



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 12:16 AM GMT

PDB ID : 2YHQ
Title : Structure of respiratory syncytial virus nucleocapsid protein, P1 crystal form
Authors : El Omari, K.; Dhaliwal, B.; Ren, J.; Abrescia, N.G.A.; Lockyer, M.; Powell, K.L.; Hawkins, A.R.; Stammers, D.K.
Deposited on : 2011-05-04
Resolution : 3.60 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

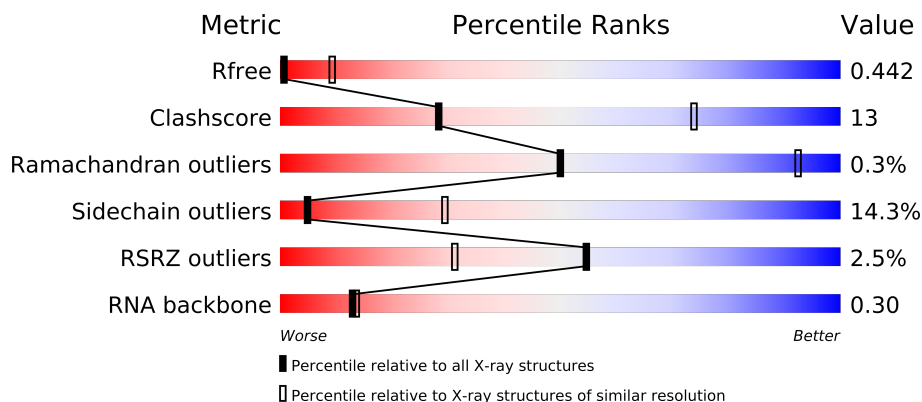
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.60 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.












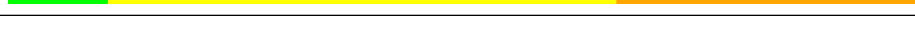
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	1020 (3.86-3.34)
Clashscore	79885	1155 (3.80-3.40)
Ramachandran outliers	78287	1109 (3.80-3.40)
Sidechain outliers	78261	1108 (3.80-3.40)
RSRZ outliers	66119	1000 (3.84-3.36)
RNA backbone	1838	1012 (4.40-2.76)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	375	
1	B	375	
1	C	375	
1	D	375	
1	E	375	
1	F	375	
1	G	375	
1	H	375	
1	I	375	
1	J	375	
1	K	375	
1	L	375	

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Mol	Chain	Length	Quality of chain
1	M	375	
1	N	375	
1	O	375	
1	P	375	
1	Q	375	
1	W	375	
1	Y	375	
1	Z	375	
2	R	70	
2	X	70	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 61200 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	B	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	C	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	D	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	E	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	F	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	G	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	H	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	I	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	J	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	K	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	L	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	M	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	N	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	O	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	P	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	W	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	Y	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			
1	Z	375	Total	C	N	O	S	0	0	0
			2920	1848	506	549	17			

- Molecule 2 is a RNA chain called RNA.

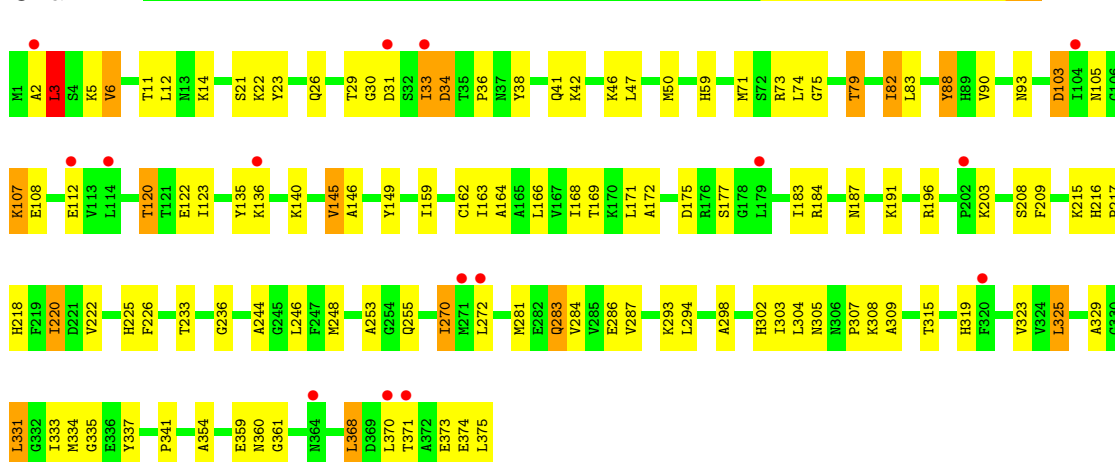
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	R	70	Total	C	N	O	P	0	0	0
			1400	630	210	490	70			
2	X	70	Total	C	N	O	P	0	0	0
			1400	630	210	490	70			

3 Residue-property plots

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

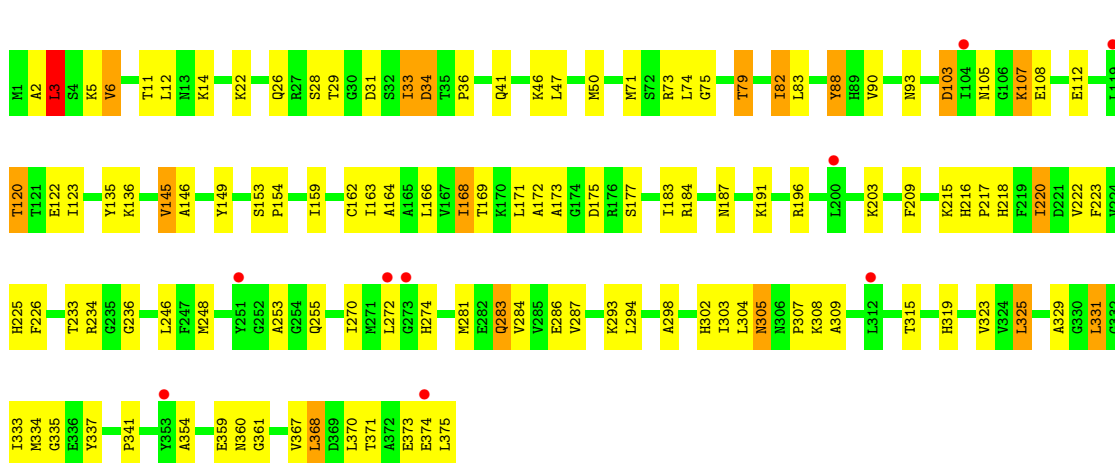
• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

Chain A:



• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

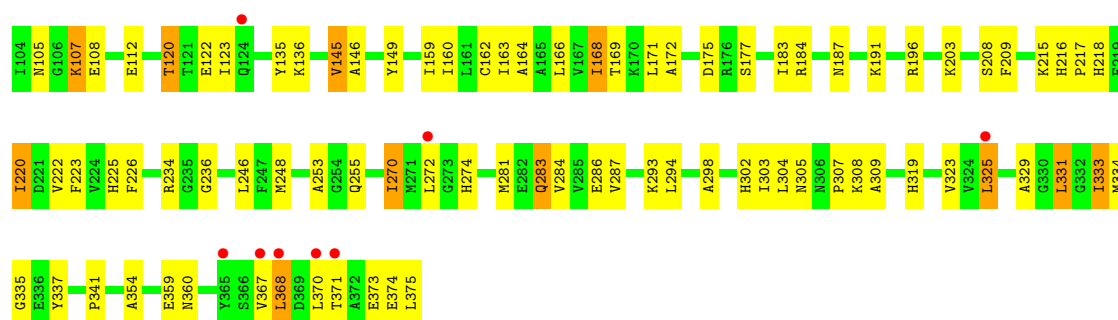
Chain B:



• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

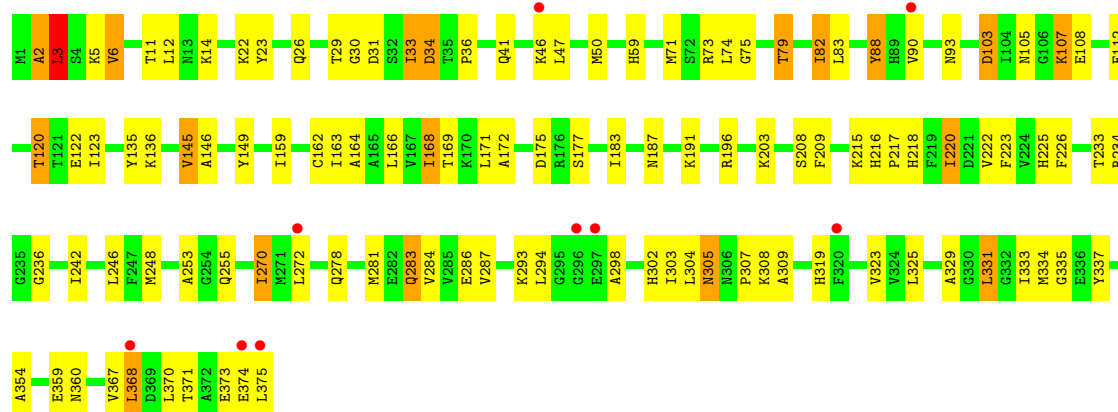
Chain C:





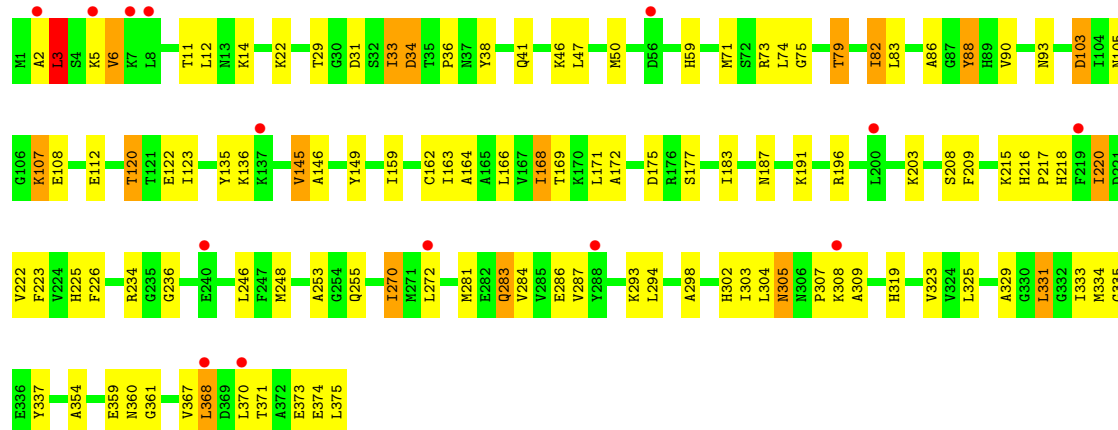
• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

Chain D:



• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

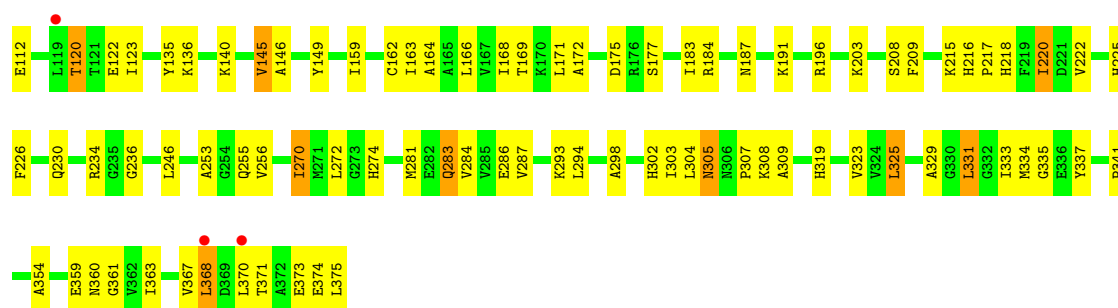
Chain E:



• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

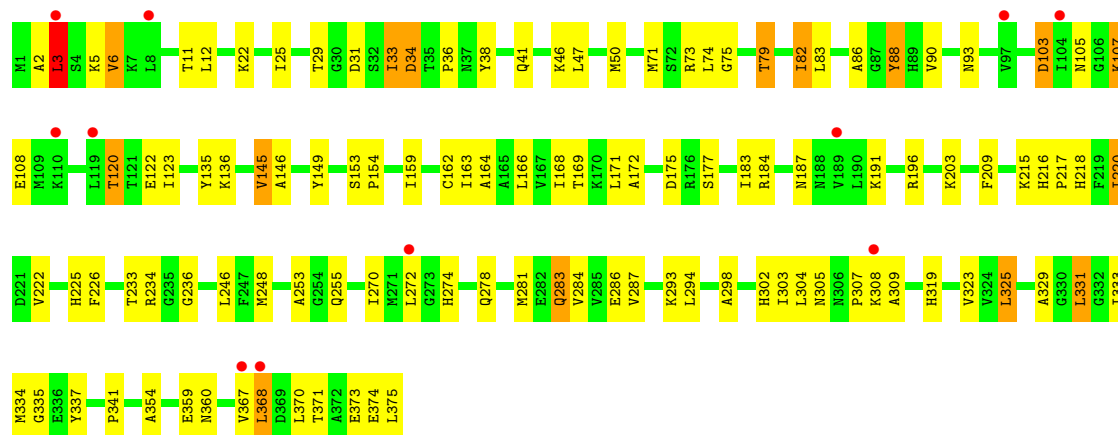
Chain F:





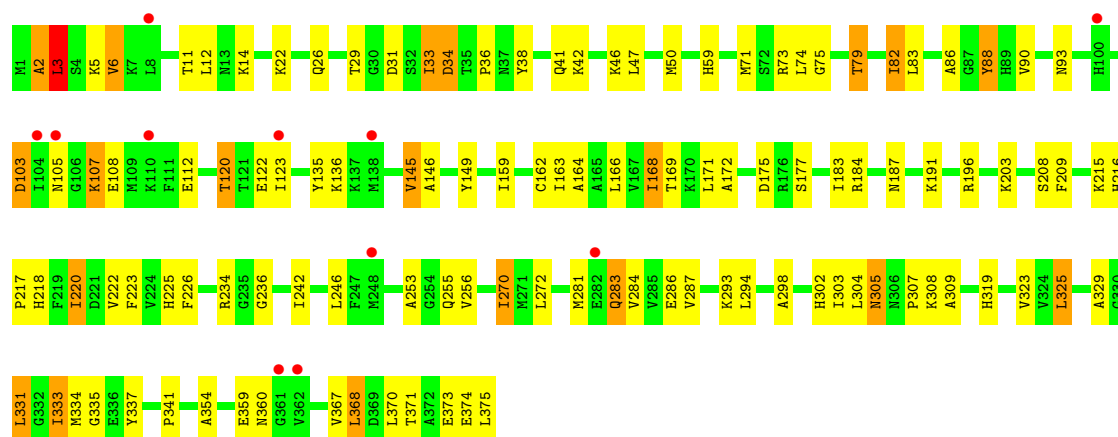
• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

Chain G:



• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

Chain H:



• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

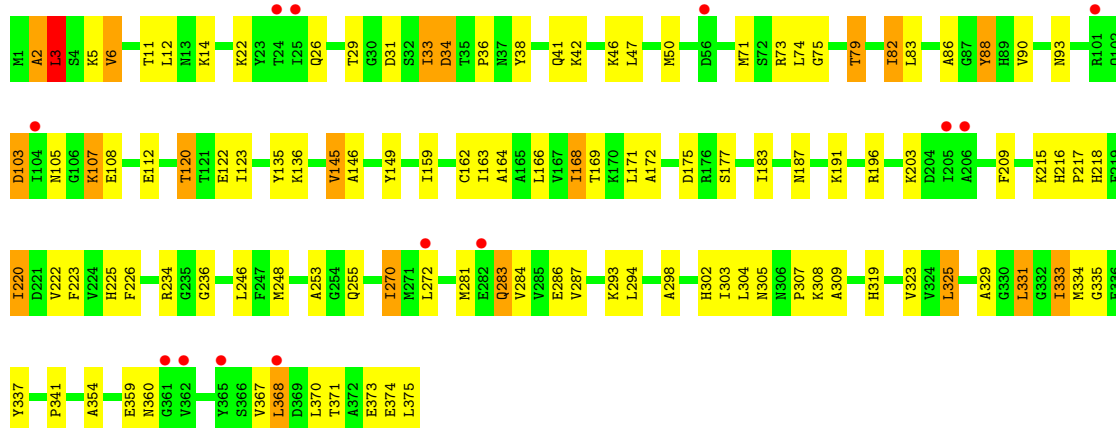
Chain I:





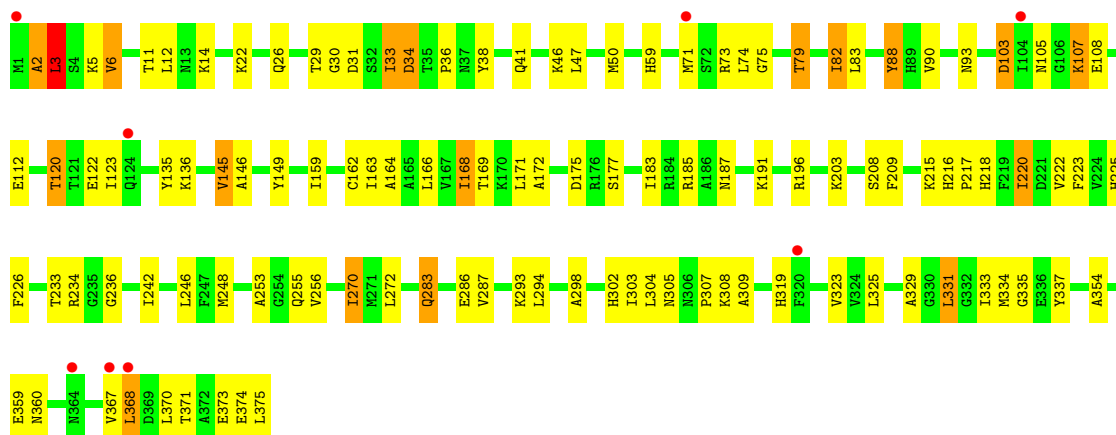
• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

Chain J:



• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

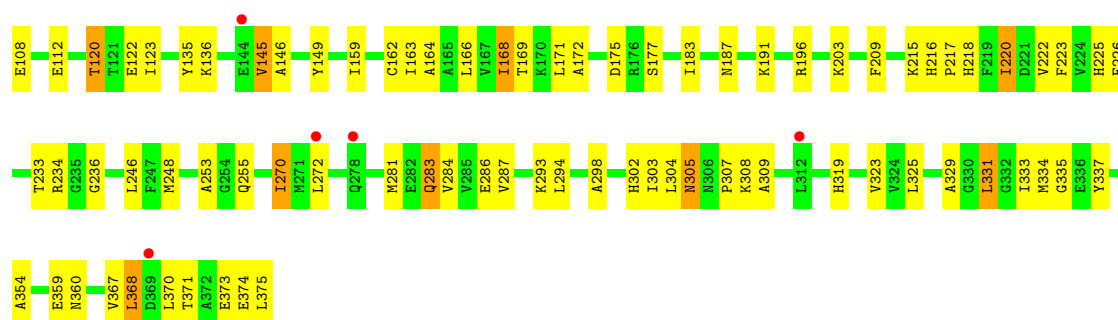
Chain K:



• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

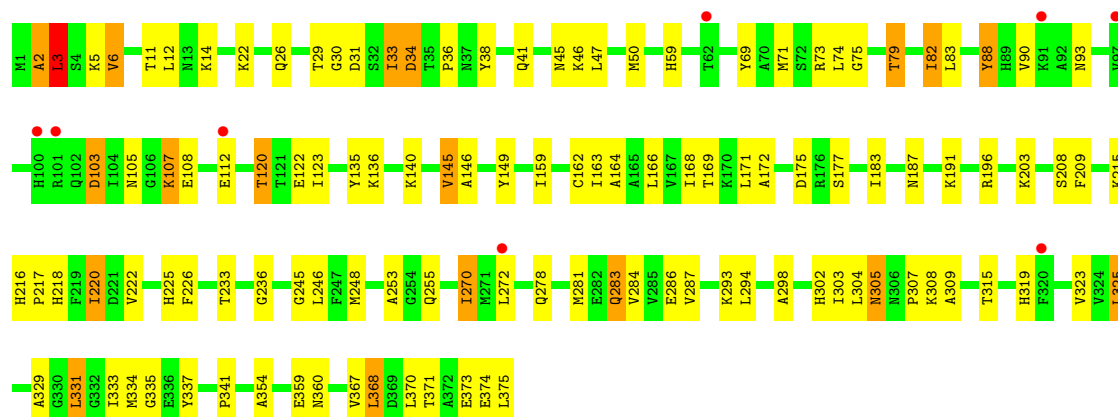
Chain L:





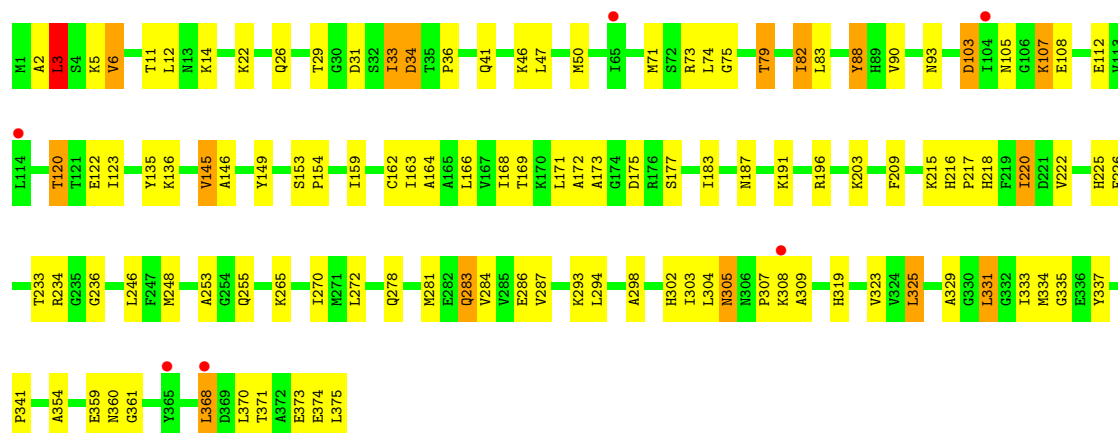
• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

Chain M:



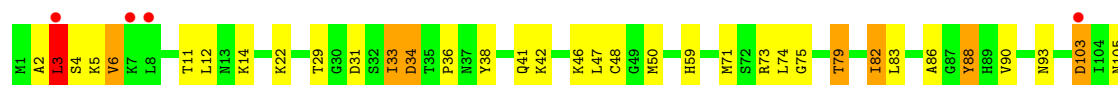
• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

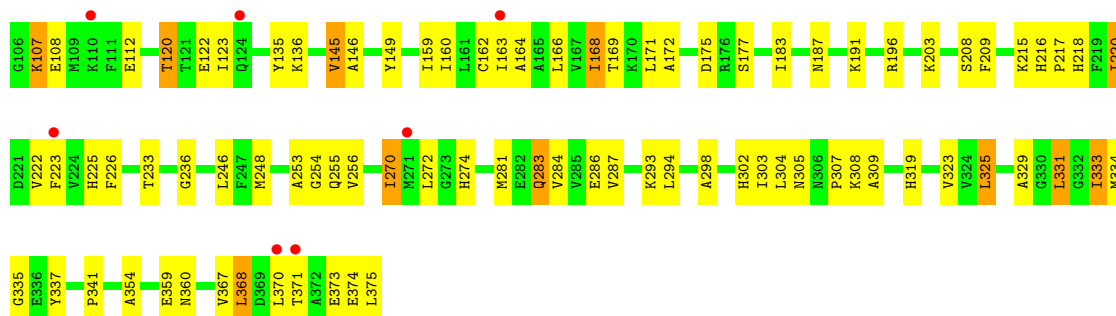
Chain N:



• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

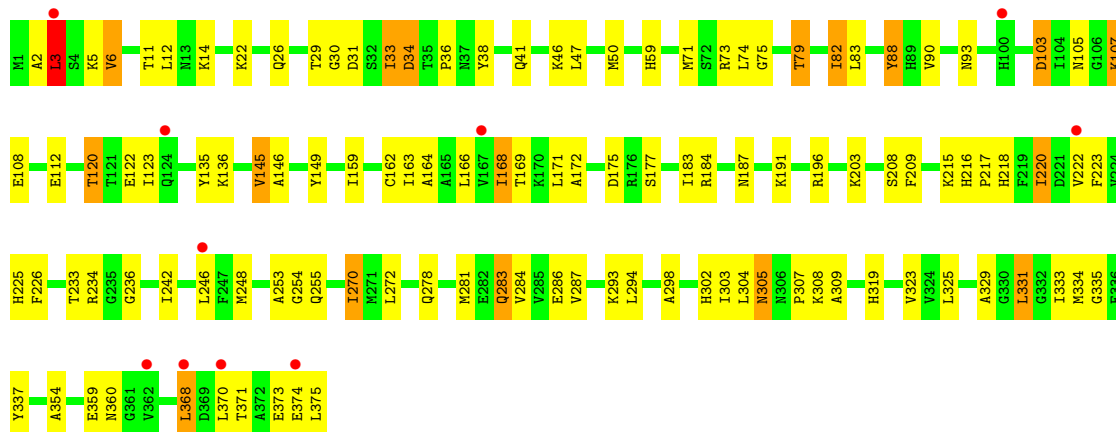
Chain O:





