:503:GLU:HB2	2.43	0.41
2:H:622:ASN:O	2:H:626:GLU:HG3	2.19	0.41
2:H:627:LYS:O	2:H:630:GLN:HB3	2.20	0.41
1:K:80:U:OP2	2:D:349:ARG:NH2	2.53	0.41
1:K:90:G:P	2:C:577:LYS:NZ	2.93	0.41
1:M:93:U:H2'	1:M:94:C:C6	2.53	0.41
2:A:325:ARG:HH11	2:A:325:ARG:CG	2.24	0.41
2:A:347:SER:HA	2:A:498:VAL:H	1.84	0.41
2:E:298:MET:CG	2:F:263:MET:HE2	2.45	0.41
2:E:313:TYR:CE1	2:E:316:ALA:HB3	2.55	0.41
2:F:501:ASP:O	2:F:502:ASN:C	2.63	0.41
2:F:594:LEU:CD2	2:F:608:VAL:HG22	2.50	0.41
2:H:253:LEU:O	2:H:270:GLY:HA3	2.20	0.41
1:L:83:U:H3	2:D:575:ASN:HD21	1.68	0.41
1:O:90:G:C8	2:G:577:LYS:HD3	2.55	0.41
2:A:342:ASN:HD22	2:A:342:ASN:HA	1.63	0.41
2:C:322:SER:O	2:C:323:GLU:C	2.63	0.41
2:C:358:PHE:N	2:C:358:PHE:CD1	2.88	0.41
2:C:545:MET:SD	2:C:581:LYS:HB3	2.61	0.41
2:C:564:ASN:C	2:C:566:GLY:N	2.78	0.41
2:D:423:ARG:HG2	2:D:423:ARG:O	2.19	0.41
2:D:569:VAL:O	2:D:570:LYS:HG2	2.19	0.41
2:D:573:LEU:O	2:D:574:ARG:C	2.63	0.41
2:E:365:GLU:N	2:E:377:ARG:HD3	2.35	0.41
2:E:551:GLN:H	2:E:551:GLN:CD	2.25	0.41
2:G:473:CYS:HB2	2:G:527:GLU:OE1	2.21	0.41
2:G:562:LEU:O	2:G:565:ALA:HB3	2.20	0.41
2:H:277:LEU:O	2:H:280:PHE:HB3	2.21	0.41
2:H:582:ILE:HA	2:H:593:MET:CE	2.50	0.41
1:J:74:A:C2	1:J:101:G:C2	3.09	0.41
1:K:99:C:O3'	2:C:249:LYS:NZ	2.54	0.41
1:M:76:G:N2	1:M:97:C:C2	2.89	0.41
2:A:430:SER:O	2:A:431:ASP:C	2.62	0.41
2:D:285:LEU:HD21	2:D:512:ARG:HH12	1.85	0.41
2:D:422:THR:HG1	2:D:456:LEU:HA	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:594:LEU:CD2	2:D:608:VAL:HG22	2.51	0.41
2:E:277:LEU:CB	2:E:518:MET:HE1	2.51	0.41
2:E:353:LEU:HB3	2:E:387:PHE:HB2	2.01	0.41
2:F:349:ARG:HE	2:F:349:ARG:HB2	1.59	0.41
2:F:536:TRP:CD1	2:F:635:ARG:HA	2.55	0.41
2:G:285:LEU:HD21	2:G:512:ARG:HH12	1.85	0.41
2:G:528:GLU:O	2:G:529:PHE:C	2.63	0.41
2:G:532:PHE:CZ	2:G:589:ARG:NH1	2.89	0.41
2:H:536:TRP:CE3	2:H:635:ARG:HD3	2.55	0.41
1:K:95:A:H2'	1:K:96:C:O4'	2.20	0.41
1:L:89:U:H5	2:D:583:ARG:NH2	2.18	0.41
2:A:270:GLY:O	2:A:271:TRP:C	2.63	0.41
2:A:455:GLN:O	2:A:456:LEU:O	2.38	0.41
2:A:526:THR:HG23	2:A:533:PHE:CZ	2.53	0.41
2:B:317:MET:HG3	2:B:331:PRO:HG3	2.03	0.41
2:B:627:LYS:CB	2:B:640:LEU:HD21	2.51	0.41
2:C:277:LEU:O	2:C:280:PHE:HB3	2.21	0.41
2:E:254:TYR:CD2	2:E:255:HIS:N	2.88	0.41
2:E:569:VAL:CG2	2:E:570:LYS:N	2.84	0.41
2:F:246:LYS:O	2:F:250:GLN:HG2	2.21	0.41
2:F:437:ALA:HB2	2:F:487:PHE:CD2	2.55	0.41
2:F:609:ARG:O	2:F:609:ARG:CG	2.69	0.41
2:G:323:GLU:O	2:G:324:ASN:O	2.38	0.41
2:G:526:THR:O	2:G:530:ALA:N	2.53	0.41
2:H:349:ARG:HE	2:H:349:ARG:HB2	1.64	0.41
2:H:465:LYS:HG3	2:H:484:GLN:CG	2.46	0.41
1:P:89:U:H2'	1:P:90:G:OP1	2.21	0.41
2:A:351:LEU:HA	2:A:352:PRO:C	2.45	0.41
2:A:501:ASP:O	2:A:502:ASN:C	2.62	0.41
2:A:542:VAL:CG1	2:A:543:VAL:N	2.84	0.41
2:C:551:GLN:HE21	2:C:597:GLY:C	2.28	0.41
2:C:555:VAL:O	2:C:559:THR:HG23	2.21	0.41
2:D:482:THR:O	2:D:512:ARG:HA	2.20	0.41
2:E:253:LEU:O	2:E:270:GLY:HA3	2.20	0.41
2:E:389:THR:HG1	2:E:392:GLN:HG3	1.84	0.41
2:G:278:GLU:HB3	2:G:358:PHE:CE2	2.54	0.41
2:G:501:ASP:O	2:G:502:ASN:C	2.62	0.41
2:G:573:LEU:O	2:G:574:ARG:C	2.63	0.41
2:G:610:THR:HG22	2:G:614:LYS:HB3	2.02	0.41
1:K:101:G:O2'	1:K:102:C:C5'	2.58	0.41
1:M:96:C:H2'	1:M:97:C:C6	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:70:G:O2'	1:O:71:C:H5'	2.21	0.41
2:A:365:GLU:N	2:A:377:ARG:HD3	2.33	0.41
2:A:519:GLU:CD	2:A:519:GLU:H	2.28	0.41
2:B:514:ILE:HG22	2:B:515:LEU:HD23	2.02	0.41
2:B:609:ARG:O	2:B:609:ARG:CG	2.69	0.41
2:C:255:HIS:HE2	2:C:257:GLN:CG	2.34	0.41
2:C:339:GLN:OE1	2:C:339:GLN:CA	2.67	0.41
2:D:274:PHE:C	2:D:274:PHE:HD1	2.29	0.41
2:D:330:LYS:HA	2:D:331:PRO:HD3	1.62	0.41
2:E:294:LYS:HA	2:E:357:GLU:HG3	2.02	0.41
2:E:562:LEU:HD12	2:E:625:ILE:HD11	2.02	0.41
2:F:351:LEU:HD13	2:F:388:CYS:HA	2.02	0.41
2:F:421:SER:HB3	2:F:458:GLU:CB	2.40	0.41
2:F:545:MET:HE3	2:F:593:MET:HB3	2.03	0.41
2:F:554:TYR:CE1	2:F:558:LEU:HD11	2.56	0.41
2:G:543:VAL:HG21	2:G:590:VAL:HG21	2.03	0.41
2:G:584:GLU:O	2:G:588:ARG:HB2	2.20	0.41
1:O:87:U:H2'	1:O:88:G:OP2	2.20	0.41
1:P:86:G:H2'	5:P:1613:HOH:O	2.20	0.41
2:A:452:PHE:HE2	2:A:454:TYR:CE1	2.23	0.41
2:B:299:MET:O	2:B:304:TRP:HZ3	2.04	0.41
2:B:365:GLU:N	2:B:377:ARG:HD3	2.35	0.41
2:C:329:ILE:N	2:C:329:ILE:CD1	2.84	0.41
2:D:288:TYR:CZ	2:D:403:LEU:HD13	2.56	0.41
2:D:526:THR:HG23	2:D:533:PHE:CZ	2.56	0.41
2:D:621:VAL:O	2:D:622:ASN:C	2.62	0.41
2:E:323:GLU:O	2:E:324:ASN:O	2.38	0.41
2:E:421:SER:HB3	2:E:458:GLU:CB	2.32	0.41
2:F:441:LEU:HD23	2:F:441:LEU:HA	1.91	0.41
2:F:548:THR:C	2:F:550:SER:H	2.29	0.41
2:G:358:PHE:N	2:G:358:PHE:CD1	2.87	0.41
2:G:550:SER:C	2:G:552:SER:N	2.78	0.41
2:H:353:LEU:HB3	2:H:387:PHE:HB2	2.02	0.41
2:H:386:ILE:O	2:H:509:MET:HA	2.20	0.41
2:H:398:ASN:O	2:H:399:GLY:C	2.64	0.41
2:H:462:TYR:CE2	2:H:511:HIS:NE2	2.89	0.41
2:H:547:ILE:HG12	2:H:578:ILE:HD12	2.02	0.41
1:I:75:U:O2'	1:I:89:U:O4	2.28	0.41
1:I:88:G:N2	5:I:111:HOH:O	2.47	0.41
1:J:93:U:N3	1:J:94:C:C5	2.89	0.41
1:L:89:U:C6	2:D:583:ARG:CZ	3.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:86:G:N3	2:F:547:ILE:HD13	2.36	0.41
1:N:89:U:O2'	2:F:580:PHE:CB	2.69	0.41
1:O:69:G:O2'	1:O:70:G:H5'	2.21	0.41
1:O:80:U:H2'	1:O:81:C:H6	1.84	0.41
1:P:78:G:H5'	5:P:1639:HOH:O	2.21	0.41
1:P:97:C:C4'	1:P:98:A:OP1	2.69	0.41
2:A:253:LEU:O	2:A:270:GLY:HA3	2.21	0.41
2:A:300:ASP:OD2	2:A:302:VAL:HB	2.20	0.41
2:A:300:ASP:HB2	2:B:259:GLU:O	2.21	0.41
2:A:459:GLY:HA2	2:A:465:LYS:HE2	2.03	0.41
2:A:460:ALA:C	2:A:462:TYR:H	2.26	0.41
2:B:280:PHE:HE2	2:B:407:MET:HG2	1.85	0.41
2:B:334:CYS:O	2:B:335:PRO:C	2.61	0.41
2:B:351:LEU:HA	2:B:352:PRO:C	2.46	0.41
2:C:452:PHE:CE2	2:C:454:TYR:HE1	2.22	0.41
2:C:480:CYS:O	2:C:516:GLY:CA	2.69	0.41
2:C:527:GLU:O	2:C:528:GLU:O	2.39	0.41
2:C:528:GLU:O	2:C:530:ALA:N	2.53	0.41
