



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:09 PM UTC

PDB ID : 4ZUX / pdb_00004zux
Title : SAGA DUB module Ubp8/Sgf11/Sus1/Sgf73 bound to ubiquitinated nucleosome
Authors : Morgan, M.; Wolberger, C.
Deposited on : 2015-05-17
Resolution : 3.82 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

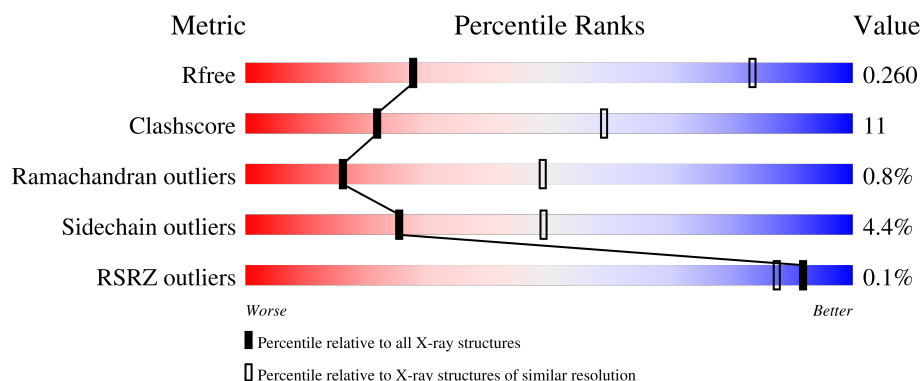
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.82 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1043 (3.98-3.66)
Clashscore	190562	1075 (3.98-3.66)
Ramachandran outliers	187476	1029 (3.98-3.66)
Sidechain outliers	187428	1024 (3.98-3.66)
RSRZ outliers	180081	1042 (3.98-3.66)

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

Mol	Chain	Length	Quality of chain
1	A	136	44% 25% • 29%
1	E	136	46% 26% • 29%
1	K	136	50% 21% • 29%
1	O	136	46% 25% • 29%
2	B	103	50% 24% 6% 19%

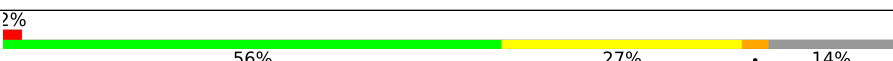
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Mol	Chain	Length	Quality of chain
2	F	103	
2	L	103	
2	P	103	
3	C	130	
3	G	130	
3	M	130	
3	Q	130	
4	D	123	
4	H	123	
4	N	123	
4	R	123	
5	I	145	
5	S	145	
6	J	145	
6	T	145	
7	U	472	
7	Z	472	
7	e	472	
7	j	472	
8	V	96	
8	a	96	
8	f	96	
8	k	96	
9	W	99	
9	b	99	

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Mol	Chain	Length	Quality of chain
9	g	99	
9	l	99	
10	X	76	
10	c	76	
10	h	76	
10	m	76	
11	Y	104	
11	d	104	
11	i	104	
11	n	104	

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 49079 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called Histone H3.2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	97	Total	C	N	O	S	0	0	0
			802	506	155	138	3			
1	E	97	Total	C	N	O	S	0	0	0
			801	504	155	139	3			
1	K	97	Total	C	N	O	S	0	0	0
			802	506	155	138	3			
1	O	97	Total	C	N	O	S	0	0	0
			801	504	155	139	3			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	102	ALA	GLY	conflict	UNP P84233
E	102	ALA	GLY	conflict	UNP P84233
K	102	ALA	GLY	conflict	UNP P84233
O	102	ALA	GLY	conflict	UNP P84233

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	83	Total	C	N	O	S	0	0	0
			662	418	129	114	1			
2	F	78	Total	C	N	O	S	0	0	0
			619	391	120	107	1			
2	L	83	Total	C	N	O	S	0	0	0
			662	418	129	114	1			
2	P	78	Total	C	N	O	S	0	0	0
			619	391	120	107	1			

- Molecule 3 is a protein called Histone H2A type 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	103	Total	C	N	O	0	0	0
			795	501	155	139			
3	G	105	Total	C	N	O	0	0	0
			809	510	158	141			
3	M	103	Total	C	N	O	0	0	0
			795	501	155	139			
3	Q	105	Total	C	N	O	0	0	0
			809	510	158	141			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	99	ARG	GLY	conflict	UNP P06897
C	123	SER	ALA	conflict	UNP P06897
G	99	ARG	GLY	conflict	UNP P06897
G	123	SER	ALA	conflict	UNP P06897
M	99	ARG	GLY	conflict	UNP P06897
M	123	SER	ALA	conflict	UNP P06897
Q	99	ARG	GLY	conflict	UNP P06897
Q	123	SER	ALA	conflict	UNP P06897

- Molecule 4 is a protein called Histone H2B 1.1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	95	Total	C	N	O	S	0	0	0
			745	469	134	140	2			
4	H	93	Total	C	N	O	S	0	0	0
			726	457	130	137	2			
4	N	95	Total	C	N	O	S	0	0	0
			745	469	134	140	2			
4	R	93	Total	C	N	O	S	0	0	0
			726	457	130	137	2			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	0	MET	-	initiating methionine	UNP P02281
D	29	THR	SER	conflict	UNP P02281
H	0	MET	-	initiating methionine	UNP P02281
H	29	THR	SER	conflict	UNP P02281
N	0	MET	-	initiating methionine	UNP P02281
N	29	THR	SER	conflict	UNP P02281
R	0	MET	-	initiating methionine	UNP P02281

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Chain	Residue	Modelled	Actual	Comment	Reference
R	29	THR	SER	conflict	UNP P02281

- Molecule 5 is a DNA chain called DNA (145-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	I	145	Total	C	N	O	P	0	0	0
			2952	1404	537	867	144			
5	S	145	Total	C	N	O	P	0	0	0
			2952	1404	537	867	144			

- Molecule 6 is a DNA chain called DNA (145-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	J	145	Total	C	N	O	P	0	0	0
			2987	1416	558	869	144			
6	T	145	Total	C	N	O	P	0	0	0
			2987	1416	558	869	144			

- Molecule 7 is a protein called Ubiquitin carboxyl-terminal hydrolase 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	U	447	Total	C	N	O	S	0	0	0
			3569	2264	609	661	35			
7	Z	451	Total	C	N	O	S	0	0	0
			3600	2283	615	667	35			
7	e	453	Total	C	N	O	S	0	0	0
			3617	2292	618	672	35			
7	j	447	Total	C	N	O	S	0	0	0
			3566	2262	609	660	35			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
U	0	ALA	-	expression tag	UNP P50102
U	146	ALA	CYS	engineered mutation	UNP P50102
Z	0	ALA	-	expression tag	UNP P50102
Z	146	ALA	CYS	engineered mutation	UNP P50102
e	0	ALA	-	expression tag	UNP P50102
e	146	ALA	CYS	engineered mutation	UNP P50102
j	0	ALA	-	expression tag	UNP P50102
j	146	ALA	CYS	engineered mutation	UNP P50102

- Molecule 8 is a protein called Transcription and mRNA export factor SUS1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	V	89	Total	C	N	O	S	0	0	0
			719	451	118	148	2			
8	a	90	Total	C	N	O	S	0	0	0
			730	457	119	152	2			
8	f	85	Total	C	N	O	S	0	0	0
			690	432	112	144	2			
8	k	86	Total	C	N	O	S	0	0	0
			695	435	115	143	2			

- Molecule 9 is a protein called SAGA-associated factor 11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	W	90	Total	C	N	O	S	0	0	0
			718	440	133	142	3			
9	b	89	Total	C	N	O	S	0	0	0
			710	434	132	141	3			
9	g	90	Total	C	N	O	S	0	0	0
			718	440	133	142	3			
9	l	93	Total	C	N	O	S	0	0	0
			739	452	136	148	3			

- Molecule 10 is a protein called Polyubiquitin-B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	X	76	Total	C	N	O	S	0	0	0
			601	378	105	117	1			
10	c	76	Total	C	N	O	S	0	0	0
			601	378	105	117	1			
10	h	76	Total	C	N	O	S	0	0	0
			601	378	105	117	1			
10	m	76	Total	C	N	O	S	0	0	0
			601	378	105	117	1			

- Molecule 11 is a protein called SAGA-associated factor 73.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	Y	82	Total	C	N	O	S	0	0	0
			663	419	112	127	5			
11	d	89	Total	C	N	O	S	0	0	0
			710	447	120	138	5			
11	i	89	Total	C	N	O	S	0	0	0
			710	447	120	138	5			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	n	86	Total	C	N	O	S	0	0	0
			693	437	117	134	5			

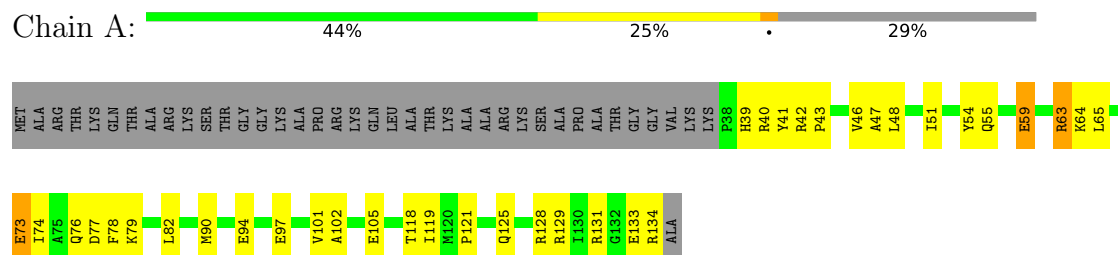
- Molecule 12 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	U	6	Total	Zn	0	0
			6	6		
12	W	1	Total	Zn	0	0
			1	1		
12	Y	1	Total	Zn	0	0
			1	1		
12	Z	6	Total	Zn	0	0
			6	6		
12	b	1	Total	Zn	0	0
			1	1		
12	d	1	Total	Zn	0	0
			1	1		
12	e	6	Total	Zn	0	0
			6	6		
12	g	1	Total	Zn	0	0
			1	1		
12	i	1	Total	Zn	0	0
			1	1		
12	j	6	Total	Zn	0	0
			6	6		
12	l	1	Total	Zn	0	0
			1	1		
12	n	1	Total	Zn	0	0
			1	1		

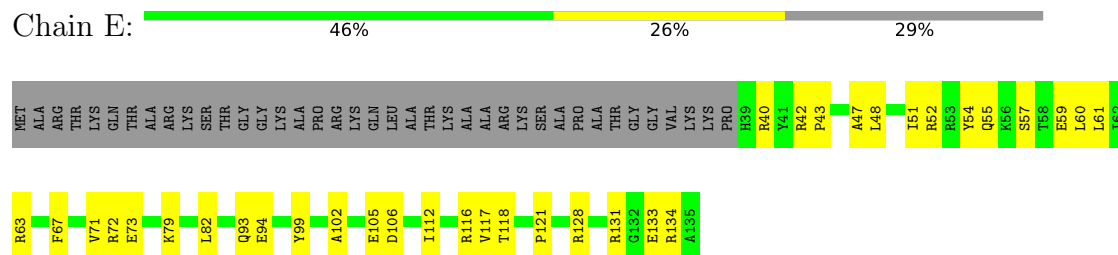
3 Residue-property plots [i](#)