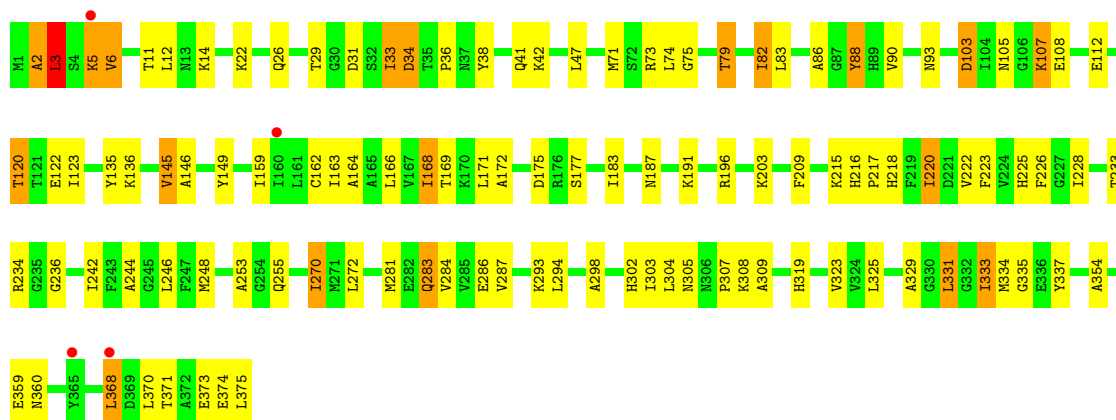
• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

Chain P:



• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

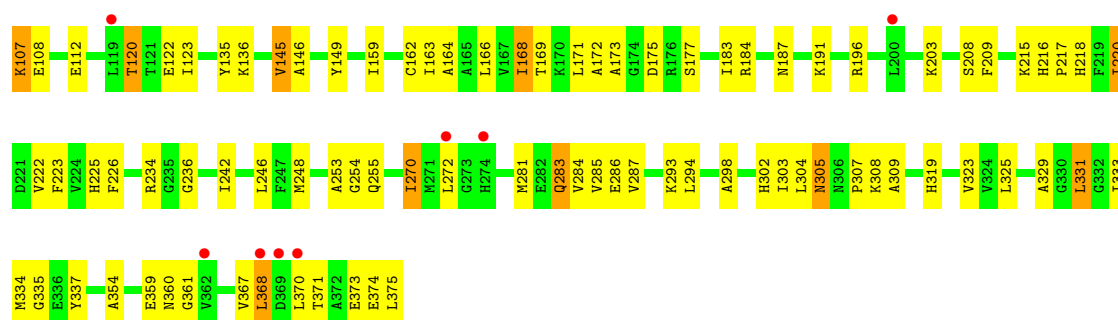
Chain Q:



• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

Chain W:





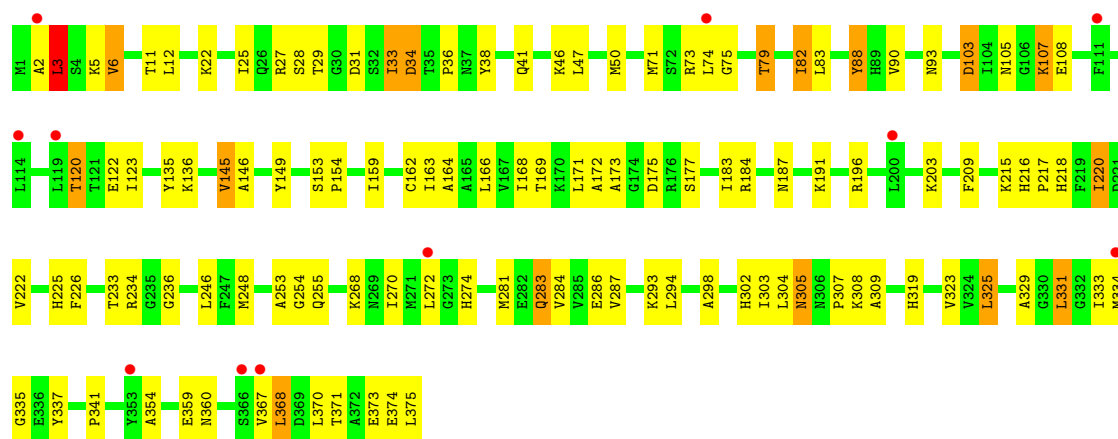
• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

Chain Y:



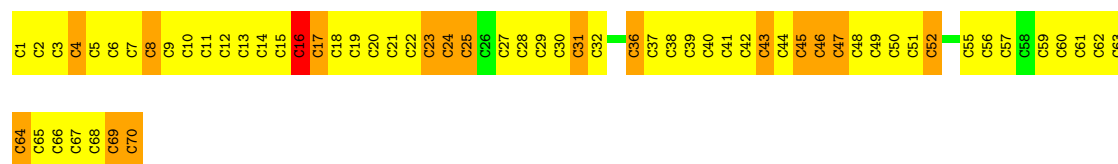
• Molecule 1: RESPIRATORY SYNCYTIAL VIRUS NUCLEOCAPSID PROTEIN

Chain Z:



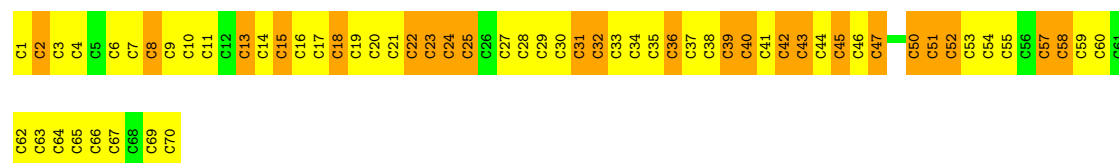
• Molecule 2: RNA

Chain R:



- Molecule 2: RNA

Chain X: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	164.35Å 175.74Å 241.98Å 90.09° 89.96° 89.92°	Depositor
Resolution (Å)	50.00 – 3.60 48.03 – 3.46	Depositor EDS
% Data completeness (in resolution range)	97.2 (50.00-3.60) 72.7 (48.03-3.46)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.28 (at 3.48Å)	Xtriage
Refinement program	REFMAC 5.6.0077	Depositor
R, R_{free}	0.229 , 0.265 0.445 , 0.442	Depositor DCC
R_{free} test set	12910 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	107.7	Xtriage
Anisotropy	0.623	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 62.0	EDS
Estimated twinning fraction	0.297 for h,-k,-l 0.058 for -h,k,-l 0.057 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	0 of 258134 reflections	Xtriage
F_o, F_c correlation	0.75	EDS
Total number of atoms	61200	wwPDB-VP
Average B, all atoms (Å ²)	139.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 23.38 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 4.7195e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

5 Model quality i

5.1 Standard geometry i

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.45	0/2968	0.63	2/3998 (0.1%)
1	B	0.45	0/2968	0.63	2/3998 (0.1%)
1	C	0.45	0/2968	0.63	2/3998 (0.1%)
1	D	0.45	0/2968	0.63	2/3998 (0.1%)
1	E	0.45	0/2968	0.63	2/3998 (0.1%)
1	F	0.45	0/2968	0.63	2/3998 (0.1%)
1	G	0.45	0/2968	0.63	2/3998 (0.1%)
1	H	0.45	0/2968	0.63	2/3998 (0.1%)
1	I	0.45	0/2968	0.63	2/3998 (0.1%)
1	J	0.45	0/2968	0.63	2/3998 (0.1%)
1	K	0.45	0/2968	0.63	2/3998 (0.1%)
1	L	0.45	0/2968	0.63	2/3998 (0.1%)
1	M	0.45	0/2968	0.63	2/3998 (0.1%)
1	N	0.45	0/2968	0.63	2/3998 (0.1%)
1	O	0.45	0/2968	0.63	2/3998 (0.1%)
1	P	0.45	0/2968	0.63	2/3998 (0.1%)
1	Q	0.45	0/2968	0.63	2/3998 (0.1%)
1	W	0.45	0/2968	0.63	2/3998 (0.1%)
1	Y	0.45	0/2968	0.63	2/3998 (0.1%)
1	Z	0.45	0/2968	0.63	2/3998 (0.1%)
2	R	0.35	0/1539	0.88	2/2376 (0.1%)
2	X	0.35	0/1539	0.91	6/2376 (0.3%)
All	All	0.45	0/62438	0.65	48/84712 (0.1%)

There are no bond length outliers.

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	R	69	C	P-O3'-C3'	-7.37	110.86	119.70
2	X	2	C	P-O3'-C3'	-7.06	111.23	119.70
2	X	16	C	P-O3'-C3'	-6.21	112.25	119.70
2	X	40	C	P-O3'-C3'	5.93	126.81	119.70
1	Y	2	ALA	N-CA-C	5.83	126.73	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	2	ALA	N-CA-C	5.82	126.72	111.00
1	M	2	ALA	N-CA-C	5.82	126.71	111.00
1	P	2	ALA	N-CA-C	5.82	126.71	111.00
1	G	2	ALA	N-CA-C	5.82	126.71	111.00
1	Q	2	ALA	N-CA-C	5.82	126.71	111.00
1	E	2	ALA	N-CA-C	5.82	126.71	111.00
1	C	2	ALA	N-CA-C	5.81	126.70	111.00
1	W	2	ALA	N-CA-C	5.81	126.69	111.00
1	N	2	ALA	N-CA-C	5.81	126.69	111.00
1	O	2	ALA	N-CA-C	5.81	126.69	111.00
1	Z	2	ALA	N-CA-C	5.81	126.69	111.00
1	A	2	ALA	N-CA-C	5.81	126.68	111.00
1	B	2	ALA	N-CA-C	5.80	126.67	111.00
1	H	2	ALA	N-CA-C	5.80	126.67	111.00
1	L	2	ALA	N-CA-C	5.80	126.67	111.00
1	D	2	ALA	N-CA-C	5.80	126.67	111.00
1	F	2	ALA	N-CA-C	5.80	126.67	111.00
1	I	2	ALA	N-CA-C	5.80	126.67	111.00
1	J	2	ALA	N-CA-C	5.80	126.66	111.00
1	C	3	LEU	CA-CB-CG	5.58	128.12	115.30
1	K	3	LEU	CA-CB-CG	5.57	128.11	115.30
1	F	3	LEU	CA-CB-CG	5.57	128.10	115.30
1	P	3	LEU	CA-CB-CG	5.57	128.10	115.30
1	O	3	LEU	CA-CB-CG	5.56	128.09	115.30
1	Z	3	LEU	CA-CB-CG	5.56	128.09	115.30
1	G	3	LEU	CA-CB-CG	5.56	128.08	115.30
1	H	3	LEU	CA-CB-CG	5.56	128.08	115.30
1	Q	3	LEU	CA-CB-CG	5.56	128.08	115.30
1	B	3	LEU	CA-CB-CG	5.55	128.07	115.30
1	Y	3	LEU	CA-CB-CG	5.55	128.07	115.30
1	N	3	LEU	CA-CB-CG	5.55	128.07	115.30
1	L	3	LEU	CA-CB-CG	5.55	128.07	115.30
1	A	3	LEU	CA-CB-CG	5.55	128.06	115.30
1	D	3	LEU	CA-CB-CG	5.55	128.06	115.30
1	J	3	LEU	CA-CB-CG	5.55	128.06	115.30
1	E	3	LEU	CA-CB-CG	5.54	128.05	115.30
1	M	3	LEU	CA-CB-CG	5.54	128.05	115.30
1	W	3	LEU	CA-CB-CG	5.54	128.05	115.30
1	I	3	LEU	CA-CB-CG	5.53	128.03	115.30
2	X	13	C	P-O3'-C3'	-5.53	113.06	119.70
2	X	23	C	P-O3'-C3'	-5.52	113.08	119.70
2	X	51	C	P-O3'-C3'	-5.28	113.36	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	R	16	C	P-O3'-C3'	-5.03	113.67	119.70

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2920	0	2955	84	1
1	B	2920	0	2955	76	0
1	C	2920	0	2955	79	0
1	D	2920	0	2955	76	0
1	E	2920	0	2955	75	0
1	F	2920	0	2955	76	1
1	G	2920	0	2955	74	0
1	H	2920	0	2955	78	0
1	I	2920	0	2955	77	0
1	J	2920	0	2955	73	0
1	K	2920	0	2955	76	0
1	L	2920	0	2955	74	0
1	M	2920	0	2955	78	1
1	N	2920	0	2955	70	0
1	O	2920	0	2955	82	0
1	P	2920	0	2955	86	0
1	Q	2920	0	2955	87	0
1	W	2920	0	2955	85	0
1	Y	2920	0	2955	77	1
1	Z	2920	0	2955	75	0
2	R	1400	0	771	87	0
2	X	1400	0	771	121	0
All	All	61200	0	60642	1533	2

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 13.

All (1533) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:K:253:ALA:HB3	1:K:303:ILE:HD11	1.27	1.17
1:I:253:ALA:HB3	1:I:303:ILE:HD11	1.27	1.16
1:F:253:ALA:HB3	1:F:303:ILE:HD11	1.27	1.16
1:O:253:ALA:HB3	1:O:303:ILE:HD11	1.27	1.16
1:Y:253:ALA:HB3	1:Y:303:ILE:HD11	1.27	1.15
1:C:253:ALA:HB3	1:C:303:ILE:HD11	1.27	1.15
1:H:253:ALA:HB3	1:H:303:ILE:HD11	1.27	1.14
1:M:253:ALA:HB3	1:M:303:ILE:HD11	1.27	1.14
1:D:253:ALA:HB3	1:D:303:ILE:HD11	1.27	1.13
1:Q:253:ALA:HB3	1:Q:303:ILE:HD11	1.27	1.13
1:J:253:ALA:HB3	1:J:303:ILE:HD11	1.27	1.13
1:P:253:ALA:HB3	1:P:303:ILE:HD11	1.27	1.13
1:A:253:ALA:HB3	1:A:303:ILE:HD11	1.27	1.12
1:N:253:ALA:HB3	1:N:303:ILE:HD11	1.27	1.12
1:L:253:ALA:HB3	1:L:303:ILE:HD11	1.27	1.12
1:B:253:ALA:HB3	1:B:303:ILE:HD11	1.27	1.12
1:W:253:ALA:HB3	1:W:303:ILE:HD11	1.27	1.09
1:G:253:ALA:HB3	1:G:303:ILE:HD11	1.27	1.08
1:Z:253:ALA:HB3	1:Z:303:ILE:HD11	1.27	1.08
1:E:253:ALA:HB3	1:E:303:ILE:HD11	1.27	1.08
1:A:305:ASN:HB2	1:Q:236:GLY:O	1.54	1.07
1:D:236:GLY:O	1:E:305:ASN:HB2	1.55	1.04
1:J:305:ASN:HB2	1:Z:236:GLY:O	1.59	1.03
1:A:236:GLY:O	1:B:305:ASN:HB2	1.59	1.02
1:K:236:GLY:O	1:L:305:ASN:HB2	1.59	1.01
1:W:82:ILE:HD11	1:W:222:VAL:HA	1.45	0.99
1:E:82:ILE:HD11	1:E:222:VAL:HA	1.45	0.98
1:A:82:ILE:HD11	1:A:222:VAL:HA	1.45	0.98
1:L:82:ILE:HD11	1:L:222:VAL:HA	1.45	0.98
1:P:236:GLY:O	1:W:305:ASN:HB2	1.62	0.98
1:L:236:GLY:O	1:M:305:ASN:HB2	1.63	0.97
1:Q:82:ILE:HD11	1:Q:222:VAL:HA	1.45	0.97
1:M:82:ILE:HD11	1:M:222:VAL:HA	1.45	0.97
1:H:82:ILE:HD11	1:H:222:VAL:HA	1.45	0.97
1:Y:82:ILE:HD11	1:Y:222:VAL:HA	1.45	0.97
1:G:82:ILE:HD11	1:G:222:VAL:HA	1.45	0.97
1:G:236:GLY:O	1:H:305:ASN:HB2	1.64	0.97
1:C:236:GLY:O	1:D:305:ASN:HB2	1.64	0.97
1:N:236:GLY:O	1:O:305:ASN:HB2	1.65	0.96
1:J:82:ILE:HD11	1:J:222:VAL:HA	1.45	0.96
1:Z:82:ILE:HD11	1:Z:222:VAL:HA	1.45	0.96
1:F:82:ILE:HD11	1:F:222:VAL:HA	1.45	0.96

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:82:ILE:HD11	1:D:222:VAL:HA	1.45	0.96
1:B:236:GLY:O	1:C:305:ASN:HB2	1.65	0.95
1:C:82:ILE:HD11	1:C:222:VAL:HA	1.45	0.95
1:O:82:ILE:HD11	1:O:222:VAL:HA	1.45	0.95
1:K:82:ILE:HD11	1:K:222:VAL:HA	1.45	0.95
1:P:82:ILE:HD11	1:P:222:VAL:HA	1.45	0.95
1:N:82:ILE:HD11	1:N:222:VAL:HA	1.45	0.95
1:J:236:GLY:O	1:K:305:ASN:HB2	1.66	0.95
1:B:82:ILE:HD11	1:B:222:VAL:HA	1.45	0.94
1:I:82:ILE:HD11	1:I:222:VAL:HA	1.45	0.94
1:M:236:GLY:O	1:N:305:ASN:HB2	1.68	0.94
2:R:45:C:C5	2:R:46:C:N4	2.35	0.94
1:I:236:GLY:O	1:Q:305:ASN:HB2	1.67	0.94
2:X:37:C:C4	2:X:38:C:C2	2.56	0.93
1:O:236:GLY:O	1:P:305:ASN:HB2	1.68	0.93
2:R:45:C:C5	2:R:46:C:C4	2.56	0.92
2:X:59:C:C5	2:X:60:C:C4	2.59	0.91
2:R:69:C:C4	2:R:70:C:C4	2.61	0.89
1:N:75:GLY:O	1:N:79:THR:HG22	1.74	0.88
1:B:75:GLY:O	1:B:79:THR:HG22	1.74	0.88
1:G:75:GLY:O	1:G:79:THR:HG22	1.74	0.87
1:Z:75:GLY:O	1:Z:79:THR:HG22	1.74	0.87
1:M:75:GLY:O	1:M:79:THR:HG22	1.74	0.87
1:A:75:GLY:O	1:A:79:THR:HG22	1.74	0.87
1:H:236:GLY:O	1:I:305:ASN:HB2	1.74	0.87
1:E:75:GLY:O	1:E:79:THR:HG22	1.74	0.87
1:W:236:GLY:O	1:Y:305:ASN:HB2	1.74	0.87
1:W:75:GLY:O	1:W:79:THR:HG22	1.74	0.86
1:K:75:GLY:O	1:K:79:THR:HG22	1.74	0.86
1:H:75:GLY:O	1:H:79:THR:HG22	1.74	0.86
1:I:75:GLY:O	1:I:79:THR:HG22	1.74	0.86
1:J:75:GLY:O	1:J:79:THR:HG22	1.74	0.86
1:D:75:GLY:O	1:D:79:THR:HG22	1.74	0.86
1:O:75:GLY:O	1:O:79:THR:HG22	1.74	0.86
2:X:37:C:N4	2:X:38:C:N3	2.23	0.86
1:C:75:GLY:O	1:C:79:THR:HG22	1.74	0.86
1:P:75:GLY:O	1:P:79:THR:HG22	1.74	0.85
1:L:75:GLY:O	1:L:79:THR:HG22	1.74	0.85
1:F:75:GLY:O	1:F:79:THR:HG22	1.74	0.85
1:Q:75:GLY:O	1:Q:79:THR:HG22	1.74	0.85
1:Y:75:GLY:O	1:Y:79:THR:HG22	1.74	0.85
2:X:59:C:C4	2:X:60:C:N3	2.45	0.84

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:X:43:C:C4	2:X:44:C:C2	2.66	0.84
2:X:65:C:C4	2:X:66:C:C2	2.67	0.82
2:X:57:C:C4	2:X:58:C:C2	2.69	0.81
2:X:9:C:H2'	2:X:10:C:O4'	1.80	0.80
1:K:2:ALA:HB3	1:Z:367:VAL:CG1	2.12	0.80
2:X:45:C:C5	2:X:46:C:N4	2.50	0.80
1:Q:29:THR:HG22	1:Q:88:TYR:HB3	1.65	0.79
1:L:29:THR:HG22	1:L:88:TYR:HB3	1.65	0.79
1:J:29:THR:HG22	1:J:88:TYR:HB3	1.65	0.79
1:H:29:THR:HG22	1:H:88:TYR:HB3	1.65	0.79
1:O:29:THR:HG22	1:O:88:TYR:HB3	1.65	0.79
1:W:29:THR:HG22	1:W:88:TYR:HB3	1.65	0.79
1:C:29:THR:HG22	1:C:88:TYR:HB3	1.65	0.79
1:E:29:THR:HG22	1:E:88:TYR:HB3	1.65	0.78
1:K:29:THR:HG22	1:K:88:TYR:HB3	1.65	0.78
1:G:29:THR:HG22	1:G:88:TYR:HB3	1.65	0.78
2:X:59:C:C4	2:X:60:C:C4	2.71	0.78
1:I:29:THR:HG22	1:I:88:TYR:HB3	1.65	0.78
1:Z:29:THR:HG22	1:Z:88:TYR:HB3	1.65	0.78
1:A:29:THR:HG22	1:A:88:TYR:HB3	1.65	0.78
1:M:29:THR:HG22	1:M:88:TYR:HB3	1.65	0.78
1:N:29:THR:HG22	1:N:88:TYR:HB3	1.65	0.78
1:B:29:THR:HG22	1:B:88:TYR:HB3	1.65	0.78
1:E:236:GLY:O	1:F:305:ASN:HB2	1.84	0.77
1:Y:29:THR:HG22	1:Y:88:TYR:HB3	1.65	0.77
1:F:29:THR:HG22	1:F:88:TYR:HB3	1.65	0.77
1:D:29:THR:HG22	1:D:88:TYR:HB3	1.65	0.77
1:P:29:THR:HG22	1:P:88:TYR:HB3	1.65	0.77
2:X:30:C:H2'	2:X:31:C:O4'	1.85	0.76
2:X:41:C:N4	2:X:42:C:N4	2.35	0.75
1:L:38:TYR:CE1	1:M:26:GLN:HB2	2.22	0.75
2:X:1:C:OP2	2:X:70:C:H3'	1.87	0.74
2:X:65:C:N3	2:X:66:C:O2	2.20	0.74
2:R:59:C:C5	2:R:60:C:N4	2.56	0.73
2:X:37:C:C5	2:X:38:C:C4	2.76	0.73
2:X:1:C:P	2:X:70:C:O3'	2.47	0.72
2:R:59:C:C5	2:R:60:C:C4	2.77	0.72
2:X:37:C:N4	2:X:38:C:C2	2.58	0.71
2:R:45:C:H5	2:R:46:C:N4	1.88	0.71
2:R:27:C:C4	2:R:28:C:C4	2.79	0.71
1:Y:361:GLY:HA3	1:Z:274:HIS:NE2	2.05	0.71
2:R:59:C:C4	2:R:60:C:C4	2.78	0.71