2:D:421:SER:HB3	2:D:459:GLY:N	2.35	0.41
2:E:247:ILE:HA	2:E:250:GLN:HB2	2.03	0.41
2:E:250:GLN:HA	2:E:250:GLN:NE2	2.36	0.41
2:E:348:TYR:CE2	2:E:500:GLU:HG3	2.56	0.41
2:E:528:GLU:O	2:E:530:ALA:N	2.53	0.41
2:F:334:CYS:O	2:F:335:PRO:C	2.62	0.41
2:F:348:TYR:CE2	2:F:500:GLU:HG3	2.55	0.41
2:F:358:PHE:N	2:F:358:PHE:CD1	2.88	0.41
2:F:462:TYR:HE2	2:F:511:HIS:CD2	2.39	0.41
2:F:493:LEU:N	2:F:493:LEU:HD23	2.35	0.41
2:G:298:MET:O	2:G:298:MET:HG2	2.21	0.41
2:G:377:ARG:HH11	2:G:377:ARG:HG3	1.86	0.41
2:G:434:TRP:O	2:G:435:ASP:C	2.64	0.41
2:G:517:SER:HB3	4:G:8002:TSB:H1'	2.00	0.41
2:G:548:THR:C	2:G:550:SER:H	2.29	0.41
2:G:594:LEU:CD2	2:G:608:VAL:HG22	2.51	0.41
2:G:602:GLU:N	2:G:602:GLU:OE1	2.54	0.41
2:H:427:ARG:HA	2:H:461:PHE:O	2.21	0.41
2:H:555:VAL:O	2:H:559:THR:HG23	2.20	0.41
2:H:564:ASN:C	2:H:566:GLY:N	2.78	0.41
1:L:72:G:H2'	1:L:73:U:C6	2.56	0.41
1:O:76:G:H21	1:O:99:C:N4	2.19	0.41
2:A:399:GLY:HA2	2:A:402:ARG:HH11	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:529:PHE:CD1	2:A:529:PHE:N	2.76	0.41
2:A:626:GLU:HG3	2:A:626:GLU:H	1.70	0.41
2:B:254:TYR:CD2	2:B:255:HIS:N	2.89	0.41
2:B:274:PHE:C	2:B:274:PHE:HD1	2.27	0.41
2:B:397:VAL:O	2:B:401:ILE:HG12	2.21	0.41
2:B:482:THR:HG21	4:B:3002:TSB:H5'2	2.02	0.41
2:C:526:THR:O	2:C:530:ALA:HA	2.20	0.41
2:D:277:LEU:O	2:D:280:PHE:HB3	2.20	0.41
2:D:320:THR:HG22	2:D:321:SER:N	2.26	0.41
2:D:390:GLU:H	2:D:390:GLU:HG2	1.71	0.41
2:D:534:PRO:O	2:D:535:THR:C	2.63	0.41
2:E:300:ASP:HB3	2:F:259:GLU:HB2	2.03	0.41
2:E:422:THR:HB	2:E:438:GLU:OE2	2.21	0.41
2:E:427:ARG:HA	2:E:461:PHE:O	2.21	0.41
2:E:558:LEU:C	2:E:560:GLN:N	2.77	0.41
2:F:423:ARG:HB3	2:F:434:TRP:CE3	2.56	0.41
2:G:255:HIS:HE2	2:G:257:GLN:CG	2.33	0.41
2:G:417:VAL:HG13	2:G:471:TYR:CE1	2.46	0.41
2:H:358:PHE:N	2:H:358:PHE:CD1	2.89	0.41
2:H:421:SER:CA	2:H:455:GLN:HB3	2.48	0.41
2:H:572:ASP:OD2	2:H:585:HIS:HE1	2.04	0.41
1:I:87:U:H2'	1:I:88:G:OP2	2.20	0.40
1:L:80:U:H2'	1:L:81:C:H6	1.85	0.40
1:M:69:G:P	1:P:105:C:C2'	3.09	0.40
2:A:298:MET:HE2	2:A:327:TYR:CD1	2.56	0.40
2:A:298:MET:HE2	2:A:327:TYR:HB2	2.03	0.40
2:A:421:SER:OG	2:A:467:GLU:OE2	2.39	0.40
2:A:441:LEU:HD23	2:A:441:LEU:HA	1.91	0.40
2:B:298:MET:CE	2:B:327:TYR:HB2	2.51	0.40
2:C:323:GLU:CD	2:C:323:GLU:N	2.79	0.40
2:C:462:TYR:HE2	2:C:511:HIS:CD2	2.38	0.40
2:D:244:HIS:ND1	2:D:374:MET:HE2	2.36	0.40
2:D:582:ILE:HA	2:D:593:MET:HE2	2.02	0.40
2:E:294:LYS:HA	2:E:294:LYS:HD2	1.85	0.40
2:F:424:PRO:HD2	2:F:434:TRP:CH2	2.55	0.40
2:G:294:LYS:HA	2:G:294:LYS:HD2	1.76	0.40
2:G:545:MET:CE	2:G:593:MET:HB3	2.51	0.40
2:H:426:LYS:HE2	2:H:426:LYS:HB3	1.90	0.40
2:H:562:LEU:HB3	2:H:569:VAL:HG11	2.02	0.40
2:A:353:LEU:HD22	2:A:353:LEU:HA	1.83	0.40
2:A:627:LYS:O	2:A:630:GLN:HB3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:380:THR:N	2:B:519:GLU:OE2	2.54	0.40
2:B:394:ARG:O	2:B:395:ASP:C	2.64	0.40
2:B:436:ARG:HG2	2:B:440:ASP:OD2	2.21	0.40
2:C:353:LEU:HD22	2:C:353:LEU:HA	1.81	0.40
2:C:499:GLY:HA3	2:C:501:ASP:OD1	2.20	0.40
2:C:567:ILE:O	2:C:569:VAL:N	2.54	0.40
2:D:313:TYR:CE1	2:D:316:ALA:HB3	2.56	0.40
2:D:386:ILE:O	2:D:509:MET:HA	2.21	0.40
2:D:625:ILE:O	2:D:626:GLU:C	2.65	0.40
2:E:362:HIS:CD2	5:E:6011:HOH:O	2.70	0.40
2:E:430:SER:O	2:E:433:MET:HB2	2.20	0.40
2:E:437:ALA:HB2	2:E:487:PHE:CD2	2.56	0.40
2:E:501:ASP:O	2:E:503:GLU:HG2	2.22	0.40
2:F:564:ASN:C	2:F:566:GLY:N	2.78	0.40
2:G:247:ILE:HD12	2:G:526:THR:HG22	2.02	0.40
2:G:567:ILE:O	2:G:569:VAL:HG12	2.22	0.40
2:H:339:GLN:OE1	2:H:339:GLN:CA	2.69	0.40
1:P:93:U:C2	1:P:94:C:C6	3.10	0.40
2:A:609:ARG:HD3	2:A:610:THR:O	2.22	0.40
2:B:538:ALA:O	2:B:539:PRO:C	2.63	0.40
2:B:602:GLU:N	2:B:602:GLU:CD	2.80	0.40
2:C:356:ALA:HB2	2:C:384:ALA:HB2	2.03	0.40
2:C:426:LYS:HE2	2:C:426:LYS:HB3	1.89	0.40
2:D:277:LEU:CB	2:D:518:MET:HE1	2.52	0.40
2:D:430:SER:O	2:D:433:MET:HB2	2.21	0.40
2:D:539:PRO:O	2:D:568:ARG:HD2	2.21	0.40
2:D:576:GLU:HG2	2:D:580:PHE:CD2	2.56	0.40
2:E:488:SER:O	2:E:489:LEU:C	2.64	0.40
2:F:400:CYS:SG	2:F:512:ARG:HG3	2.61	0.40
2:G:544:ILE:CG2	2:G:555:VAL:HG13	2.49	0.40
2:H:281:VAL:O	2:H:282:ARG:C	2.64	0.40
2:H:288:TYR:CZ	2:H:403:LEU:HD13	2.57	0.40
2:H:437:ALA:HA	2:H:485:LEU:HD23	2.03	0.40
1:K:79:A:OP1	2:D:349:ARG:NH2	2.54	0.40
1:M:84:U:O2'	1:M:85:C:P	2.79	0.40
1:O:92:G:H2'	1:O:93:U:C6	2.55	0.40
1:P:95:A:H2'	1:P:96:C:O4'	2.20	0.40
2:A:393:ILE:HD11	2:A:508:VAL:HG11	2.02	0.40
2:A:545:MET:SD	2:A:581:LYS:HB3	2.61	0.40
2:B:249:LYS:HE2	2:B:249:LYS:HA	2.03	0.40
2:B:356:ALA:HA	2:B:384:ALA:HA	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:358:PHE:N	2:B:358:PHE:HD1	2.19	0.40
2:B:583:ARG:O	2:B:586:THR:N	2.54	0.40
2:C:423:ARG:HG2	2:C:423:ARG:O	2.21	0.40
2:C:554:TYR:CE1	2:C:558:LEU:HD11	2.56	0.40
2:D:356:ALA:HA	2:D:384:ALA:CB	2.52	0.40
2:D:427:ARG:HA	2:D:461:PHE:O	2.22	0.40
2:E:608:VAL:CB	2:E:619:MET:HE3	2.51	0.40
2:F:526:THR:O	2:F:527:GLU:C	2.65	0.40
2:G:349:ARG:HE	2:G:349:ARG:HB2	1.64	0.40
2:G:554:TYR:CE1	2:G:558:LEU:HD11	2.56	0.40
2:H:251:LEU:CA	2:H:588:ARG:HH12	2.27	0.40
2:H:348:TYR:CE2	2:H:500:GLU:HG3	2.56	0.40
2:H:488:SER:O	2:H:489:LEU:C	2.65	0.40
2:A:415:LYS:HG2	2:A:471:TYR:CD1	2.57	0.40
2:A:545:MET:HE3	2:A:593:MET:HB3	2.04	0.40
2:C:334:CYS:O	2:C:335:PRO:C	2.63	0.40
2:D:323:GLU:O	2:D:324:ASN:O	2.39	0.40
2:D:480:CYS:O	2:D:516:GLY:HA3	2.22	0.40
2:D:550:SER:C	2:D:552:SER:N	2.80	0.40
2:E:490:PRO:HD3	2:E:509:MET:CE	2.51	0.40
2:E:548:THR:HG22	2:E:551:GLN:HE22	1.86	0.40
2:F:254:TYR:CE2	2:F:256:MET:HE2	2.57	0.40
2:F:294:LYS:HA	2:F:294:LYS:HD2	1.83	0.40
2:F:356:ALA:CB	2:F:384:ALA:HB2	2.51	0.40
2:G:482:THR:O	2:G:512:ARG:HA	2.21	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
2	A	399/401 (100%)	327 (82%)	53 (13%)	19 (5%)	2 16