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

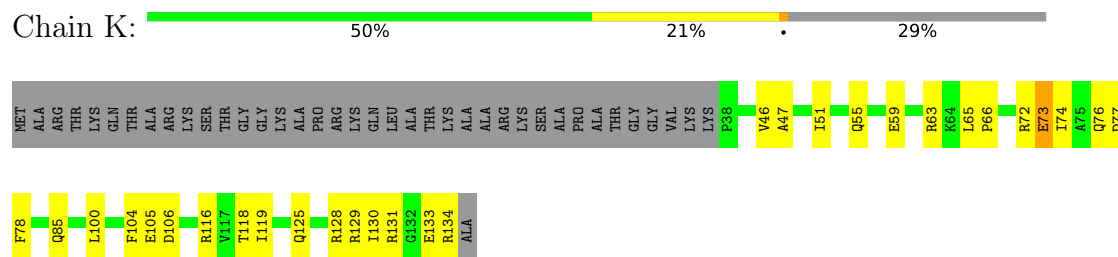
• Molecule 1: Histone H3.2



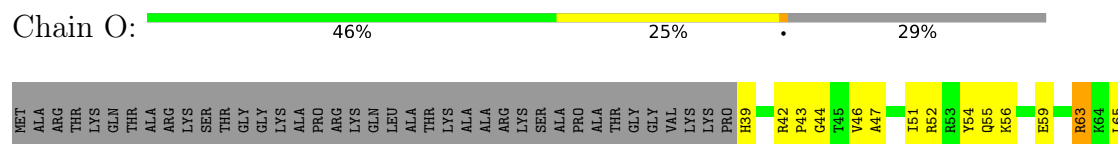
• Molecule 1: Histone H3.2



• Molecule 1: Histone H3.2

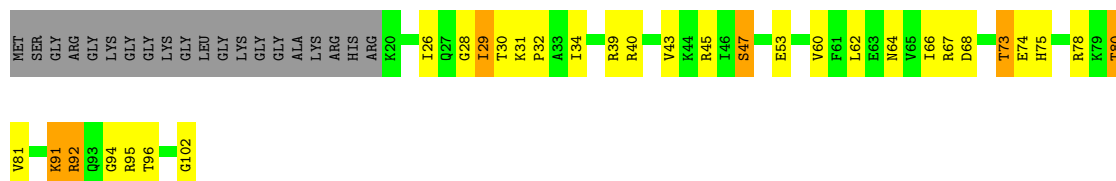


• Molecule 1: Histone H3.2

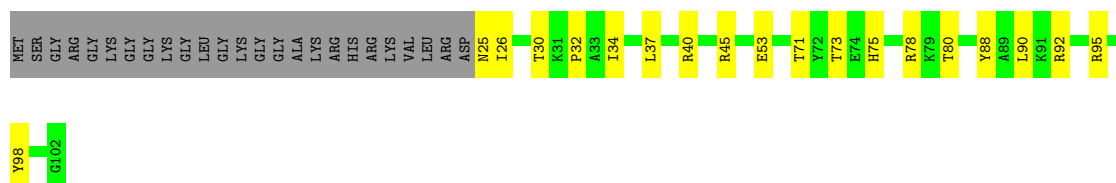




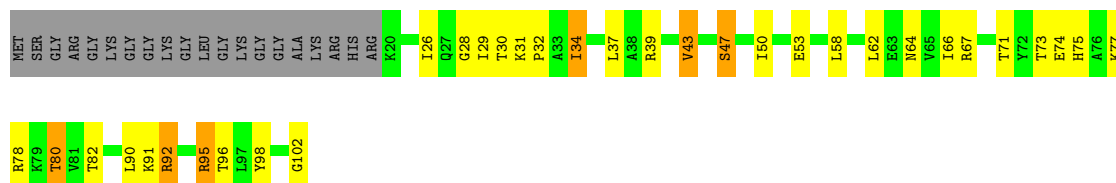
• Molecule 2: Histone H4



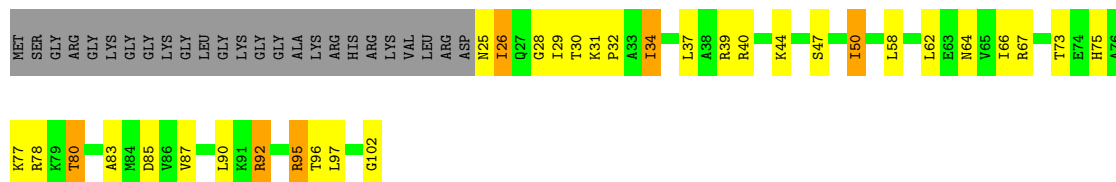
• Molecule 2: Histone H4



• Molecule 2: Histone H4



• Molecule 2: Histone H4

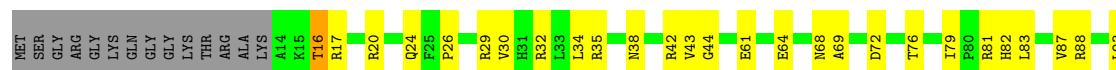


• Molecule 3: Histone H2A type 1

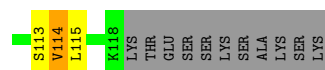




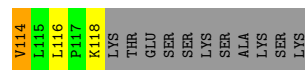
- Molecule 3: Histone H2A type 1



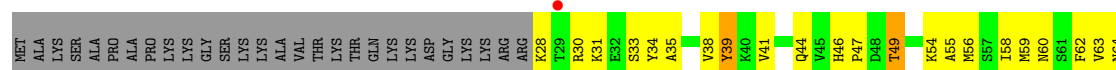
- Molecule 3: Histone H2A type 1



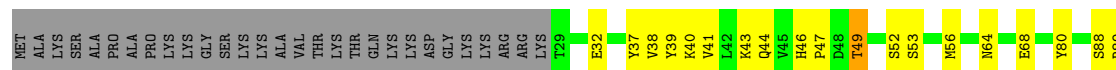
- Molecule 3: Histone H2A type 1



- Molecule 4: Histone H2B 1.1



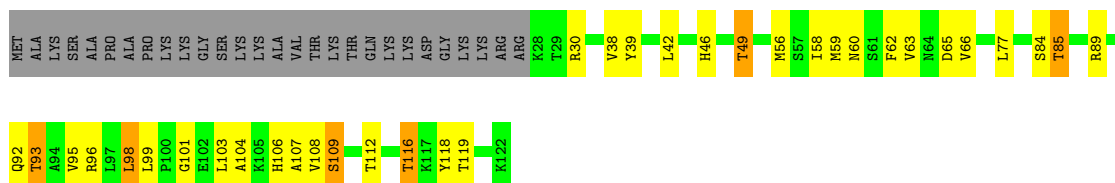
- Molecule 4: Histone H2B 1.1





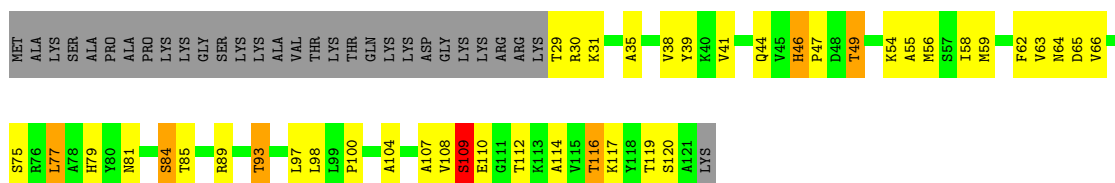
• Molecule 4: Histone H2B 1.1

Chain N: 49% 24% 5% 23%



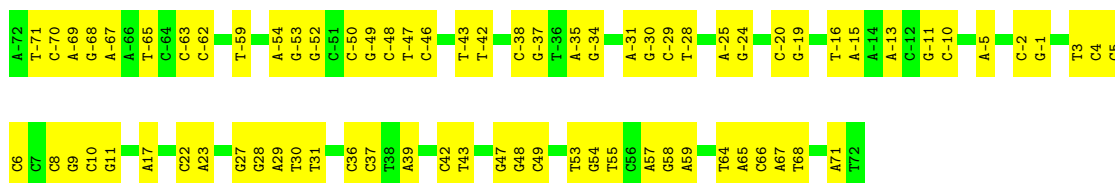
• Molecule 4: Histone H2B 1.1

Chain R: 41% 29% 5% 24%



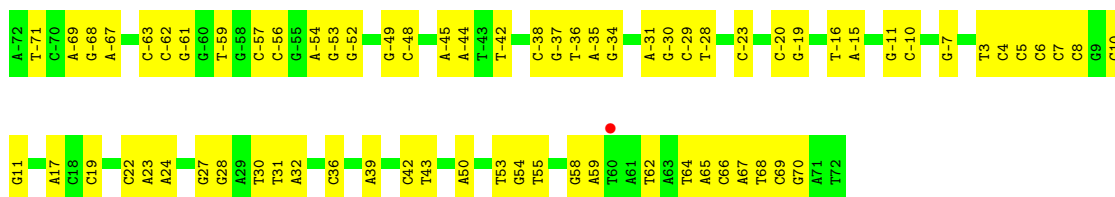
• Molecule 5: DNA (145-MER)

Chain I: 48% 52%



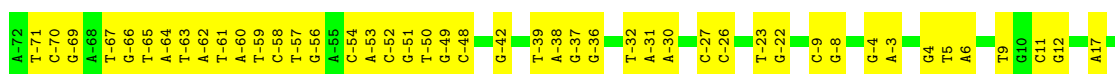
• Molecule 5: DNA (145-MER)

Chain S: 51% 49%



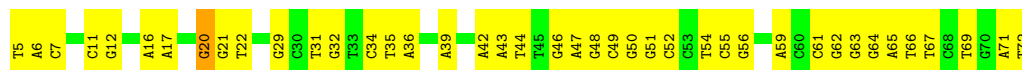
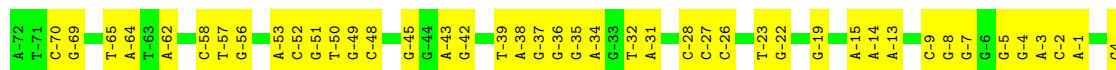
• Molecule 6: DNA (145-MER)

Chain J: 49% 51%

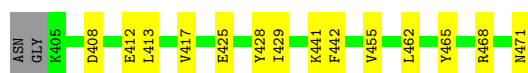
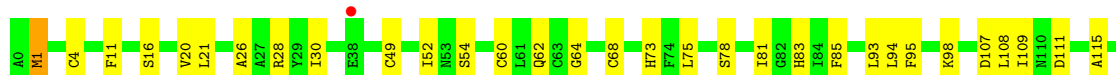




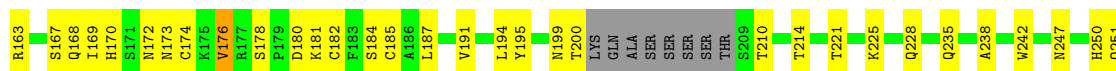
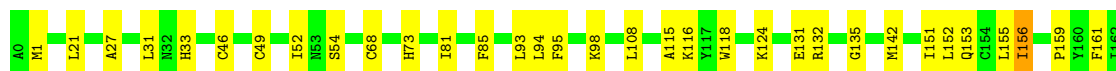
• Molecule 6: DNA (145-MER)