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:J:38:TYR:CE1	1:K:26:GLN:HB2	2.26	0.71
1:Y:236:GLY:O	1:Z:305:ASN:HB2	1.89	0.70
1:I:331:LEU:HD13	1:I:354:ALA:HB1	1.74	0.70
1:Y:331:LEU:HD13	1:Y:354:ALA:HB1	1.74	0.70
1:A:331:LEU:HD13	1:A:354:ALA:HB1	1.74	0.70
1:K:331:LEU:HD13	1:K:354:ALA:HB1	1.74	0.70
1:M:331:LEU:HD13	1:M:354:ALA:HB1	1.74	0.70
1:B:331:LEU:HD13	1:B:354:ALA:HB1	1.74	0.70
1:E:331:LEU:HD13	1:E:354:ALA:HB1	1.74	0.70
1:F:331:LEU:HD13	1:F:354:ALA:HB1	1.74	0.70
1:N:331:LEU:HD13	1:N:354:ALA:HB1	1.74	0.70
1:W:331:LEU:HD13	1:W:354:ALA:HB1	1.74	0.70
2:R:65:C:H2'	2:R:66:C:O4'	1.92	0.69
1:P:331:LEU:HD13	1:P:354:ALA:HB1	1.74	0.69
1:D:331:LEU:HD13	1:D:354:ALA:HB1	1.74	0.69
2:X:45:C:C4	2:X:46:C:N4	2.60	0.69
1:Q:331:LEU:HD13	1:Q:354:ALA:HB1	1.74	0.69
2:X:59:C:H41	2:X:60:C:H42	1.38	0.69
1:C:29:THR:CG2	1:C:88:TYR:HB3	2.23	0.69
1:L:331:LEU:HD13	1:L:354:ALA:HB1	1.74	0.69
1:O:29:THR:CG2	1:O:88:TYR:HB3	2.23	0.69
1:J:29:THR:CG2	1:J:88:TYR:HB3	2.23	0.69
1:A:29:THR:CG2	1:A:88:TYR:HB3	2.23	0.69
1:Y:29:THR:CG2	1:Y:88:TYR:HB3	2.23	0.69
1:F:29:THR:CG2	1:F:88:TYR:HB3	2.23	0.69
1:O:331:LEU:HD13	1:O:354:ALA:HB1	1.74	0.69
1:H:29:THR:CG2	1:H:88:TYR:HB3	2.23	0.69
1:K:29:THR:CG2	1:K:88:TYR:HB3	2.23	0.69
2:X:51:C:N4	2:X:52:C:N3	2.41	0.69
1:C:331:LEU:HD13	1:C:354:ALA:HB1	1.74	0.69
1:M:29:THR:CG2	1:M:88:TYR:HB3	2.23	0.69
1:I:29:THR:CG2	1:I:88:TYR:HB3	2.23	0.68
1:W:242:ILE:HG12	2:X:8:C:C4	2.29	0.68
2:X:59:C:C5	2:X:60:C:N4	2.61	0.68
1:D:29:THR:CG2	1:D:88:TYR:HB3	2.23	0.68
1:G:331:LEU:HD13	1:G:354:ALA:HB1	1.74	0.68
1:Z:331:LEU:HD13	1:Z:354:ALA:HB1	1.74	0.68
2:R:37:C:C4	2:R:38:C:N3	2.62	0.68
2:X:65:C:C4	2:X:66:C:N3	2.62	0.68
1:B:29:THR:CG2	1:B:88:TYR:HB3	2.23	0.68
1:P:29:THR:CG2	1:P:88:TYR:HB3	2.23	0.68
2:R:30:C:H2'	2:R:31:C:O4'	1.93	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:29:THR:CG2	1:L:88:TYR:HB3	2.23	0.68
1:Z:29:THR:CG2	1:Z:88:TYR:HB3	2.23	0.68
1:Y:335:GLY:HA2	1:Y:337:TYR:H	1.59	0.68
1:C:38:TYR:CE1	1:D:26:GLN:HB2	2.29	0.68
1:H:331:LEU:HD13	1:H:354:ALA:HB1	1.74	0.68
1:J:331:LEU:HD13	1:J:354:ALA:HB1	1.74	0.68
1:Q:29:THR:CG2	1:Q:88:TYR:HB3	2.23	0.67
1:G:29:THR:CG2	1:G:88:TYR:HB3	2.23	0.67
1:N:29:THR:CG2	1:N:88:TYR:HB3	2.23	0.67
1:F:335:GLY:HA2	1:F:337:TYR:H	1.60	0.67
1:A:335:GLY:HA2	1:A:337:TYR:H	1.59	0.67
1:J:335:GLY:HA2	1:J:337:TYR:H	1.60	0.67
2:R:69:C:N4	2:R:70:C:C4	2.63	0.67
2:R:59:C:C4	2:R:60:C:N4	2.63	0.67
2:R:37:C:C4	2:R:38:C:C2	2.82	0.67
1:H:335:GLY:HA2	1:H:337:TYR:H	1.60	0.67
1:M:335:GLY:HA2	1:M:337:TYR:H	1.60	0.67
2:R:69:C:C4	2:R:70:C:N3	2.63	0.67
1:B:335:GLY:HA2	1:B:337:TYR:H	1.60	0.67
1:E:29:THR:CG2	1:E:88:TYR:HB3	2.23	0.67
1:C:215:LYS:HE3	1:C:216:HIS:HE1	1.59	0.67
1:N:335:GLY:HA2	1:N:337:TYR:H	1.60	0.67
1:E:3:LEU:O	1:E:6:VAL:HG13	1.95	0.67
1:J:215:LYS:HE3	1:J:216:HIS:HE1	1.59	0.67
1:W:335:GLY:HA2	1:W:337:TYR:H	1.60	0.67
1:W:29:THR:CG2	1:W:88:TYR:HB3	2.23	0.67
1:J:3:LEU:O	1:J:6:VAL:HG13	1.95	0.67
1:Z:3:LEU:O	1:Z:6:VAL:HG13	1.95	0.67
1:H:3:LEU:O	1:H:6:VAL:HG13	1.95	0.67
1:F:335:GLY:CA	1:F:337:TYR:H	2.08	0.67
1:C:335:GLY:CA	1:C:337:TYR:H	2.08	0.67
1:W:3:LEU:O	1:W:6:VAL:HG13	1.95	0.67
1:G:3:LEU:O	1:G:6:VAL:HG13	1.95	0.67
1:Y:215:LYS:HE3	1:Y:216:HIS:HE1	1.59	0.67
1:E:335:GLY:HA2	1:E:337:TYR:H	1.59	0.67
1:J:335:GLY:CA	1:J:337:TYR:H	2.08	0.67
1:H:335:GLY:CA	1:H:337:TYR:H	2.08	0.67
1:C:335:GLY:HA2	1:C:337:TYR:H	1.60	0.67
1:O:215:LYS:HE3	1:O:216:HIS:HE1	1.59	0.67
1:P:335:GLY:HA2	1:P:337:TYR:H	1.59	0.67
1:Y:335:GLY:CA	1:Y:337:TYR:H	2.08	0.67
1:C:3:LEU:O	1:C:6:VAL:HG13	1.95	0.67