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	399/401 (100%)	325 (82%)	57 (14%)	17 (4%)	2	18
2	C	399/401 (100%)	327 (82%)	52 (13%)	20 (5%)	1	15
2	D	399/401 (100%)	327 (82%)	54 (14%)	18 (4%)	2	17
2	E	399/401 (100%)	327 (82%)	55 (14%)	17 (4%)	2	18
2	F	399/401 (100%)	327 (82%)	52 (13%)	20 (5%)	1	15
2	G	399/401 (100%)	323 (81%)	55 (14%)	21 (5%)	1	14
2	H	399/401 (100%)	323 (81%)	55 (14%)	21 (5%)	1	14
All	All	3192/3208 (100%)	2606 (82%)	433 (14%)	153 (5%)	2	16

All (153) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	313	TYR
2	A	324	ASN
2	A	456	LEU
2	A	528	GLU
2	A	574	ARG
2	B	313	TYR
2	B	324	ASN
2	B	456	LEU
2	B	528	GLU
2	C	313	TYR
2	C	324	ASN
2	C	456	LEU
2	C	528	GLU
2	D	313	TYR
2	D	324	ASN
2	D	456	LEU
2	D	528	GLU
2	E	313	TYR
2	E	324	ASN
2	E	456	LEU
2	E	528	GLU
2	E	574	ARG
2	F	313	TYR
2	F	324	ASN
2	F	456	LEU
2	F	528	GLU
2	G	313	TYR
2	G	324	ASN

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Mol	Chain	Res	Type
2	G	456	LEU
2	G	528	GLU
2	G	529	PHE
2	H	313	TYR
2	H	324	ASN
2	H	456	LEU
2	H	528	GLU
2	A	394	ARG
2	A	494	SER
2	A	529	PHE
2	B	394	ARG
2	B	551	GLN
2	B	574	ARG
2	C	394	ARG
2	C	494	SER
2	C	529	PHE
2	C	574	ARG
2	C	617	GLY
2	D	394	ARG
2	D	529	PHE
2	D	574	ARG
2	D	614	LYS
2	E	394	ARG
2	F	394	ARG
2	F	494	SER
2	F	529	PHE
2	F	574	ARG
2	F	614	LYS
2	G	394	ARG
2	G	574	ARG
2	G	617	GLY
2	H	394	ARG
2	H	494	SER
2	H	529	PHE
2	H	574	ARG
2	H	614	LYS
2	A	254	TYR
2	A	614	LYS
2	B	254	TYR
2	B	494	SER
2	C	614	LYS
2	D	254	TYR

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Mol	Chain	Res	Type
2	D	494	SER
2	D	617	GLY
2	E	254	TYR
2	E	494	SER
2	E	529	PHE
2	E	551	GLN
2	E	614	LYS
2	E	617	GLY
2	F	617	GLY
2	G	254	TYR
2	G	494	SER
2	G	551	GLN
2	G	568	ARG
2	G	614	LYS
2	H	294	LYS
2	H	551	GLN
2	A	551	GLN
2	A	617	GLY
2	B	583	ARG
2	B	617	GLY
2	C	254	TYR
2	C	551	GLN
2	D	551	GLN
2	F	254	TYR
2	F	551	GLN
2	G	331	PRO
2	G	372	GLY
2	G	583	ARG
2	H	254	TYR
2	H	331	PRO
2	H	617	GLY
2	A	294	LYS
2	A	331	PRO
2	A	501	ASP
2	B	331	PRO
2	B	401	ILE
2	B	614	LYS
2	C	331	PRO
2	C	445	LEU
2	D	331	PRO
2	D	501	ASP
2	D	583	ARG

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Mol	Chain	Res	Type
2	E	331	PRO
2	E	501	ASP
2	E	583	ARG
2	F	445	LEU
2	G	298	MET
2	G	501	ASP
2	H	568	ARG
2	B	501	ASP
2	C	568	ARG
2	C	583	ARG
2	F	331	PRO
2	F	372	GLY
2	F	501	ASP
2	H	293	VAL
2	H	445	LEU
2	A	423	ARG
2	A	489	LEU
2	B	293	VAL
2	C	293	VAL
2	C	372	GLY
2	C	423	ARG
2	D	372	GLY
2	A	293	VAL
2	A	625	ILE
2	D	293	VAL
2	D	423	ARG
2	E	293	VAL
2	E	423	ARG
2	F	423	ARG
2	H	423	ARG
2	H	489	LEU
2	B	372	GLY
2	F	293	VAL
2	F	489	LEU
2	G	293	VAL
2	G	625	ILE
2	H	372	GLY
2	H	625	ILE
2	C	625	ILE
2	F	625	ILE
2	G	489	LEU

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	356/356 (100%)	313 (88%)	43 (12%)	5	22
2	B	356/356 (100%)	314 (88%)	42 (12%)	5	23
2	C	356/356 (100%)	317 (89%)	39 (11%)	6	26
2	D	356/356 (100%)	314 (88%)	42 (12%)	5	23
2	E	356/356 (100%)	316 (89%)	40 (11%)	6	25
2	F	356/356 (100%)	314 (88%)	42 (12%)	5	23
2	G	356/356 (100%)	315 (88%)	41 (12%)	5	24
2	H	356/356 (100%)	314 (88%)	42 (12%)	5	23
All	All	2848/2848 (100%)	2517 (88%)	331 (12%)	5	24

All (331) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	258	GLU
2	A	259	GLU
2	A	279	VAL
2	A	288	TYR
2	A	296	PRO
2	A	297	PHE
2	A	303	LEU
2	A	312	ASN
2	A	315	ASP
2	A	317	MET
2	A	323	GLU
2	A	325	ARG
2	A	327	TYR
2	A	333	ASN
2	A	334	CYS
2	A	353	LEU
2	A	390	GLU
2	A	391	GLU
2	A	395	ASP

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Mol	Chain	Res	Type
2	A	396	GLU
2	A	406	ASP
2	A	411	PHE
2	A	417	VAL
2	A	432	GLU
2	A	445	LEU
2	A	458	GLU
2	A	491	SER
2	A	494	SER
2	A	507	PRO
2	A	512	ARG
2	A	519	GLU
2	A	522	ILE
2	A	526	THR
2	A	528	GLU
2	A	537	LEU
2	A	539	PRO
2	A	577	LYS
2	A	587	LEU
2	A	602	GLU
2	A	606	VAL
2	A	609	ARG
2	A	638	LYS
2	A	641	GLU
2	B	258	GLU
2	B	259	GLU
2	B	279	VAL
2	B	288	TYR
2	B	296	PRO
2	B	297	PHE
2	B	298	MET
2	B	303	LEU
2	B	312	ASN
2	B	315	ASP
2	B	320	THR
2	B	323	GLU
2	B	327	TYR
2	B	333	ASN
2	B	353	LEU
2	B	390	GLU
2	B	391	GLU
2	B	395	ASP

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Mol	Chain	Res	Type
2	B	396	GLU
2	B	406	ASP
2	B	411	PHE
2	B	417	VAL
2	B	432	GLU
2	B	445	LEU
2	B	458	GLU
2	B	491	SER
2	B	494	SER
2	B	512	ARG
2	B	519	GLU
2	B	522	ILE
2	B	526	THR
2	B	528	GLU
2	B	537	LEU
2	B	539	PRO
2	B	577	LYS
2	B	587	LEU
2	B	602	GLU
2	B	606	VAL
2	B	609	ARG
2	B	624	VAL
2	B	638	LYS
2	B	641	GLU
2	C	258	GLU
2	C	259	GLU
2	C	279	VAL
2	C	288	TYR
2	C	296	PRO
2	C	297	PHE
2	C	303	LEU
2	C	312	ASN
2	C	315	ASP
2	C	323	GLU
2	C	327	TYR
2	C	333	ASN
2	C	353	LEU
2	C	390	GLU
2	C	391	GLU
2	C	395	ASP
2	C	396	GLU
2	C	406	ASP

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Mol	Chain	Res	Type
2	C	411	PHE
2	C	417	VAL
2	C	432	GLU
2	C	445	LEU
2	C	458	GLU
2	C	491	SER
2	C	494	SER
2	C	512	ARG
2	C	519	GLU
2	C	522	ILE
2	C	526	THR
2	C	537	LEU
2	C	539	PRO
2	C	551	GLN
2	C	577	LYS
2	C	587	LEU
2	C	602	GLU
2	C	606	VAL
2	C	609	ARG
2	C	638	LYS
2	C	641	GLU
2	D	258	GLU
2	D	259	GLU
2	D	279	VAL
2	D	288	TYR
2	D	296	PRO
2	D	297	PHE
2	D	303	LEU
2	D	312	ASN
2	D	315	ASP
2	D	317	MET
2	D	320	THR
2	D	323	GLU
2	D	327	TYR
2	D	333	ASN
2	D	334	CYS
2	D	353	LEU
2	D	390	GLU
2	D	391	GLU
2	D	395	ASP
2	D	396	GLU
2	D	406	ASP

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Mol	Chain	Res	Type
2	D	411	PHE
2	D	417	VAL
2	D	432	GLU
2	D	445	LEU
2	D	458	GLU
2	D	491	SER
2	D	494	SER
2	D	512	ARG
2	D	519	GLU
2	D	522	ILE
2	D	526	THR
2	D	528	GLU
2	D	537	LEU
2	D	539	PRO
2	D	577	LYS
2	D	587	LEU
2	D	602	GLU
2	D	606	VAL
2	D	609	ARG
2	D	638	LYS
2	D	641	GLU
2	E	258	GLU
2	E	259	GLU
2	E	279	VAL
2	E	288	TYR
2	E	297	PHE
2	E	303	LEU
2	E	312	ASN
2	E	315	ASP
2	E	320	THR
2	E	323	GLU
2	E	327	TYR
2	E	333	ASN
2	E	334	CYS
2	E	353	LEU
2	E	390	GLU
2	E	391	GLU
2	E	395	ASP
2	E	396	GLU
2	E	406	ASP
2	E	411	PHE
2	E	417	VAL

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Mol	Chain	Res	Type
2	E	432	GLU
2	E	445	LEU
2	E	458	GLU
2	E	491	SER
2	E	494	SER
2	E	512	ARG
2	E	519	GLU
2	E	522	ILE
2	E	526	THR
2	E	528	GLU
2	E	537	LEU
2	E	539	PRO
2	E	577	LYS
2	E	587	LEU
2	E	602	GLU
2	E	606	VAL
2	E	609	ARG
2	E	638	LYS
2	E	641	GLU
2	F	258	GLU
2	F	259	GLU
2	F	279	VAL
2	F	288	TYR
2	F	297	PHE
2	F	303	LEU
2	F	312	ASN
2	F	315	ASP
2	F	317	MET
2	F	323	GLU
2	F	327	TYR
2	F	333	ASN
2	F	334	CYS
2	F	353	LEU
2	F	390	GLU
2	F	391	GLU
2	F	395	ASP
2	F	396	GLU
2	F	406	ASP
2	F	411	PHE
2	F	417	VAL
2	F	432	GLU
2	F	445	LEU