• Molecule 7: Ubiquitin carboxyl-terminal hydrolase 8

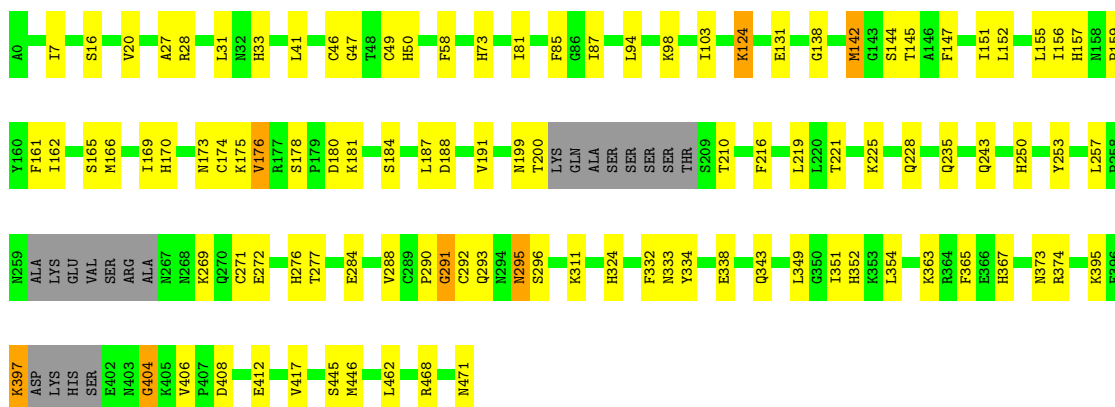


• Molecule 7: Ubiquitin carboxyl-terminal hydrolase 8



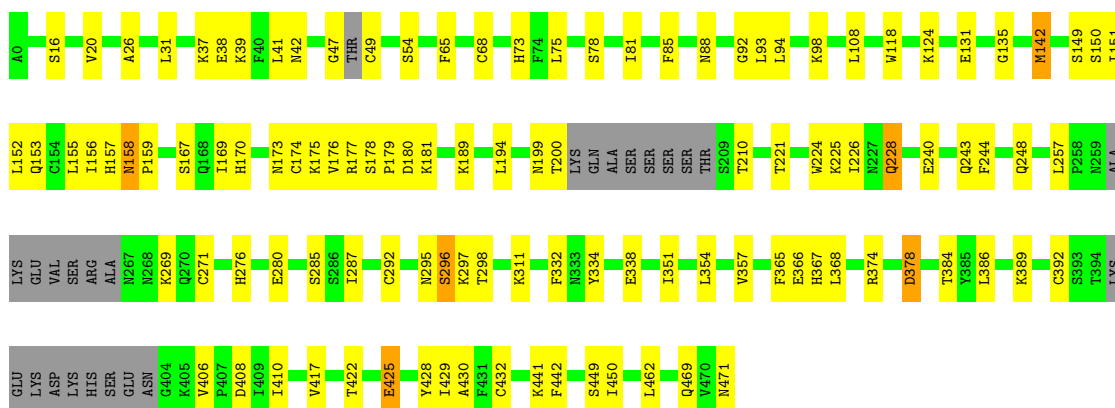
• Molecule 7: Ubiquitin carboxyl-terminal hydrolase 8





• Molecule 7: Ubiquitin carboxyl-terminal hydrolase 8

Chain j: 71% 22% • 5%



• Molecule 8: Transcription and mRNA export factor SUS1

Chain V: 69% 24% 7%



• Molecule 8: Transcription and mRNA export factor SUS1

Chain a: 77% 16% • 6%



• Molecule 8: Transcription and mRNA export factor SUS1

Chain f: 67% 21% • 11%



- Molecule 8: Transcription and mRNA export factor SUS1

Chain k:  72% 18% 10%



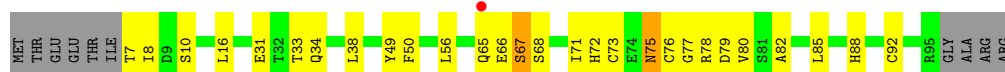
- Molecule 9: SAGA-associated factor 11

Chain W:  69% 19% 9%



- Molecule 9: SAGA-associated factor 11

Chain b:  62% 26% 10%



- Molecule 9: SAGA-associated factor 11

Chain g:  67% 23% 9%



- Molecule 9: SAGA-associated factor 11

Chain l:  71% 19% 6%



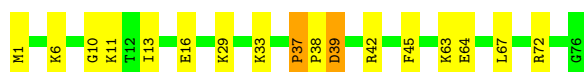
- Molecule 10: Polyubiquitin-B

Chain X:  80% 18% 2%



- Molecule 10: Polyubiquitin-B

Chain c:  78% 20% 2%



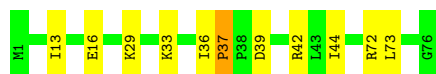
- Molecule 10: Polyubiquitin-B

Chain h:  76% 22% .



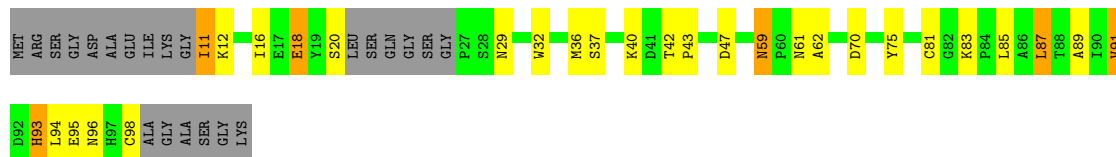
• Molecule 10: Polyubiquitin-B

Chain m:  86% 13% .



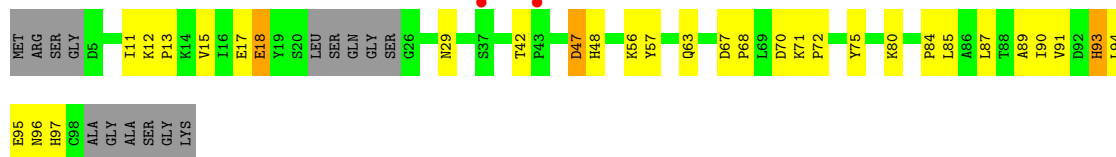
• Molecule 11: SAGA-associated factor 73

Chain Y:  51% 22% 6% 21%



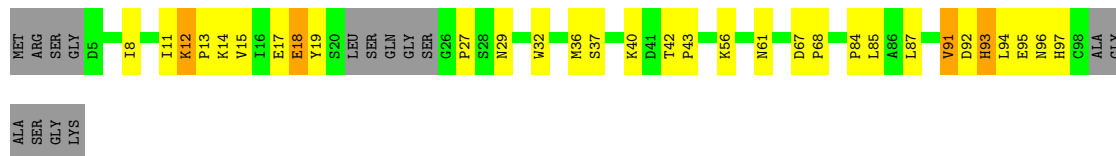
• Molecule 11: SAGA-associated factor 73

Chain d:  2% 56% 27% 14%



• Molecule 11: SAGA-associated factor 73

Chain i:  56% 26% 14%



• Molecule 11: SAGA-associated factor 73

Chain n:  59% 20% 17%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	178.80Å 179.17Å 353.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.50 – 3.82 49.50 – 3.82	Depositor EDS
% Data completeness (in resolution range)	98.9 (49.50-3.82) 99.4 (49.50-3.82)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 3.67Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.234 , 0.256 0.232 , 0.260	Depositor DCC
R_{free} test set	1997 reflections (1.64%)	wwPDB-VP
Wilson B-factor (Å ²)	168.1	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 368.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	0.228 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	49079	wwPDB-VP
Average B, all atoms (Å ²)	197.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.69% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.38	0/814	0.88	0/1092
1	E	0.34	0/812	0.88	2/1088 (0.2%)
1	K	0.37	0/814	0.86	0/1092
1	O	0.39	0/812	0.94	0/1088
2	B	0.41	0/669	0.99	0/894
2	F	0.45	0/626	0.96	0/837
2	L	0.43	0/669	0.96	1/894 (0.1%)
2	P	0.42	0/626	1.02	2/837 (0.2%)
3	C	0.36	0/805	1.00	4/1088 (0.4%)
3	G	0.38	0/819	1.08	3/1106 (0.3%)
3	M	0.36	0/805	0.92	2/1088 (0.2%)
3	Q	0.36	0/819	1.06	2/1106 (0.2%)
4	D	0.40	0/756	0.96	2/1015 (0.2%)
4	H	0.37	0/737	0.96	0/993
4	N	0.37	0/756	0.94	0/1015
4	R	0.40	0/737	1.01	2/993 (0.2%)
5	I	0.28	0/3308	0.61	0/5099
5	S	0.28	0/3308	0.61	0/5099
6	J	0.27	0/3354	0.62	0/5180
6	T	0.26	0/3354	0.63	1/5180 (0.0%)
7	U	0.32	0/3647	0.83	6/4916 (0.1%)
7	Z	0.31	0/3678	0.83	3/4955 (0.1%)
7	e	0.32	0/3695	0.85	8/4978 (0.2%)
7	j	0.33	0/3643	0.85	9/4908 (0.2%)
8	V	0.31	0/725	0.81	0/977
8	a	0.32	0/735	0.86	1/990 (0.1%)
8	f	0.37	0/695	1.06	4/936 (0.4%)
8	k	0.34	0/701	0.77	0/944
9	W	0.34	0/726	0.92	2/982 (0.2%)
9	b	0.32	0/718	0.83	0/971
9	g	0.37	0/726	0.87	0/982
9	l	0.34	0/747	0.91	1/1011 (0.1%)

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
10	X	0.29	0/607	0.83	0/816
10	c	0.30	0/607	0.84	2/816 (0.2%)
10	h	0.29	0/607	0.82	0/816
10	m	0.28	0/607	0.83	2/816 (0.2%)
11	Y	0.37	0/678	1.18	7/915 (0.8%)
11	d	0.36	0/725	1.09	3/978 (0.3%)
11	i	0.37	0/725	1.12	5/978 (0.5%)
11	n	0.35	0/707	1.05	7/953 (0.7%)
All	All	0.33	0/51099	0.84	81/71422 (0.1%)

There are no bond length outliers.