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:215:LYS:HE3	1:H:216:HIS:HE1	1.59	0.67
1:P:3:LEU:O	1:P:6:VAL:HG13	1.95	0.67
1:O:335:GLY:HA2	1:O:337:TYR:H	1.60	0.67
1:O:335:GLY:CA	1:O:337:TYR:H	2.08	0.67
1:L:3:LEU:O	1:L:6:VAL:HG13	1.95	0.67
1:O:3:LEU:O	1:O:6:VAL:HG13	1.95	0.67
1:Z:335:GLY:HA2	1:Z:337:TYR:H	1.59	0.66
1:Q:3:LEU:O	1:Q:6:VAL:HG13	1.95	0.66
1:D:335:GLY:HA2	1:D:337:TYR:H	1.60	0.66
1:F:215:LYS:HE3	1:F:216:HIS:HE1	1.59	0.66
1:B:3:LEU:O	1:B:6:VAL:HG13	1.95	0.66
1:D:3:LEU:O	1:D:6:VAL:HG13	1.95	0.66
1:P:215:LYS:HE3	1:P:216:HIS:HE1	1.59	0.66
1:W:215:LYS:HE3	1:W:216:HIS:HE1	1.59	0.66
1:N:3:LEU:O	1:N:6:VAL:HG13	1.95	0.66
1:P:248:MET:CE	1:W:14:LYS:HD3	2.26	0.66
1:G:335:GLY:HA2	1:G:337:TYR:H	1.60	0.66
1:A:335:GLY:CA	1:A:337:TYR:H	2.08	0.66
1:A:215:LYS:HE3	1:A:216:HIS:HE1	1.59	0.66
1:M:335:GLY:CA	1:M:337:TYR:H	2.08	0.66
1:I:3:LEU:O	1:I:6:VAL:HG13	1.95	0.66
1:A:26:GLN:HB2	1:Q:38:TYR:CE1	2.31	0.66
1:I:215:LYS:HE3	1:I:216:HIS:HE1	1.59	0.66
1:D:215:LYS:HE3	1:D:216:HIS:HE1	1.59	0.66
1:Z:335:GLY:CA	1:Z:337:TYR:H	2.08	0.66
1:K:3:LEU:O	1:K:6:VAL:HG13	1.95	0.66
1:E:215:LYS:HE3	1:E:216:HIS:HE1	1.59	0.66
1:K:335:GLY:HA2	1:K:337:TYR:H	1.59	0.66
1:L:335:GLY:HA2	1:L:337:TYR:H	1.59	0.66
1:G:335:GLY:CA	1:G:337:TYR:H	2.08	0.66
1:I:335:GLY:HA2	1:I:337:TYR:H	1.60	0.66
1:A:3:LEU:O	1:A:6:VAL:HG13	1.95	0.66
1:G:217:PRO:O	1:G:220:ILE:HG23	1.96	0.66
1:Z:217:PRO:O	1:Z:220:ILE:HG23	1.96	0.66
1:Q:335:GLY:HA2	1:Q:337:TYR:H	1.60	0.66
1:G:215:LYS:HE3	1:G:216:HIS:HE1	1.59	0.66
1:C:217:PRO:O	1:C:220:ILE:HG23	1.96	0.66
1:K:215:LYS:HE3	1:K:216:HIS:HE1	1.60	0.66
1:Y:3:LEU:O	1:Y:6:VAL:HG13	1.95	0.66
1:B:335:GLY:CA	1:B:337:TYR:H	2.08	0.65
1:E:335:GLY:CA	1:E:337:TYR:H	2.08	0.65
1:P:335:GLY:CA	1:P:337:TYR:H	2.08	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:Q:335:GLY:CA	1:Q:337:TYR:H	2.08	0.65
1:O:217:PRO:O	1:O:220:ILE:HG23	1.96	0.65
1:I:217:PRO:O	1:I:220:ILE:HG23	1.96	0.65
1:F:3:LEU:O	1:F:6:VAL:HG13	1.95	0.65
1:M:215:LYS:HE3	1:M:216:HIS:HE1	1.59	0.65
1:M:3:LEU:O	1:M:6:VAL:HG13	1.95	0.65
1:Z:215:LYS:HE3	1:Z:216:HIS:HE1	1.59	0.65
2:X:43:C:H2'	2:X:44:C:O4'	1.96	0.65
1:N:335:GLY:CA	1:N:337:TYR:H	2.08	0.65
1:W:335:GLY:CA	1:W:337:TYR:H	2.08	0.65
1:K:335:GLY:CA	1:K:337:TYR:H	2.08	0.65
1:H:217:PRO:O	1:H:220:ILE:HG23	1.96	0.65
1:Q:215:LYS:HE3	1:Q:216:HIS:HE1	1.59	0.65
1:K:217:PRO:O	1:K:220:ILE:HG23	1.96	0.65
1:L:335:GLY:CA	1:L:337:TYR:H	2.08	0.65
1:N:217:PRO:O	1:N:220:ILE:HG23	1.96	0.65
1:B:215:LYS:HE3	1:B:216:HIS:HE1	1.59	0.65
1:L:215:LYS:HE3	1:L:216:HIS:HE1	1.60	0.65
2:X:37:C:C4	2:X:38:C:N3	2.62	0.65
1:D:335:GLY:CA	1:D:337:TYR:H	2.08	0.65
1:B:217:PRO:O	1:B:220:ILE:HG23	1.96	0.65
2:X:24:C:N4	2:X:25:C:H42	1.95	0.65
2:X:24:C:C4	2:X:25:C:N3	2.64	0.65
1:J:217:PRO:O	1:J:220:ILE:HG23	1.96	0.65
1:D:217:PRO:O	1:D:220:ILE:HG23	1.96	0.65
1:E:217:PRO:O	1:E:220:ILE:HG23	1.96	0.65
1:P:217:PRO:O	1:P:220:ILE:HG23	1.96	0.65
1:N:215:LYS:HE3	1:N:216:HIS:HE1	1.59	0.65
1:A:217:PRO:O	1:A:220:ILE:HG23	1.96	0.65
1:W:217:PRO:O	1:W:220:ILE:HG23	1.96	0.65
1:M:217:PRO:O	1:M:220:ILE:HG23	1.96	0.65
1:I:335:GLY:CA	1:I:337:TYR:H	2.08	0.65
1:Y:217:PRO:O	1:Y:220:ILE:HG23	1.96	0.65
1:O:38:TYR:CE1	1:P:26:GLN:HB2	2.32	0.65
1:C:74:LEU:HD21	1:C:226:PHE:HA	1.79	0.65
1:F:217:PRO:O	1:F:220:ILE:HG23	1.96	0.65
1:Y:74:LEU:HD21	1:Y:226:PHE:HA	1.80	0.64
1:F:74:LEU:HD21	1:F:226:PHE:HA	1.80	0.64
1:L:217:PRO:O	1:L:220:ILE:HG23	1.96	0.64
1:O:74:LEU:HD21	1:O:226:PHE:HA	1.80	0.64
1:A:248:MET:CE	1:B:14:LYS:HD3	2.28	0.64
1:P:74:LEU:HD21	1:P:226:PHE:HA	1.80	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:74:LEU:HD21	1:D:226:PHE:HA	1.80	0.64
1:E:74:LEU:HD21	1:E:226:PHE:HA	1.79	0.64
1:Q:217:PRO:O	1:Q:220:ILE:HG23	1.96	0.64
1:H:74:LEU:HD21	1:H:226:PHE:HA	1.79	0.64
1:J:74:LEU:HD21	1:J:226:PHE:HA	1.79	0.64
2:X:45:C:C4	2:X:46:C:C4	2.86	0.64
2:X:13:C:C5	2:X:14:C:C5	2.86	0.64
2:X:18:C:OP1	1:Z:254:GLY:HA2	1.98	0.64
1:W:74:LEU:HD21	1:W:226:PHE:HA	1.80	0.64
1:M:74:LEU:HD21	1:M:226:PHE:HA	1.80	0.64
1:A:74:LEU:HD21	1:A:226:PHE:HA	1.80	0.64
2:X:57:C:C4	2:X:58:C:O2	2.51	0.64
2:R:43:C:C4	2:R:44:C:C2	2.86	0.64
1:H:38:TYR:CE1	1:I:26:GLN:HB2	2.33	0.64
1:L:74:LEU:HD21	1:L:226:PHE:HA	1.80	0.63
1:I:74:LEU:HD21	1:I:226:PHE:HA	1.80	0.63
1:K:74:LEU:HD21	1:K:226:PHE:HA	1.80	0.63
1:Q:74:LEU:HD21	1:Q:226:PHE:HA	1.80	0.63
1:B:74:LEU:HD21	1:B:226:PHE:HA	1.80	0.63
1:Z:74:LEU:HD21	1:Z:226:PHE:HA	1.80	0.63
1:Y:361:GLY:HA3	1:Z:274:HIS:HE2	1.64	0.63
1:G:74:LEU:HD21	1:G:226:PHE:HA	1.80	0.63
2:X:24:C:C5	2:X:25:C:C4	2.87	0.63
1:N:74:LEU:HD21	1:N:226:PHE:HA	1.80	0.63
2:R:69:C:N4	2:R:70:C:N4	2.47	0.63
1:I:38:TYR:CE1	1:Q:26:GLN:HB2	2.34	0.62
2:R:43:C:N4	2:R:44:C:N3	2.47	0.62
1:I:146:ALA:HB3	1:I:149:TYR:HD2	1.64	0.62
2:X:24:C:C5	2:X:25:C:N4	2.68	0.62
1:G:166:LEU:O	1:G:169:THR:HB	2.00	0.62
1:Y:166:LEU:O	1:Y:169:THR:HB	2.00	0.62
1:I:166:LEU:O	1:I:169:THR:HB	2.00	0.62
1:I:248:MET:CE	1:Q:14:LYS:HD3	2.30	0.62
1:K:146:ALA:HB3	1:K:149:TYR:HD2	1.64	0.62
2:X:65:C:N3	2:X:66:C:C2	2.68	0.62
1:F:166:LEU:O	1:F:169:THR:HB	2.00	0.62
1:L:166:LEU:O	1:L:169:THR:HB	2.00	0.62
1:B:234:ARG:HH21	1:C:86:ALA:HB2	1.65	0.62
1:Z:166:LEU:O	1:Z:169:THR:HB	2.00	0.62
1:Q:166:LEU:O	1:Q:169:THR:HB	2.00	0.62
1:P:146:ALA:HB3	1:P:149:TYR:HD2	1.64	0.62
1:J:368:LEU:H	1:J:368:LEU:HD12	1.65	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:K:166:LEU:O	1:K:169:THR:HB	2.00	0.62
1:H:368:LEU:H	1:H:368:LEU:HD12	1.65	0.62
2:X:1:C:P	2:X:70:C:C3'	2.88	0.62
1:D:146:ALA:HB3	1:D:149:TYR:HD2	1.64	0.62
1:F:368:LEU:HD12	1:F:368:LEU:H	1.65	0.62
2:R:37:C:H2'	2:R:38:C:O4'	2.00	0.61
1:Q:146:ALA:HB3	1:Q:149:TYR:HD2	1.64	0.61
1:M:368:LEU:H	1:M:368:LEU:HD12	1.65	0.61
1:Y:146:ALA:HB3	1:Y:149:TYR:HD2	1.64	0.61
1:W:166:LEU:O	1:W:169:THR:HB	2.00	0.61
1:E:166:LEU:O	1:E:169:THR:HB	2.00	0.61
1:P:368:LEU:HD12	1:P:368:LEU:H	1.65	0.61
1:Y:368:LEU:H	1:Y:368:LEU:HD12	1.65	0.61
1:A:368:LEU:HD12	1:A:368:LEU:H	1.65	0.61
1:A:166:LEU:O	1:A:169:THR:HB	2.00	0.61
1:W:368:LEU:HD12	1:W:368:LEU:H	1.65	0.61
1:B:368:LEU:H	1:B:368:LEU:HD12	1.65	0.61
1:O:368:LEU:HD12	1:O:368:LEU:H	1.65	0.61
2:X:1:C:P	2:X:70:C:H3'	2.40	0.61
1:G:367:VAL:CG1	1:I:2:ALA:HB3	2.30	0.61
1:M:146:ALA:HB3	1:M:149:TYR:HD2	1.64	0.61
1:L:146:ALA:HB3	1:L:149:TYR:HD2	1.64	0.61
1:W:146:ALA:HB3	1:W:149:TYR:HD2	1.64	0.61
1:D:166:LEU:O	1:D:169:THR:HB	2.00	0.61
1:E:368:LEU:HD12	1:E:368:LEU:H	1.65	0.61
1:N:368:LEU:HD12	1:N:368:LEU:H	1.65	0.61
1:Z:146:ALA:HB3	1:Z:149:TYR:HD2	1.64	0.61
1:N:146:ALA:HB3	1:N:149:TYR:HD2	1.64	0.61
1:M:166:LEU:O	1:M:169:THR:HB	2.00	0.61
1:H:166:LEU:O	1:H:169:THR:HB	2.00	0.61
1:O:166:LEU:O	1:O:169:THR:HB	2.00	0.61
1:P:166:LEU:O	1:P:169:THR:HB	2.00	0.61
1:C:368:LEU:H	1:C:368:LEU:HD12	1.65	0.61
1:C:166:LEU:O	1:C:169:THR:HB	2.00	0.61
1:F:236:GLY:O	1:G:305:ASN:HB2	1.99	0.61
1:J:166:LEU:O	1:J:169:THR:HB	2.00	0.61
1:D:368:LEU:H	1:D:368:LEU:HD12	1.65	0.61
1:G:368:LEU:HD12	1:G:368:LEU:H	1.65	0.61
1:Z:368:LEU:HD12	1:Z:368:LEU:H	1.65	0.61
1:G:146:ALA:HB3	1:G:149:TYR:HD2	1.64	0.61
1:B:146:ALA:HB3	1:B:149:TYR:HD2	1.64	0.61
1:F:146:ALA:HB3	1:F:149:TYR:HD2	1.64	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:146:ALA:HB3	1:E:149:TYR:HD2	1.64	0.61
1:N:234:ARG:HH21	1:O:86:ALA:HB2	1.66	0.61
1:N:166:LEU:O	1:N:169:THR:HB	2.00	0.61
1:L:88:TYR:CD1	1:L:88:TYR:N	2.69	0.61
1:E:337:TYR:CE1	2:R:39:C:H2'	2.35	0.61
1:A:146:ALA:HB3	1:A:149:TYR:HD2	1.64	0.61
1:E:234:ARG:HH21	1:F:86:ALA:HB2	1.66	0.61
1:O:256:VAL:HG12	2:X:60:C:OP1	2.01	0.60
1:Q:88:TYR:CD1	1:Q:88:TYR:N	2.69	0.60
1:A:88:TYR:CD1	1:A:88:TYR:N	2.69	0.60
1:O:146:ALA:HB3	1:O:149:TYR:HD2	1.64	0.60
1:C:146:ALA:HB3	1:C:149:TYR:HD2	1.64	0.60
1:B:166:LEU:O	1:B:169:THR:HB	2.00	0.60
1:J:88:TYR:N	1:J:88:TYR:CD1	2.69	0.60
1:H:88:TYR:N	1:H:88:TYR:CD1	2.69	0.60
1:W:88:TYR:N	1:W:88:TYR:CD1	2.69	0.60
1:E:88:TYR:N	1:E:88:TYR:CD1	2.69	0.60
1:M:88:TYR:CD1	1:M:88:TYR:N	2.69	0.60
1:Q:368:LEU:H	1:Q:368:LEU:HD12	1.65	0.60
2:R:59:C:N4	2:R:60:C:N4	2.49	0.60
1:H:146:ALA:HB3	1:H:149:TYR:HD2	1.64	0.60
1:L:368:LEU:HD12	1:L:368:LEU:H	1.65	0.60
1:N:88:TYR:CD1	1:N:88:TYR:N	2.69	0.60
1:D:88:TYR:CD1	1:D:88:TYR:N	2.69	0.60
1:F:367:VAL:HG11	1:G:278:GLN:OE1	2.00	0.60
1:I:368:LEU:H	1:I:368:LEU:HD12	1.65	0.60
1:Q:88:TYR:CE2	1:Q:218:HIS:HB3	2.37	0.60
1:L:88:TYR:CE2	1:L:218:HIS:HB3	2.37	0.60
1:W:88:TYR:CE2	1:W:218:HIS:HB3	2.37	0.60
1:B:88:TYR:CD1	1:B:88:TYR:N	2.69	0.60
1:F:88:TYR:CE2	1:F:218:HIS:HB3	2.37	0.60
1:P:88:TYR:CE2	1:P:218:HIS:HB3	2.37	0.60
1:J:146:ALA:HB3	1:J:149:TYR:HD2	1.64	0.60
1:E:88:TYR:CE2	1:E:218:HIS:HB3	2.37	0.60
1:K:88:TYR:N	1:K:88:TYR:CD1	2.69	0.60
1:G:88:TYR:CE2	1:G:218:HIS:HB3	2.37	0.60
1:G:88:TYR:N	1:G:88:TYR:CD1	2.69	0.60
1:Z:88:TYR:N	1:Z:88:TYR:CD1	2.69	0.60
1:Y:88:TYR:CD1	1:Y:88:TYR:N	2.69	0.60
1:D:88:TYR:CE2	1:D:218:HIS:HB3	2.37	0.60
1:P:88:TYR:N	1:P:88:TYR:CD1	2.69	0.60
1:M:248:MET:CE	1:N:14:LYS:HD3	2.31	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:Z:88:TYR:CE2	1:Z:218:HIS:HB3	2.37	0.60
1:G:234:ARG:HH21	1:H:86:ALA:HB2	1.66	0.60
1:K:368:LEU:HD12	1:K:368:LEU:H	1.65	0.60
2:X:65:C:H2'	2:X:66:C:O4'	2.02	0.60
1:O:88:TYR:CE2	1:O:218:HIS:HB3	2.37	0.60
1:C:88:TYR:CE2	1:C:218:HIS:HB3	2.37	0.60
1:I:88:TYR:CD1	1:I:88:TYR:N	2.69	0.60
1:A:14:LYS:HD3	1:Q:248:MET:CE	2.32	0.60
1:K:88:TYR:CE2	1:K:218:HIS:HB3	2.37	0.60
1:I:88:TYR:CE2	1:I:218:HIS:HB3	2.37	0.60
1:Y:88:TYR:CE2	1:Y:218:HIS:HB3	2.37	0.60
2:X:41:C:C4	2:X:42:C:C4	2.90	0.60
2:X:24:C:C4	2:X:25:C:C4	2.90	0.60
1:M:88:TYR:CE2	1:M:218:HIS:HB3	2.37	0.59
1:N:88:TYR:CE2	1:N:218:HIS:HB3	2.37	0.59
1:F:88:TYR:N	1:F:88:TYR:CD1	2.69	0.59
2:X:59:C:N4	2:X:60:C:N3	2.50	0.59
1:H:88:TYR:CE2	1:H:218:HIS:HB3	2.37	0.59
1:A:88:TYR:CE2	1:A:218:HIS:HB3	2.37	0.59
1:B:88:TYR:CE2	1:B:218:HIS:HB3	2.37	0.59
1:E:34:ASP:HA	1:E:93:ASN:HB3	1.85	0.59
1:G:34:ASP:HA	1:G:93:ASN:HB3	1.85	0.59
1:Z:34:ASP:HA	1:Z:93:ASN:HB3	1.85	0.59
1:Y:34:ASP:HA	1:Y:93:ASN:HB3	1.85	0.59
1:O:88:TYR:CD1	1:O:88:TYR:N	2.69	0.59
1:F:34:ASP:HA	1:F:93:ASN:HB3	1.85	0.59
2:R:20:C:C4	2:R:21:C:C4	2.90	0.59
2:R:23:C:N4	2:R:24:C:N3	2.50	0.59
2:R:55:C:C5	2:R:56:C:C4	2.91	0.59
1:W:34:ASP:HA	1:W:93:ASN:HB3	1.85	0.59
1:J:88:TYR:CE2	1:J:218:HIS:HB3	2.37	0.59
1:N:34:ASP:HA	1:N:93:ASN:HB3	1.85	0.59
1:B:34:ASP:HA	1:B:93:ASN:HB3	1.85	0.59
1:D:248:MET:CE	1:E:14:LYS:HD3	2.33	0.59
1:J:86:ALA:HB2	1:Z:234:ARG:HH21	1.67	0.59
1:K:34:ASP:HA	1:K:93:ASN:HB3	1.85	0.59
1:P:38:TYR:CE1	1:W:26:GLN:HB2	2.38	0.59
1:C:34:ASP:HA	1:C:93:ASN:HB3	1.85	0.59
1:C:88:TYR:N	1:C:88:TYR:CD1	2.69	0.59
2:R:55:C:C4	2:R:56:C:C4	2.90	0.59
1:G:184:ARG:NH1	2:R:56:C:OP2	2.36	0.59
2:R:69:C:H2'	2:R:70:C:O4'	2.02	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:367:VAL:HG12	1:G:3:LEU:HD12	1.85	0.59
1:I:34:ASP:HA	1:I:93:ASN:HB3	1.85	0.59
1:O:34:ASP:HA	1:O:93:ASN:HB3	1.85	0.59
1:L:34:ASP:HA	1:L:93:ASN:HB3	1.85	0.59
1:J:34:ASP:HA	1:J:93:ASN:HB3	1.85	0.59
2:R:1:C:P	2:R:70:C:O3'	2.61	0.58
2:X:9:C:N3	2:X:10:C:C2	2.71	0.58
2:R:37:C:N4	2:R:38:C:N3	2.51	0.58
1:Q:34:ASP:HA	1:Q:93:ASN:HB3	1.85	0.58
1:H:34:ASP:HA	1:H:93:ASN:HB3	1.85	0.58
1:M:34:ASP:HA	1:M:93:ASN:HB3	1.85	0.58
1:A:34:ASP:HA	1:A:93:ASN:HB3	1.85	0.58
2:R:45:C:C4	2:R:46:C:C4	2.91	0.58
1:D:34:ASP:HA	1:D:93:ASN:HB3	1.85	0.58
1:F:361:GLY:HA3	1:G:274:HIS:HE2	1.69	0.58
1:E:234:ARG:NH2	1:F:86:ALA:HB2	2.19	0.58
1:P:34:ASP:HA	1:P:93:ASN:HB3	1.85	0.58
2:R:59:C:N4	2:R:60:C:H42	2.02	0.58
2:X:51:C:C4	2:X:52:C:C2	2.92	0.58
2:X:37:C:H5	2:X:38:C:C4	2.21	0.57
1:Y:135:TYR:OH	1:Y:145:VAL:HG21	2.05	0.57
1:Z:135:TYR:OH	1:Z:145:VAL:HG21	2.05	0.57
1:F:135:TYR:OH	1:F:145:VAL:HG21	2.05	0.57
1:G:135:TYR:OH	1:G:145:VAL:HG21	2.05	0.57
2:X:13:C:C4	2:X:14:C:C4	2.93	0.57
1:D:234:ARG:HH21	1:E:86:ALA:HB2	1.69	0.57
2:X:59:C:N4	2:X:60:C:H42	2.01	0.57
1:K:234:ARG:HH21	1:L:86:ALA:HB2	1.70	0.57
1:M:135:TYR:OH	1:M:145:VAL:HG21	2.05	0.57
1:C:337:TYR:CE1	2:R:25:C:H2'	2.39	0.56
1:N:135:TYR:OH	1:N:145:VAL:HG21	2.05	0.56
1:O:135:TYR:OH	1:O:145:VAL:HG21	2.05	0.56
1:J:135:TYR:OH	1:J:145:VAL:HG21	2.05	0.56
1:B:135:TYR:OH	1:B:145:VAL:HG21	2.05	0.56
2:X:45:C:N4	2:X:46:C:H42	2.03	0.56
2:X:69:C:C4	2:X:70:C:C4	2.93	0.56
1:A:135:TYR:OH	1:A:145:VAL:HG21	2.05	0.56
1:H:135:TYR:OH	1:H:145:VAL:HG21	2.05	0.56
1:C:135:TYR:OH	1:C:145:VAL:HG21	2.05	0.56
1:P:135:TYR:OH	1:P:145:VAL:HG21	2.05	0.56
1:E:135:TYR:OH	1:E:145:VAL:HG21	2.05	0.56
1:J:2:ALA:HB3	1:Y:367:VAL:CG1	2.35	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:X:33:C:H2'	2:X:34:C:O4'	2.05	0.56
1:W:135:TYR:OH	1:W:145:VAL:HG21	2.05	0.56
1:Q:135:TYR:OH	1:Q:145:VAL:HG21	2.05	0.56
2:R:69:C:N4	2:R:70:C:N3	2.53	0.56
2:R:27:C:C4	2:R:28:C:N3	2.74	0.56
1:M:245:GLY:HA3	2:X:50:C:O2	2.06	0.56
1:D:135:TYR:OH	1:D:145:VAL:HG21	2.05	0.56
2:R:41:C:C4	2:R:42:C:C4	2.94	0.56
1:L:135:TYR:OH	1:L:145:VAL:HG21	2.05	0.56
1:K:135:TYR:OH	1:K:145:VAL:HG21	2.05	0.56
2:X:30:C:C4	2:X:31:C:N3	2.74	0.56
2:R:30:C:C4	2:R:31:C:C2	2.94	0.56
1:E:367:VAL:HG12	1:G:3:LEU:CD1	2.36	0.56
1:P:233:THR:HG22	1:W:307:PRO:HG2	1.86	0.56
1:I:135:TYR:OH	1:I:145:VAL:HG21	2.05	0.56
1:N:248:MET:CE	1:O:14:LYS:HD3	2.35	0.56
1:D:367:VAL:CG1	1:F:2:ALA:HB3	2.36	0.56
1:Y:82:ILE:HD12	1:Y:225:HIS:HB2	1.89	0.55
1:F:82:ILE:HD12	1:F:225:HIS:HB2	1.89	0.55
1:I:82:ILE:HD12	1:I:225:HIS:HB2	1.88	0.55
1:H:82:ILE:HD12	1:H:225:HIS:HB2	1.88	0.55
1:J:82:ILE:HD12	1:J:225:HIS:HB2	1.89	0.55
1:D:82:ILE:HD12	1:D:225:HIS:HB2	1.89	0.55
1:P:82:ILE:HD12	1:P:225:HIS:HB2	1.88	0.55
1:W:82:ILE:HD12	1:W:225:HIS:HB2	1.89	0.55
1:E:82:ILE:HD12	1:E:225:HIS:HB2	1.89	0.55
1:C:82:ILE:HD12	1:C:225:HIS:HB2	1.89	0.55
1:K:82:ILE:HD12	1:K:225:HIS:HB2	1.88	0.55
1:J:367:VAL:CG1	1:L:2:ALA:HB3	2.37	0.55
2:X:65:C:C4	2:X:66:C:O2	2.57	0.55
1:L:82:ILE:HD12	1:L:225:HIS:HB2	1.88	0.55
1:M:82:ILE:HD12	1:M:225:HIS:HB2	1.88	0.55
1:O:82:ILE:HD12	1:O:225:HIS:HB2	1.89	0.55
1:K:3:LEU:HD11	1:Z:368:LEU:HG	1.89	0.55
1:Q:82:ILE:HD12	1:Q:225:HIS:HB2	1.89	0.55
1:A:82:ILE:HD12	1:A:225:HIS:HB2	1.89	0.55
2:X:59:C:N4	2:X:60:C:N4	2.54	0.55
2:R:30:C:C4	2:R:31:C:N3	2.74	0.55
1:F:234:ARG:HH21	1:G:86:ALA:HB2	1.72	0.55
1:G:82:ILE:HD12	1:G:225:HIS:HB2	1.88	0.55
1:Z:82:ILE:HD12	1:Z:225:HIS:HB2	1.89	0.55
1:F:361:GLY:HA3	1:G:274:HIS:NE2	2.21	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:X:6:C:C4	2:X:7:C:C4	2.95	0.55
2:X:37:C:C5	2:X:38:C:C2	2.94	0.55
2:X:45:C:C5	2:X:46:C:C4	2.94	0.54
2:R:9:C:C5	2:R:10:C:C4	2.95	0.54
1:Y:164:ALA:HB2	1:Y:209:PHE:CE2	2.42	0.54
1:F:164:ALA:HB2	1:F:209:PHE:CE2	2.42	0.54
1:W:164:ALA:HB2	1:W:209:PHE:CE2	2.42	0.54
1:Q:164:ALA:HB2	1:Q:209:PHE:CE2	2.43	0.54
1:E:164:ALA:HB2	1:E:209:PHE:CE2	2.42	0.54
1:I:164:ALA:HB2	1:I:209:PHE:CE2	2.43	0.54
1:A:164:ALA:HB2	1:A:209:PHE:CE2	2.42	0.54
1:N:82:ILE:HD12	1:N:225:HIS:HB2	1.89	0.54
1:B:82:ILE:HD12	1:B:225:HIS:HB2	1.89	0.54
1:L:42:LYS:NZ	1:M:30:GLY:C	2.60	0.54
1:L:164:ALA:HB2	1:L:209:PHE:CE2	2.43	0.54
1:Z:164:ALA:HB2	1:Z:209:PHE:CE2	2.43	0.54
1:M:164:ALA:HB2	1:M:209:PHE:CE2	2.42	0.54
1:G:164:ALA:HB2	1:G:209:PHE:CE2	2.42	0.54
1:I:146:ALA:HB3	1:I:149:TYR:CD2	2.43	0.54
2:R:16:C:C4	2:R:17:C:C2	2.95	0.54
1:O:164:ALA:HB2	1:O:209:PHE:CE2	2.42	0.54
1:O:42:LYS:NZ	1:P:30:GLY:C	2.61	0.54
1:P:164:ALA:HB2	1:P:209:PHE:CE2	2.42	0.54
2:X:41:C:N4	2:X:42:C:C4	2.75	0.54
1:N:146:ALA:HB3	1:N:149:TYR:CD2	2.43	0.54
1:C:164:ALA:HB2	1:C:209:PHE:CE2	2.42	0.54
1:K:164:ALA:HB2	1:K:209:PHE:CE2	2.42	0.54
2:R:2:C:N4	2:R:3:C:N3	2.55	0.54
2:X:62:C:C4	2:X:63:C:C4	2.96	0.54
2:X:43:C:N4	2:X:44:C:C2	2.75	0.54
2:X:57:C:N3	2:X:58:C:O2	2.40	0.54
2:X:18:C:H4'	1:Z:173:ALA:HB2	1.89	0.54
1:K:146:ALA:HB3	1:K:149:TYR:CD2	2.43	0.54
1:P:146:ALA:HB3	1:P:149:TYR:CD2	2.43	0.54
1:B:146:ALA:HB3	1:B:149:TYR:CD2	2.43	0.54
1:B:164:ALA:HB2	1:B:209:PHE:CE2	2.42	0.54
1:I:234:ARG:NH2	1:Q:86:ALA:HB2	2.23	0.54
2:X:13:C:C5	2:X:14:C:C4	2.96	0.54
2:R:23:C:C4	2:R:24:C:C2	2.96	0.54
1:D:164:ALA:HB2	1:D:209:PHE:CE2	2.42	0.54
1:N:164:ALA:HB2	1:N:209:PHE:CE2	2.42	0.54
1:D:146:ALA:HB3	1:D:149:TYR:CD2	2.43	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:X:11:C:H4'	1:Y:173:ALA:HB2	1.90	0.54
1:J:164:ALA:HB2	1:J:209:PHE:CE2	2.42	0.54
1:G:234:ARG:NH2	1:H:86:ALA:HB2	2.22	0.54
1:J:294:LEU:HB3	1:J:298:ALA:HB2	1.90	0.54
2:R:69:C:C5	2:R:70:C:C4	2.96	0.53
1:G:146:ALA:HB3	1:G:149:TYR:CD2	2.43	0.53
1:F:146:ALA:HB3	1:F:149:TYR:CD2	2.43	0.53
1:H:294:LEU:HB3	1:H:298:ALA:HB2	1.90	0.53
2:R:27:C:N3	2:R:28:C:C2	2.76	0.53
1:B:173:ALA:HB2	2:R:18:C:H4'	1.89	0.53
1:J:14:LYS:HD3	1:Z:248:MET:CE	2.37	0.53
1:H:164:ALA:HB2	1:H:209:PHE:CE2	2.43	0.53
1:M:38:TYR:CE1	1:N:26:GLN:HB2	2.43	0.53
1:M:146:ALA:HB3	1:M:149:TYR:CD2	2.43	0.53
2:X:21:C:OP2	1:Z:184:ARG:NH1	2.42	0.53
2:X:37:C:C5	2:X:38:C:C5	2.97	0.53
1:Z:146:ALA:HB3	1:Z:149:TYR:CD2	2.43	0.53
1:C:294:LEU:HB3	1:C:298:ALA:HB2	1.90	0.53
1:K:248:MET:CE	1:L:14:LYS:HD3	2.39	0.53
1:L:88:TYR:HD1	1:L:88:TYR:N	2.07	0.53
1:C:146:ALA:HB3	1:C:149:TYR:CD2	2.43	0.53
1:B:248:MET:CE	1:C:14:LYS:HD3	2.37	0.53
1:B:367:VAL:CG1	1:D:2:ALA:HB3	2.38	0.53
1:J:234:ARG:CZ	1:K:225:HIS:CE1	2.92	0.53
1:Q:88:TYR:HD1	1:Q:88:TYR:N	2.07	0.53
1:J:88:TYR:N	1:J:88:TYR:HD1	2.07	0.53
1:B:88:TYR:N	1:B:88:TYR:HD1	2.07	0.53
1:Y:88:TYR:HD1	1:Y:88:TYR:N	2.07	0.53
1:P:335:GLY:HA2	1:P:337:TYR:N	2.24	0.53
1:Y:146:ALA:HB3	1:Y:149:TYR:CD2	2.43	0.53
1:J:146:ALA:HB3	1:J:149:TYR:CD2	2.43	0.53
1:H:367:VAL:CG1	1:Q:2:ALA:HB3	2.38	0.53
1:O:294:LEU:HB3	1:O:298:ALA:HB2	1.90	0.53
1:H:88:TYR:N	1:H:88:TYR:HD1	2.07	0.53
1:E:88:TYR:N	1:E:88:TYR:HD1	2.07	0.53
1:N:88:TYR:N	1:N:88:TYR:HD1	2.07	0.53
1:Y:88:TYR:HD1	1:Y:88:TYR:H	1.57	0.53
1:A:335:GLY:HA2	1:A:337:TYR:N	2.24	0.53
1:A:146:ALA:HB3	1:A:149:TYR:CD2	2.43	0.53
1:H:146:ALA:HB3	1:H:149:TYR:CD2	2.43	0.53
1:E:294:LEU:HB3	1:E:298:ALA:HB2	1.90	0.53
1:G:294:LEU:HB3	1:G:298:ALA:HB2	1.90	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:W:88:TYR:N	1:W:88:TYR:HD1	2.07	0.53
1:F:88:TYR:N	1:F:88:TYR:HD1	2.07	0.53
1:F:294:LEU:HB3	1:F:298:ALA:HB2	1.90	0.53
1:M:294:LEU:HB3	1:M:298:ALA:HB2	1.90	0.53
1:Z:294:LEU:HB3	1:Z:298:ALA:HB2	1.90	0.53
1:O:88:TYR:HD1	1:O:88:TYR:N	2.07	0.53
1:C:88:TYR:HD1	1:C:88:TYR:N	2.07	0.53
1:D:335:GLY:HA2	1:D:337:TYR:N	2.24	0.53
1:W:294:LEU:HB3	1:W:298:ALA:HB2	1.90	0.53
1:L:248:MET:CE	1:M:14:LYS:HD3	2.39	0.53
1:N:294:LEU:HB3	1:N:298:ALA:HB2	1.90	0.53
1:M:88:TYR:H	1:M:88:TYR:HD1	1.57	0.53
1:F:88:TYR:H	1:F:88:TYR:HD1	1.57	0.53
1:M:335:GLY:HA2	1:M:337:TYR:N	2.24	0.53
1:W:36:PRO:HD2	1:W:71:MET:HB3	1.91	0.53
1:P:294:LEU:HB3	1:P:298:ALA:HB2	1.90	0.53
1:L:294:LEU:HB3	1:L:298:ALA:HB2	1.90	0.53
1:C:234:ARG:CZ	1:D:225:HIS:CE1	2.93	0.52
1:A:88:TYR:HD1	1:A:88:TYR:H	1.57	0.52
2:X:69:C:H2'	2:X:70:C:O4'	2.09	0.52
1:E:36:PRO:HD2	1:E:71:MET:HB3	1.92	0.52
1:Y:294:LEU:HB3	1:Y:298:ALA:HB2	1.90	0.52
1:J:88:TYR:H	1:J:88:TYR:HD1	1.57	0.52
1:H:88:TYR:HD1	1:H:88:TYR:H	1.57	0.52
1:Q:36:PRO:HD2	1:Q:71:MET:HB3	1.91	0.52
1:Q:294:LEU:HB3	1:Q:298:ALA:HB2	1.90	0.52
1:J:36:PRO:HD2	1:J:71:MET:HB3	1.91	0.52
1:B:294:LEU:HB3	1:B:298:ALA:HB2	1.90	0.52
1:Z:88:TYR:H	1:Z:88:TYR:HD1	1.57	0.52
1:P:242:ILE:HG12	2:X:1:C:C4	2.45	0.52
1:Y:335:GLY:HA2	1:Y:337:TYR:N	2.24	0.52
1:O:146:ALA:HB3	1:O:149:TYR:CD2	2.43	0.52
1:H:36:PRO:HD2	1:H:71:MET:HB3	1.92	0.52
1:L:234:ARG:CZ	1:M:225:HIS:CE1	2.92	0.52
1:G:88:TYR:HD1	1:G:88:TYR:H	1.57	0.52
1:M:88:TYR:HD1	1:M:88:TYR:N	2.07	0.52
1:P:234:ARG:NH2	1:W:86:ALA:HB2	2.24	0.52
1:L:36:PRO:HD2	1:L:71:MET:HB3	1.92	0.52
1:A:294:LEU:HB3	1:A:298:ALA:HB2	1.90	0.52
1:L:88:TYR:HD1	1:L:88:TYR:H	1.57	0.52
1:L:335:GLY:HA2	1:L:337:TYR:N	2.24	0.52
2:X:24:C:C4	2:X:25:C:N4	2.78	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:J:86:ALA:HB2	1:Z:234:ARG:NH2	2.24	0.52
1:K:294:LEU:HB3	1:K:298:ALA:HB2	1.90	0.52
1:Q:88:TYR:H	1:Q:88:TYR:HD1	1.57	0.52
1:A:88:TYR:HD1	1:A:88:TYR:N	2.07	0.52
1:F:335:GLY:HA2	1:F:337:TYR:N	2.24	0.52
1:B:36:PRO:HD2	1:B:71:MET:HB3	1.92	0.52
1:D:294:LEU:HB3	1:D:298:ALA:HB2	1.90	0.52
1:A:30:GLY:C	1:Q:42:LYS:NZ	2.63	0.52
1:O:367:VAL:HG11	1:P:278:GLN:NE2	2.24	0.52
1:Q:335:GLY:HA2	1:Q:337:TYR:N	2.24	0.52
1:N:36:PRO:HD2	1:N:71:MET:HB3	1.92	0.52
1:F:36:PRO:HD2	1:F:71:MET:HB3	1.91	0.52
1:Y:36:PRO:HD2	1:Y:71:MET:HB3	1.92	0.52
1:O:335:GLY:HA2	1:O:337:TYR:N	2.24	0.52
1:O:367:VAL:HG11	1:P:278:GLN:OE1	2.09	0.52
1:W:361:GLY:HA3	1:Y:274:HIS:NE2	2.24	0.52
1:H:159:ILE:O	1:H:162:CYS:HB2	2.10	0.52
1:G:253:ALA:O	1:G:302:HIS:HB2	2.10	0.52
1:Z:253:ALA:O	1:Z:302:HIS:HB2	2.10	0.52
2:X:43:C:N4	2:X:44:C:N3	2.57	0.52
1:O:88:TYR:HD1	1:O:88:TYR:H	1.57	0.52
1:C:36:PRO:HD2	1:C:71:MET:HB3	1.91	0.52
1:I:159:ILE:O	1:I:162:CYS:HB2	2.10	0.52
1:J:159:ILE:O	1:J:162:CYS:HB2	2.10	0.52
1:C:335:GLY:HA2	1:C:337:TYR:N	2.24	0.52
1:B:253:ALA:O	1:B:302:HIS:HB2	2.10	0.51
1:A:225:HIS:CE1	1:Q:234:ARG:CZ	2.93	0.51
1:C:88:TYR:HD1	1:C:88:TYR:H	1.57	0.51
1:E:88:TYR:HD1	1:E:88:TYR:H	1.57	0.51
1:P:88:TYR:HD1	1:P:88:TYR:H	1.57	0.51
2:R:9:C:N4	2:R:10:C:N3	2.58	0.51
1:O:36:PRO:HD2	1:O:71:MET:HB3	1.92	0.51
1:Z:159:ILE:O	1:Z:162:CYS:HB2	2.10	0.51
1:O:159:ILE:O	1:O:162:CYS:HB2	2.10	0.51
1:P:159:ILE:O	1:P:162:CYS:HB2	2.10	0.51
1:D:159:ILE:O	1:D:162:CYS:HB2	2.10	0.51
1:I:294:LEU:HB3	1:I:298:ALA:HB2	1.90	0.51
1:C:253:ALA:O	1:C:302:HIS:HB2	2.10	0.51
1:P:253:ALA:O	1:P:302:HIS:HB2	2.10	0.51
1:N:253:ALA:O	1:N:302:HIS:HB2	2.10	0.51
1:N:88:TYR:H	1:N:88:TYR:HD1	1.57	0.51
1:B:88:TYR:H	1:B:88:TYR:HD1	1.57	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:88:TYR:H	1:D:88:TYR:HD1	1.57	0.51
1:H:335:GLY:HA2	1:H:337:TYR:N	2.24	0.51
1:Q:146:ALA:HB3	1:Q:149:TYR:CD2	2.43	0.51
1:L:146:ALA:HB3	1:L:149:TYR:CD2	2.43	0.51
1:C:159:ILE:O	1:C:162:CYS:HB2	2.10	0.51
1:K:159:ILE:O	1:K:162:CYS:HB2	2.10	0.51
1:P:236:GLY:O	1:W:305:ASN:CB	2.49	0.51
1:W:88:TYR:H	1:W:88:TYR:HD1	1.57	0.51
1:G:88:TYR:HD1	1:G:88:TYR:N	2.07	0.51
1:Z:88:TYR:N	1:Z:88:TYR:HD1	2.07	0.51
1:D:88:TYR:HD1	1:D:88:TYR:N	2.07	0.51
1:Y:159:ILE:O	1:Y:162:CYS:HB2	2.10	0.51
1:G:159:ILE:O	1:G:162:CYS:HB2	2.10	0.51
1:D:253:ALA:O	1:D:302:HIS:HB2	2.10	0.51
1:Q:253:ALA:O	1:Q:302:HIS:HB2	2.10	0.51
2:X:37:C:C5	2:X:38:C:N3	2.79	0.51
1:J:335:GLY:HA2	1:J:337:TYR:N	2.24	0.51
1:F:159:ILE:O	1:F:162:CYS:HB2	2.10	0.51
1:Z:36:PRO:HD2	1:Z:71:MET:HB3	1.91	0.51
1:O:253:ALA:O	1:O:302:HIS:HB2	2.10	0.51
1:A:253:ALA:O	1:A:302:HIS:HB2	2.10	0.51
1:L:253:ALA:O	1:L:302:HIS:HB2	2.10	0.51
1:M:159:ILE:O	1:M:162:CYS:HB2	2.10	0.51
1:A:159:ILE:O	1:A:162:CYS:HB2	2.10	0.51
1:G:36:PRO:HD2	1:G:71:MET:HB3	1.92	0.51
1:M:253:ALA:O	1:M:302:HIS:HB2	2.10	0.51
1:A:23:TYR:CD1	1:Q:82:ILE:HG22	2.46	0.51
1:W:335:GLY:HA2	1:W:337:TYR:N	2.24	0.51
1:L:337:TYR:CE1	2:X:39:C:H2'	2.45	0.51
1:W:146:ALA:HB3	1:W:149:TYR:CD2	2.43	0.51
1:H:329:ALA:HA	1:H:334:MET:HG2	1.93	0.51
1:I:88:TYR:N	1:I:88:TYR:HD1	2.07	0.51
1:A:26:GLN:HB2	1:Q:38:TYR:CD1	2.46	0.51
2:X:20:C:C4	2:X:21:C:C4	2.99	0.51
2:R:12:C:C4	2:R:13:C:C5	2.98	0.51
1:K:36:PRO:HD2	1:K:71:MET:HB3	1.92	0.51
1:J:329:ALA:HA	1:J:334:MET:HG2	1.93	0.51
1:H:42:LYS:NZ	1:I:30:GLY:C	2.64	0.51
1:N:335:GLY:HA2	1:N:337:TYR:N	2.24	0.51
1:E:335:GLY:HA2	1:E:337:TYR:N	2.24	0.51
1:C:248:MET:CE	1:D:14:LYS:HD3	2.41	0.51
1:I:253:ALA:O	1:I:302:HIS:HB2	2.10	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:335:GLY:HA2	1:B:337:TYR:N	2.24	0.51
1:K:335:GLY:HA2	1:K:337:TYR:N	2.24	0.51
1:I:36:PRO:HD2	1:I:71:MET:HB3	1.92	0.51
1:A:36:PRO:HD2	1:A:71:MET:HB3	1.92	0.51
1:J:253:ALA:O	1:J:302:HIS:HB2	2.10	0.51
1:E:253:ALA:O	1:E:302:HIS:HB2	2.10	0.51
1:K:88:TYR:H	1:K:88:TYR:HD1	1.57	0.51
2:R:43:C:N4	2:R:44:C:C2	2.79	0.51
1:P:283:GLN:O	1:P:286:GLU:HB2	2.11	0.51
1:E:159:ILE:O	1:E:162:CYS:HB2	2.10	0.51
1:B:159:ILE:O	1:B:162:CYS:HB2	2.10	0.51
1:Q:329:ALA:HA	1:Q:334:MET:HG2	1.93	0.51
1:M:36:PRO:HD2	1:M:71:MET:HB3	1.92	0.51
1:K:253:ALA:O	1:K:302:HIS:HB2	2.10	0.50
1:Y:253:ALA:O	1:Y:302:HIS:HB2	2.10	0.50
1:H:253:ALA:O	1:H:302:HIS:HB2	2.10	0.50
1:P:88:TYR:HD1	1:P:88:TYR:N	2.07	0.50
1:P:242:ILE:HG12	2:X:1:C:N4	2.26	0.50
1:D:283:GLN:O	1:D:286:GLU:HB2	2.11	0.50
1:I:283:GLN:O	1:I:286:GLU:HB2	2.11	0.50
1:N:159:ILE:O	1:N:162:CYS:HB2	2.10	0.50
1:W:159:ILE:O	1:W:162:CYS:HB2	2.10	0.50
1:W:253:ALA:O	1:W:302:HIS:HB2	2.10	0.50
1:E:172:ALA:HB2	1:E:253:ALA:HB1	1.93	0.50
1:I:88:TYR:H	1:I:88:TYR:HD1	1.57	0.50
1:E:146:ALA:HB3	1:E:149:TYR:CD2	2.43	0.50
2:R:16:C:N4	2:R:17:C:C2	2.79	0.50
1:Y:329:ALA:HA	1:Y:334:MET:HG2	1.93	0.50
1:N:307:PRO:O	1:N:309:ALA:N	2.45	0.50
1:B:307:PRO:O	1:B:309:ALA:N	2.45	0.50
1:I:307:PRO:O	1:I:309:ALA:N	2.45	0.50
1:P:307:PRO:O	1:P:309:ALA:N	2.45	0.50
1:D:307:PRO:O	1:D:309:ALA:N	2.45	0.50
1:D:233:THR:HG22	1:E:307:PRO:HG2	1.93	0.50
1:B:283:GLN:O	1:B:286:GLU:HB2	2.11	0.50
1:Y:307:PRO:O	1:Y:309:ALA:N	2.45	0.50
1:Q:159:ILE:O	1:Q:162:CYS:HB2	2.10	0.50
1:C:172:ALA:HB2	1:C:253:ALA:HB1	1.93	0.50
1:K:88:TYR:N	1:K:88:TYR:HD1	2.07	0.50
2:X:41:C:C5	2:X:42:C:C5	2.99	0.50
1:L:159:ILE:O	1:L:162:CYS:HB2	2.10	0.50
1:H:307:PRO:O	1:H:309:ALA:N	2.45	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:329:ALA:HA	1:L:334:MET:HG2	1.93	0.50
1:N:329:ALA:HA	1:N:334:MET:HG2	1.93	0.50
1:I:233:THR:HG22	1:Q:307:PRO:HG2	1.93	0.50
1:C:42:LYS:NZ	1:D:30:GLY:C	2.65	0.50
1:K:307:PRO:O	1:K:309:ALA:N	2.45	0.50
1:F:329:ALA:HA	1:F:334:MET:HG2	1.93	0.50
1:N:283:GLN:O	1:N:286:GLU:HB2	2.11	0.50
1:F:253:ALA:O	1:F:302:HIS:HB2	2.10	0.50
1:I:335:GLY:HA2	1:I:337:TYR:N	2.24	0.50
1:J:187:ASN:O	1:J:191:LYS:HB3	2.12	0.50
1:K:283:GLN:O	1:K:286:GLU:HB2	2.12	0.50
1:P:36:PRO:HD2	1:P:71:MET:HB3	1.92	0.50
1:J:307:PRO:O	1:J:309:ALA:N	2.45	0.50
1:F:307:PRO:O	1:F:309:ALA:N	2.45	0.50
1:O:172:ALA:HB2	1:O:253:ALA:HB1	1.93	0.50
1:M:172:ALA:HB2	1:M:253:ALA:HB1	1.93	0.50
1:J:172:ALA:HB2	1:J:253:ALA:HB1	1.93	0.50
2:R:2:C:N4	2:R:3:C:C4	2.80	0.50
1:Q:307:PRO:O	1:Q:309:ALA:N	2.45	0.50
1:O:307:PRO:O	1:O:309:ALA:N	2.45	0.50
1:B:329:ALA:HA	1:B:334:MET:HG2	1.93	0.50
1:I:329:ALA:HA	1:I:334:MET:HG2	1.93	0.50
1:H:187:ASN:O	1:H:191:LYS:HB3	2.12	0.50
1:H:172:ALA:HB2	1:H:253:ALA:HB1	1.93	0.50
1:A:172:ALA:HB2	1:A:253:ALA:HB1	1.93	0.50
1:W:172:ALA:HB2	1:W:253:ALA:HB1	1.93	0.50
1:Z:172:ALA:HB2	1:Z:253:ALA:HB1	1.93	0.50
1:G:335:GLY:HA2	1:G:337:TYR:N	2.24	0.50
1:M:307:PRO:O	1:M:309:ALA:N	2.45	0.50
1:W:187:ASN:O	1:W:191:LYS:HB3	2.12	0.50
1:E:187:ASN:O	1:E:191:LYS:HB3	2.12	0.50
1:C:307:PRO:O	1:C:309:ALA:N	2.45	0.50
1:D:36:PRO:HD2	1:D:71:MET:HB3	1.92	0.50
1:O:329:ALA:HA	1:O:334:MET:HG2	1.93	0.50
1:C:329:ALA:HA	1:C:334:MET:HG2	1.93	0.50
1:L:307:PRO:O	1:L:309:ALA:N	2.45	0.50
1:A:307:PRO:O	1:A:309:ALA:N	2.45	0.50
1:G:172:ALA:HB2	1:G:253:ALA:HB1	1.93	0.50
1:Z:335:GLY:HA2	1:Z:337:TYR:N	2.24	0.50
1:G:329:ALA:HA	1:G:334:MET:HG2	1.93	0.50
1:W:283:GLN:O	1:W:286:GLU:HB2	2.11	0.50
1:G:307:PRO:O	1:G:309:ALA:N	2.45	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:Q:283:GLN:O	1:Q:286:GLU:HB2	2.11	0.50
1:C:187:ASN:O	1:C:191:LYS:HB3	2.12	0.50
2:R:6:C:C5	2:R:7:C:C5	3.00	0.50
1:K:329:ALA:HA	1:K:334:MET:HG2	1.93	0.50
1:F:172:ALA:HB2	1:F:253:ALA:HB1	1.93	0.50
1:B:172:ALA:HB2	1:B:253:ALA:HB1	1.93	0.50
1:F:187:ASN:O	1:F:191:LYS:HB3	2.12	0.50
1:F:283:GLN:O	1:F:286:GLU:HB2	2.11	0.50
1:O:187:ASN:O	1:O:191:LYS:HB3	2.12	0.50
1:Z:307:PRO:O	1:Z:309:ALA:N	2.45	0.50
1:P:172:ALA:HB2	1:P:253:ALA:HB1	1.93	0.50
1:N:172:ALA:HB2	1:N:253:ALA:HB1	1.93	0.50
2:X:45:C:N4	2:X:46:C:N4	2.60	0.50
1:Z:329:ALA:HA	1:Z:334:MET:HG2	1.93	0.50
1:L:187:ASN:O	1:L:191:LYS:HB3	2.12	0.50
1:Y:283:GLN:O	1:Y:286:GLU:HB2	2.11	0.50
1:P:329:ALA:HA	1:P:334:MET:HG2	1.93	0.50
1:H:283:GLN:O	1:H:286:GLU:HB2	2.11	0.50
1:E:307:PRO:O	1:E:309:ALA:N	2.45	0.49
1:Q:187:ASN:O	1:Q:191:LYS:HB3	2.12	0.49
1:A:329:ALA:HA	1:A:334:MET:HG2	1.93	0.49
1:L:283:GLN:O	1:L:286:GLU:HB2	2.11	0.49
1:Z:283:GLN:O	1:Z:286:GLU:HB2	2.11	0.49
1:D:329:ALA:HA	1:D:334:MET:HG2	1.93	0.49
1:I:187:ASN:O	1:I:191:LYS:HB3	2.12	0.49
1:D:172:ALA:HB2	1:D:253:ALA:HB1	1.93	0.49
1:L:172:ALA:HB2	1:L:253:ALA:HB1	1.93	0.49
1:W:307:PRO:O	1:W:309:ALA:N	2.45	0.49
1:G:283:GLN:O	1:G:286:GLU:HB2	2.11	0.49
1:P:187:ASN:O	1:P:191:LYS:HB3	2.12	0.49
1:Y:172:ALA:HB2	1:Y:253:ALA:HB1	1.93	0.49
2:R:27:C:H2'	2:R:28:C:O4'	2.11	0.49
1:A:248:MET:HE2	1:B:14:LYS:HD3	1.94	0.49
1:J:283:GLN:O	1:J:286:GLU:HB2	2.12	0.49
1:J:42:LYS:NZ	1:K:30:GLY:C	2.66	0.49
1:K:187:ASN:O	1:K:191:LYS:HB3	2.12	0.49
1:M:329:ALA:HA	1:M:334:MET:HG2	1.93	0.49
1:E:283:GLN:O	1:E:286:GLU:HB2	2.12	0.49
1:Y:187:ASN:O	1:Y:191:LYS:HB3	2.12	0.49