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Mol	Chain	Res	Type
2	F	458	GLU
2	F	491	SER
2	F	494	SER
2	F	507	PRO
2	F	512	ARG
2	F	519	GLU
2	F	522	ILE
2	F	526	THR
2	F	528	GLU
2	F	537	LEU
2	F	539	PRO
2	F	551	GLN
2	F	577	LYS
2	F	587	LEU
2	F	602	GLU
2	F	606	VAL
2	F	609	ARG
2	F	638	LYS
2	F	641	GLU
2	G	258	GLU
2	G	259	GLU
2	G	288	TYR
2	G	297	PHE
2	G	298	MET
2	G	303	LEU
2	G	312	ASN
2	G	315	ASP
2	G	317	MET
2	G	323	GLU
2	G	327	TYR
2	G	333	ASN
2	G	334	CYS
2	G	353	LEU
2	G	390	GLU
2	G	391	GLU
2	G	395	ASP
2	G	396	GLU
2	G	406	ASP
2	G	411	PHE
2	G	417	VAL
2	G	432	GLU
2	G	445	LEU

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Mol	Chain	Res	Type
2	G	458	GLU
2	G	491	SER
2	G	494	SER
2	G	507	PRO
2	G	512	ARG
2	G	519	GLU
2	G	522	ILE
2	G	526	THR
2	G	528	GLU
2	G	537	LEU
2	G	539	PRO
2	G	577	LYS
2	G	587	LEU
2	G	602	GLU
2	G	606	VAL
2	G	609	ARG
2	G	638	LYS
2	G	641	GLU
2	H	242	ARG
2	H	258	GLU
2	H	259	GLU
2	H	279	VAL
2	H	288	TYR
2	H	297	PHE
2	H	298	MET
2	H	303	LEU
2	H	312	ASN
2	H	315	ASP
2	H	317	MET
2	H	323	GLU
2	H	327	TYR
2	H	333	ASN
2	H	334	CYS
2	H	353	LEU
2	H	390	GLU
2	H	391	GLU
2	H	395	ASP
2	H	396	GLU
2	H	406	ASP
2	H	411	PHE
2	H	417	VAL
2	H	432	GLU

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Mol	Chain	Res	Type
2	H	445	LEU
2	H	458	GLU
2	H	491	SER
2	H	494	SER
2	H	512	ARG
2	H	519	GLU
2	H	522	ILE
2	H	526	THR
2	H	528	GLU
2	H	537	LEU
2	H	539	PRO
2	H	577	LYS
2	H	587	LEU
2	H	602	GLU
2	H	606	VAL
2	H	609	ARG
2	H	638	LYS
2	H	641	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (112) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	250	GLN
2	A	267	HIS
2	A	268	ASN
2	A	342	ASN
2	A	362	HIS
2	A	371	HIS
2	A	449	ASN
2	A	479	GLN
2	A	484	GLN
2	A	556	ASN
2	A	560	GLN
2	A	575	ASN
2	A	585	HIS
2	B	250	GLN
2	B	267	HIS
2	B	268	ASN
2	B	291	GLN
2	B	342	ASN
2	B	362	HIS
2	B	371	HIS

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Mol	Chain	Res	Type
2	B	448	ASN
2	B	479	GLN
2	B	556	ASN
2	B	560	GLN
2	B	575	ASN
2	B	585	HIS
2	B	629	GLN
2	B	639	GLN
2	C	250	GLN
2	C	267	HIS
2	C	291	GLN
2	C	342	ASN
2	C	362	HIS
2	C	371	HIS
2	C	479	GLN
2	C	484	GLN
2	C	556	ASN
2	C	560	GLN
2	C	575	ASN
2	C	585	HIS
2	D	250	GLN
2	D	267	HIS
2	D	268	ASN
2	D	291	GLN
2	D	342	ASN
2	D	362	HIS
2	D	371	HIS
2	D	448	ASN
2	D	479	GLN
2	D	484	GLN
2	D	556	ASN
2	D	560	GLN
2	D	575	ASN
2	D	585	HIS
2	D	629	GLN
2	E	250	GLN
2	E	257	GLN
2	E	267	HIS
2	E	333	ASN
2	E	342	ASN
2	E	362	HIS
2	E	371	HIS

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Mol	Chain	Res	Type
2	E	479	GLN
2	E	484	GLN
2	E	556	ASN
2	E	560	GLN
2	E	575	ASN
2	E	585	HIS
2	E	629	GLN
2	F	250	GLN
2	F	267	HIS
2	F	268	ASN
2	F	291	GLN
2	F	342	ASN
2	F	362	HIS
2	F	371	HIS
2	F	479	GLN
2	F	484	GLN
2	F	556	ASN
2	F	560	GLN
2	F	575	ASN
2	F	585	HIS
2	F	629	GLN
2	G	250	GLN
2	G	257	GLN
2	G	267	HIS
2	G	291	GLN
2	G	342	ASN
2	G	362	HIS
2	G	371	HIS
2	G	448	ASN
2	G	479	GLN
2	G	484	GLN
2	G	556	ASN
2	G	560	GLN
2	G	575	ASN
2	G	585	HIS
2	G	629	GLN
2	H	250	GLN
2	H	257	GLN
2	H	267	HIS
2	H	291	GLN
2	H	342	ASN
2	H	362	HIS

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Mol	Chain	Res	Type
2	H	371	HIS
2	H	479	GLN
2	H	484	GLN
2	H	556	ASN
2	H	560	GLN
2	H	575	ASN
2	H	585	HIS
2	H	639	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	I	36/37 (97%)	6 (16%)	4 (11%)
1	J	36/37 (97%)	9 (25%)	5 (13%)
1	K	36/37 (97%)	11 (30%)	4 (11%)
1	L	36/37 (97%)	7 (19%)	4 (11%)
1	M	36/37 (97%)	8 (22%)	5 (13%)
1	N	36/37 (97%)	8 (22%)	4 (11%)
1	O	36/37 (97%)	9 (25%)	5 (13%)
1	P	36/37 (97%)	8 (22%)	5 (13%)
All	All	288/296 (97%)	66 (22%)	36 (12%)

All (66) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	I	76	G
1	I	85	C
1	I	86	G
1	I	89	U
1	I	90	G
1	I	98	A
1	J	74	A
1	J	75	U
1	J	76	G
1	J	85	C
1	J	86	G
1	J	88	G
1	J	89	U
1	J	90	G
1	J	98	A
1	K	74	A

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Mol	Chain	Res	Type
1	K	76	G
1	K	85	C
1	K	86	G
1	K	87	U
1	K	88	G
1	K	89	U
1	K	90	G
1	K	91	G
1	K	98	A
1	K	105	C
1	L	76	G
1	L	85	C
1	L	86	G
1	L	88	G
1	L	89	U
1	L	90	G
1	L	98	A
1	M	75	U
1	M	76	G
1	M	85	C
1	M	86	G
1	M	88	G
1	M	89	U
1	M	90	G
1	M	98	A
1	N	74	A
1	N	76	G
1	N	85	C
1	N	86	G
1	N	88	G
1	N	89	U
1	N	90	G
1	N	98	A
1	O	74	A
1	O	76	G
1	O	85	C
1	O	86	G
1	O	88	G
1	O	89	U
1	O	90	G
1	O	96	C
1	O	98	A

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Mol	Chain	Res	Type
1	P	75	U
1	P	76	G
1	P	85	C
1	P	86	G
1	P	88	G
1	P	89	U
1	P	90	G
1	P	98	A

All (36) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	I	75	U
1	I	84	U
1	I	88	G
1	I	97	C
1	J	74	A
1	J	75	U
1	J	84	U
1	J	88	G
1	J	97	C
1	K	84	U
1	K	86	G
1	K	88	G
1	K	97	C
1	L	75	U
1	L	84	U
1	L	88	G
1	L	97	C
1	M	74	A
1	M	75	U
1	M	84	U
1	M	88	G
1	M	97	C
1	N	75	U
1	N	84	U
1	N	88	G
1	N	97	C
1	O	74	A
1	O	75	U
1	O	84	U
1	O	88	G

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Mol	Chain	Res	Type
1	O	97	C
1	P	74	A
1	P	75	U
1	P	84	U
1	P	88	G
1	P	97	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 8 are monoatomic - leaving 8 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	TSB	F	7002	3	32,32,32	1.83	3 (9%)	46,48,48	0.93	5 (10%)
4	TSB	D	5002	3	32,32,32	2.30	4 (12%)	46,48,48	0.75	1 (2%)
4	TSB	E	6002	3	32,32,32	2.43	5 (15%)	46,48,48	0.87	2 (4%)
4	TSB	H	9002	3	32,32,32	2.03	3 (9%)	46,48,48	0.96	3 (6%)
4	TSB	G	8002	3	32,32,32	2.24	3 (9%)	46,48,48	1.18	3 (6%)
4	TSB	B	3002	3	32,32,32	2.69	4 (12%)	46,48,48	0.81	3 (6%)
4	TSB	A	2002	3	32,32,32	1.77	4 (12%)	46,48,48	0.93	3 (6%)
4	TSB	C	4002	3	32,32,32	2.06	4 (12%)	46,48,48	0.87	3 (6%)

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	TSB	F	7002	3	-	3/22/39/39	0/3/3/3
4	TSB	D	5002	3	-	4/22/39/39	0/3/3/3
4	TSB	E	6002	3	-	5/22/39/39	0/3/3/3
4	TSB	H	9002	3	-	3/22/39/39	0/3/3/3
4	TSB	G	8002	3	-	3/22/39/39	0/3/3/3
4	TSB	B	3002	3	-	5/22/39/39	0/3/3/3
4	TSB	A	2002	3	-	4/22/39/39	0/3/3/3
4	TSB	C	4002	3	-	3/22/39/39	0/3/3/3