All (81) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	n	93	HIS	N-CA-C	-9.45	100.86	111.82
8	f	15	TYR	N-CA-C	9.41	121.14	111.07
8	f	16	LEU	N-CA-C	9.39	122.69	111.33
11	d	93	HIS	N-CA-C	-9.36	100.96	111.82
3	Q	116	LEU	CA-C-N	9.35	129.29	120.03
3	Q	116	LEU	C-N-CA	9.35	129.29	120.03
11	Y	12	LYS	CA-C-N	8.94	129.11	119.28
11	Y	12	LYS	C-N-CA	8.94	129.11	119.28
11	i	93	HIS	N-CA-C	-8.82	101.75	111.36
11	Y	93	HIS	N-CA-C	-8.72	101.71	111.82
3	G	116	LEU	CA-C-N	8.47	128.50	119.78
3	G	116	LEU	C-N-CA	8.47	128.50	119.78
7	U	157	HIS	N-CA-C	8.41	122.07	112.57
11	i	12	LYS	CA-C-N	8.21	129.00	119.47
11	i	12	LYS	C-N-CA	8.21	129.00	119.47
7	Z	436	GLY	N-CA-C	-8.03	105.13	114.69
11	d	67	ASP	CA-C-N	7.66	127.38	119.56
11	d	67	ASP	C-N-CA	7.66	127.38	119.56
11	n	36	MET	N-CA-C	-7.66	102.93	111.28
7	e	404	GLY	N-CA-C	7.25	120.55	111.85
1	E	42	ARG	CA-C-N	6.74	128.26	119.84
1	E	42	ARG	C-N-CA	6.74	128.26	119.84
7	j	157	HIS	N-CA-C	6.64	120.08	112.57
11	n	12	LYS	CA-C-N	6.60	127.12	119.47
11	n	12	LYS	C-N-CA	6.60	127.12	119.47
11	Y	47	ASP	N-CA-C	6.45	118.39	111.36
7	j	384	THR	N-CA-C	6.43	119.28	111.82
7	Z	214	THR	N-CA-C	6.37	117.88	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	l	82	ALA	N-CA-C	6.34	118.19	111.28
7	Z	152	LEU	N-CA-C	6.33	118.99	111.71
11	Y	59	ASN	CA-C-N	6.32	126.23	119.28
11	Y	59	ASN	C-N-CA	6.32	126.23	119.28
3	C	116	LEU	CA-C-N	6.28	126.25	120.03
3	C	116	LEU	C-N-CA	6.28	126.25	120.03
7	e	291	GLY	N-CA-C	-6.03	104.50	112.82
4	R	46	HIS	CA-C-N	5.99	126.15	119.32
4	R	46	HIS	C-N-CA	5.99	126.15	119.32
2	P	83	ALA	N-CA-C	5.95	118.53	111.33
7	e	157	HIS	N-CA-C	5.87	119.20	112.57
4	D	99	LEU	CA-C-N	5.82	127.39	120.23
4	D	99	LEU	C-N-CA	5.82	127.39	120.23
10	c	37	PRO	CA-C-N	5.78	125.64	119.28
10	c	37	PRO	C-N-CA	5.78	125.64	119.28
7	U	4	CYS	CA-C-N	5.74	125.36	119.56
7	U	4	CYS	C-N-CA	5.74	125.36	119.56
7	e	406	VAL	CA-C-N	5.67	125.64	120.03
7	e	406	VAL	C-N-CA	5.67	125.64	120.03
8	a	65	VAL	N-CA-C	5.64	115.83	110.53
2	P	50	ILE	N-CA-C	5.58	115.78	110.53
6	T	20	DG	O3'-P-O5'	-5.55	95.68	104.00
7	j	158	ASN	CA-C-N	5.49	125.32	119.28
7	j	158	ASN	C-N-CA	5.49	125.32	119.28
3	M	100	VAL	N-CA-C	5.42	115.66	107.75
11	i	91	VAL	N-CA-C	5.41	116.14	110.62
7	j	389	LYS	N-CA-C	5.40	117.92	111.71
8	f	14	GLN	CA-C-N	5.37	127.42	120.44
8	f	14	GLN	C-N-CA	5.37	127.42	120.44
11	Y	91	VAL	N-CA-C	5.34	116.07	110.62
7	e	124	LYS	N-CA-C	5.31	118.92	112.23
7	j	93	LEU	N-CA-C	5.30	118.57	110.20
9	W	42	TYR	CA-C-N	5.25	125.55	119.47
9	W	42	TYR	C-N-CA	5.25	125.55	119.47
3	C	25	PHE	CA-C-N	5.23	126.38	119.84
3	C	25	PHE	C-N-CA	5.23	126.38	119.84
11	n	59	ASN	CA-C-N	5.22	125.02	119.28
11	n	59	ASN	C-N-CA	5.22	125.02	119.28
7	j	386	LEU	N-CA-C	5.20	116.89	108.41
11	i	32	TRP	N-CA-C	5.19	117.68	111.71
11	n	39	ALA	N-CA-C	5.19	116.94	111.28
3	M	81	ARG	NE-CZ-NH1	-5.11	116.39	121.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	j	226	ILE	N-CA-C	-5.11	106.17	111.58
3	G	16	THR	N-CA-C	5.11	117.96	110.30
2	L	43	VAL	N-CA-C	5.10	115.62	108.89
7	j	410	ILE	N-CA-C	5.07	115.27	108.17
10	m	37	PRO	CA-C-N	5.07	124.85	119.28
10	m	37	PRO	C-N-CA	5.07	124.85	119.28
7	U	158	ASN	CA-C-N	5.04	124.83	119.28
7	U	158	ASN	C-N-CA	5.04	124.83	119.28
7	e	295	ASN	CA-C-N	5.01	130.73	121.70
7	e	295	ASN	C-N-CA	5.01	130.73	121.70
7	U	455	VAL	N-CA-C	-5.01	105.51	110.62