1:W:248:MET:CE	1:Y:14:LYS:HD3	2.41	0.49
1:K:234:ARG:NH2	1:L:86:ALA:HB2	2.27	0.49
1:W:329:ALA:HA	1:W:334:MET:HG2	1.93	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:Z:187:ASN:O	1:Z:191:LYS:HB3	2.12	0.49
1:O:283:GLN:O	1:O:286:GLU:HB2	2.11	0.49
1:D:187:ASN:O	1:D:191:LYS:HB3	2.12	0.49
1:Q:172:ALA:HB2	1:Q:253:ALA:HB1	1.93	0.49
1:A:23:TYR:CZ	1:Q:82:ILE:HB	2.47	0.49
1:W:285:VAL:CG1	1:Y:5:LYS:HD2	2.43	0.49
1:G:248:MET:CE	1:H:14:LYS:HD3	2.42	0.49
1:K:256:VAL:HG12	2:X:32:C:OP1	2.13	0.49
1:G:187:ASN:O	1:G:191:LYS:HB3	2.12	0.49
1:N:187:ASN:O	1:N:191:LYS:HB3	2.12	0.49
1:E:329:ALA:HA	1:E:334:MET:HG2	1.93	0.49
1:P:175:ASP:OD2	1:P:177:SER:HB3	2.13	0.49
1:H:175:ASP:OD2	1:H:177:SER:HB3	2.13	0.49
1:B:187:ASN:O	1:B:191:LYS:HB3	2.12	0.49
1:I:172:ALA:HB2	1:I:253:ALA:HB1	1.93	0.49
2:R:45:C:C4	2:R:46:C:N3	2.81	0.49
2:X:57:C:N4	2:X:58:C:C2	2.80	0.49
1:D:175:ASP:OD2	1:D:177:SER:HB3	2.13	0.49
1:J:175:ASP:OD2	1:J:177:SER:HB3	2.13	0.49
1:L:367:VAL:HG11	1:M:278:GLN:OE1	2.13	0.49
1:M:283:GLN:O	1:M:286:GLU:HB2	2.11	0.49
1:M:187:ASN:O	1:M:191:LYS:HB3	2.12	0.49
1:B:361:GLY:HA3	1:C:274:HIS:NE2	2.27	0.49
1:W:184:ARG:HD2	2:X:7:C:OP1	2.12	0.49
1:C:283:GLN:O	1:C:286:GLU:HB2	2.11	0.49
1:K:172:ALA:HB2	1:K:253:ALA:HB1	1.93	0.49
1:M:315:THR:OG1	2:X:45:C:OP2	2.21	0.49
1:P:248:MET:HE3	1:W:14:LYS:HD3	1.95	0.49
1:D:234:ARG:NH2	1:E:86:ALA:HB2	2.28	0.49
1:A:283:GLN:O	1:A:286:GLU:HB2	2.11	0.49
1:O:175:ASP:OD2	1:O:177:SER:HB3	2.13	0.49
1:I:175:ASP:OD2	1:I:177:SER:HB3	2.13	0.49
1:C:175:ASP:OD2	1:C:177:SER:HB3	2.13	0.49
2:R:9:C:C5	2:R:10:C:C5	3.01	0.48
1:F:175:ASP:OD2	1:F:177:SER:HB3	2.13	0.48
2:R:37:C:C5	2:R:38:C:C4	3.01	0.48
1:K:175:ASP:OD2	1:K:177:SER:HB3	2.13	0.48
1:A:187:ASN:O	1:A:191:LYS:HB3	2.12	0.48
1:E:175:ASP:OD2	1:E:177:SER:HB3	2.13	0.48
1:B:175:ASP:OD2	1:B:177:SER:HB3	2.13	0.48
1:L:175:ASP:OD2	1:L:177:SER:HB3	2.13	0.48
1:M:175:ASP:OD2	1:M:177:SER:HB3	2.13	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:W:337:TYR:CE1	2:X:4:C:H2'	2.49	0.48
2:R:6:C:C4	2:R:7:C:C4	3.02	0.48
1:Y:175:ASP:OD2	1:Y:177:SER:HB3	2.13	0.48
1:N:175:ASP:OD2	1:N:177:SER:HB3	2.13	0.48
1:W:175:ASP:OD2	1:W:177:SER:HB3	2.13	0.48
1:K:367:VAL:CG1	1:M:2:ALA:HB3	2.44	0.48
1:M:335:GLY:O	2:X:47:C:O2'	2.31	0.48
1:W:254:GLY:HA2	2:X:4:C:OP1	2.13	0.48
1:Q:175:ASP:OD2	1:Q:177:SER:HB3	2.13	0.48
1:A:175:ASP:OD2	1:A:177:SER:HB3	2.13	0.48
1:Y:42:LYS:HD2	1:Z:28:SER:HB3	1.96	0.48
1:W:242:ILE:HG12	2:X:8:C:N4	2.28	0.48
1:A:368:LEU:N	1:A:368:LEU:HD12	2.29	0.48
1:N:234:ARG:NH2	1:O:86:ALA:HB2	2.29	0.48
1:L:368:LEU:HD12	1:L:368:LEU:N	2.29	0.48
1:N:248:MET:HE3	1:O:14:LYS:HD3	1.95	0.48
1:G:175:ASP:OD2	1:G:177:SER:HB3	2.13	0.48
1:Z:175:ASP:OD2	1:Z:177:SER:HB3	2.13	0.48
2:X:59:C:C4	2:X:60:C:N4	2.82	0.48
1:B:234:ARG:NH2	1:C:86:ALA:HB2	2.29	0.48
1:Q:368:LEU:N	1:Q:368:LEU:HD12	2.29	0.48
2:R:18:C:O2'	2:R:19:C:H5'	2.14	0.48
1:P:234:ARG:HH21	1:W:86:ALA:HB2	1.79	0.48
2:X:24:C:H41	2:X:25:C:H42	1.62	0.47
1:J:368:LEU:N	1:J:368:LEU:HD12	2.29	0.47
1:M:368:LEU:N	1:M:368:LEU:HD12	2.29	0.47
1:Y:368:LEU:N	1:Y:368:LEU:HD12	2.29	0.47
1:B:315:THR:OG1	2:R:17:C:OP2	2.24	0.47
1:Y:303:ILE:HD12	1:Y:303:ILE:N	2.30	0.47
1:H:368:LEU:HD12	1:H:368:LEU:N	2.29	0.47
1:W:184:ARG:HH11	2:X:7:C:P	2.37	0.47
1:F:368:LEU:HD12	1:F:368:LEU:N	2.29	0.47
1:I:234:ARG:HH21	1:Q:86:ALA:HB2	1.78	0.47
1:B:368:LEU:HD12	1:B:368:LEU:N	2.29	0.47
1:N:368:LEU:HD12	1:N:368:LEU:N	2.29	0.47
1:C:368:LEU:N	1:C:368:LEU:HD12	2.29	0.47
2:R:23:C:C4	2:R:24:C:N3	2.83	0.47
1:F:303:ILE:N	1:F:303:ILE:HD12	2.30	0.47
1:G:303:ILE:N	1:G:303:ILE:HD12	2.30	0.47
1:Z:303:ILE:HD12	1:Z:303:ILE:N	2.30	0.47
2:X:51:C:N4	2:X:52:C:C2	2.82	0.47
1:W:367:VAL:HG12	1:Z:3:LEU:CD1	2.44	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:O:368:LEU:HD12	1:O:368:LEU:N	2.29	0.47
1:G:368:LEU:HD12	1:G:368:LEU:N	2.29	0.47
1:M:303:ILE:HD12	1:M:303:ILE:N	2.30	0.47
1:A:303:ILE:HD12	1:A:303:ILE:N	2.30	0.47
1:N:303:ILE:HD12	1:N:303:ILE:N	2.30	0.47
1:B:303:ILE:N	1:B:303:ILE:HD12	2.30	0.47
1:W:303:ILE:N	1:W:303:ILE:HD12	2.30	0.47
1:Z:368:LEU:N	1:Z:368:LEU:HD12	2.29	0.47
1:W:368:LEU:HD12	1:W:368:LEU:N	2.29	0.47
1:E:368:LEU:HD12	1:E:368:LEU:N	2.29	0.47
1:H:303:ILE:HD12	1:H:303:ILE:N	2.30	0.47
1:L:303:ILE:N	1:L:303:ILE:HD12	2.30	0.47
1:E:303:ILE:HD12	1:E:303:ILE:N	2.30	0.47
1:F:367:VAL:HG21	1:G:278:GLN:CD	2.34	0.47
1:E:38:TYR:CE1	1:F:26:GLN:HB2	2.50	0.47
1:J:248:MET:CE	1:K:14:LYS:HD3	2.45	0.47
1:K:368:LEU:N	1:K:368:LEU:HD12	2.29	0.47
1:O:303:ILE:HD12	1:O:303:ILE:N	2.30	0.47
1:Q:303:ILE:N	1:Q:303:ILE:HD12	2.30	0.47
1:J:303:ILE:HD12	1:J:303:ILE:N	2.30	0.46
2:R:9:C:H2'	2:R:10:C:O4'	2.14	0.46
2:X:15:C:C4	1:Y:242:ILE:HG12	2.50	0.46
1:E:361:GLY:HA3	1:F:274:HIS:NE2	2.30	0.46
1:A:361:GLY:HA3	1:B:274:HIS:NE2	2.30	0.46
1:A:42:LYS:HD2	1:B:28:SER:HB3	1.97	0.46
1:N:105:ASN:HB3	1:N:107:LYS:HE2	1.97	0.46
1:C:303:ILE:N	1:C:303:ILE:HD12	2.30	0.46
1:A:14:LYS:HD3	1:Q:248:MET:HE2	1.98	0.46
1:B:105:ASN:HB3	1:B:107:LYS:HE2	1.98	0.46
1:P:368:LEU:HD12	1:P:368:LEU:N	2.29	0.46
1:I:368:LEU:N	1:I:368:LEU:HD12	2.29	0.46
1:I:103:ASP:HA	1:I:108:GLU:HA	1.98	0.46
1:P:105:ASN:HB3	1:P:107:LYS:HE2	1.98	0.46
1:F:105:ASN:HB3	1:F:107:LYS:HE2	1.98	0.46
1:I:105:ASN:HB3	1:I:107:LYS:HE2	1.98	0.46
1:D:105:ASN:HB3	1:D:107:LYS:HE2	1.98	0.46
1:K:103:ASP:HA	1:K:108:GLU:HA	1.98	0.46
1:F:103:ASP:HA	1:F:108:GLU:HA	1.98	0.46
1:Y:103:ASP:HA	1:Y:108:GLU:HA	1.98	0.46
1:D:242:ILE:HG12	2:R:36:C:C4	2.51	0.46
1:P:270:ILE:HA	1:P:270:ILE:HD13	1.83	0.46
2:X:59:C:N4	2:X:60:C:C4	2.83	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:O:367:VAL:HG21	1:P:278:GLN:CD	2.36	0.46
1:K:105:ASN:HB3	1:K:107:LYS:HE2	1.98	0.46
1:M:103:ASP:HA	1:M:108:GLU:HA	1.98	0.46
1:K:303:ILE:HD12	1:K:303:ILE:N	2.30	0.46
1:P:303:ILE:N	1:P:303:ILE:HD12	2.30	0.46
2:R:23:C:H2'	2:R:24:C:O4'	2.16	0.46
1:F:107:LYS:HG3	1:F:107:LYS:H	1.58	0.46
1:Q:103:ASP:HA	1:Q:108:GLU:HA	1.98	0.46
1:M:120:THR:C	1:M:122:GLU:H	2.19	0.46
1:L:103:ASP:HA	1:L:108:GLU:HA	1.98	0.46
1:Y:105:ASN:HB3	1:Y:107:LYS:HE2	1.98	0.46
1:Z:107:LYS:H	1:Z:107:LYS:HG3	1.58	0.46
1:P:233:THR:HG22	1:W:307:PRO:CG	2.46	0.46
1:M:233:THR:HG22	1:N:307:PRO:HG2	1.97	0.46
1:H:105:ASN:HB3	1:H:107:LYS:HE2	1.98	0.46
1:C:103:ASP:HA	1:C:108:GLU:HA	1.98	0.46
1:P:103:ASP:HA	1:P:108:GLU:HA	1.98	0.46
1:J:105:ASN:HB3	1:J:107:LYS:HE2	1.98	0.46
1:A:103:ASP:HA	1:A:108:GLU:HA	1.98	0.46
1:D:103:ASP:HA	1:D:108:GLU:HA	1.98	0.46
1:G:107:LYS:HG3	1:G:107:LYS:H	1.58	0.46
2:X:65:C:N4	2:X:66:C:N3	2.64	0.46
1:D:368:LEU:N	1:D:368:LEU:HD12	2.29	0.46
1:G:105:ASN:HB3	1:G:107:LYS:HE2	1.98	0.46
1:O:103:ASP:HA	1:O:108:GLU:HA	1.98	0.46
1:M:105:ASN:HB3	1:M:107:LYS:HE2	1.98	0.46
1:I:303:ILE:N	1:I:303:ILE:HD12	2.30	0.46
1:Z:105:ASN:HB3	1:Z:107:LYS:HE2	1.98	0.46
1:Q:120:THR:C	1:Q:122:GLU:H	2.19	0.46
1:K:120:THR:C	1:K:122:GLU:H	2.19	0.46
1:L:120:THR:C	1:L:122:GLU:H	2.19	0.46
1:Q:270:ILE:HA	1:Q:270:ILE:HD13	1.83	0.46
1:L:270:ILE:HA	1:L:270:ILE:HD13	1.83	0.46
1:D:303:ILE:N	1:D:303:ILE:HD12	2.30	0.45
1:K:183:ILE:HD11	1:K:203:LYS:HG3	1.98	0.45
1:I:120:THR:C	1:I:122:GLU:H	2.19	0.45
1:C:120:THR:C	1:C:122:GLU:H	2.19	0.45
1:E:105:ASN:HB3	1:E:107:LYS:HE2	1.98	0.45
1:Z:120:THR:C	1:Z:122:GLU:H	2.19	0.45
2:X:59:C:H41	2:X:60:C:N4	2.05	0.45
1:J:103:ASP:HA	1:J:108:GLU:HA	1.98	0.45
1:O:120:THR:C	1:O:122:GLU:H	2.19	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:120:THR:C	1:G:122:GLU:H	2.19	0.45
1:Y:370:LEU:HG	1:Z:268:LYS:HD3	1.98	0.45
1:C:105:ASN:HB3	1:C:107:LYS:HE2	1.98	0.45
1:W:103:ASP:HA	1:W:108:GLU:HA	1.98	0.45
1:W:105:ASN:HB3	1:W:107:LYS:HE2	1.98	0.45
1:P:120:THR:C	1:P:122:GLU:H	2.19	0.45
1:E:103:ASP:HA	1:E:108:GLU:HA	1.98	0.45
1:A:105:ASN:HB3	1:A:107:LYS:HE2	1.98	0.45
1:N:361:GLY:HA3	1:O:274:HIS:NE2	2.31	0.45
1:I:183:ILE:HD11	1:I:203:LYS:HG3	1.98	0.45
1:D:270:ILE:HA	1:D:270:ILE:HD13	1.83	0.45
1:C:270:ILE:HA	1:C:270:ILE:HD13	1.83	0.45
1:O:367:VAL:HG11	1:P:278:GLN:CD	2.36	0.45
1:J:107:LYS:H	1:J:107:LYS:HG3	1.58	0.45
1:O:105:ASN:HB3	1:O:107:LYS:HE2	1.98	0.45
1:Q:183:ILE:HD11	1:Q:203:LYS:HG3	1.98	0.45
1:L:183:ILE:HD11	1:L:203:LYS:HG3	1.98	0.45
1:H:103:ASP:HA	1:H:108:GLU:HA	1.98	0.45
2:R:55:C:C4	2:R:56:C:N3	2.84	0.45
2:R:9:C:C4	2:R:10:C:C2	3.04	0.45
1:D:183:ILE:HD11	1:D:203:LYS:HG3	1.98	0.45
1:B:120:THR:C	1:B:122:GLU:H	2.19	0.45
1:B:103:ASP:HA	1:B:108:GLU:HA	1.98	0.45
1:D:120:THR:C	1:D:122:GLU:H	2.19	0.45
1:H:183:ILE:HD11	1:H:203:LYS:HG3	1.98	0.45
1:M:183:ILE:HD11	1:M:203:LYS:HG3	1.98	0.45
1:A:183:ILE:HD11	1:A:203:LYS:HG3	1.98	0.45
1:A:30:GLY:C	1:Q:42:LYS:HZ1	2.20	0.45
1:J:183:ILE:HD11	1:J:203:LYS:HG3	1.98	0.45
2:X:15:C:O2	1:Y:245:GLY:HA3	2.17	0.45
1:N:120:THR:C	1:N:122:GLU:H	2.19	0.45
1:Z:103:ASP:HA	1:Z:108:GLU:HA	1.98	0.45
1:N:103:ASP:HA	1:N:108:GLU:HA	1.98	0.45
1:O:270:ILE:HD13	1:O:270:ILE:HA	1.83	0.45
1:P:183:ILE:HD11	1:P:203:LYS:HG3	1.98	0.45
1:N:183:ILE:HD11	1:N:203:LYS:HG3	1.98	0.45
1:B:183:ILE:HD11	1:B:203:LYS:HG3	1.98	0.45
1:W:173:ALA:HB2	2:X:4:C:H4'	1.99	0.45
1:F:234:ARG:NH2	1:G:86:ALA:HB2	2.31	0.45
1:Y:107:LYS:HG3	1:Y:107:LYS:H	1.58	0.45
1:Q:105:ASN:HB3	1:Q:107:LYS:HE2	1.98	0.45
1:L:105:ASN:HB3	1:L:107:LYS:HE2	1.98	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:367:VAL:HG11	1:D:278:GLN:OE1	2.17	0.45
1:B:107:LYS:HG3	1:B:107:LYS:H	1.58	0.44
1:B:120:THR:CG2	1:B:122:GLU:HG2	2.47	0.44
1:N:120:THR:CG2	1:N:122:GLU:HG2	2.47	0.44
1:O:183:ILE:HD11	1:O:203:LYS:HG3	1.98	0.44
1:W:120:THR:C	1:W:122:GLU:H	2.19	0.44
1:G:103:ASP:HA	1:G:108:GLU:HA	1.98	0.44
2:X:57:C:N4	2:X:58:C:O2	2.50	0.44
2:X:24:C:N4	2:X:25:C:N4	2.63	0.44
1:I:285:VAL:CG1	1:Q:5:LYS:HD2	2.48	0.44
1:M:248:MET:HE2	1:N:14:LYS:HD3	1.99	0.44
1:M:120:THR:CG2	1:M:122:GLU:HG2	2.48	0.44
1:W:183:ILE:HD11	1:W:203:LYS:HG3	1.98	0.44
1:A:120:THR:C	1:A:122:GLU:H	2.19	0.44
1:J:33:ILE:HD11	1:J:218:HIS:NE2	2.33	0.44
1:P:254:GLY:HA2	2:X:67:C:OP1	2.18	0.44
1:O:368:LEU:CD1	1:O:368:LEU:H	2.30	0.44
1:B:184:ARG:NH1	2:R:21:C:OP2	2.50	0.44
1:J:307:PRO:HG2	1:Z:233:THR:HG22	1.98	0.44
1:N:107:LYS:H	1:N:107:LYS:HG3	1.58	0.44
1:A:120:THR:CG2	1:A:122:GLU:HG2	2.48	0.44
1:E:183:ILE:HD11	1:E:203:LYS:HG3	1.98	0.44
1:F:183:ILE:HD11	1:F:203:LYS:HG3	1.98	0.44
1:G:183:ILE:HD11	1:G:203:LYS:HG3	1.98	0.44
1:C:183:ILE:HD11	1:C:203:LYS:HG3	1.98	0.44
1:A:248:MET:HE3	1:B:14:LYS:HD3	2.00	0.44
1:H:74:LEU:HD23	1:I:25:ILE:HG22	1.99	0.44
1:C:368:LEU:H	1:C:368:LEU:CD1	2.30	0.44
1:Z:183:ILE:HD11	1:Z:203:LYS:HG3	1.99	0.44
1:F:120:THR:C	1:F:122:GLU:H	2.19	0.44
2:X:57:C:N4	2:X:58:C:N3	2.66	0.44
1:A:33:ILE:HD11	1:A:218:HIS:NE2	2.33	0.44
1:N:368:LEU:CD1	1:N:368:LEU:H	2.30	0.44
1:H:107:LYS:HG3	1:H:107:LYS:H	1.58	0.44
1:C:120:THR:CG2	1:C:122:GLU:HG2	2.47	0.44
1:O:120:THR:CG2	1:O:122:GLU:HG2	2.47	0.44
1:Y:183:ILE:HD11	1:Y:203:LYS:HG3	1.98	0.44
1:E:120:THR:C	1:E:122:GLU:H	2.19	0.44
1:K:319:HIS:O	1:K:323:VAL:HG12	2.18	0.44
1:A:23:TYR:CE1	1:Q:82:ILE:HG22	2.53	0.44
2:X:43:C:C4	2:X:44:C:N3	2.85	0.44
1:Q:33:ILE:HD11	1:Q:218:HIS:NE2	2.33	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:33:ILE:HD11	1:H:218:HIS:NE2	2.33	0.44
1:M:33:ILE:HD11	1:M:218:HIS:NE2	2.33	0.44
1:I:248:MET:HE2	1:Q:14:LYS:HD3	2.00	0.44
1:N:319:HIS:O	1:N:323:VAL:HG12	2.18	0.44
1:J:120:THR:C	1:J:122:GLU:H	2.19	0.44
1:I:319:HIS:O	1:I:323:VAL:HG12	2.18	0.44
1:Y:240:GLU:OE2	1:Z:27:ARG:HD3	2.17	0.44
1:D:319:HIS:O	1:D:323:VAL:HG12	2.18	0.44
1:C:33:ILE:HD11	1:C:218:HIS:NE2	2.33	0.44
1:C:38:TYR:CD1	1:D:26:GLN:HB2	2.52	0.44
1:B:368:LEU:H	1:B:368:LEU:CD1	2.30	0.44
1:K:120:THR:CG2	1:K:122:GLU:HG2	2.48	0.44
1:P:120:THR:CG2	1:P:122:GLU:HG2	2.48	0.44
1:F:120:THR:CG2	1:F:122:GLU:HG2	2.48	0.44
1:K:2:ALA:HB3	1:Z:367:VAL:HG12	1.97	0.44
1:O:33:ILE:HD11	1:O:218:HIS:NE2	2.33	0.44
1:E:33:ILE:HD11	1:E:218:HIS:NE2	2.33	0.44
1:F:367:VAL:CG1	1:H:2:ALA:HB3	2.48	0.44
1:H:42:LYS:HZ2	1:I:30:GLY:C	2.21	0.44
1:H:120:THR:C	1:H:122:GLU:H	2.19	0.44
1:B:319:HIS:O	1:B:323:VAL:HG12	2.18	0.44
1:F:270:ILE:HA	1:F:270:ILE:HD13	1.83	0.44
1:Y:270:ILE:HD13	1:Y:270:ILE:HA	1.83	0.44
1:L:33:ILE:HD11	1:L:218:HIS:NE2	2.33	0.43
2:R:6:C:C5	2:R:7:C:C4	3.06	0.43
1:M:107:LYS:H	1:M:107:LYS:HG3	1.58	0.43
1:D:120:THR:CG2	1:D:122:GLU:HG2	2.48	0.43
1:Y:120:THR:CG2	1:Y:122:GLU:HG2	2.48	0.43
2:X:37:C:H2'	2:X:38:C:O4'	2.17	0.43
1:W:33:ILE:HD11	1:W:218:HIS:NE2	2.33	0.43
1:F:33:ILE:HD11	1:F:218:HIS:NE2	2.33	0.43
2:X:41:C:C4	2:X:42:C:C5	3.06	0.43
2:X:14:C:OP2	1:Y:184:ARG:NH1	2.51	0.43
1:A:368:LEU:CD1	1:A:368:LEU:H	2.30	0.43
2:X:20:C:N4	2:X:21:C:N4	2.66	0.43
1:L:120:THR:CG2	1:L:122:GLU:HG2	2.47	0.43
1:I:120:THR:CG2	1:I:122:GLU:HG2	2.48	0.43
1:G:120:THR:CG2	1:G:122:GLU:HG2	2.48	0.43
1:C:107:LYS:HG3	1:C:107:LYS:H	1.58	0.43
1:H:120:THR:CG2	1:H:122:GLU:HG2	2.47	0.43
1:W:234:ARG:HH21	1:Y:86:ALA:HB2	1.83	0.43
1:Z:153:SER:HA	1:Z:154:PRO:HD2	1.86	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:Y:319:HIS:O	1:Y:323:VAL:HG12	2.18	0.43
1:P:319:HIS:O	1:P:323:VAL:HG12	2.18	0.43
1:F:319:HIS:O	1:F:323:VAL:HG12	2.18	0.43
1:K:33:ILE:HD11	1:K:218:HIS:NE2	2.33	0.43
1:I:33:ILE:HD11	1:I:218:HIS:NE2	2.33	0.43
1:Z:120:THR:CG2	1:Z:122:GLU:HG2	2.48	0.43
1:Y:120:THR:C	1:Y:122:GLU:H	2.19	0.43
1:D:41:GLN:NE2	1:D:73:ARG:HA	2.34	0.43
1:O:41:GLN:NE2	1:O:73:ARG:HA	2.34	0.43
1:D:33:ILE:HD11	1:D:218:HIS:NE2	2.33	0.43
1:P:33:ILE:HD11	1:P:218:HIS:NE2	2.33	0.43
1:P:248:MET:HE2	1:W:14:LYS:HD3	2.00	0.43
1:J:368:LEU:H	1:J:368:LEU:CD1	2.30	0.43
1:H:368:LEU:CD1	1:H:368:LEU:H	2.30	0.43
1:A:244:ALA:HB1	1:B:307:PRO:HB3	2.00	0.43
1:A:307:PRO:HB3	1:Q:244:ALA:HB1	1.99	0.43
1:W:120:THR:CG2	1:W:122:GLU:HG2	2.48	0.43
1:J:120:THR:CG2	1:J:122:GLU:HG2	2.48	0.43
1:C:41:GLN:NE2	1:C:73:ARG:HA	2.34	0.43
1:Y:41:GLN:NE2	1:Y:73:ARG:HA	2.34	0.43
1:F:41:GLN:NE2	1:F:73:ARG:HA	2.34	0.43
1:F:256:VAL:HG12	2:R:46:C:OP1	2.18	0.43
1:Y:33:ILE:HD11	1:Y:218:HIS:NE2	2.33	0.43
1:M:368:LEU:H	1:M:368:LEU:CD1	2.30	0.43
1:Y:368:LEU:H	1:Y:368:LEU:CD1	2.30	0.43
2:X:6:C:C5	2:X:7:C:C4	3.06	0.43
1:Q:120:THR:CG2	1:Q:122:GLU:HG2	2.48	0.43
1:E:120:THR:CG2	1:E:122:GLU:HG2	2.48	0.43
1:P:41:GLN:NE2	1:P:73:ARG:HA	2.34	0.43
1:G:319:HIS:O	1:G:323:VAL:HG12	2.18	0.43
1:Z:319:HIS:O	1:Z:323:VAL:HG12	2.18	0.43
1:H:319:HIS:O	1:H:323:VAL:HG12	2.18	0.43
1:Q:242:ILE:HG12	2:R:8:C:C4	2.53	0.43
1:J:319:HIS:O	1:J:323:VAL:HG12	2.18	0.43
1:E:367:VAL:CG1	1:G:3:LEU:HD12	2.47	0.43
1:K:368:LEU:CD1	1:K:368:LEU:H	2.30	0.43
1:I:153:SER:HA	1:I:154:PRO:HD2	1.85	0.43
1:E:319:HIS:O	1:E:323:VAL:HG12	2.18	0.43
1:A:270:ILE:HA	1:A:270:ILE:HD13	1.83	0.43
1:Z:33:ILE:HD11	1:Z:218:HIS:NE2	2.33	0.43
1:N:33:ILE:HD11	1:N:218:HIS:NE2	2.33	0.43
2:R:20:C:C5	2:R:21:C:C4	3.06	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:K:38:TYR:CE1	1:L:26:GLN:HB2	2.54	0.43
1:G:41:GLN:NE2	1:G:73:ARG:HA	2.34	0.43
2:R:62:C:C4	2:R:63:C:C4	3.07	0.43
1:N:41:GLN:NE2	1:N:73:ARG:HA	2.34	0.43
1:B:41:GLN:NE2	1:B:73:ARG:HA	2.34	0.43
2:R:69:C:N3	2:R:70:C:C2	2.86	0.43
1:G:33:ILE:HD11	1:G:218:HIS:NE2	2.33	0.43
1:C:184:ARG:HD2	2:R:28:C:OP1	2.19	0.43
1:F:368:LEU:CD1	1:F:368:LEU:H	2.30	0.43
1:G:153:SER:HA	1:G:154:PRO:HD2	1.85	0.43
1:I:41:GLN:NE2	1:I:73:ARG:HA	2.34	0.43
1:K:41:GLN:NE2	1:K:73:ARG:HA	2.34	0.43
1:O:319:HIS:O	1:O:323:VAL:HG12	2.18	0.43
1:Z:41:GLN:NE2	1:Z:73:ARG:HA	2.34	0.43
1:D:236:GLY:O	1:E:305:ASN:CB	2.46	0.43
1:B:33:ILE:HD11	1:B:218:HIS:NE2	2.33	0.43
1:L:38:TYR:CD1	1:M:26:GLN:HB2	2.53	0.43
1:L:368:LEU:CD1	1:L:368:LEU:H	2.30	0.43
1:L:42:LYS:HZ2	1:M:30:GLY:C	2.21	0.43
1:B:153:SER:HA	1:B:154:PRO:HD2	1.85	0.43
1:W:319:HIS:O	1:W:323:VAL:HG12	2.18	0.43
1:I:234:ARG:CZ	1:Q:225:HIS:CE1	3.02	0.43
1:Q:337:TYR:CE1	2:R:4:C:H2'	2.54	0.43
1:Q:368:LEU:CD1	1:Q:368:LEU:H	2.30	0.43
1:P:233:THR:O	1:W:307:PRO:HD2	2.19	0.43
1:W:285:VAL:HG11	1:Y:5:LYS:HD2	1.99	0.43
1:A:319:HIS:O	1:A:323:VAL:HG12	2.18	0.43
1:L:319:HIS:O	1:L:323:VAL:HG12	2.18	0.43
1:C:319:HIS:O	1:C:323:VAL:HG12	2.18	0.43
1:Q:319:HIS:O	1:Q:323:VAL:HG12	2.18	0.43
1:H:41:GLN:NE2	1:H:73:ARG:HA	2.34	0.43
1:M:319:HIS:O	1:M:323:VAL:HG12	2.18	0.43
1:I:368:LEU:CD1	1:I:368:LEU:H	2.30	0.42
1:B:248:MET:HE3	1:C:14:LYS:HD3	2.00	0.42
1:J:41:GLN:NE2	1:J:73:ARG:HA	2.34	0.42
1:A:21:SER:HB2	1:Q:228:ILE:HG22	2.00	0.42
1:K:233:THR:HG22	1:L:307:PRO:HG2	2.00	0.42
1:Q:41:GLN:NE2	1:Q:73:ARG:HA	2.34	0.42
1:M:41:GLN:NE2	1:M:73:ARG:HA	2.34	0.42
1:A:41:GLN:NE2	1:A:73:ARG:HA	2.34	0.42
1:M:270:ILE:HD13	1:M:270:ILE:HA	1.83	0.42
1:D:368:LEU:H	1:D:368:LEU:CD1	2.30	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:W:184:ARG:NH1	2:X:7:C:P	2.92	0.42
1:O:107:LYS:HG3	1:O:107:LYS:H	1.58	0.42
1:W:41:GLN:NE2	1:W:73:ARG:HA	2.34	0.42
1:L:41:GLN:NE2	1:L:73:ARG:HA	2.34	0.42
1:C:215:LYS:HE3	1:C:216:HIS:CE1	2.48	0.42
2:X:33:C:C4	2:X:34:C:C4	3.07	0.42
1:D:164:ALA:HB2	1:D:209:PHE:CZ	2.55	0.42
1:Y:230:GLN:HA	1:Z:25:ILE:HG12	2.01	0.42
1:E:41:GLN:NE2	1:E:73:ARG:HA	2.34	0.42
1:A:253:ALA:HB3	1:A:303:ILE:CD1	2.21	0.42
1:E:368:LEU:CD1	1:E:368:LEU:H	2.30	0.42
2:X:33:C:N4	2:X:34:C:C4	2.88	0.42
1:P:164:ALA:HB2	1:P:209:PHE:CZ	2.55	0.42
1:A:107:LYS:HG3	1:A:107:LYS:H	1.58	0.42
2:R:68:C:H2'	2:R:69:C:O4'	2.18	0.42
2:R:65:C:N3	2:R:66:C:C2	2.88	0.42
1:D:215:LYS:HE3	1:D:216:HIS:CE1	2.48	0.42
1:Y:164:ALA:HB2	1:Y:209:PHE:CZ	2.55	0.42
1:A:164:ALA:HB2	1:A:209:PHE:CZ	2.55	0.42
1:C:42:LYS:HZ1	1:D:30:GLY:C	2.22	0.42
1:L:233:THR:HG22	1:M:307:PRO:HG2	2.00	0.42
1:P:107:LYS:H	1:P:107:LYS:HG3	1.58	0.42
1:N:153:SER:HA	1:N:154:PRO:HD2	1.85	0.42
1:A:184:ARG:HD2	2:R:14:C:OP1	2.19	0.42
1:I:270:ILE:HA	1:I:270:ILE:HD13	1.83	0.42
1:W:368:LEU:CD1	1:W:368:LEU:H	2.30	0.42
2:R:55:C:H2'	2:R:56:C:O4'	2.19	0.42
1:F:164:ALA:HB2	1:F:209:PHE:CZ	2.55	0.42
1:M:164:ALA:HB2	1:M:209:PHE:CZ	2.55	0.42
1:N:164:ALA:HB2	1:N:209:PHE:CZ	2.55	0.42
1:J:270:ILE:HA	1:J:270:ILE:HD13	1.83	0.42
2:X:2:C:H2'	2:X:2:C:O2	2.19	0.42
1:O:215:LYS:HE3	1:O:216:HIS:CE1	2.48	0.42
1:I:164:ALA:HB2	1:I:209:PHE:CZ	2.55	0.42
1:B:164:ALA:HB2	1:B:209:PHE:CZ	2.55	0.42
1:J:333:ILE:H	1:J:333:ILE:HG13	1.70	0.42
2:X:41:C:H41	2:X:42:C:N4	2.16	0.42
1:P:337:TYR:CE1	2:X:67:C:H2'	2.55	0.42
1:N:6:VAL:HG22	1:N:6:VAL:O	2.20	0.42
1:P:368:LEU:CD1	1:P:368:LEU:H	2.30	0.42
1:O:164:ALA:HB2	1:O:209:PHE:CZ	2.55	0.42
1:C:164:ALA:HB2	1:C:209:PHE:CZ	2.55	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:K:164:ALA:HB2	1:K:209:PHE:CZ	2.55	0.42
1:J:164:ALA:HB2	1:J:209:PHE:CZ	2.55	0.42
1:A:307:PRO:HG2	1:Q:233:THR:HG22	2.02	0.42
1:B:6:VAL:O	1:B:6:VAL:HG22	2.20	0.42
1:K:6:VAL:O	1:K:6:VAL:HG22	2.20	0.42
1:C:166:LEU:HD21	1:C:246:LEU:HD22	2.02	0.42
1:D:248:MET:HE3	1:E:14:LYS:HD3	2.02	0.42
2:X:62:C:N4	2:X:63:C:N4	2.68	0.42
1:E:107:LYS:HG3	1:E:107:LYS:H	1.58	0.42
1:F:184:ARG:NH1	2:R:49:C:OP2	2.53	0.42
1:Q:333:ILE:O	2:R:5:C:O2	2.38	0.42
1:Y:281:MET:HA	1:Y:284:VAL:HG22	2.02	0.42
1:M:253:ALA:HB3	1:M:303:ILE:CD1	2.21	0.41
1:L:234:ARG:CZ	1:M:225:HIS:HE1	2.33	0.41
2:R:27:C:N4	2:R:28:C:N3	2.68	0.41
1:F:215:LYS:HE3	1:F:216:HIS:CE1	2.48	0.41
1:P:215:LYS:HE3	1:P:216:HIS:CE1	2.48	0.41
1:I:6:VAL:O	1:I:6:VAL:HG22	2.20	0.41
1:G:166:LEU:HD21	1:G:246:LEU:HD22	2.02	0.41
1:I:166:LEU:HD21	1:I:246:LEU:HD22	2.02	0.41
1:L:166:LEU:HD21	1:L:246:LEU:HD22	2.02	0.41
1:Z:166:LEU:HD21	1:Z:246:LEU:HD22	2.02	0.41
1:W:166:LEU:HD21	1:W:246:LEU:HD22	2.02	0.41
1:E:166:LEU:HD21	1:E:246:LEU:HD22	2.02	0.41
1:M:166:LEU:HD21	1:M:246:LEU:HD22	2.02	0.41
1:H:166:LEU:HD21	1:H:246:LEU:HD22	2.02	0.41
1:O:166:LEU:HD21	1:O:246:LEU:HD22	2.02	0.41
1:J:166:LEU:HD21	1:J:246:LEU:HD22	2.02	0.41
1:O:42:LYS:HZ1	1:P:30:GLY:C	2.22	0.41
1:H:164:ALA:HB2	1:H:209:PHE:CZ	2.55	0.41
2:X:19:C:H2'	2:X:20:C:O4'	2.20	0.41
1:J:42:LYS:HZ1	1:K:30:GLY:C	2.22	0.41
1:I:285:VAL:HG11	1:Q:5:LYS:HD2	2.01	0.41
1:W:38:TYR:CE1	1:Y:26:GLN:HB2	2.54	0.41
1:F:281:MET:HA	1:F:284:VAL:HG22	2.02	0.41
1:H:281:MET:HA	1:H:284:VAL:HG22	2.02	0.41
1:D:6:VAL:HG22	1:D:6:VAL:O	2.20	0.41
1:F:166:LEU:HD21	1:F:246:LEU:HD22	2.02	0.41
1:Q:166:LEU:HD21	1:Q:246:LEU:HD22	2.02	0.41
1:N:166:LEU:HD21	1:N:246:LEU:HD22	2.02	0.41
1:B:166:LEU:HD21	1:B:246:LEU:HD22	2.02	0.41
2:X:20:C:N4	2:X:21:C:C4	2.88	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:367:VAL:HG21	1:I:278:GLN:CD	2.40	0.41
1:N:265:LYS:HE3	1:O:4:SER:HA	2.03	0.41
1:K:270:ILE:HA	1:K:270:ILE:HD13	1.83	0.41
1:H:270:ILE:HD13	1:H:270:ILE:HA	1.83	0.41
1:C:82:ILE:HG22	1:D:23:TYR:CD1	2.55	0.41
1:P:6:VAL:O	1:P:6:VAL:HG22	2.20	0.41
1:A:166:LEU:HD21	1:A:246:LEU:HD22	2.02	0.41
1:E:164:ALA:HB2	1:E:209:PHE:CZ	2.55	0.41
1:J:281:MET:HA	1:J:284:VAL:HG22	2.03	0.41
1:M:325:LEU:HD12	1:M:341:PRO:HD3	2.03	0.41
1:M:367:VAL:HG11	1:N:278:GLN:OE1	2.20	0.41
1:I:112:GLU:HA	1:I:112:GLU:OE1	2.21	0.41
1:L:112:GLU:HA	1:L:112:GLU:OE1	2.21	0.41
1:A:23:TYR:CE2	1:Q:82:ILE:HB	2.55	0.41
1:Y:215:LYS:HE3	1:Y:216:HIS:CE1	2.48	0.41
2:R:39:C:H6	2:R:39:C:H2'	1.73	0.41
1:Y:166:LEU:HD21	1:Y:246:LEU:HD22	2.02	0.41
1:K:166:LEU:HD21	1:K:246:LEU:HD22	2.03	0.41
1:A:315:THR:OG1	2:R:10:C:OP2	2.23	0.41
1:W:164:ALA:HB2	1:W:209:PHE:CZ	2.55	0.41
1:N:233:THR:HG22	1:O:307:PRO:HG2	2.01	0.41
1:K:185:ARG:NH1	2:X:35:C:C5	2.88	0.41
1:H:234:ARG:HH21	1:I:86:ALA:HB2	1.84	0.41
1:P:112:GLU:HA	1:P:112:GLU:OE1	2.21	0.41
1:W:112:GLU:HA	1:W:112:GLU:OE1	2.21	0.41
1:K:112:GLU:HA	1:K:112:GLU:OE1	2.21	0.41
1:F:230:GLN:HA	1:G:25:ILE:HG12	2.03	0.41
1:C:281:MET:HA	1:C:284:VAL:HG22	2.02	0.41
1:F:325:LEU:HD12	1:F:341:PRO:HD3	2.03	0.41
1:A:281:MET:HA	1:A:284:VAL:HG22	2.02	0.41
1:H:333:ILE:HG13	1:H:333:ILE:H	1.70	0.41
1:Q:112:GLU:OE1	1:Q:112:GLU:HA	2.21	0.41
1:E:112:GLU:OE1	1:E:112:GLU:HA	2.21	0.41
1:C:333:ILE:H	1:C:333:ILE:HG13	1.70	0.41
1:A:112:GLU:HA	1:A:112:GLU:OE1	2.21	0.41
1:W:367:VAL:HG12	1:Z:3:LEU:HD12	2.02	0.41
1:W:215:LYS:HE3	1:W:216:HIS:CE1	2.48	0.41
1:D:166:LEU:HD21	1:D:246:LEU:HD22	2.02	0.41
1:P:166:LEU:HD21	1:P:246:LEU:HD22	2.02	0.41
1:F:363:ILE:HA	1:G:274:HIS:HA	2.02	0.41
1:L:164:ALA:HB2	1:L:209:PHE:CZ	2.55	0.41
1:O:367:VAL:HG11	1:P:278:GLN:HE22	1.84	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:233:THR:HG22	1:B:307:PRO:HG2	2.02	0.41
1:Y:325:LEU:HD12	1:Y:341:PRO:HD3	2.03	0.41
1:A:325:LEU:HD12	1:A:341:PRO:HD3	2.03	0.41
1:E:248:MET:CE	1:F:14:LYS:HD3	2.51	0.41
1:N:325:LEU:HD12	1:N:341:PRO:HD3	2.03	0.41
1:B:325:LEU:HD12	1:B:341:PRO:HD3	2.03	0.41
1:M:281:MET:HA	1:M:284:VAL:HG22	2.02	0.41
1:O:281:MET:HA	1:O:284:VAL:HG22	2.03	0.41
1:D:112:GLU:HA	1:D:112:GLU:OE1	2.21	0.41
1:O:333:ILE:HG13	1:O:333:ILE:H	1.70	0.41
1:I:236:GLY:O	1:Q:305:ASN:CB	2.54	0.41
1:P:184:ARG:HD2	2:X:70:C:OP1	2.20	0.41
1:E:215:LYS:HE3	1:E:216:HIS:CE1	2.48	0.41
1:Q:164:ALA:HB2	1:Q:209:PHE:CZ	2.55	0.41
1:P:307:PRO:C	1:P:309:ALA:H	2.24	0.41
1:I:233:THR:HG22	1:Q:307:PRO:CG	2.51	0.41
1:Z:325:LEU:HD12	1:Z:341:PRO:HD3	2.03	0.41
1:G:38:TYR:CE1	1:H:26:GLN:HB2	2.56	0.41
1:B:281:MET:HA	1:B:284:VAL:HG22	2.02	0.41
1:K:242:ILE:HG12	2:X:36:C:C4	2.56	0.41
1:F:112:GLU:HA	1:F:112:GLU:OE1	2.21	0.41
1:E:6:VAL:O	1:E:6:VAL:HG22	2.20	0.41
1:D:307:PRO:C	1:D:309:ALA:H	2.24	0.41
1:O:307:PRO:C	1:O:309:ALA:H	2.24	0.41
1:G:325:LEU:HD12	1:G:341:PRO:HD3	2.03	0.41
1:J:46:LYS:O	1:J:50:MET:HG3	2.21	0.41
1:A:46:LYS:O	1:A:50:MET:HG3	2.21	0.41
1:Q:168:ILE:HG12	1:Q:223:PHE:CE2	2.56	0.41
1:G:6:VAL:O	1:G:6:VAL:HG22	2.20	0.41
1:Y:6:VAL:O	1:Y:6:VAL:HG22	2.20	0.41
1:G:164:ALA:HB2	1:G:209:PHE:CZ	2.55	0.41
1:Z:164:ALA:HB2	1:Z:209:PHE:CZ	2.55	0.41
1:I:307:PRO:C	1:I:309:ALA:H	2.24	0.41
1:H:307:PRO:C	1:H:309:ALA:H	2.24	0.41
1:K:307:PRO:C	1:K:309:ALA:H	2.24	0.41
1:J:307:PRO:C	1:J:309:ALA:H	2.24	0.41
1:C:307:PRO:C	1:C:309:ALA:H	2.25	0.41
1:P:46:LYS:O	1:P:50:MET:HG3	2.21	0.41
1:O:46:LYS:O	1:O:50:MET:HG3	2.21	0.41
1:D:46:LYS:O	1:D:50:MET:HG3	2.21	0.41
2:X:22:C:H2'	2:X:23:C:O4'	2.21	0.41
2:X:27:C:O2'	2:X:28:C:H5'	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:K:46:LYS:O	1:K:50:MET:HG3	2.21	0.41
1:I:46:LYS:O	1:I:50:MET:HG3	2.21	0.41
1:L:281:MET:HA	1:L:284:VAL:HG22	2.02	0.41
2:R:51:C:H2'	2:R:52:C:O4'	2.21	0.41
1:M:46:LYS:O	1:M:50:MET:HG3	2.21	0.41
1:G:281:MET:HA	1:G:284:VAL:HG22	2.02	0.41
1:I:168:ILE:HG12	1:I:223:PHE:CE2	2.56	0.41
1:K:168:ILE:HG12	1:K:223:PHE:CE2	2.56	0.41
1:N:281:MET:HA	1:N:284:VAL:HG22	2.02	0.41
1:L:168:ILE:HG12	1:L:223:PHE:CE2	2.56	0.41
1:H:46:LYS:O	1:H:50:MET:HG3	2.21	0.41
1:E:270:ILE:HA	1:E:270:ILE:HD13	1.83	0.41
1:Q:281:MET:HA	1:Q:284:VAL:HG22	2.03	0.41
1:F:46:LYS:O	1:F:50:MET:HG3	2.21	0.41
1:Y:46:LYS:O	1:Y:50:MET:HG3	2.21	0.41
1:O:325:LEU:HD12	1:O:341:PRO:HD3	2.03	0.41
1:Z:281:MET:HA	1:Z:284:VAL:HG22	2.02	0.41
1:Y:112:GLU:OE1	1:Y:112:GLU:HA	2.21	0.41
1:M:112:GLU:OE1	1:M:112:GLU:HA	2.21	0.41
1:K:253:ALA:HB3	1:K:303:ILE:CD1	2.20	0.41
1:J:215:LYS:HE3	1:J:216:HIS:CE1	2.48	0.41
1:E:337:TYR:CD1	2:R:39:C:H2'	2.56	0.41
1:O:233:THR:HG22	1:P:307:PRO:HG2	2.02	0.41
1:W:107:LYS:H	1:W:107:LYS:HG3	1.58	0.41
1:C:46:LYS:O	1:C:50:MET:HG3	2.21	0.41
1:J:26:GLN:HB2	1:Z:38:TYR:CE1	2.56	0.41
1:C:325:LEU:HD12	1:C:341:PRO:HD3	2.03	0.41
1:H:256:VAL:HG21	2:R:61:C:C5	2.56	0.41
1:N:173:ALA:HB2	2:X:53:C:H4'	2.03	0.41
1:J:325:LEU:HD12	1:J:341:PRO:HD3	2.03	0.41
1:E:168:ILE:HG12	1:E:223:PHE:CE2	2.56	0.41
1:P:168:ILE:HG12	1:P:223:PHE:CE2	2.56	0.41
1:O:112:GLU:OE1	1:O:112:GLU:HA	2.21	0.41
1:C:112:GLU:OE1	1:C:112:GLU:HA	2.21	0.41
1:J:38:TYR:HE1	1:K:26:GLN:HB2	1.83	0.40
1:Z:6:VAL:O	1:Z:6:VAL:HG22	2.20	0.40
1:W:6:VAL:HG22	1:W:6:VAL:O	2.20	0.40
1:Q:307:PRO:C	1:Q:309:ALA:H	2.24	0.40
1:B:233:THR:HG22	1:C:307:PRO:HG2	2.01	0.40
1:L:307:PRO:C	1:L:309:ALA:H	2.24	0.40
1:I:107:LYS:H	1:I:107:LYS:HG3	1.58	0.40
1:H:184:ARG:NH1	2:R:63:C:OP2	2.54	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:P:281:MET:HA	1:P:284:VAL:HG22	2.02	0.40
1:D:281:MET:HA	1:D:284:VAL:HG22	2.02	0.40
1:I:59:HIS:O	1:I:208:SER:HB2	2.22	0.40
1:H:168:ILE:HG12	1:H:223:PHE:CE2	2.56	0.40
1:W:59:HIS:O	1:W:208:SER:HB2	2.22	0.40
1:K:59:HIS:O	1:K:208:SER:HB2	2.22	0.40
1:J:168:ILE:HG12	1:J:223:PHE:CE2	2.56	0.40
1:A:38:TYR:CE1	1:B:26:GLN:HB2	2.56	0.40
1:W:168:ILE:HG12	1:W:223:PHE:CE2	2.56	0.40
1:B:46:LYS:O	1:B:50:MET:HG3	2.21	0.40
1:D:168:ILE:HG12	1:D:223:PHE:CE2	2.56	0.40
1:O:135:TYR:CZ	1:O:145:VAL:HG21	2.57	0.40
1:C:135:TYR:CZ	1:C:145:VAL:HG21	2.57	0.40
1:P:135:TYR:CZ	1:P:145:VAL:HG21	2.57	0.40
1:L:135:TYR:CZ	1:L:145:VAL:HG21	2.57	0.40
1:F:59:HIS:O	1:F:208:SER:HB2	2.22	0.40
1:O:168:ILE:HG12	1:O:223:PHE:CE2	2.56	0.40
1:C:168:ILE:HG12	1:C:223:PHE:CE2	2.56	0.40
1:N:46:LYS:O	1:N:50:MET:HG3	2.21	0.40
1:Z:46:LYS:O	1:Z:50:MET:HG3	2.21	0.40
1:W:270:ILE:HD13	1:W:270:ILE:HA	1.83	0.40
1:O:254:GLY:HA2	2:X:59:C:O3'	2.21	0.40
1:F:6:VAL:HG22	1:F:6:VAL:O	2.20	0.40
1:Q:135:TYR:CZ	1:Q:145:VAL:HG21	2.57	0.40
1:L:42:LYS:HZ1	1:M:30:GLY:C	2.25	0.40
1:A:307:PRO:C	1:A:309:ALA:H	2.24	0.40
1:P:59:HIS:O	1:P:208:SER:HB2	2.22	0.40
1:H:325:LEU:HD12	1:H:341:PRO:HD3	2.03	0.40
1:H:242:ILE:HG12	2:R:64:C:C4	2.56	0.40
2:X:54:C:N4	2:X:55:C:N4	2.70	0.40
1:Y:59:HIS:O	1:Y:208:SER:HB2	2.22	0.40
1:O:248:MET:CE	1:P:14:LYS:HD3	2.52	0.40
1:N:112:GLU:OE1	1:N:112:GLU:HA	2.21	0.40
1:Y:135:TYR:CZ	1:Y:145:VAL:HG21	2.57	0.40
1:C:59:HIS:O	1:C:208:SER:HB2	2.22	0.40
1:D:59:HIS:O	1:D:208:SER:HB2	2.22	0.40
1:E:59:HIS:O	1:E:208:SER:HB2	2.22	0.40
1:O:48:CYS:SG	1:O:160:ILE:HD12	2.62	0.40
1:G:46:LYS:O	1:G:50:MET:HG3	2.21	0.40
1:H:112:GLU:HA	1:H:112:GLU:OE1	2.21	0.40
1:B:112:GLU:OE1	1:B:112:GLU:HA	2.21	0.40
1:O:59:HIS:O	1:O:208:SER:HB2	2.22	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:J:112:GLU:OE1	1:J:112:GLU:HA	2.21	0.40
2:X:51:C:N3	2:X:52:C:C2	2.90	0.40
1:H:215:LYS:HE3	1:H:216:HIS:CE1	2.48	0.40
1:L:6:VAL:O	1:L:6:VAL:HG22	2.20	0.40
1:Q:6:VAL:HG22	1:Q:6:VAL:O	2.20	0.40
1:Z:368:LEU:CD1	1:Z:368:LEU:H	2.30	0.40
1:A:14:LYS:HD3	1:Q:248:MET:HE3	2.03	0.40
1:H:135:TYR:CZ	1:H:145:VAL:HG21	2.57	0.40
1:G:233:THR:HG22	1:H:307:PRO:HG2	2.02	0.40
1:G:307:PRO:C	1:G:309:ALA:H	2.24	0.40
1:W:281:MET:HA	1:W:284:VAL:HG22	2.02	0.40
1:E:281:MET:HA	1:E:284:VAL:HG22	2.02	0.40
1:W:46:LYS:O	1:W:50:MET:HG3	2.21	0.40
1:E:46:LYS:O	1:E:50:MET:HG3	2.21	0.40
2:R:47:C:N4	2:R:48:C:N4	2.70	0.40
1:B:168:ILE:HG12	1:B:223:PHE:CE2	2.56	0.40
1:C:48:CYS:SG	1:C:160:ILE:HD12	2.62	0.40
1:M:59:HIS:O	1:M:208:SER:HB2	2.22	0.40
1:A:59:HIS:O	1:A:208:SER:HB2	2.22	0.40
1:M:45:ASN:HB2	1:M:69:TYR:CE1	2.57	0.40
1:H:59:HIS:O	1:H:208:SER:HB2	2.22	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:M:140:LYS:NZ	1:Y:140:LYS:NZ[1_545]	1.76	0.44
1:A:140:LYS:NZ	1:F:140:LYS:NZ[1_565]	1.84	0.36