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	3002	TSB	O1S-S1	10.94	1.52	1.42
4	E	6002	TSB	O1S-S1	9.90	1.51	1.42
4	G	8002	TSB	O2S-S1	9.44	1.50	1.42
4	B	3002	TSB	O2S-S1	9.17	1.50	1.42
4	D	5002	TSB	O1S-S1	8.94	1.50	1.42
4	E	6002	TSB	O2S-S1	8.51	1.49	1.42
4	H	9002	TSB	O2S-S1	8.25	1.49	1.42
4	C	4002	TSB	O2S-S1	8.12	1.49	1.42
4	D	5002	TSB	O2S-S1	7.77	1.49	1.42
4	F	7002	TSB	O2S-S1	7.24	1.48	1.42
4	A	2002	TSB	O2S-S1	6.87	1.48	1.42
4	H	9002	TSB	O1S-S1	6.72	1.48	1.42
4	C	4002	TSB	O1S-S1	6.60	1.48	1.42
4	G	8002	TSB	O1S-S1	6.58	1.48	1.42
4	F	7002	TSB	O1S-S1	6.07	1.47	1.42
4	A	2002	TSB	O1S-S1	5.49	1.47	1.42
4	C	4002	TSB	S1-N8	4.30	1.66	1.59
4	B	3002	TSB	S1-N8	3.74	1.65	1.59
4	F	7002	TSB	S1-N8	3.59	1.65	1.59
4	D	5002	TSB	S1-N8	3.53	1.65	1.59
4	D	5002	TSB	CA-C	3.53	1.56	1.53
4	A	2002	TSB	S1-N8	3.27	1.65	1.59
4	H	9002	TSB	S1-N8	3.16	1.65	1.59
4	G	8002	TSB	S1-N8	3.08	1.64	1.59
4	B	3002	TSB	CA-C	2.39	1.55	1.53
4	E	6002	TSB	S1-N8	2.32	1.63	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	6002	TSB	CA-C	2.28	1.55	1.53
4	A	2002	TSB	C-N8	-2.18	1.33	1.37
4	C	4002	TSB	CA-C	2.10	1.55	1.53
4	E	6002	TSB	CB-CA	2.02	1.60	1.53

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	8002	TSB	O-C-CA	-4.64	116.99	120.66
4	G	8002	TSB	C-CA-N	-4.14	104.23	110.31
4	E	6002	TSB	C-CA-N	-4.03	104.40	110.31
4	H	9002	TSB	O-C-CA	-3.69	117.75	120.66
4	B	3002	TSB	C-CA-N	-3.31	105.45	110.31
4	A	2002	TSB	C-CA-N	-3.17	105.66	110.31
4	F	7002	TSB	O-C-CA	-3.12	118.20	120.66
4	A	2002	TSB	O-C-CA	-3.01	118.28	120.66
4	C	4002	TSB	O-C-CA	-2.92	118.35	120.66
4	G	8002	TSB	CA-C-N8	2.75	120.68	114.20
4	C	4002	TSB	C-CA-N	-2.66	106.41	110.31
4	D	5002	TSB	C-CA-N	-2.66	106.41	110.31
4	H	9002	TSB	C-CA-N	-2.63	106.45	110.31
4	B	3002	TSB	O-C-CA	-2.41	118.76	120.66
4	H	9002	TSB	CA-C-N8	2.29	119.59	114.20
4	F	7002	TSB	O5'-C5'-C4'	2.28	111.64	107.57
4	F	7002	TSB	C5'-O5'-S1	2.27	121.74	116.97
4	A	2002	TSB	CA-C-N8	2.22	119.43	114.20
4	F	7002	TSB	C-CA-N	-2.19	107.10	110.31
4	B	3002	TSB	CA-C-N8	2.17	119.31	114.20
4	F	7002	TSB	CA-C-N8	2.11	119.16	114.20
4	C	4002	TSB	C5'-O5'-S1	2.05	121.27	116.97
4	E	6002	TSB	O-C-CA	-2.03	119.06	120.66

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	2002	TSB	C5'-O5'-S1-N8
4	B	3002	TSB	C5'-O5'-S1-N8
4	C	4002	TSB	C5'-O5'-S1-N8
4	D	5002	TSB	C5'-O5'-S1-N8
4	E	6002	TSB	C5'-O5'-S1-N8
4	F	7002	TSB	C5'-O5'-S1-N8

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Mol	Chain	Res	Type	Atoms
4	G	8002	TSB	C5'-O5'-S1-N8
4	H	9002	TSB	C5'-O5'-S1-N8
4	A	2002	TSB	C5'-O5'-S1-O1S
4	B	3002	TSB	C5'-O5'-S1-O1S
4	C	4002	TSB	C5'-O5'-S1-O1S
4	D	5002	TSB	C5'-O5'-S1-O1S
4	E	6002	TSB	C5'-O5'-S1-O1S
4	F	7002	TSB	C5'-O5'-S1-O1S
4	H	9002	TSB	C5'-O5'-S1-O1S
4	B	3002	TSB	C-N8-S1-O2S
4	E	6002	TSB	C-N8-S1-O2S
4	F	7002	TSB	C5'-O5'-S1-O2S
4	G	8002	TSB	C5'-O5'-S1-O1S
4	A	2002	TSB	C-N8-S1-O2S
4	B	3002	TSB	C-N8-S1-O1S
4	D	5002	TSB	C-N8-S1-O2S
4	E	6002	TSB	C-N8-S1-O1S
4	G	8002	TSB	C-N8-S1-O2S
4	A	2002	TSB	C5'-O5'-S1-O2S
4	C	4002	TSB	C5'-O5'-S1-O2S
4	B	3002	TSB	C5'-O5'-S1-O2S
4	D	5002	TSB	C5'-O5'-S1-O2S
4	E	6002	TSB	C5'-O5'-S1-O2S
4	H	9002	TSB	C5'-O5'-S1-O2S

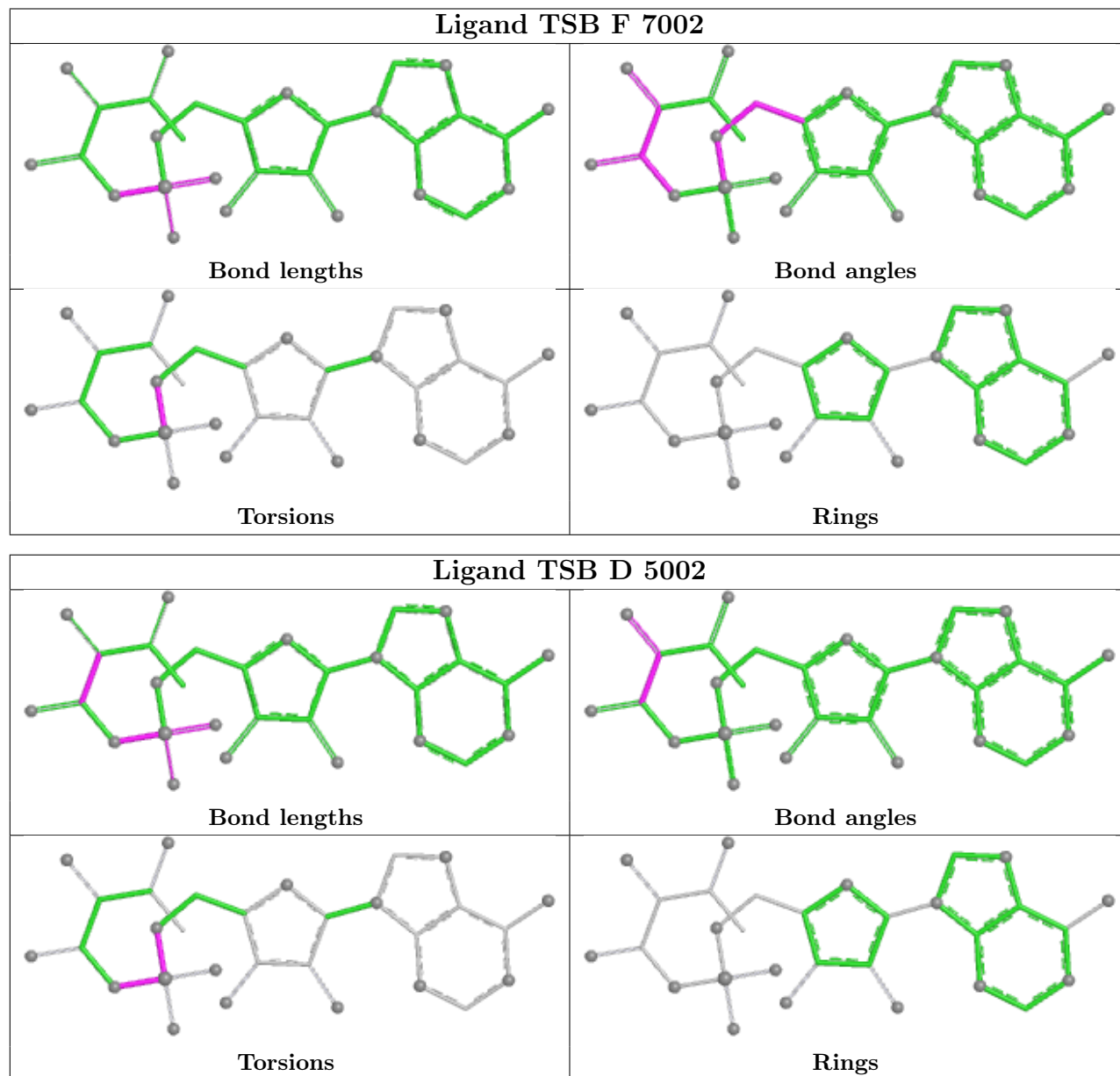
There are no ring outliers.