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	802	0	841	25	0
1	E	801	0	838	24	0
1	K	802	0	841	19	0
1	O	801	0	838	27	0
2	B	662	0	709	23	0
2	F	619	0	659	18	0
2	L	662	0	709	25	0
2	P	619	0	659	27	0
3	C	795	0	846	25	0
3	G	809	0	864	28	0
3	M	795	0	846	27	0
3	Q	809	0	864	25	0
4	D	745	0	773	38	0
4	H	726	0	747	29	0
4	N	745	0	773	28	0
4	R	726	0	747	33	0
5	I	2952	0	1629	81	0
5	S	2952	0	1629	88	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	J	2987	0	1630	84	0
6	T	2987	0	1630	96	0
7	U	3569	0	3476	70	0
7	Z	3600	0	3514	64	0
7	e	3617	0	3524	71	0
7	j	3566	0	3472	68	0
8	V	719	0	730	14	0
8	a	730	0	737	11	0
8	f	690	0	691	18	0
8	k	695	0	704	12	0
9	W	718	0	710	20	0
9	b	710	0	699	21	0
9	g	718	0	709	20	0
9	l	739	0	725	21	0
10	X	601	0	629	12	0
10	c	601	0	629	11	0
10	h	601	0	629	17	0
10	m	601	0	629	8	0
11	Y	663	0	653	23	0
11	d	710	0	697	21	0
11	i	710	0	697	24	0
11	n	693	0	682	18	0
12	U	6	0	0	0	0
12	W	1	0	0	0	0
12	Y	1	0	0	0	0
12	Z	6	0	0	0	0
12	b	1	0	0	0	0
12	d	1	0	0	0	0
12	e	6	0	0	0	0
12	g	1	0	0	0	0
12	i	1	0	0	0	0
12	j	6	0	0	0	0
12	l	1	0	0	0	0
12	n	1	0	0	0	0
All	All	49079	0	44008	986	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (986) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:47:PRO:HG3	9:I:79:ASP:HB2	1.52	0.88
7:Z:170:HIS:CE1	7:Z:182:CYS:SG	2.64	0.85
5:S:19:DC:H42	6:T:-19:DG:H1	1.21	0.84
5:S:70:DG:N2	6:T:-70:DC:O2	2.13	0.82
3:C:20:ARG:NH2	5:I:-42:DT:OP1	2.11	0.82
10:h:39:ASP:OD1	10:h:39:ASP:N	2.14	0.80
7:Z:46:CYS:SG	7:Z:73:HIS:ND1	2.54	0.78
3:M:97:LEU:HB3	3:M:100:VAL:HB	1.66	0.77
4:N:38:VAL:HG21	4:N:56:MET:HE1	1.65	0.77
4:N:95:VAL:HG13	4:N:99:LEU:HD12	1.65	0.77
4:D:95:VAL:HG13	4:D:99:LEU:HD12	1.66	0.77
9:b:75:ASN:ND2	9:b:92:CYS:SG	2.59	0.76
1:A:125:GLN:NE2	2:B:53:GLU:OE2	2.18	0.75
5:I:-70:DC:O2	6:J:70:DG:N2	2.18	0.75
10:c:39:ASP:OD1	10:c:39:ASP:N	2.14	0.74
2:F:75:HIS:NE2	4:H:90:GLU:OE2	2.20	0.74
1:A:94:GLU:OE1	3:G:104:GLN:NE2	2.19	0.74
2:P:78:ARG:NH1	2:P:80:THR:O	2.20	0.74
7:e:33:HIS:ND1	11:i:42:THR:O	2.20	0.74
7:e:295:ASN:HB3	7:e:296:SER:HA	1.70	0.74
7:Z:295:ASN:ND2	7:Z:338:GLU:OE2	2.21	0.73
7:j:295:ASN:HB3	7:j:296:SER:HA	1.71	0.73
7:U:295:ASN:HB3	7:U:296:SER:HA	1.71	0.73
2:P:92:ARG:HH21	4:R:98:LEU:HD23	1.53	0.73
7:Z:295:ASN:HB3	7:Z:296:SER:HA	1.69	0.72
10:m:16:GLU:O	10:m:29:LYS:NZ	2.22	0.72
3:C:42:ARG:HD2	5:I:-35:DA:H4'	1.69	0.72
7:e:288:VAL:HG22	7:e:296:SER:HB3	1.72	0.71
3:G:42:ARG:HG2	5:I:39:DA:H5''	1.72	0.71
3:M:71:ARG:HH12	9:g:78:ARG:HE	1.36	0.71
3:G:68:ASN:OD1	9:b:78:ARG:NH2	2.23	0.71
3:Q:20:ARG:NH2	6:T:-42:DG:OP2	2.25	0.70
7:j:392:CYS:HB2	7:j:406:VAL:HG21	1.72	0.70
3:Q:97:LEU:HB3	3:Q:100:VAL:HB	1.74	0.70
3:M:42:ARG:HD2	5:S:-35:DA:H4'	1.74	0.70
5:I:-53:DG:H2''	5:I:-52:DG:OP2	1.92	0.70
3:M:21:ALA:HB2	4:N:118:TYR:HD1	1.56	0.70
5:I:23:DA:H61	6:J:-23:DT:H3	1.39	0.69
8:V:38:TRP:HA	8:V:41:LYS:HE2	1.74	0.69
9:l:80:VAL:HG21	9:l:88:HIS:CG	2.26	0.69
4:D:38:VAL:HG21	4:D:56:MET:HE1	1.74	0.69
7:Z:33:HIS:ND1	11:d:42:THR:O	2.24	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Q:90:ASP:HB3	3:Q:93:LEU:HB2	1.75	0.69
9:W:34:GLN:NE2	11:Y:70:ASP:O	2.24	0.69
8:k:47:LYS:NZ	9:l:31:GLU:OE1	2.26	0.69
7:e:124:LYS:HD3	11:i:94:LEU:HD22	1.75	0.68
7:e:250:HIS:ND1	7:e:276:HIS:HE1	1.90	0.68
9:W:66:GLU:HG3	9:W:83:ASN:HB3	1.74	0.68
1:E:60:LEU:HD13	1:E:93:GLN:HG2	1.76	0.68
2:P:75:HIS:HD2	4:R:93:THR:HG21	1.59	0.67
5:S:-67:DA:H61	6:T:67:DT:H3	1.41	0.67
5:S:8:DC:H42	6:T:-8:DG:H1	1.42	0.67
8:a:46:THR:HG22	8:a:50:MET:HE2	1.75	0.67
7:j:169:ILE:O	7:j:173:ASN:ND2	2.27	0.67
9:W:66:GLU:CD	9:W:67:SER:HA	2.20	0.67
9:g:46:ARG:NH2	9:g:58:ILE:O	2.28	0.67
1:E:116:ARG:NH1	1:E:118:THR:O	2.28	0.67
3:M:104:GLN:NE2	1:O:94:GLU:OE2	2.27	0.67
7:e:169:ILE:O	7:e:173:ASN:ND2	2.28	0.67
5:S:-53:DG:H2"	5:S:-52:DG:OP2	1.94	0.66
7:j:295:ASN:ND2	7:j:338:GLU:OE2	2.27	0.66
4:D:112:THR:O	4:D:116:THR:OG1	2.13	0.66
2:L:74:GLU:O	4:N:89:ARG:NH2	2.28	0.66
8:V:70:LEU:O	8:V:78:ARG:NH2	2.27	0.66
4:H:116:THR:OG1	7:Z:425:GLU:OE2	2.12	0.66
7:U:280:GLU:O	11:Y:59:ASN:ND2	2.29	0.66
10:h:16:GLU:O	10:h:29:LYS:NZ	2.29	0.66
4:D:84:SER:N	5:I:-34:DG:OP1	2.26	0.65
5:S:69:DC:O2	6:T:-69:DG:N2	2.29	0.65
6:J:-32:DT:H2"	6:J:-31:DA:C8	2.32	0.65
8:V:47:LYS:NZ	9:W:31:GLU:OE1	2.17	0.65
5:S:-28:DT:O2	6:T:29:DG:N2	2.29	0.65
3:Q:61:GLU:OE1	9:l:91:ARG:NH2	2.30	0.65
1:O:100:LEU:HD11	2:P:58:LEU:HD13	1.79	0.65
7:j:287:ILE:HB	7:j:297:LYS:HB2	1.78	0.65
7:j:408:ASP:HB2	7:j:471:ASN:HB2	1.78	0.65
7:j:26:ALA:HB1	11:n:36:MET:HE1	1.78	0.65
10:X:39:ASP:N	10:X:39:ASP:OD1	2.27	0.64
6:T:31:DT:H2"	6:T:32:DG:C8	2.31	0.64
9:W:80:VAL:HG21	9:W:88:HIS:CG	2.31	0.64
10:X:42:ARG:HD3	10:X:72:ARG:HG3	1.77	0.64
4:R:38:VAL:HG21	4:R:56:MET:HE1	1.79	0.64
4:D:47:PRO:HG2	9:W:79:ASP:HB2	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:j:295:ASN:CB	7:j:296:SER:HA	2.27	0.64
2:B:60:VAL:O	2:B:64:ASN:ND2	2.30	0.64
2:F:75:HIS:CD2	4:H:93:THR:HG21	2.32	0.64
10:X:45:PHE:HB3	10:X:50:LEU:HD21	1.80	0.64
7:Z:295:ASN:CB	7:Z:296:SER:HA	2.27	0.64
8:k:38:TRP:HA	8:k:41:LYS:HE2	1.80	0.64
5:S:-34:DG:H1	6:T:34:DC:H42	1.43	0.63
8:f:12:ILE:HA	11:i:11:ILE:HD11	1.81	0.63
2:L:78:ARG:NH1	2:L:80:THR:O	2.31	0.63
4:D:39:TYR:OH	5:I:-53:DG:OP2	2.14	0.63
7:e:295:ASN:ND2	7:e:338:GLU:OE2	2.27	0.63
5:S:58:DG:H1	6:T:-58:DC:H42	1.46	0.63
9:b:66:GLU:CD	9:b:67:SER:HA	2.24	0.63
5:S:7:DC:H42	6:T:-7:DG:H1	1.47	0.63
2:L:75:HIS:CG	4:N:93:THR:HG21	2.33	0.62
7:U:292:CYS:SG	7:U:339:CYS:HB3	2.39	0.62
7:U:295:ASN:CB	7:U:296:SER:HA	2.29	0.62
9:l:66:GLU:CD	9:l:67:SER:HA	2.24	0.62
7:e:295:ASN:CB	7:e:296:SER:HA	2.28	0.62
6:T:-53:DA:H1'	6:T:-52:DC:H5'	1.82	0.62
10:c:16:GLU:O	10:c:29:LYS:NZ	2.32	0.62
2:B:78:ARG:NH1	2:B:80:THR:O	2.32	0.62
5:S:-59:DT:H3	6:T:59:DA:H61	1.47	0.62
5:S:54:DG:H2''	5:S:55:DT:OP2	2.00	0.62
7:U:54:SER:HB3	7:U:68:CYS:HB2	1.81	0.62
7:j:124:LYS:HD3	11:n:94:LEU:HD22	1.82	0.62
10:X:16:GLU:O	10:X:29:LYS:NZ	2.33	0.62
5:I:-46:DC:O2	6:J:46:DG:N2	2.27	0.62
9:g:75:ASN:ND2	9:g:92:CYS:SG	2.71	0.62
2:F:78:ARG:NH1	2:F:80:THR:O	2.33	0.62
7:Z:124:LYS:HD3	11:d:94:LEU:HD22	1.81	0.62
7:U:64:GLY:HA2	11:Y:32:TRP:CD1	2.35	0.61
7:e:161:PHE:HZ	7:e:187:LEU:HD13	1.64	0.61
1:K:125:GLN:NE2	2:L:53:GLU:OE2	2.22	0.61
9:g:66:GLU:CD	9:g:67:SER:HA	2.26	0.61
3:G:97:LEU:HB3	3:G:100:VAL:HB	1.83	0.61
3:C:104:GLN:NE2	1:E:94:GLU:OE2	2.33	0.61
4:H:46:HIS:HB3	4:H:49:THR:OG1	2.00	0.61
1:E:117:VAL:N	6:J:-3:DA:OP1	2.34	0.61
2:P:77:LYS:HG3	4:R:89:ARG:HH22	1.65	0.61
7:Z:169:ILE:O	7:Z:173:ASN:ND2	2.33	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:S:66:DC:H1'	5:S:67:DA:H5'	1.84	0.60
7:Z:155:LEU:HD21	7:Z:242:TRP:HZ3	1.66	0.60
7:e:221:THR:HG22	7:e:225:LYS:HE3	1.84	0.60
8:k:70:LEU:O	8:k:78:ARG:NH2	2.30	0.60
1:A:54:TYR:O	2:B:40:ARG:NE	2.33	0.60
7:U:124:LYS:HD3	11:Y:94:LEU:HD13	1.82	0.60