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	373/375 (100%)	361 (97%)	11 (3%)	1 (0%)	50	92
1	B	373/375 (100%)	361 (97%)	11 (3%)	1 (0%)	50	92

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	373/375 (100%)	361 (97%)	11 (3%)	1 (0%)	50	92
1	D	373/375 (100%)	361 (97%)	11 (3%)	1 (0%)	50	92
1	E	373/375 (100%)	360 (96%)	12 (3%)	1 (0%)	50	92
1	F	373/375 (100%)	360 (96%)	12 (3%)	1 (0%)	50	92
1	G	373/375 (100%)	361 (97%)	11 (3%)	1 (0%)	50	92
1	H	373/375 (100%)	361 (97%)	11 (3%)	1 (0%)	50	92
1	I	373/375 (100%)	360 (96%)	12 (3%)	1 (0%)	50	92
1	J	373/375 (100%)	361 (97%)	11 (3%)	1 (0%)	50	92
1	K	373/375 (100%)	360 (96%)	12 (3%)	1 (0%)	50	92
1	L	373/375 (100%)	360 (96%)	12 (3%)	1 (0%)	50	92
1	M	373/375 (100%)	360 (96%)	12 (3%)	1 (0%)	50	92
1	N	373/375 (100%)	361 (97%)	11 (3%)	1 (0%)	50	92
1	O	373/375 (100%)	360 (96%)	12 (3%)	1 (0%)	50	92
1	P	373/375 (100%)	360 (96%)	12 (3%)	1 (0%)	50	92
1	Q	373/375 (100%)	361 (97%)	11 (3%)	1 (0%)	50	92
1	W	373/375 (100%)	361 (97%)	11 (3%)	1 (0%)	50	92
1	Y	373/375 (100%)	360 (96%)	12 (3%)	1 (0%)	50	92
1	Z	373/375 (100%)	361 (97%)	11 (3%)	1 (0%)	50	92
All	All	7460/7500 (100%)	7211 (97%)	229 (3%)	20 (0%)	50	92