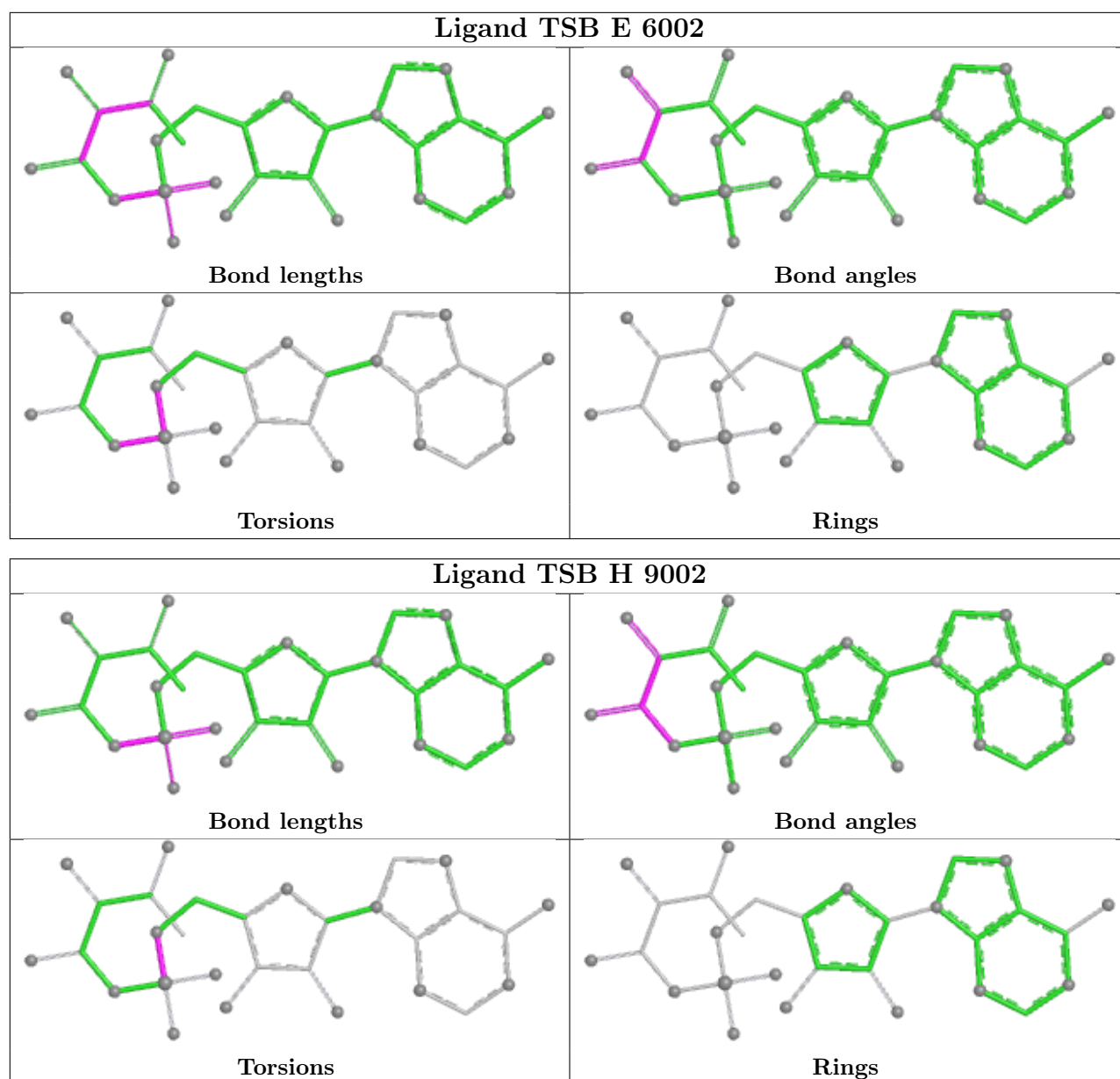
8 monomers are involved in 31 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	F	7002	TSB	5	0
4	D	5002	TSB	4	0
4	E	6002	TSB	4	0
4	H	9002	TSB	4	0
4	G	8002	TSB	4	0
4	B	3002	TSB	4	0
4	A	2002	TSB	3	0
4	C	4002	TSB	3	0

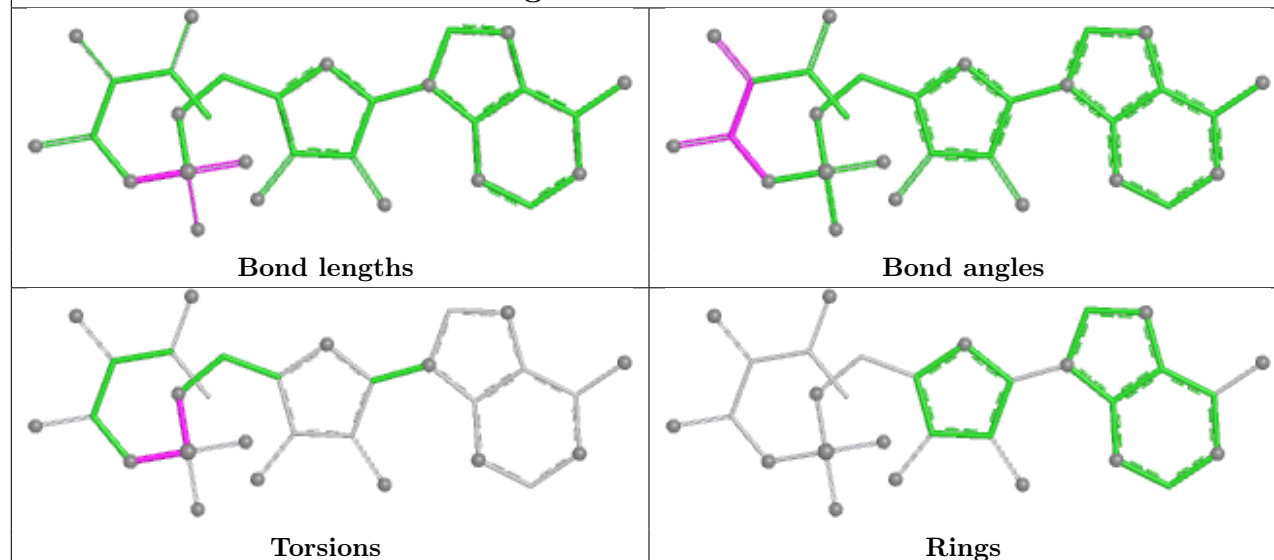
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is