7:e:277:THR:HG22	11:i:61:ASN:HD22	1.66	0.60
2:B:74:GLU:O	4:D:89:ARG:NH2	2.35	0.60
3:C:55:LEU:O	3:C:59:THR:HG22	2.01	0.60
7:Z:221:THR:HG22	7:Z:225:LYS:HE3	1.82	0.60
7:e:243:GLN:HE22	10:h:44:ILE:HD12	1.64	0.60
8:f:38:TRP:HA	8:f:41:LYS:HE2	1.84	0.60
7:U:169:ILE:O	7:U:173:ASN:ND2	2.35	0.60
7:Z:250:HIS:NE2	7:Z:271:CYS:SG	2.74	0.60
8:a:38:TRP:HA	8:a:41:LYS:HE2	1.83	0.60
1:E:61:LEU:HD12	2:F:37:LEU:HD23	1.82	0.60
10:m:13:ILE:HD12	10:m:33:LYS:HD3	1.83	0.60
3:C:69:ALA:O	3:C:73:ASN:ND2	2.35	0.59
7:e:367:HIS:CG	10:h:36:ILE:HG12	2.37	0.59
7:e:412:GLU:OE1	7:e:468:ARG:NH1	2.35	0.59
4:R:112:THR:O	4:R:116:THR:OG1	2.19	0.59
5:S:-31:DA:OP1	7:e:374:ARG:NH2	2.30	0.59
6:T:21:DG:H2''	6:T:22:DT:OP2	2.03	0.59
7:U:21:LEU:HD13	7:U:115:ALA:HB1	1.83	0.59
5:S:64:DT:H3	6:T:-64:DA:H61	1.49	0.59
11:i:93:HIS:HA	11:i:97:HIS:HB2	1.83	0.59
3:G:29:ARG:NH1	5:I:49:DC:OP1	2.35	0.59
7:U:221:THR:HG22	7:U:225:LYS:HE3	1.84	0.59
1:E:67:PHE:O	1:E:71:VAL:HG23	2.02	0.59
2:P:90:LEU:HB3	2:P:95:ARG:O	2.02	0.59
4:R:109:SER:OG	4:R:110:GLU:N	2.33	0.59
10:m:42:ARG:HD3	10:m:72:ARG:HG3	1.84	0.59
7:Z:163:ARG:NH2	11:d:63:GLN:OE1	2.35	0.59
5:I:8:DC:H42	6:J:-8:DG:H1	1.49	0.58
6:T:46:DG:H4'	6:T:47:DA:OP1	2.02	0.58
6:T:63:DG:H2''	6:T:64:DG:OP2	2.02	0.58
3:C:97:LEU:HB3	3:C:100:VAL:HB	1.86	0.58
7:e:152:LEU:HD23	7:e:155:LEU:HD12	1.85	0.58
7:Z:287:ILE:HB	7:Z:297:LYS:HB2	1.84	0.58
6:T:-9:DC:H2''	6:T:-8:DG:C8	2.39	0.58
2:B:96:THR:HB	3:G:100:VAL:HG22	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:71:ARG:NH1	9:W:78:ARG:HE	2.00	0.58
6:J:63:DG:H2''	6:J:64:DG:OP2	2.03	0.58
1:K:106:ASP:HB3	1:O:130:ILE:HG12	1.86	0.58
4:R:117:LYS:O	4:R:120:SER:OG	2.18	0.58
7:e:291:GLY:O	7:e:293:GLN:N	2.33	0.58
2:P:75:HIS:CD2	4:R:93:THR:HG21	2.37	0.58
7:U:124:LYS:HE3	11:Y:94:LEU:HB3	1.85	0.58
7:Z:288:VAL:HG22	7:Z:296:SER:HB3	1.86	0.58
9:l:67:SER:OG	9:l:83:ASN:ND2	2.35	0.58
6:J:-63:DT:H2''	6:J:-62:DA:C8	2.39	0.58
5:I:64:DT:H3	6:J:-64:DA:H61	1.52	0.58
5:S:66:DC:H2''	5:S:67:DA:OP2	2.03	0.58
7:j:280:GLU:O	11:n:59:ASN:ND2	2.37	0.58
3:C:71:ARG:HH12	9:W:78:ARG:HG2	1.69	0.57
1:O:63:ARG:HG2	5:S:17:DA:O3'	2.04	0.57
5:I:-34:DG:H1	6:J:34:DC:H42	1.52	0.57
5:S:-20:DC:H2''	5:S:-19:DG:C8	2.39	0.57
7:U:146:ALA:HB3	10:X:76:GLY:HA2	1.86	0.57
9:W:76:CYS:SG	9:W:78:ARG:HB2	2.45	0.57
10:c:42:ARG:HD3	10:c:72:ARG:HG3	1.86	0.57
6:T:-57:DT:H2''	6:T:-56:DG:C8	2.39	0.57
6:J:46:DG:H4'	6:J:47:DA:OP1	2.04	0.57
1:O:116:ARG:NH1	1:O:118:THR:O	2.37	0.57
5:S:-31:DA:H1'	5:S:-30:DG:H5'	1.87	0.57
7:j:180:ASP:OD1	7:j:181:LYS:NZ	2.38	0.57
3:G:20:ARG:NH1	6:J:-42:DG:OP2	2.29	0.57
5:I:54:DG:H2''	5:I:55:DT:OP2	2.04	0.57
2:L:90:LEU:HB3	2:L:95:ARG:O	2.05	0.57
7:j:221:THR:HG22	7:j:225:LYS:HE3	1.86	0.57
3:C:20:ARG:O	4:D:118:TYR:HA	2.05	0.57
3:G:20:ARG:NH2	6:J:-42:DG:OP1	2.38	0.57
6:J:21:DG:H2''	6:J:22:DT:OP2	2.04	0.57
7:j:37:LYS:HG2	8:k:17:VAL:HG12	1.87	0.57
1:O:69:ARG:NH1	2:P:25:ASN:OD1	2.38	0.57
8:f:93:VAL:HG11	11:i:8:ILE:HD12	1.87	0.57
7:U:271:CYS:HB3	7:U:276:HIS:CG	2.40	0.57
5:S:69:DC:H42	6:T:-70:DC:H42	1.52	0.56
6:T:61:DC:H2''	6:T:62:DG:C8	2.40	0.56
1:O:69:ARG:HD2	2:P:25:ASN:HD21	1.71	0.56
5:I:66:DC:H2''	5:I:67:DA:OP2	2.04	0.56
1:K:119:ILE:HG13	2:L:50:ILE:HG13	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:30:ARG:HD2	5:S:50:DA:H5'	1.88	0.56
7:U:250:HIS:CE1	7:U:271:CYS:SG	2.99	0.56
5:I:-50:DC:N3	6:J:51:DG:N2	2.54	0.56
7:e:408:ASP:HB2	7:e:471:ASN:HB2	1.88	0.56
4:R:31:LYS:N	5:S:50:DA:OP1	2.39	0.56
5:I:23:DA:N6	6:J:-23:DT:H3	2.04	0.56
1:K:116:ARG:NH1	1:K:118:THR:O	2.39	0.56
3:Q:87:VAL:HG13	3:Q:93:LEU:HB3	1.88	0.56
7:Z:178:SER:OG	7:Z:180:ASP:OD2	2.23	0.56
7:e:85:PHE:HB3	7:e:94:LEU:HD11	1.88	0.56
1:O:84:PHE:N	6:T:-23:DT:OP1	2.31	0.55
7:j:75:LEU:O	7:j:78:SER:OG	2.22	0.55
1:E:54:TYR:HB3	2:F:40:ARG:HB2	1.88	0.55
7:j:378:ASP:OD1	7:j:378:ASP:N	2.30	0.55
5:I:42:DC:H2''	5:I:43:DT:OP2	2.07	0.55
4:D:30:ARG:NH2	6:J:49:DC:O4'	2.39	0.55
6:T:-49:DG:H2''	6:T:-48:DC:C6	2.41	0.55
10:c:13:ILE:HD12	10:c:33:LYS:HD3	1.89	0.55
6:J:-39:DT:H2''	6:J:-38:DA:C8	2.42	0.55
1:O:54:TYR:O	2:P:40:ARG:NE	2.38	0.55
3:G:24:GLN:HE21	4:H:40:LYS:HD3	1.71	0.55
5:S:-36:DT:H3	6:T:36:DA:H61	1.55	0.55
7:Z:378:ASP:OD1	7:Z:378:ASP:N	2.30	0.55
4:H:39:TYR:OH	6:J:-53:DA:H3'	2.07	0.55
7:e:152:LEU:HD11	7:e:219:LEU:HD22	1.88	0.55
1:E:43:PRO:HA	5:I:9:DG:H5'	1.89	0.54
5:S:28:DG:N2	6:T:-28:DC:O2	2.39	0.54
5:S:62:DT:H3	6:T:-62:DA:H61	1.53	0.54
5:I:27:DG:H2''	5:I:28:DG:C8	2.41	0.54
4:R:59:MET:O	4:R:63:VAL:HG23	2.08	0.54
7:U:367:HIS:CE1	10:X:36:ILE:HG12	2.42	0.54
5:I:-29:DC:H2''	5:I:-28:DT:OP2	2.08	0.54
8:a:47:LYS:NZ	9:b:31:GLU:OE1	2.41	0.54
2:B:73:THR:HG21	2:B:81:VAL:HG22	1.88	0.54
5:I:-31:DA:H1'	5:I:-30:DG:H5'	1.89	0.54
2:L:39:ARG:NH1	2:L:43:VAL:O	2.41	0.54
7:j:152:LEU:HD23	7:j:155:LEU:HD12	1.90	0.54
7:j:271:CYS:HB3	7:j:276:HIS:CG	2.42	0.54
1:E:47:ALA:N	5:I:9:DG:OP1	2.38	0.54
5:I:66:DC:H1'	5:I:67:DA:H5'	1.89	0.54
6:T:51:DG:H2''	6:T:52:DC:OP2	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:U:128:PRO:O	9:W:42:TYR:OH	2.15	0.54
7:e:180:ASP:OD1	7:e:181:LYS:NZ	2.41	0.54
4:H:102:GLU:OE2	4:H:105:LYS:NZ	2.41	0.54
5:S:27:DG:H2''	5:S:28:DG:C8	2.41	0.54
6:T:66:DT:H2''	6:T:67:DT:C5	2.43	0.54
8:V:8:LEU:HD12	11:Y:11:ILE:HG12	1.90	0.54
7:Z:85:PHE:HB3	7:Z:94:LEU:HD11	1.90	0.54
8:a:93:VAL:HG21	9:b:8:ILE:HD11	1.90	0.54
8:f:86:ARG:NE	9:g:9:ASP:OD1	2.41	0.54
11:i:85:LEU:HD11	11:i:93:HIS:CD2	2.43	0.54
11:n:85:LEU:HD11	11:n:93:HIS:CD2	2.43	0.54
4:D:35:ALA:HA	4:D:56:MET:SD	2.48	0.53
5:I:5:DC:O2	6:J:-4:DG:N2	2.42	0.53
6:J:51:DG:H2''	6:J:52:DC:OP2	2.07	0.53
6:T:-32:DT:H2''	6:T:-31:DA:C8	2.43	0.53
7:Z:54:SER:HB3	7:Z:68:CYS:HB2	1.89	0.53
7:Z:351:ILE:HG21	7:Z:354:LEU:HG	1.91	0.53
7:j:367:HIS:CD2	10:m:36:ILE:HG12	2.43	0.53
3:Q:24:GLN:HG3	4:R:44:GLN:NE2	2.22	0.53
5:S:-29:DC:H2''	5:S:-28:DT:OP2	2.08	0.53
7:U:271:CYS:HB3	7:U:276:HIS:ND1	2.22	0.53
7:U:295:ASN:ND2	7:U:338:GLU:OE2	2.42	0.53
7:j:31:LEU:HD22	7:j:39:LYS:HE3	1.89	0.53
4:H:39:TYR:CE1	4:H:43:LYS:HE3	2.43	0.53
2:L:31:LYS:N	2:L:32:PRO:HD2	2.24	0.53
7:e:253:TYR:OH	7:e:272:GLU:HB2	2.09	0.53
3:G:83:LEU:O	3:G:87:VAL:HG23	2.08	0.53
5:I:-54:DA:H61	6:J:53:DC:H42	1.56	0.53
3:M:87:VAL:HG11	3:M:97:LEU:HD12	1.91	0.53
5:S:-11:DG:H2''	5:S:-10:DC:OP2	2.08	0.53
7:U:60:CYS:SG	7:U:83:HIS:HD2	2.32	0.53
7:Z:417:VAL:HB	7:Z:462:LEU:HB2	1.91	0.53
9:g:67:SER:HB3	9:g:83:ASN:ND2	2.22	0.53
6:J:61:DC:H2''	6:J:62:DG:C8	2.44	0.53
7:U:167:SER:HA	11:Y:89:ALA:HB2	1.91	0.53
10:h:42:ARG:HB2	10:h:70:VAL:HG23	1.90	0.53
6:T:5:DT:H2''	6:T:6:DA:C8	2.43	0.53
7:e:178:SER:OG	7:e:180:ASP:OD2	2.25	0.53
11:n:12:LYS:O	11:n:15:VAL:HG12	2.09	0.53
5:I:-49:DG:H2''	5:I:-48:DC:C5	2.43	0.53
5:S:42:DC:H2''	5:S:43:DT:OP2	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:Z:52:ILE:HG13	8:a:17:VAL:HG21	1.91	0.53