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	308	LYS
1	B	308	LYS
1	C	308	LYS
1	D	308	LYS
1	E	308	LYS
1	F	308	LYS
1	G	308	LYS
1	H	308	LYS
1	I	308	LYS
1	J	308	LYS
1	K	308	LYS
1	L	308	LYS
1	M	308	LYS

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Mol	Chain	Res	Type
1	N	308	LYS
1	O	308	LYS
1	P	308	LYS
1	Q	308	LYS
1	W	308	LYS
1	Y	308	LYS
1	Z	308	LYS

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	313/313 (100%)	269 (86%)	44 (14%)	5	30
1	B	313/313 (100%)	268 (86%)	45 (14%)	5	28
1	C	313/313 (100%)	269 (86%)	44 (14%)	5	30
1	D	313/313 (100%)	268 (86%)	45 (14%)	5	28
1	E	313/313 (100%)	268 (86%)	45 (14%)	5	28
1	F	313/313 (100%)	268 (86%)	45 (14%)	5	28
1	G	313/313 (100%)	269 (86%)	44 (14%)	5	30
1	H	313/313 (100%)	268 (86%)	45 (14%)	5	28
1	I	313/313 (100%)	268 (86%)	45 (14%)	5	28
1	J	313/313 (100%)	269 (86%)	44 (14%)	5	30
1	K	313/313 (100%)	269 (86%)	44 (14%)	5	30
1	L	313/313 (100%)	268 (86%)	45 (14%)	5	28
1	M	313/313 (100%)	268 (86%)	45 (14%)	5	28
1	N	313/313 (100%)	268 (86%)	45 (14%)	5	28
1	O	313/313 (100%)	269 (86%)	44 (14%)	5	30
1	P	313/313 (100%)	268 (86%)	45 (14%)	5	28
1	Q	313/313 (100%)	269 (86%)	44 (14%)	5	30
1	W	313/313 (100%)	268 (86%)	45 (14%)	5	28
1	Y	313/313 (100%)	268 (86%)	45 (14%)	5	28

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	Z	313/313 (100%)	268 (86%)	45 (14%)	5	28
All	All	6260/6260 (100%)	5367 (86%)	893 (14%)	5	29

All (893) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	5	LYS
1	A	6	VAL
1	A	11	THR
1	A	12	LEU
1	A	22	LYS
1	A	31	ASP
1	A	33	ILE
1	A	34	ASP
1	A	47	LEU
1	A	79	THR
1	A	82	ILE
1	A	83	LEU
1	A	88	TYR
1	A	90	VAL
1	A	103	ASP
1	A	107	LYS
1	A	120	THR
1	A	123	ILE
1	A	136	LYS
1	A	145	VAL
1	A	163	ILE
1	A	168	ILE
1	A	171	LEU
1	A	196	ARG
1	A	220	ILE
1	A	255	GLN
1	A	270	ILE
1	A	272	LEU
1	A	283	GLN
1	A	287	VAL
1	A	293	LYS
1	A	304	LEU
1	A	325	LEU
1	A	331	LEU
1	A	333	ILE

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Mol	Chain	Res	Type
1	A	359	GLU
1	A	360	ASN
1	A	368	LEU
1	A	370	LEU
1	A	371	THR
1	A	373	GLU
1	A	374	GLU
1	A	375	LEU
1	B	3	LEU
1	B	5	LYS
1	B	6	VAL
1	B	11	THR
1	B	12	LEU
1	B	22	LYS
1	B	31	ASP
1	B	33	ILE
1	B	34	ASP
1	B	47	LEU
1	B	79	THR
1	B	82	ILE
1	B	83	LEU
1	B	88	TYR
1	B	90	VAL
1	B	103	ASP
1	B	107	LYS
1	B	120	THR
1	B	123	ILE
1	B	136	LYS
1	B	145	VAL
1	B	163	ILE
1	B	168	ILE
1	B	171	LEU
1	B	196	ARG
1	B	220	ILE
1	B	255	GLN
1	B	270	ILE
1	B	272	LEU
1	B	283	GLN
1	B	287	VAL
1	B	293	LYS
1	B	304	LEU
1	B	305	ASN

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Mol	Chain	Res	Type
1	B	325	LEU
1	B	331	LEU
1	B	333	ILE
1	B	359	GLU
1	B	360	ASN
1	B	368	LEU
1	B	370	LEU
1	B	371	THR
1	B	373	GLU
1	B	374	GLU
1	B	375	LEU
1	C	3	LEU
1	C	5	LYS
1	C	6	VAL
1	C	11	THR
1	C	12	LEU
1	C	22	LYS
1	C	31	ASP
1	C	33	ILE
1	C	34	ASP
1	C	47	LEU
1	C	79	THR
1	C	82	ILE
1	C	83	LEU
1	C	88	TYR
1	C	90	VAL
1	C	103	ASP
1	C	107	LYS
1	C	120	THR
1	C	123	ILE
1	C	136	LYS
1	C	145	VAL
1	C	163	ILE
1	C	168	ILE
1	C	171	LEU
1	C	196	ARG
1	C	220	ILE
1	C	255	GLN
1	C	270	ILE
1	C	272	LEU
1	C	283	GLN
1	C	287	VAL

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Mol	Chain	Res	Type
1	C	293	LYS
1	C	304	LEU
1	C	325	LEU
1	C	331	LEU
1	C	333	ILE
1	C	359	GLU
1	C	360	ASN
1	C	368	LEU
1	C	370	LEU
1	C	371	THR
1	C	373	GLU
1	C	374	GLU
1	C	375	LEU
1	D	3	LEU
1	D	5	LYS
1	D	6	VAL
1	D	11	THR
1	D	12	LEU
1	D	22	LYS
1	D	31	ASP
1	D	33	ILE
1	D	34	ASP
1	D	47	LEU
1	D	79	THR
1	D	82	ILE
1	D	83	LEU
1	D	88	TYR
1	D	90	VAL
1	D	103	ASP
1	D	107	LYS
1	D	120	THR
1	D	123	ILE
1	D	136	LYS
1	D	145	VAL
1	D	163	ILE
1	D	168	ILE
1	D	171	LEU
1	D	196	ARG
1	D	220	ILE
1	D	255	GLN
1	D	270	ILE
1	D	272	LEU

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Mol	Chain	Res	Type
1	D	283	GLN
1	D	287	VAL
1	D	293	LYS
1	D	304	LEU
1	D	305	ASN
1	D	325	LEU
1	D	331	LEU
1	D	333	ILE
1	D	359	GLU
1	D	360	ASN
1	D	368	LEU
1	D	370	LEU
1	D	371	THR
1	D	373	GLU
1	D	374	GLU
1	D	375	LEU
1	E	3	LEU
1	E	5	LYS
1	E	6	VAL
1	E	11	THR
1	E	12	LEU
1	E	22	LYS
1	E	31	ASP
1	E	33	ILE
1	E	34	ASP
1	E	47	LEU
1	E	79	THR
1	E	82	ILE
1	E	83	LEU
1	E	88	TYR
1	E	90	VAL
1	E	103	ASP
1	E	107	LYS
1	E	120	THR
1	E	123	ILE
1	E	136	LYS
1	E	145	VAL
1	E	163	ILE
1	E	168	ILE
1	E	171	LEU
1	E	196	ARG
1	E	220	ILE

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Mol	Chain	Res	Type
1	E	255	GLN
1	E	270	ILE
1	E	272	LEU
1	E	283	GLN
1	E	287	VAL
1	E	293	LYS
1	E	304	LEU
1	E	305	ASN
1	E	325	LEU
1	E	331	LEU
1	E	333	ILE
1	E	359	GLU
1	E	360	ASN
1	E	368	LEU
1	E	370	LEU
1	E	371	THR
1	E	373	GLU
1	E	374	GLU
1	E	375	LEU
1	F	3	LEU
1	F	5	LYS
1	F	6	VAL
1	F	11	THR
1	F	12	LEU
1	F	22	LYS
1	F	31	ASP
1	F	33	ILE
1	F	34	ASP
1	F	47	LEU
1	F	79	THR
1	F	82	ILE
1	F	83	LEU
1	F	88	TYR
1	F	90	VAL
1	F	103	ASP
1	F	107	LYS
1	F	120	THR
1	F	123	ILE
1	F	136	LYS
1	F	145	VAL
1	F	163	ILE
1	F	168	ILE

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Mol	Chain	Res	Type
1	F	171	LEU
1	F	196	ARG
1	F	220	ILE
1	F	255	GLN
1	F	270	ILE
1	F	272	LEU
1	F	283	GLN
1	F	287	VAL
1	F	293	LYS
1	F	304	LEU
1	F	305	ASN
1	F	325	LEU
1	F	331	LEU
1	F	333	ILE
1	F	359	GLU
1	F	360	ASN
1	F	368	LEU
1	F	370	LEU
1	F	371	THR
1	F	373	GLU
1	F	374	GLU
1	F	375	LEU
1	G	3	LEU
1	G	5	LYS
1	G	6	VAL
1	G	11	THR
1	G	12	LEU
1	G	22	LYS
1	G	31	ASP
1	G	33	ILE
1	G	34	ASP
1	G	47	LEU
1	G	79	THR
1	G	82	ILE
1	G	83	LEU
1	G	88	TYR
1	G	90	VAL
1	G	103	ASP
1	G	107	LYS
1	G	120	THR
1	G	123	ILE
1	G	136	LYS

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Mol	Chain	Res	Type
1	G	145	VAL
1	G	163	ILE
1	G	168	ILE
1	G	171	LEU
1	G	196	ARG
1	G	220	ILE
1	G	255	GLN
1	G	270	ILE
1	G	272	LEU
1	G	283	GLN
1	G	287	VAL
1	G	293	LYS
1	G	304	LEU
1	G	325	LEU
1	G	331	LEU
1	G	333	ILE
1	G	359	GLU
1	G	360	ASN
1	G	368	LEU
1	G	370	LEU
1	G	371	THR
1	G	373	GLU
1	G	374	GLU
1	G	375	LEU
1	H	3	LEU
1	H	5	LYS
1	H	6	VAL
1	H	11	THR
1	H	12	LEU
1	H	22	LYS
1	H	31	ASP
1	H	33	ILE
1	H	34	ASP
1	H	47	LEU
1	H	79	THR
1	H	82	ILE
1	H	83	LEU
1	H	88	TYR
1	H	90	VAL
1	H	103	ASP
1	H	107	LYS
1	H	120	THR