within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

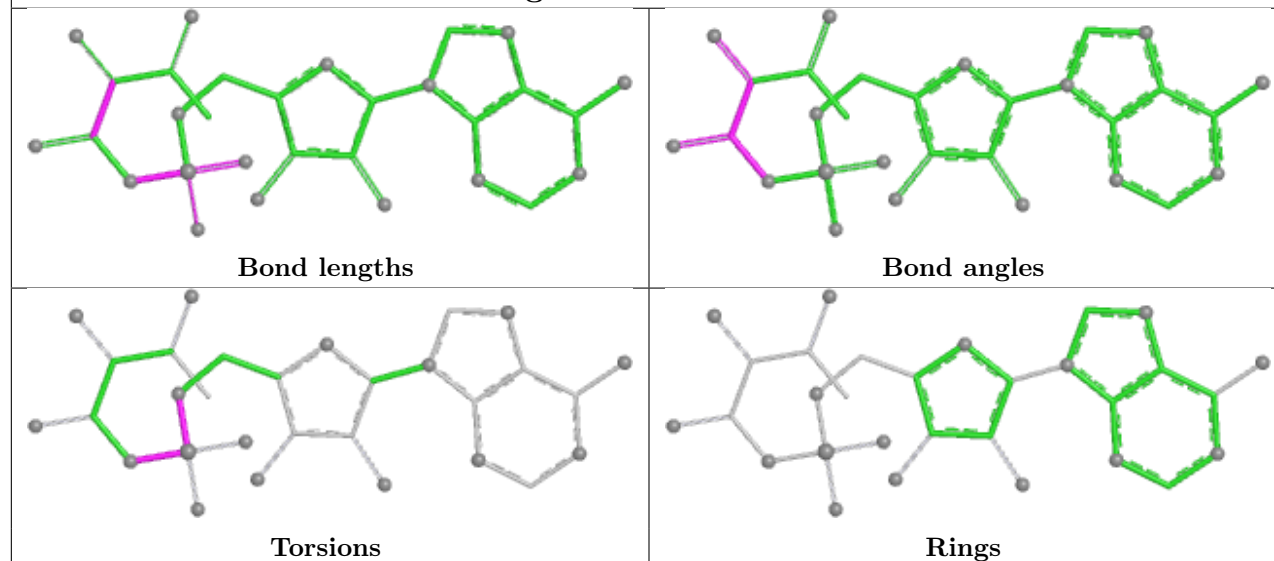


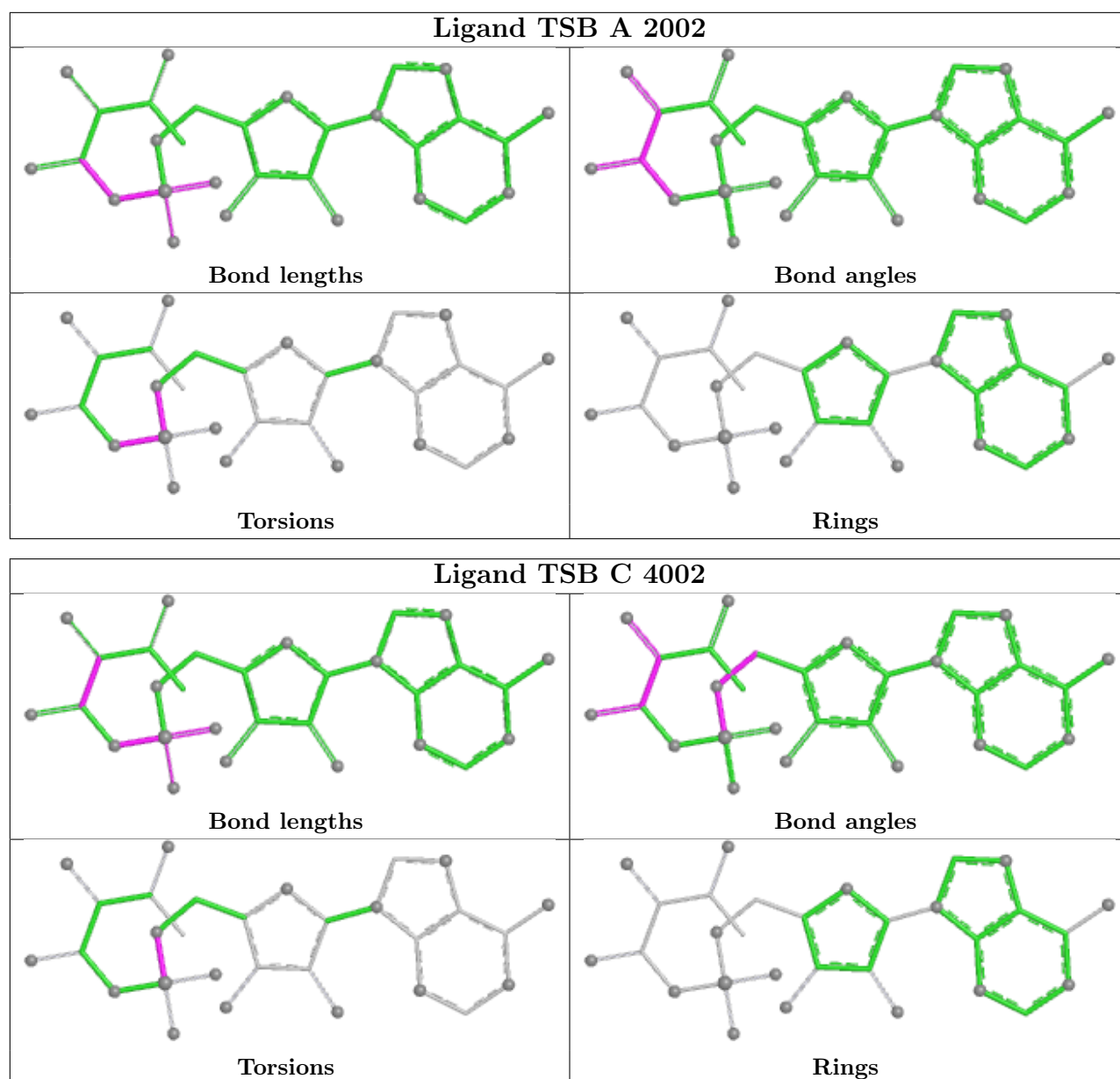


Ligand TSB G 8002



Ligand TSB B 3002





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	I	37/37 (100%)	1.02	4 (10%) 11 8	82, 115, 150, 153	0
1	J	37/37 (100%)	0.45	2 (5%) 31 18	55, 88, 135, 152	0
1	K	37/37 (100%)	1.62	13 (35%) 1 1	119, 148, 168, 184	0
1	L	37/37 (100%)	0.76	1 (2%) 56 33	110, 134, 170, 180	0
1	M	37/37 (100%)	0.49	3 (8%) 18 12	60, 86, 153, 166	0
1	N	37/37 (100%)	1.44	6 (16%) 4 4	140, 163, 181, 184	0
1	O	37/37 (100%)	0.39	1 (2%) 56 33	45, 67, 115, 124	0
1	P	37/37 (100%)	0.49	0 100 100	46, 104, 155, 159	0
2	A	401/401 (100%)	-0.09	1 (0%) 91 82	15, 68, 109, 136	0
2	B	401/401 (100%)	-0.16	1 (0%) 91 82	17, 62, 105, 137	0
2	C	401/401 (100%)	0.38	15 (3%) 45 25	55, 101, 146, 160	0
2	D	401/401 (100%)	0.52	19 (4%) 36 20	80, 123, 147, 158	0
2	E	401/401 (100%)	0.80	31 (7%) 19 13	75, 126, 150, 168	0
2	F	401/401 (100%)	0.51	21 (5%) 33 19	47, 105, 148, 168	0
2	G	401/401 (100%)	-0.31	0 100 100	7, 48, 98, 132	0
2	H	401/401 (100%)	-0.16	0 100 100	14, 68, 105, 135	0
All	All	3504/3504 (100%)	0.24	118 (3%) 48 27	7, 91, 147, 184	0

All (118) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	482	THR	5.5
2	F	260	ALA	5.2
2	E	356	ALA	4.9
2	C	595	VAL	4.7
2	C	578	ILE	4.6

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Mol	Chain	Res	Type	RSRZ
2	C	582	ILE	4.6
2	D	462	TYR	4.4
2	E	345	LEU	4.1
2	E	510	ILE	4.1
1	J	69	G	3.9
2	D	510	ILE	3.8
1	N	97	C	3.8
2	C	586	THR	3.6
2	F	461	PHE	3.6
2	E	485	LEU	3.6
2	E	484	GLN	3.5
2	D	338	VAL	3.4
2	E	309	HIS	3.4
2	D	386	ILE	3.3
2	F	595	VAL	3.3
2	D	485	LEU	3.3
2	D	464	PRO	3.2
2	F	376	VAL	3.2
2	C	596	CYS	3.2
2	E	483	VAL	3.2
2	E	397	VAL	3.1
1	K	76	G	3.1
2	C	543	VAL	3.1
2	F	596	CYS	3.0
2	D	356	ALA	3.0
1	L	85	C	3.0
2	F	621	VAL	3.0
2	E	512	ARG	3.0
2	A	499	GLY	2.9
2	F	372	GLY	2.9
1	K	86	G	2.9
1	K	81	C	2.9
2	D	640	LEU	2.9
2	F	318	PHE	2.9
1	N	98	A	2.8
2	E	468	PHE	2.8
2	F	299	MET	2.8
2	E	335	PRO	2.8
2	F	363	ARG	2.8
1	I	70	G	2.8
1	K	97	C	2.7
1	K	85	C	2.7

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Mol	Chain	Res	Type	RSRZ
2	B	499	GLY	2.7
2	E	493	LEU	2.7
2	E	316	ALA	2.7
2	C	355	MET	2.7
2	F	250	GLN	2.7
2	E	401	ILE	2.6
2	E	469	THR	2.6
2	D	504	ARG	2.6
2	E	504	ARG	2.6
2	E	481	GLY	2.6
1	K	92	G	2.6
1	M	88	G	2.6
2	D	468	PHE	2.6
2	F	259	GLU	2.6
2	F	546	ASN	2.6
2	D	299	MET	2.6
2	C	266	TRP	2.5
2	C	558	LEU	2.5
2	E	299	MET	2.5
2	C	547	ILE	2.5
2	D	328	CYS	2.5
1	K	82	U	2.5
1	N	87	U	2.5
2	D	340	ILE	2.5
2	E	333	ASN	2.4
1	K	94	C	2.4
1	N	80	U	2.4
2	E	373	LEU	2.4
2	E	507	PRO	2.4
2	E	613	GLY	2.4
1	N	85	C	2.4
2	C	576	GLU	2.4
1	I	87	U	2.4
1	K	96	C	2.4
1	K	99	C	2.4
2	D	487	PHE	2.3
2	E	355	MET	2.3
1	J	85	C	2.3
2	F	571	ALA	2.3
2	C	463	GLY	2.3
2	E	388	CYS	2.3
2	D	388	CYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	K	100	U	2.2
1	N	88	G	2.2
2	F	460	ALA	2.2
2	C	592	TYR	2.2
1	M	85	C	2.2
1	K	74	A	2.2
2	C	478	TRP	2.2
2	D	493	LEU	2.2
1	O	105	C	2.1
2	D	482	THR	2.1
2	F	364	ASN	2.1
2	F	543	VAL	2.1
2	E	296	PRO	2.1
2	D	460	ALA	2.1
2	F	578	ILE	2.1
2	E	341	PHE	2.1
1	I	71	C	2.1
1	I	102	C	2.1
2	F	262	GLY	2.1
2	F	545	MET	2.1
2	E	487	PHE	2.1
2	E	566	GLY	2.0
2	E	460	ALA	2.0
2	C	546	ASN	2.0
2	E	534	PRO	2.0
2	D	291	GLN	2.0
2	F	329	ILE	2.0
1	M	105	C	2.0
1	K	93	U	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 oligosaccharides in this entry.

6.4 Ligands ⓘ

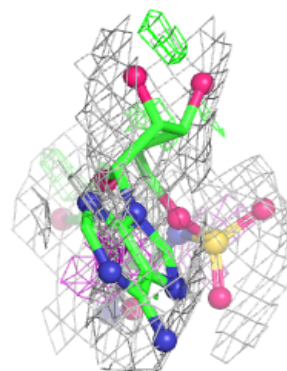
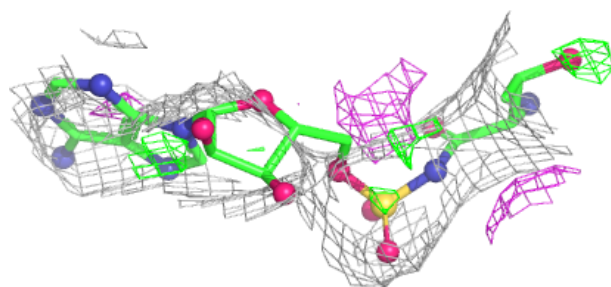
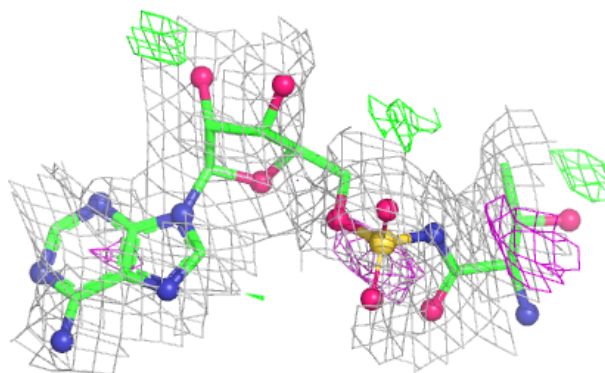
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	TSB	E	6002	30/30	0.71	0.14	86,122,126,127	0
4	TSB	D	5002	30/30	0.81	0.13	94,112,119,120	0
4	TSB	F	7002	30/30	0.86	0.14	89,103,111,112	0
4	TSB	C	4002	30/30	0.91	0.12	73,88,100,102	0
4	TSB	H	9002	30/30	0.93	0.11	30,48,71,72	0
3	ZN	E	1	1/1	0.94	0.06	185,185,185,185	0
3	ZN	D	1	1/1	0.95	0.05	135,135,135,135	0
4	TSB	A	2002	30/30	0.96	0.09	31,44,58,65	0
4	TSB	G	8002	30/30	0.96	0.09	0,18,49,50	0
3	ZN	F	1	1/1	0.96	0.04	101,101,101,101	0
4	TSB	B	3002	30/30	0.97	0.08	32,50,57,58	0
3	ZN	C	1	1/1	0.98	0.03	95,95,95,95	0
3	ZN	H	1	1/1	0.99	0.04	54,54,54,54	0
3	ZN	B	1	1/1	0.99	0.06	52,52,52,52	0
3	ZN	G	1	1/1	0.99	0.02	27,27,27,27	0
3	ZN	A	1	1/1	1.00	0.03	42,42,42,42	0

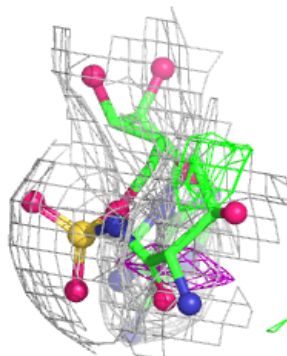
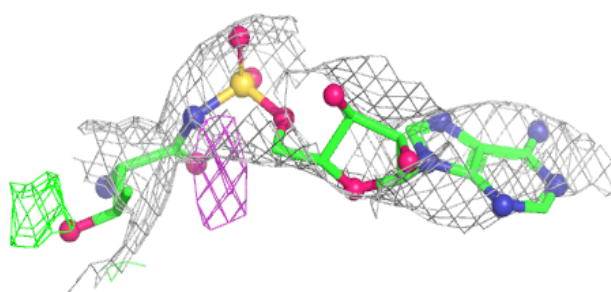
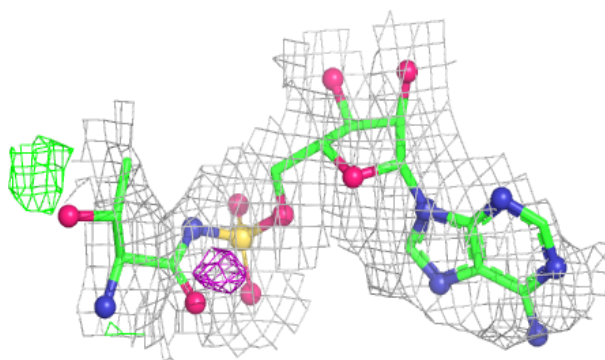
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around TSB E 6002:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