7:j:175:LYS:NZ	7:j:257:LEU:O	2.39	0.53
4:D:46:HIS:HB3	4:D:49:THR:OG1	2.09	0.53
5:I:36:DC:H42	6:J:-36:DG:H1	1.56	0.53
1:K:131:ARG:O	3:M:99:ARG:NH1	2.41	0.53
6:T:-39:DT:H2''	6:T:-38:DA:C8	2.43	0.53
1:K:85:GLN:OE1	2:L:82:THR:HG22	2.09	0.53
9:g:76:CYS:SG	9:g:88:HIS:CE1	2.98	0.53
7:j:167:SER:HA	11:n:89:ALA:HB2	1.91	0.53
2:P:31:LYS:N	2:P:32:PRO:HD2	2.23	0.52
5:S:36:DC:H42	6:T:-36:DG:H1	1.56	0.52
6:T:55:DC:H2'	6:T:55:DC:OP2	2.09	0.52
7:U:412:GLU:OE1	7:U:468:ARG:NH1	2.42	0.52
7:Z:161:PHE:HZ	7:Z:187:LEU:HD13	1.72	0.52
1:E:106:ASP:OD2	1:E:131:ARG:NH2	2.39	0.52
3:M:21:ALA:HB2	4:N:118:TYR:CD1	2.40	0.52
7:e:468:ARG:NE	11:i:67:ASP:OD1	2.30	0.52
4:H:41:VAL:O	4:H:44:GLN:HB2	2.09	0.52
7:U:152:LEU:HD23	7:U:155:LEU:HD12	1.90	0.52
7:U:180:ASP:OD1	7:U:181:LYS:NZ	2.42	0.52
8:f:82:LEU:O	8:f:86:ARG:HG3	2.09	0.52
5:I:-43:DT:O4	6:J:42:DA:N6	2.42	0.52
7:U:178:SER:OG	7:U:180:ASP:OD2	2.28	0.52
3:M:100:VAL:HG22	2:P:96:THR:HB	1.92	0.52
9:g:65:GLN:O	9:g:68:SER:OG	2.21	0.52
1:E:57:SER:OG	2:F:40:ARG:NH2	2.43	0.52
6:J:-49:DG:H2''	6:J:-48:DC:C6	2.44	0.52
3:Q:15:LYS:HB3	6:T:-43:DA:H4'	1.92	0.52
5:I:-71:DT:H3	6:J:71:DA:H61	1.56	0.52
3:M:50:TYR:O	3:M:54:VAL:HG23	2.10	0.52
7:U:277:THR:HG22	11:Y:61:ASN:OD1	2.10	0.52
6:T:-2:DC:H2''	6:T:-1:DA:H8	1.75	0.52
7:Z:392:CYS:HB2	7:Z:406:VAL:HG11	1.91	0.52
2:B:31:LYS:N	2:B:32:PRO:HD2	2.25	0.52
4:D:58:ILE:HG23	2:F:98:TYR:HB3	1.92	0.52
4:H:38:VAL:HG21	4:H:56:MET:HE1	1.91	0.52
7:Z:170:HIS:CD2	7:Z:185:CYS:SG	3.03	0.52
7:e:446:MET:SD	9:g:63:LYS:HA	2.50	0.52
11:n:93:HIS:CD2	11:n:98:CYS:SG	3.03	0.52
1:A:63:ARG:HG2	6:J:17:DA:O3'	2.10	0.51
4:R:41:VAL:HA	4:R:44:GLN:HB2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:I:-38:DC:H2''	5:I:-37:DG:OP2	2.10	0.51
6:T:-65:DT:H2''	6:T:-64:DA:H8	1.75	0.51
7:U:471:ASN:OD1	8:V:27:ASN:HB3	2.10	0.51
3:Q:54:VAL:HG13	4:R:107:ALA:HB1	1.91	0.51
5:S:50:DA:H61	6:T:-50:DT:H3	1.57	0.51
7:j:449:SER:HB2	9:l:50:PHE:HA	1.91	0.51
3:G:26:PRO:HG3	4:H:37:TYR:CZ	2.45	0.51
11:d:18:GLU:O	11:d:18:GLU:HG2	2.11	0.51
7:e:351:ILE:HG21	7:e:354:LEU:HG	1.92	0.51
4:R:44:GLN:O	9:l:81:SER:N	2.42	0.51
5:I:28:DG:H2''	5:I:29:DA:C8	2.46	0.51
6:J:4:DG:H2''	6:J:5:DT:OP2	2.10	0.51
7:U:378:ASP:OD1	7:U:378:ASP:N	2.28	0.51
7:j:368:LEU:HD11	7:j:374:ARG:HD2	1.92	0.51
1:E:72:ARG:HH22	6:J:-23:DT:P	2.34	0.51
4:H:43:LYS:HD2	4:H:47:PRO:HA	1.93	0.51
5:I:22:DC:H2''	5:I:23:DA:C8	2.46	0.51
5:I:65:DA:H2''	5:I:66:DC:OP2	2.11	0.51
1:O:63:ARG:NH1	6:T:-13:DA:OP1	2.43	0.51
7:U:26:ALA:HB1	11:Y:36:MET:HE1	1.92	0.51
6:J:31:DT:H2''	6:J:32:DG:C8	2.45	0.51
1:O:39:HIS:HE1	5:S:-67:DA:H4'	1.76	0.51
3:Q:23:LEU:HD11	4:R:114:ALA:HB1	1.93	0.51
3:Q:50:TYR:O	3:Q:54:VAL:HG23	2.11	0.51
11:Y:81:CYS:HB3	11:Y:98:CYS:SG	2.50	0.51
11:d:13:PRO:O	11:d:17:GLU:HB2	2.11	0.51
3:M:90:ASP:HB3	3:M:93:LEU:HB2	1.91	0.51
5:S:8:DC:N4	6:T:-8:DG:H1	2.09	0.51
5:S:65:DA:H2''	5:S:66:DC:OP2	2.10	0.51
8:f:15:TYR:CE1	11:i:15:VAL:HG21	2.47	0.51
4:N:59:MET:O	4:N:63:VAL:HG23	2.11	0.50
3:Q:77:ARG:HB2	5:S:58:DG:OP1	2.11	0.50
2:B:96:THR:O	3:G:101:THR:N	2.32	0.50
5:I:-11:DG:H2''	5:I:-10:DC:OP2	2.11	0.50
5:I:37:DC:H42	6:J:-37:DG:H1	1.59	0.50
5:I:58:DG:H2''	5:I:59:DA:C8	2.47	0.50
3:M:71:ARG:NH1	9:g:78:ARG:HE	2.08	0.50
7:U:180:ASP:O	7:U:225:LYS:NZ	2.36	0.50
9:W:65:GLN:O	9:W:68:SER:OG	2.27	0.50
1:A:102:ALA:HB1	1:A:131:ARG:HH22	1.77	0.50
5:I:28:DG:H2''	5:I:29:DA:H8	1.75	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:54:TYR:HB3	2:P:40:ARG:HB2	1.93	0.50
3:Q:32:ARG:HA	3:Q:35:ARG:HH12	1.76	0.50
5:S:-7:DG:H1	6:T:7:DC:H42	1.59	0.50
7:e:175:LYS:NZ	7:e:257:LEU:O	2.36	0.50
1:E:40:ARG:NH2	5:I:9:DG:N3	2.58	0.50
5:S:22:DC:H2''	5:S:23:DA:C8	2.47	0.50
7:j:42:ASN:HB2	11:n:35:LEU:HD11	1.94	0.50
7:j:88:ASN:O	7:j:92:GLY:N	2.44	0.50
4:D:34:TYR:CZ	4:D:63:VAL:HG11	2.47	0.50
4:D:59:MET:O	4:D:63:VAL:HG23	2.12	0.50
4:H:46:HIS:CE1	9:b:78:ARG:HE	2.29	0.50
5:S:-49:DG:H1	6:T:49:DC:N4	2.09	0.50
7:j:178:SER:OG	7:j:180:ASP:OD2	2.29	0.50
11:n:16:ILE:HA	11:n:20:SER:HA	1.92	0.50
3:Q:87:VAL:HG11	3:Q:97:LEU:HD12	1.93	0.50
5:S:-38:DC:H2''	5:S:-37:DG:OP2	2.12	0.50
7:Z:159:PRO:HG3	11:d:68:PRO:HB2	1.93	0.50
7:j:49:CYS:HB3	7:j:73:HIS:HE1	1.77	0.50
5:I:-48:DC:H2''	5:I:-47:DT:C6	2.47	0.50
6:T:-2:DC:H2''	6:T:-1:DA:C8	2.46	0.50
7:e:235:GLN:HB3	10:h:74:ARG:HB3	1.93	0.50
9:l:65:GLN:O	9:l:68:SER:OG	2.19	0.50
4:D:79:HIS:NE2	3:G:38:ASN:HB2	2.27	0.50
2:L:92:ARG:HH21	4:N:98:LEU:HD12	1.76	0.50
2:P:34:ILE:HA	2:P:37:LEU:HD12	1.93	0.50
7:Z:191:VAL:HG12	11:d:84:PRO:HG2	1.94	0.50
9:l:76:CYS:SG	9:l:78:ARG:HB2	2.51	0.50
3:Q:68:ASN:HA	9:l:78:ARG:HH21	1.77	0.50
6:T:-58:DC:H2''	6:T:-57:DT:OP2	2.12	0.50
4:N:112:THR:O	4:N:116:THR:OG1	2.30	0.49
6:T:4:DG:H2''	6:T:5:DT:OP2	2.11	0.49
11:d:91:VAL:O	11:d:95:GLU:HG3	2.12	0.49
2:L:98:TYR:OH	4:R:65:ASP:OD2	2.28	0.49
7:e:41:LEU:HG	8:f:17:VAL:HG11	1.94	0.49
2:B:75:HIS:HD2	4:D:93:THR:OG1	1.95	0.49
4:H:53:SER:HB2	6:J:-54:DC:OP2	2.13	0.49
3:M:100:VAL:HA	2:P:96:THR:O	2.12	0.49
4:R:29:THR:O	5:S:50:DA:H5''	2.13	0.49
11:d:85:LEU:HD11	11:d:93:HIS:CD2	2.47	0.49
7:U:93:LEU:HD13	9:W:33:THR:HG21	1.94	0.49
7:Z:180:ASP:OD1	7:Z:181:LYS:NZ	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:U:417:VAL:HB	7:U:462:LEU:HB2	1.95	0.49
7:Z:412:GLU:OE1	7:Z:468:ARG:NH1	2.46	0.49
11:n:91:VAL:O	11:n:95:GLU:HG3	2.12	0.49
3:C:68:ASN:CG	9:W:78:ARG:HH22	2.21	0.49
4:D:104:ALA:O	4:D:108:VAL:HG23	2.13	0.49
6:J:-38:DA:H2''	6:J:-37:DG:OP2	2.12	0.49
3:M:77:ARG:HD2	5:S:-54:DA:H4'	1.94	0.49
7:e:170:HIS:O	7:e:174:CYS:N	2.46	0.49
4:H:52:SER:HA	6:J:-54:DC:H5''	1.95	0.49
5:I:53:DT:H2''	5:I:54:DG:C8	2.47	0.49
7:e:363:LYS:NZ	10:h:7:THR:O	2.45	0.49
8:k:46:THR:HG22	8:k:50:MET:HE2	1.95	0.49
2:F:75:HIS:O	4:H:89:ARG:NH1	2.42	0.49
6:J:-23:DT:H2''	6:J:-22:DG:H5'	1.94	0.49
7:j:85:PHE:HB3	7:j:94:LEU:HD11	1.95	0.49
5:I:58:DG:H1	6:J:-58:DC:H42	1.59	0.48
5:S:-34:DG:H1	6:T:34:DC:N4	2.11	0.48
6:T:-38:DA:H2''	6:T:-37:DG:OP2	2.13	0.48
7:e:324:HIS:ND1	7:e:349:LEU:O	2.46	0.48
8:f:86:ARG:HG2	9:g:8:ILE:HG22	1.94	0.48
7:j:240:GLU:OE1	10:m:72:ARG:NE	2.42	0.48
6:J:-65:DT:H2''	6:J:-64:DA:H8	1.79	0.48
3:M:68:ASN:OD1	9:g:78:ARG:NH2	2.39	0.48
9:W:6:ILE:O	9:W:6:ILE:HG13	2.13	0.48
5:S:-20:DC:H42	6:T:20:DG:H1	1.59	0.48
11:Y:18:GLU:O	11:Y:18:GLU:HG2	2.13	0.48
7:j:332:PHE:CE2	7:j:334:TYR:HB2	2.48	0.48
6:J:43:DA:H2''	6:J:44:DT:OP2	2.13	0.48
8:a:61:ILE:O	8:a:65:VAL:HG23	2.14	0.48
4:D:28:LYS:HG3	6:J:51:DG:OP1	2.14	0.48
2:L:96:THR:O	3:Q:101:THR:N	2.42	0.48
1:O:39:HIS:CE1	5:S:-67:DA:H4'	2.49	0.48
11:i:18:GLU:O	11:i:18:GLU:HG2	2.12	0.48
7:j:108:LEU:HD13	7:j:118:TRP:CG	2.49	0.48
1:A:119:ILE:O	2:B:47:SER:HB3	2.14	0.48
3:G:24:GLN:NE2	4:H:40:LYS:HD3	2.28	0.48
4:H:98:LEU:HB3	4:H:99:LEU:HD23	1.94	0.48
6:J:50:DG:H1'	6:J:51:DG:H5'	1.94	0.48
7:U:161:PHE:HZ	7:U:187:LEU:HD13	1.79	0.48