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Mol	Chain	Res	Type
1	H	123	ILE
1	H	136	LYS
1	H	145	VAL
1	H	163	ILE
1	H	168	ILE
1	H	171	LEU
1	H	196	ARG
1	H	220	ILE
1	H	255	GLN
1	H	270	ILE
1	H	272	LEU
1	H	283	GLN
1	H	287	VAL
1	H	293	LYS
1	H	304	LEU
1	H	305	ASN
1	H	325	LEU
1	H	331	LEU
1	H	333	ILE
1	H	359	GLU
1	H	360	ASN
1	H	368	LEU
1	H	370	LEU
1	H	371	THR
1	H	373	GLU
1	H	374	GLU
1	H	375	LEU
1	I	3	LEU
1	I	5	LYS
1	I	6	VAL
1	I	11	THR
1	I	12	LEU
1	I	22	LYS
1	I	31	ASP
1	I	33	ILE
1	I	34	ASP
1	I	47	LEU
1	I	79	THR
1	I	82	ILE
1	I	83	LEU
1	I	88	TYR
1	I	90	VAL

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Mol	Chain	Res	Type
1	I	103	ASP
1	I	107	LYS
1	I	120	THR
1	I	123	ILE
1	I	136	LYS
1	I	145	VAL
1	I	163	ILE
1	I	168	ILE
1	I	171	LEU
1	I	196	ARG
1	I	220	ILE
1	I	255	GLN
1	I	270	ILE
1	I	272	LEU
1	I	283	GLN
1	I	287	VAL
1	I	293	LYS
1	I	304	LEU
1	I	305	ASN
1	I	325	LEU
1	I	331	LEU
1	I	333	ILE
1	I	359	GLU
1	I	360	ASN
1	I	368	LEU
1	I	370	LEU
1	I	371	THR
1	I	373	GLU
1	I	374	GLU
1	I	375	LEU
1	J	3	LEU
1	J	5	LYS
1	J	6	VAL
1	J	11	THR
1	J	12	LEU
1	J	22	LYS
1	J	31	ASP
1	J	33	ILE
1	J	34	ASP
1	J	47	LEU
1	J	79	THR
1	J	82	ILE

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Mol	Chain	Res	Type
1	J	83	LEU
1	J	88	TYR
1	J	90	VAL
1	J	103	ASP
1	J	107	LYS
1	J	120	THR
1	J	123	ILE
1	J	136	LYS
1	J	145	VAL
1	J	163	ILE
1	J	168	ILE
1	J	171	LEU
1	J	196	ARG
1	J	220	ILE
1	J	255	GLN
1	J	270	ILE
1	J	272	LEU
1	J	283	GLN
1	J	287	VAL
1	J	293	LYS
1	J	304	LEU
1	J	325	LEU
1	J	331	LEU
1	J	333	ILE
1	J	359	GLU
1	J	360	ASN
1	J	368	LEU
1	J	370	LEU
1	J	371	THR
1	J	373	GLU
1	J	374	GLU
1	J	375	LEU
1	K	3	LEU
1	K	5	LYS
1	K	6	VAL
1	K	11	THR
1	K	12	LEU
1	K	22	LYS
1	K	31	ASP
1	K	33	ILE
1	K	34	ASP
1	K	47	LEU

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Mol	Chain	Res	Type
1	K	79	THR
1	K	82	ILE
1	K	83	LEU
1	K	88	TYR
1	K	90	VAL
1	K	103	ASP
1	K	107	LYS
1	K	120	THR
1	K	123	ILE
1	K	136	LYS
1	K	145	VAL
1	K	163	ILE
1	K	168	ILE
1	K	171	LEU
1	K	196	ARG
1	K	220	ILE
1	K	255	GLN
1	K	270	ILE
1	K	272	LEU
1	K	283	GLN
1	K	287	VAL
1	K	293	LYS
1	K	304	LEU
1	K	325	LEU
1	K	331	LEU
1	K	333	ILE
1	K	359	GLU
1	K	360	ASN
1	K	368	LEU
1	K	370	LEU
1	K	371	THR
1	K	373	GLU
1	K	374	GLU
1	K	375	LEU
1	L	3	LEU
1	L	5	LYS
1	L	6	VAL
1	L	11	THR
1	L	12	LEU
1	L	22	LYS
1	L	31	ASP
1	L	33	ILE

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Mol	Chain	Res	Type
1	L	34	ASP
1	L	47	LEU
1	L	79	THR
1	L	82	ILE
1	L	83	LEU
1	L	88	TYR
1	L	90	VAL
1	L	103	ASP
1	L	107	LYS
1	L	120	THR
1	L	123	ILE
1	L	136	LYS
1	L	145	VAL
1	L	163	ILE
1	L	168	ILE
1	L	171	LEU
1	L	196	ARG
1	L	220	ILE
1	L	255	GLN
1	L	270	ILE
1	L	272	LEU
1	L	283	GLN
1	L	287	VAL
1	L	293	LYS
1	L	304	LEU
1	L	305	ASN
1	L	325	LEU
1	L	331	LEU
1	L	333	ILE
1	L	359	GLU
1	L	360	ASN
1	L	368	LEU
1	L	370	LEU
1	L	371	THR
1	L	373	GLU
1	L	374	GLU
1	L	375	LEU
1	M	3	LEU
1	M	5	LYS
1	M	6	VAL
1	M	11	THR
1	M	12	LEU

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Mol	Chain	Res	Type
1	M	22	LYS
1	M	31	ASP
1	M	33	ILE
1	M	34	ASP
1	M	47	LEU
1	M	79	THR
1	M	82	ILE
1	M	83	LEU
1	M	88	TYR
1	M	90	VAL
1	M	103	ASP
1	M	107	LYS
1	M	120	THR
1	M	123	ILE
1	M	136	LYS
1	M	145	VAL
1	M	163	ILE
1	M	168	ILE
1	M	171	LEU
1	M	196	ARG
1	M	220	ILE
1	M	255	GLN
1	M	270	ILE
1	M	272	LEU
1	M	283	GLN
1	M	287	VAL
1	M	293	LYS
1	M	304	LEU
1	M	305	ASN
1	M	325	LEU
1	M	331	LEU
1	M	333	ILE
1	M	359	GLU
1	M	360	ASN
1	M	368	LEU
1	M	370	LEU
1	M	371	THR
1	M	373	GLU
1	M	374	GLU
1	M	375	LEU
1	N	3	LEU
1	N	5	LYS

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Mol	Chain	Res	Type
1	N	6	VAL
1	N	11	THR
1	N	12	LEU
1	N	22	LYS
1	N	31	ASP
1	N	33	ILE
1	N	34	ASP
1	N	47	LEU
1	N	79	THR
1	N	82	ILE
1	N	83	LEU
1	N	88	TYR
1	N	90	VAL
1	N	103	ASP
1	N	107	LYS
1	N	120	THR
1	N	123	ILE
1	N	136	LYS
1	N	145	VAL
1	N	163	ILE
1	N	168	ILE
1	N	171	LEU
1	N	196	ARG
1	N	220	ILE
1	N	255	GLN
1	N	270	ILE
1	N	272	LEU
1	N	283	GLN
1	N	287	VAL
1	N	293	LYS
1	N	304	LEU
1	N	305	ASN
1	N	325	LEU
1	N	331	LEU
1	N	333	ILE
1	N	359	GLU
1	N	360	ASN
1	N	368	LEU
1	N	370	LEU
1	N	371	THR
1	N	373	GLU
1	N	374	GLU

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Mol	Chain	Res	Type
1	N	375	LEU
1	O	3	LEU
1	O	5	LYS
1	O	6	VAL
1	O	11	THR
1	O	12	LEU
1	O	22	LYS
1	O	31	ASP
1	O	33	ILE
1	O	34	ASP
1	O	47	LEU
1	O	79	THR
1	O	82	ILE
1	O	83	LEU
1	O	88	TYR
1	O	90	VAL
1	O	103	ASP
1	O	107	LYS
1	O	120	THR
1	O	123	ILE
1	O	136	LYS
1	O	145	VAL
1	O	163	ILE
1	O	168	ILE
1	O	171	LEU
1	O	196	ARG
1	O	220	ILE
1	O	255	GLN
1	O	270	ILE
1	O	272	LEU
1	O	283	GLN
1	O	287	VAL
1	O	293	LYS
1	O	304	LEU
1	O	325	LEU
1	O	331	LEU
1	O	333	ILE
1	O	359	GLU
1	O	360	ASN
1	O	368	LEU
1	O	370	LEU
1	O	371	THR

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Mol	Chain	Res	Type
1	O	373	GLU
1	O	374	GLU
1	O	375	LEU
1	P	3	LEU
1	P	5	LYS
1	P	6	VAL
1	P	11	THR
1	P	12	LEU
1	P	22	LYS
1	P	31	ASP
1	P	33	ILE
1	P	34	ASP
1	P	47	LEU
1	P	79	THR
1	P	82	ILE
1	P	83	LEU
1	P	88	TYR
1	P	90	VAL
1	P	103	ASP
1	P	107	LYS
1	P	120	THR
1	P	123	ILE
1	P	136	LYS
1	P	145	VAL
1	P	163	ILE
1	P	168	ILE
1	P	171	LEU
1	P	196	ARG
1	P	220	ILE
1	P	255	GLN
1	P	270	ILE
1	P	272	LEU
1	P	283	GLN
1	P	287	VAL
1	P	293	LYS
1	P	304	LEU
1	P	305	ASN
1	P	325	LEU
1	P	331	LEU
1	P	333	ILE
1	P	359	GLU
1	P	360	ASN

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Mol	Chain	Res	Type
1	P	368	LEU
1	P	370	LEU
1	P	371	THR
1	P	373	GLU
1	P	374	GLU
1	P	375	LEU
1	Q	3	LEU
1	Q	5	LYS
1	Q	6	VAL
1	Q	11	THR
1	Q	12	LEU
1	Q	22	LYS
1	Q	31	ASP
1	Q	33	ILE
1	Q	34	ASP
1	Q	47	LEU
1	Q	79	THR
1	Q	82	ILE
1	Q	83	LEU
1	Q	88	TYR
1	Q	90	VAL
1	Q	103	ASP
1	Q	107	LYS
1	Q	120	THR
1	Q	123	ILE
1	Q	136	LYS
1	Q	145	VAL
1	Q	163	ILE
1	Q	168	ILE
1	Q	171	LEU
1	Q	196	ARG
1	Q	220	ILE
1	Q	255	GLN
1	Q	270	ILE
1	Q	272	LEU
1	Q	283	GLN
1	Q	287	VAL
1	Q	293	LYS
1	Q	304	LEU
1	Q	325	LEU
1	Q	331	LEU
1	Q	333	ILE

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Mol	Chain	Res	Type
1	Q	359	GLU
1	Q	360	ASN
1	Q	368	LEU
1	Q	370	LEU
1	Q	371	THR
1	Q	373	GLU
1	Q	374	GLU
1	Q	375	LEU
1	W	3	LEU
1	W	5	LYS
1	W	6	VAL
1	W	11	THR
1	W	12	LEU
1	W	22	LYS
1	W	31	ASP
1	W	33	ILE
1	W	34	ASP
1	W	47	LEU
1	W	79	THR
1	W	82	ILE
1	W	83	LEU
1	W	88	TYR
1	W	90	VAL
1	W	103	ASP
1	W	107	LYS
1	W	120	THR
1	W	123	ILE
1	W	136	LYS
1	W	145	VAL
1	W	163	ILE
1	W	168	ILE
1	W	171	LEU
1	W	196	ARG
1	W	220	ILE
1	W	255	GLN
1	W	270	ILE
1	W	272	LEU
1	W	283	GLN
1	W	287	VAL
1	W	293	LYS
1	W	304	LEU
1	W	305	ASN

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Mol	Chain	Res	Type
1	W	325	LEU
1	W	331	LEU
1	W	333	ILE
1	W	359	GLU
1	W	360	ASN
1	W	368	LEU
1	W	370	LEU
1	W	371	THR
1	W	373	GLU
1	W	374	GLU
1	W	375	LEU
1	Y	3	LEU
1	Y	5	LYS
1	Y	6	VAL
1	Y	11	THR
1	Y	12	LEU
1	Y	22	LYS
1	Y	31	ASP
1	Y	33	ILE
1	Y	34	ASP
1	Y	47	LEU
1	Y	79	THR
1	Y	82	ILE
1	Y	83	LEU
1	Y	88	TYR
1	Y	90	VAL
1	Y	103	ASP
1	Y	107	LYS
1	Y	120	THR
1	Y	123	ILE
1	Y	136	LYS
1	Y	145	VAL
1	Y	163	ILE
1	Y	168	ILE
1	Y	171	LEU
1	Y	196	ARG
1	Y	220	ILE
1	Y	255	GLN
1	Y	270	ILE
1	Y	272	LEU
1	Y	283	GLN
1	Y	287	VAL

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Mol	Chain	Res	Type
1	Y	293	LYS
1	Y	304	LEU
1	Y	305	ASN
1	Y	325	LEU
1	Y	331	LEU
1	Y	333	ILE
1	Y	359	GLU
1	Y	360	ASN
1	Y	368	LEU
1	Y	370	LEU
1	Y	371	THR
1	Y	373	GLU
1	Y	374	GLU
1	Y	375	LEU
1	Z	3	LEU
1	Z	5	LYS
1	Z	6	VAL
1	Z	11	THR
1	Z	12	LEU
1	Z	22	LYS
1	Z	31	ASP
1	Z	33	ILE
1	Z	34	ASP
1	Z	47	LEU
1	Z	79	THR
1	Z	82	ILE
1	Z	83	LEU
1	Z	88	TYR
1	Z	90	VAL
1	Z	103	ASP
1	Z	107	LYS
1	Z	120	THR
1	Z	123	ILE
1	Z	136	LYS
1	Z	145	VAL
1	Z	163	ILE
1	Z	168	ILE
1	Z	171	LEU
1	Z	196	ARG
1	Z	220	ILE
1	Z	255	GLN
1	Z	270	ILE

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Mol	Chain	Res	Type
1	Z	272	LEU
1	Z	283	GLN
1	Z	287	VAL
1	Z	293	LYS
1	Z	304	LEU
1	Z	305	ASN
1	Z	325	LEU
1	Z	331	LEU
1	Z	333	ILE
1	Z	359	GLU
1	Z	360	ASN
1	Z	368	LEU
1	Z	370	LEU
1	Z	371	THR
1	Z	373	GLU
1	Z	374	GLU
1	Z	375	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (72) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	151	HIS
1	A	187	ASN
1	A	216	HIS
1	B	41	GLN
1	B	151	HIS
1	B	187	ASN
1	B	216	HIS
1	C	151	HIS
1	C	187	ASN
1	C	216	HIS
1	D	41	GLN
1	D	151	HIS
1	D	187	ASN
1	D	216	HIS
1	E	41	GLN
1	E	151	HIS
1	E	187	ASN
1	E	216	HIS
1	E	292	GLN
1	F	151	HIS
1	F	187	ASN

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Mol	Chain	Res	Type
1	F	216	HIS
1	F	292	GLN
1	G	41	GLN
1	G	151	HIS
1	G	187	ASN
1	G	216	HIS
1	H	151	HIS
1	H	187	ASN
1	H	216	HIS
1	I	41	GLN
1	I	151	HIS
1	I	187	ASN
1	I	216	HIS
1	J	151	HIS
1	J	187	ASN
1	J	216	HIS
1	K	41	GLN
1	K	151	HIS
1	K	187	ASN
1	K	216	HIS
1	L	151	HIS
1	L	187	ASN
1	L	216	HIS
1	M	151	HIS
1	M	187	ASN
1	M	216	HIS
1	M	292	GLN
1	N	41	GLN
1	N	151	HIS
1	N	187	ASN
1	N	216	HIS
1	O	151	HIS
1	O	187	ASN
1	O	216	HIS
1	P	41	GLN
1	P	151	HIS
1	P	187	ASN
1	P	216	HIS
1	Q	151	HIS
1	Q	187	ASN
1	Q	216	HIS
1	W	151	HIS

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Mol	Chain	Res	Type
1	W	187	ASN
1	W	216	HIS
1	Y	151	HIS
1	Y	187	ASN
1	Y	216	HIS
1	Z	41	GLN
1	Z	151	HIS
1	Z	187	ASN
1	Z	216	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	R	69/70 (98%)	19 (27%)	7 (10%)
2	X	69/70 (98%)	19 (27%)	4 (5%)
All	All	138/140 (98%)	38 (27%)	11 (7%)

All (38) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	R	8	C
2	R	15	C
2	R	16	C
2	R	17	C
2	R	22	C
2	R	24	C
2	R	29	C
2	R	31	C
2	R	36	C
2	R	40	C
2	R	43	C
2	R	45	C
2	R	47	C
2	R	50	C
2	R	52	C
2	R	57	C
2	R	64	C
2	R	67	C
2	R	70	C
2	X	3	C
2	X	8	C

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Mol	Chain	Res	Type
2	X	15	C
2	X	17	C
2	X	22	C
2	X	24	C
2	X	29	C
2	X	31	C
2	X	36	C
2	X	40	C
2	X	42	C
2	X	43	C
2	X	45	C
2	X	47	C
2	X	50	C
2	X	52	C
2	X	57	C
2	X	58	C
2	X	64	C

All (11) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	R	4	C
2	R	11	C
2	R	23	C
2	R	25	C
2	R	32	C
2	R	46	C
2	R	67	C
2	X	18	C
2	X	25	C
2	X	32	C
2	X	39	C

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	A	375/375 (100%)	0.31	15 (4%)	36	20	77, 135, 189, 218	0
1	B	375/375 (100%)	0.28	9 (2%)	56	32	77, 135, 189, 218	0
1	C	375/375 (100%)	0.31	10 (2%)	52	29	77, 135, 189, 218	0
1	D	375/375 (100%)	0.25	9 (2%)	56	32	77, 135, 189, 218	0
1	E	375/375 (100%)	0.29	14 (3%)	39	22	77, 135, 189, 218	0
1	F	375/375 (100%)	0.18	4 (1%)	77	50	77, 135, 189, 218	0
1	G	375/375 (100%)	0.26	11 (2%)	49	28	77, 135, 189, 218	0
1	H	375/375 (100%)	0.27	11 (2%)	49	28	77, 135, 189, 218	0
1	I	375/375 (100%)	0.27	10 (2%)	52	29	77, 135, 189, 218	0
1	J	375/375 (100%)	0.33	13 (3%)	42	23	77, 135, 189, 218	0
1	K	375/375 (100%)	0.23	8 (2%)	60	35	77, 135, 189, 218	0
1	L	375/375 (100%)	0.25	8 (2%)	60	35	77, 135, 189, 218	0
1	M	375/375 (100%)	0.23	8 (2%)	60	35	77, 135, 189, 218	0
1	N	375/375 (100%)	0.23	6 (1%)	68	41	77, 135, 189, 218	0
1	O	375/375 (100%)	0.27	11 (2%)	49	28	77, 135, 189, 218	0
1	P	375/375 (100%)	0.28	10 (2%)	52	29	77, 135, 189, 218	0
1	Q	375/375 (100%)	0.19	4 (1%)	77	50	77, 135, 189, 218	0
1	W	375/375 (100%)	0.24	9 (2%)	56	32	77, 135, 189, 218	0
1	Y	375/375 (100%)	0.24	7 (1%)	64	37	77, 135, 189, 218	0
1	Z	375/375 (100%)	0.31	11 (2%)	49	28	77, 135, 189, 218	0
2	R	70/70 (100%)	0.09	0	100	100	129, 145, 158, 162	0
2	X	70/70 (100%)	0.09	0	100	100	119, 145, 156, 159	0
All	All	7640/7640 (100%)	0.26	188 (2%)	54	31	77, 136, 189, 218	0

All (188) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	368	LEU	6.7
1	P	368	LEU	6.3
1	W	102	GLN	6.2
1	J	362	VAL	5.7
1	E	272	LEU	5.2
1	A	31	ASP	4.8
1	M	272	LEU	4.8
1	I	104	ILE	4.8
1	N	368	LEU	4.7
1	E	240	GLU	4.6
1	Z	272	LEU	4.6
1	M	100	HIS	4.6
1	C	365	TYR	4.5
1	Z	2	ALA	4.4
1	Z	367	VAL	4.4
1	A	272	LEU	4.3
1	G	367	VAL	4.3
1	I	364	ASN	4.2
1	W	369	ASP	4.2
1	O	110	LYS	4.1
1	I	361	GLY	4.1
1	G	8	LEU	4.1
1	O	371	THR	4.0
1	D	374	GLU	4.0
1	L	105	ASN	3.9
1	H	362	VAL	3.9
1	J	272	LEU	3.8
1	L	272	LEU	3.8
1	G	104	ILE	3.8
1	K	367	VAL	3.7
1	E	5	LYS	3.6
1	G	119	LEU	3.6
1	M	101	ARG	3.6
1	A	370	LEU	3.5
1	L	144	GLU	3.4
1	B	119	LEU	3.4
1	P	167	VAL	3.4
1	Z	366	SER	3.4
1	A	320	PHE	3.4
1	Y	257	MET	3.3
1	J	56	ASP	3.2
1	E	370	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
1	W	272	LEU	3.1
1	I	367	VAL	3.1
1	M	62	THR	3.1
1	A	114	LEU	3.1
1	K	124	GLN	3.1
1	I	124	GLN	3.1
1	A	371	THR	3.1
1	B	312	LEU	3.1
1	G	272	LEU	3.1
1	C	370	LEU	3.0
1	B	104	ILE	3.0
1	C	371	THR	3.0
1	J	368	LEU	3.0
1	C	124	GLN	3.0
1	E	200	LEU	3.0
1	J	25	ILE	3.0
1	A	136	LYS	2.9
1	D	375	LEU	2.9
1	O	103	ASP	2.9
1	Q	5	LYS	2.9
1	C	272	LEU	2.9
1	H	248	MET	2.9
1	G	368	LEU	2.9
1	F	370	LEU	2.9
1	W	200	LEU	2.9
1	D	297	GLU	2.8
1	I	362	VAL	2.8
1	B	353	TYR	2.8
1	K	1	MET	2.8
1	P	3	LEU	2.8
1	I	334	MET	2.8
1	K	320	PHE	2.8
1	P	100	HIS	2.8
1	C	367	VAL	2.8
1	E	219	PHE	2.8
1	G	3	LEU	2.7
1	K	104	ILE	2.7
1	D	296	GLY	2.7
1	M	112	GLU	2.7
1	Z	74	LEU	2.7
1	P	124	GLN	2.7
1	J	104	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	104	ILE	2.6
1	H	104	ILE	2.6
1	M	91	LYS	2.6
1	H	138	MET	2.6
1	Y	272	LEU	2.6
1	J	101	ARG	2.6
1	H	100	HIS	2.6
1	O	8	LEU	2.6
1	W	362	VAL	2.6
1	E	8	LEU	2.6
1	J	205	ILE	2.5
1	M	320	PHE	2.5
1	E	137	LYS	2.5
1	D	272	LEU	2.5
1	Y	145	VAL	2.5
1	J	206	ALA	2.5
1	E	7	LYS	2.5
1	F	119	LEU	2.5
1	A	33	ILE	2.5
1	L	312	LEU	2.4
1	O	7	LYS	2.4
1	N	65	ILE	2.4
1	L	278	GLN	2.4
1	H	123	ILE	2.4
1	A	202	PRO	2.3
1	J	282	GLU	2.3
1	A	271	MET	2.3
1	K	364	ASN	2.3
1	F	3	LEU	2.3
1	H	105	ASN	2.3
1	O	271	MET	2.3
1	C	53	ILE	2.3
1	W	370	LEU	2.3
1	I	100	HIS	2.3
1	E	308	LYS	2.3
1	E	2	ALA	2.3
1	Z	111	PHE	2.3
1	C	368	LEU	2.3
1	A	112	GLU	2.3
1	C	325	LEU	2.3
1	B	251	TYR	2.2
1	O	124	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	P	370	LEU	2.2
1	B	374	GLU	2.2
1	Z	334	MET	2.2
1	Z	200	LEU	2.2
1	Z	119	LEU	2.2
1	P	362	VAL	2.2
1	Y	202	PRO	2.2
1	D	46	LYS	2.2
1	N	308	LYS	2.2
1	H	361	GLY	2.2
1	W	274	HIS	2.2
1	H	282	GLU	2.2
1	G	110	LYS	2.2
1	Y	114	LEU	2.2
1	J	24	THR	2.2
1	L	56	ASP	2.2
1	J	361	GLY	2.2
1	Z	353	TYR	2.2
1	N	104	ILE	2.2
1	O	223	PHE	2.2
1	D	320	PHE	2.2
1	K	368	LEU	2.2
1	P	246	LEU	2.2
1	B	272	LEU	2.2
1	N	365	TYR	2.2
1	J	365	TYR	2.1
1	A	179	LEU	2.1
1	E	288	TYR	2.1
1	A	2	ALA	2.1
1	B	273	GLY	2.1
1	B	200	LEU	2.1
1	A	364	ASN	2.1
1	L	100	HIS	2.1
1	L	369	ASP	2.1
1	K	71	MET	2.1
1	Q	365	TYR	2.1
1	E	56	ASP	2.1
1	F	368	LEU	2.1
1	W	368	LEU	2.1
1	G	308	LYS	2.1
1	Y	320	PHE	2.1
1	P	374	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	M	97	VAL	2.1
1	O	163	ILE	2.1
1	O	3	LEU	2.1
1	C	56	ASP	2.1
1	P	222	VAL	2.1
1	N	114	LEU	2.1
1	H	8	LEU	2.0
1	Y	26	GLN	2.0
1	I	144	GLU	2.0
1	E	368	LEU	2.0
1	Q	160	ILE	2.0
1	D	90	VAL	2.0
1	W	119	LEU	2.0
1	H	110	LYS	2.0
1	I	33	ILE	2.0
1	O	370	LEU	2.0
1	Q	368	LEU	2.0
1	G	97	VAL	2.0
1	G	189	VAL	2.0
1	Z	114	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.