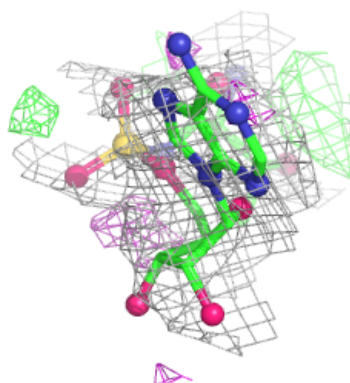
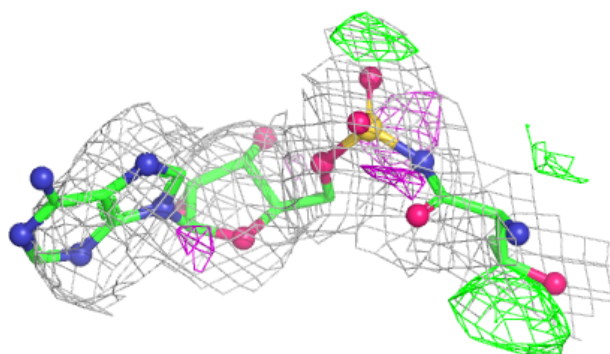
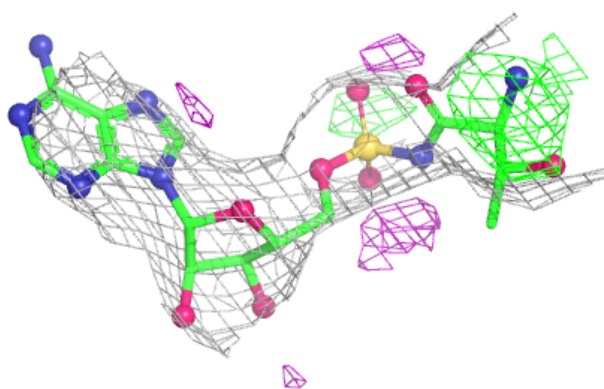
**Electron density around TSB D 5002:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

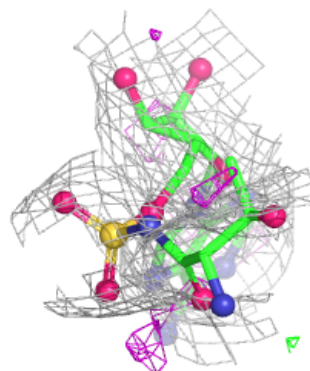
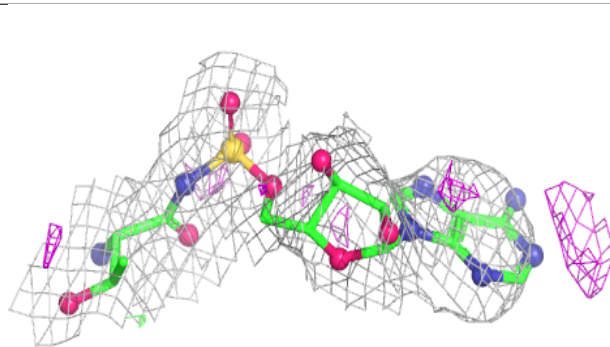
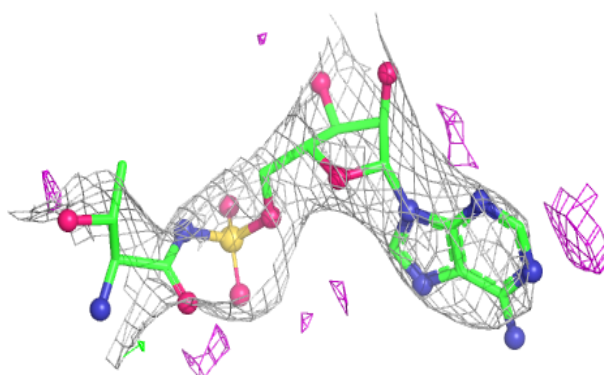


Electron density around TSB F 7002:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

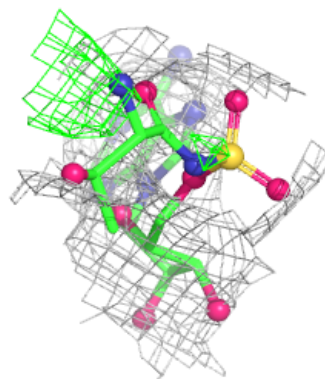
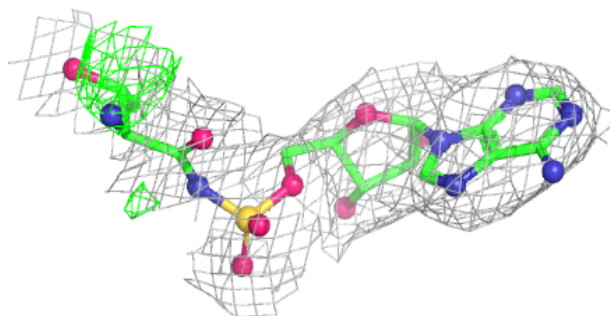
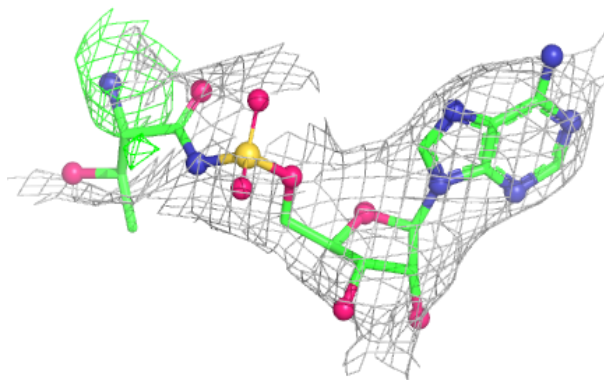
**Electron density around TSB C 4002:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

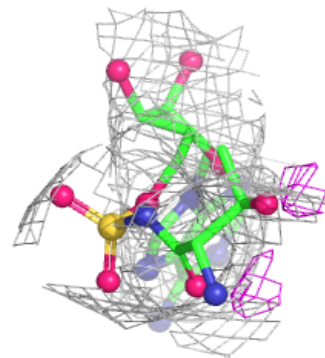
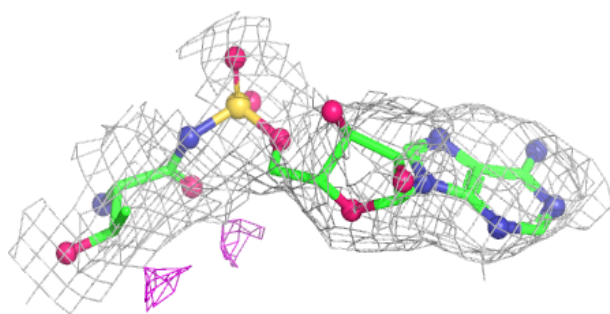
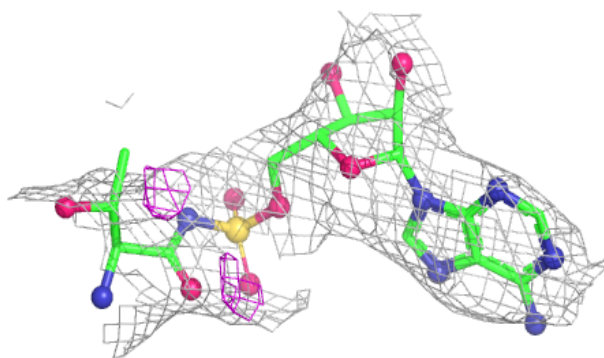


Electron density around TSB H 9002:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

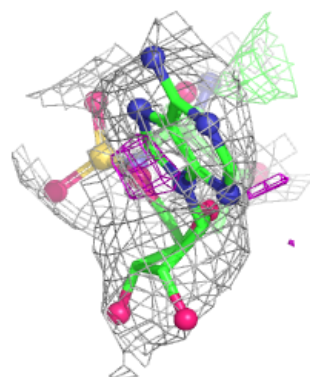
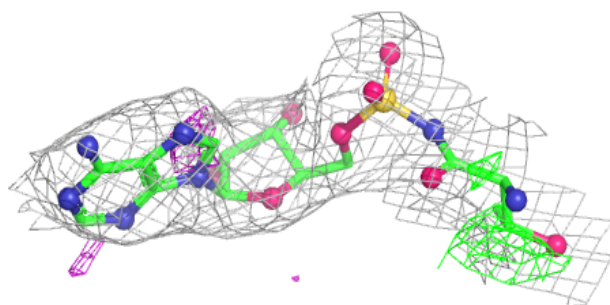
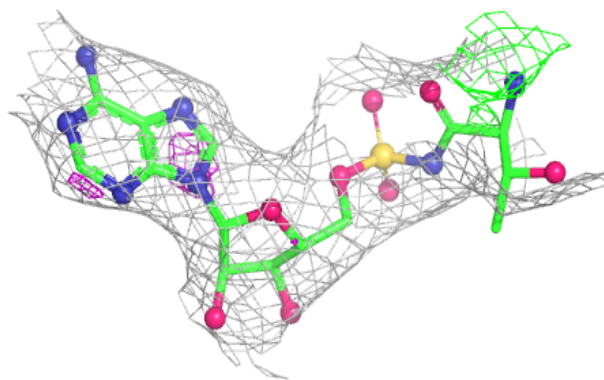
**Electron density around TSB A 2002:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

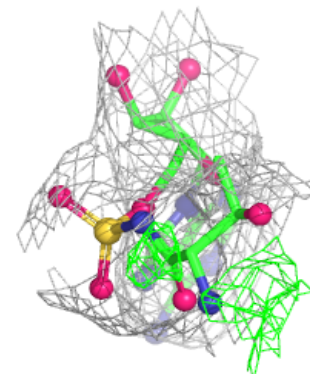
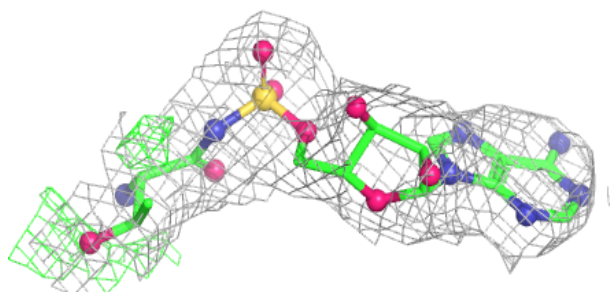
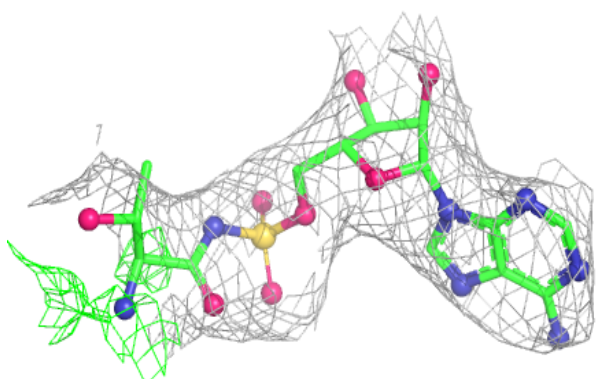


Electron density around TSB G 8002:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around TSB B 3002:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.