7:e:142:MET:SD	9:g:82:ALA:HB1	2.54	0.48
2:F:88:TYR:CE2	4:H:80:TYR:HB3	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:S:-30:DG:H2''	5:S:-29:DC:OP2	2.13	0.48
7:Z:49:CYS:HB3	7:Z:73:HIS:HE1	1.79	0.48
7:j:174:CYS:HB3	7:j:177:ARG:HG3	1.95	0.48
2:L:96:THR:HB	3:Q:100:VAL:HG22	1.95	0.48
11:Y:85:LEU:HD11	11:Y:93:HIS:CD2	2.48	0.48
7:e:47:GLY:O	11:i:27:PRO:HA	2.14	0.48
7:e:165:SER:OG	7:e:188:ASP:OD1	2.23	0.48
7:j:224:TRP:HA	7:j:228:GLN:HA	1.95	0.48
7:j:417:VAL:HB	7:j:462:LEU:HB2	1.96	0.48
5:I:-20:DC:H2''	5:I:-19:DG:C8	2.49	0.48
6:J:5:DT:H2''	6:J:6:DA:C8	2.49	0.48
6:J:55:DC:OP2	6:J:55:DC:H2'	2.14	0.48
6:T:-35:DG:H2''	6:T:-34:DA:C8	2.49	0.48
6:T:61:DC:H2''	6:T:62:DG:N7	2.28	0.48
7:e:333:ASN:H	10:h:2:GLN:HE22	1.62	0.48
4:D:62:PHE:HA	2:F:98:TYR:CZ	2.49	0.47
5:S:-57:DC:H2''	5:S:-56:DC:C5	2.48	0.47
7:U:85:PHE:HB3	7:U:94:LEU:HD11	1.96	0.47
7:j:159:PRO:HG3	11:n:68:PRO:HB2	1.96	0.47
3:M:80:PRO:HG3	4:N:58:ILE:HD12	1.96	0.47
5:S:69:DC:N4	6:T:-70:DC:H42	2.11	0.47
7:Z:168:GLN:O	7:Z:172:ASN:N	2.41	0.47
1:A:79:LYS:HB3	1:A:82:LEU:HD11	1.96	0.47
6:J:-53:DA:H1'	6:J:-52:DC:H5'	1.96	0.47
1:K:47:ALA:O	1:K:51:ILE:HG13	2.14	0.47
7:Z:93:LEU:HD13	9:b:33:THR:HG21	1.96	0.47
8:a:15:TYR:CZ	8:a:92:ILE:HG21	2.48	0.47
7:e:373:ASN:ND2	10:h:40:GLN:OE1	2.47	0.47
1:A:118:THR:HA	2:B:45:ARG:HB3	1.96	0.47
5:S:30:DT:H2''	5:S:31:DT:OP2	2.15	0.47
10:X:45:PHE:HB2	10:X:67:LEU:HD22	1.95	0.47
7:Z:199:ASN:HA	7:Z:200:THR:HA	1.58	0.47
7:e:151:ILE:O	7:e:155:LEU:HG	2.14	0.47
8:f:47:LYS:NZ	9:g:31:GLU:OE1	2.24	0.47
1:A:76:GLN:C	1:A:78:PHE:H	2.23	0.47
3:C:50:TYR:O	3:C:54:VAL:HG23	2.14	0.47
5:I:-63:DC:H2''	5:I:-62:DC:OP2	2.15	0.47
1:K:76:GLN:C	1:K:78:PHE:H	2.22	0.47
4:N:30:ARG:NH2	6:T:48:DG:H21	2.13	0.47
8:V:46:THR:HG22	8:V:50:MET:HE2	1.96	0.47
2:B:28:GLY:O	2:B:30:THR:N	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:I:3:DT:H2''	5:I:4:DC:C6	2.49	0.47
3:M:42:ARG:HG3	4:N:85:THR:HG23	1.97	0.47
1:O:44:GLY:HA3	2:P:44:LYS:HE2	1.97	0.47
2:P:28:GLY:O	2:P:30:THR:N	2.46	0.47
2:P:95:ARG:HB3	2:P:95:ARG:CZ	2.42	0.47
7:U:11:PHE:HE2	7:U:108:LEU:HD23	1.80	0.47
7:U:111:ASP:HB3	11:Y:87:LEU:HD11	1.96	0.47
7:Z:199:ASN:ND2	11:d:80:LYS:O	2.34	0.47
7:Z:271:CYS:HB3	7:Z:276:HIS:ND1	2.29	0.47
7:e:27:ALA:O	7:e:31:LEU:HG	2.15	0.47
2:B:64:ASN:HA	2:B:67:ARG:NH1	2.29	0.47
5:I:-30:DG:H2''	5:I:-29:DC:OP2	2.13	0.47
6:J:61:DC:H2''	6:J:62:DG:N7	2.29	0.47
1:K:51:ILE:O	1:K:55:GLN:HG3	2.15	0.47
1:O:42:ARG:HH21	6:T:-5:DG:P	2.38	0.47
2:P:62:LEU:O	2:P:66:ILE:HG13	2.14	0.47
4:R:84:SER:N	6:T:-34:DA:OP1	2.40	0.47
5:S:-71:DT:H3	6:T:71:DA:H61	1.61	0.47
7:U:75:LEU:O	7:U:78:SER:OG	2.28	0.47
7:j:41:LEU:HD11	8:k:14:GLN:OE1	2.14	0.47
1:A:43:PRO:HA	6:J:9:DT:H5'	1.97	0.47
6:J:-61:DT:H2'	6:J:-60:DA:C8	2.50	0.47
5:S:39:DA:H61	6:T:-39:DT:H3	1.63	0.47
6:T:43:DA:H2''	6:T:44:DT:OP2	2.15	0.47
7:Z:395:LYS:O	11:d:56:LYS:NZ	2.47	0.47
11:d:47:ASP:OD2	11:d:48:HIS:ND1	2.45	0.47
7:e:180:ASP:O	7:e:225:LYS:NZ	2.42	0.47
1:E:79:LYS:HB3	1:E:82:LEU:HD11	1.95	0.47
5:I:-65:DT:H3	6:J:65:DA:H61	1.63	0.47
6:J:31:DT:H2''	6:J:32:DG:H8	1.80	0.47
4:N:46:HIS:HB3	4:N:49:THR:OG1	2.14	0.47
2:B:94:GLY:O	3:G:99:ARG:HB3	2.15	0.47
5:I:-49:DG:H1	6:J:49:DC:H42	1.63	0.47
5:S:3:DT:H2''	5:S:4:DC:C6	2.50	0.47
5:S:10:DC:H2''	5:S:11:DG:C8	2.50	0.47
9:W:51:ASP:HB2	9:W:56:LEU:HB2	1.96	0.47
5:I:8:DC:N4	6:J:-8:DG:H1	2.13	0.46
6:J:-57:DT:H2''	6:J:-56:DG:C8	2.50	0.46
3:M:20:ARG:NH2	5:S:-42:DT:OP1	2.32	0.46
4:N:65:ASP:HB2	2:P:102:GLY:HA3	1.97	0.46
7:U:62:GLN:HB2	7:U:83:HIS:CD2	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:U:192:HIS:CE1	7:U:198:LEU:HD23	2.50	0.46
8:f:92:ILE:HA	11:i:12:LYS:HD2	1.97	0.46
7:j:151:ILE:O	7:j:155:LEU:HG	2.15	0.46
3:M:115:LEU:HD11	1:O:112:ILE:HD11	1.96	0.46
5:S:-63:DC:H1'	5:S:-62:DC:O5'	2.15	0.46
5:S:-49:DG:H1	6:T:49:DC:H42	1.62	0.46
6:T:-51:DG:H2''	6:T:-50:DT:OP2	2.16	0.46
7:U:332:PHE:CE2	7:U:334:TYR:HB2	2.50	0.46
8:V:32:ARG:NH2	8:V:80:THR:HG21	2.31	0.46
7:Z:167:SER:HA	11:d:89:ALA:HB2	1.98	0.46
7:Z:353:LYS:HA	11:d:57:TYR:O	2.16	0.46
9:l:6:ILE:HG13	9:l:6:ILE:O	2.14	0.46
5:S:53:DT:H2''	5:S:54:DG:C8	2.51	0.46
5:S:64:DT:H3	6:T:-64:DA:N6	2.12	0.46
7:Z:311:LYS:HB3	7:Z:311:LYS:HE2	1.79	0.46
3:C:68:ASN:OD1	9:W:78:ARG:NH2	2.48	0.46
3:C:118:LYS:HE3	3:C:118:LYS:HA	1.97	0.46
6:J:-58:DC:H2''	6:J:-57:DT:OP2	2.16	0.46
1:O:76:GLN:C	1:O:78:PHE:H	2.24	0.46
6:T:-23:DT:H2''	6:T:-22:DG:H5'	1.96	0.46
7:e:161:PHE:CZ	7:e:187:LEU:HD13	2.49	0.46
7:e:284:GLU:HB2	7:e:352:HIS:HB2	1.98	0.46
10:c:42:ARG:HD3	10:c:72:ARG:CG	2.44	0.46
7:e:395:LYS:HD3	7:e:404:GLY:O	2.15	0.46
3:G:30:VAL:O	3:G:34:LEU:HD12	2.16	0.46
3:G:69:ALA:HB3	3:G:82:HIS:HB3	1.97	0.46
6:J:-9:DC:H2''	6:J:-8:DG:C8	2.51	0.46
4:R:54:LYS:O	4:R:58:ILE:HG13	2.16	0.46
7:U:428:TYR:OH	10:X:73:LEU:O	2.32	0.46
9:b:72:HIS:HD1	9:b:79:ASP:CG	2.23	0.46
7:j:38:GLU:HB3	11:n:35:LEU:HD22	1.97	0.46
3:G:26:PRO:HB3	5:I:48:DG:OP1	2.15	0.46
1:O:99:TYR:HD1	2:P:95:ARG:CZ	2.29	0.46
3:Q:24:GLN:HG3	4:R:44:GLN:HE22	1.80	0.46
3:C:53:ALA:HB2	4:D:114:ALA:HB3	1.98	0.46
7:Z:108:LEU:HD13	7:Z:118:TRP:CD1	2.50	0.46
11:i:91:VAL:O	11:i:95:GLU:HG3	2.16	0.46
1:A:59:GLU:H	1:A:59:GLU:HG3	1.56	0.46
1:E:73:GLU:OE2	2:F:25:ASN:HB3	2.16	0.46
1:K:100:LEU:HD11	2:L:58:LEU:HD13	1.97	0.46
3:Q:32:ARG:HA	3:Q:35:ARG:NH1	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:T:11:DC:H2''	6:T:12:DG:C8	2.51	0.46
7:U:163:ARG:O	7:U:167:SER:OG	2.29	0.46
7:Z:170:HIS:O	7:Z:174:CYS:N	2.48	0.46
1:A:121:PRO:HB3	2:B:53:GLU:CD	2.41	0.45
3:C:17:ARG:NH2	5:I:-43:DT:OP2	2.49	0.45
4:D:99:LEU:HA	4:D:100:PRO:HD3	1.74	0.45
7:j:450:ILE:HA	9:l:50:PHE:HE1	1.82	0.45
9:l:66:GLU:HG3	9:l:83:ASN:HB3	1.98	0.45
1:A:97:GLU:O	1:A:101:VAL:HG23	2.15	0.45
3:C:69:ALA:HB3	3:C:82:HIS:HB3	1.97	0.45
3:G:32:ARG:HA	3:G:35:ARG:NH1	2.32	0.45
1:K:66:PRO:HB3	2:L:28:GLY:HA3	1.99	0.45
2:P:87:VAL:HG22	2:P:97:LEU:HD23	1.97	0.45
3:Q:32:ARG:HH11	6:T:-45:DG:H3'	1.81	0.45
4:R:46:HIS:HB3	4:R:49:THR:OG1	2.16	0.45
3:Q:31:HIS:HA	3:Q:34:LEU:HD12	1.98	0.45
6:T:66:DT:H2''	6:T:67:DT:C6	2.51	0.45
9:b:75:ASN:O	9:b:77:GLY:N	2.49	0.45
5:I:-70:DC:H2'	5:I:-69:DA:C8	2.52	0.45
6:J:-61:DT:H2''	6:J:-60:DA:O4'	2.16	0.45
2:L:71:THR:HG23	4:N:96:ARG:NH1	2.32	0.45
4:N:39:TYR:OH	5:S:-53:DG:H3'	2.16	0.45
6:T:31:DT:H2''	6:T:32:DG:H8	1.80	0.45
8:f:15:TYR:OH	8:f:92:ILE:HG12	2.17	0.45
1:A:54:TYR:HB3	2:B:40:ARG:HB2	1.97	0.45
2:L:102:GLY:OXT	4:R:64:ASN:HB2	2.17	0.45
6:T:-8:DG:H2''	6:T:-7:DG:C8	2.51	0.45
7:U:408:ASP:HB2	7:U:471:ASN:ND2	2.32	0.45
7:j:170:HIS:O	7:j:174:CYS:N	2.49	0.45
4:D:60:ASN:OD1	4:D:64:ASN:ND2	2.45	0.45
5:I:23:DA:H5''	6:T:35:DT:C7	2.47	0.45
6:J:-60:DA:H2''	6:J:-59:DT:OP2	2.16	0.45
1:K:104:PHE:HE2	2:L:37:LEU:HB3	1.82	0.45
7:U:199:ASN:HA	7:U:200:THR:HA	1.57	0.45
7:U:288:VAL:HG22	7:U:296:SER:HB2	1.98	0.45
7:Z:132:ARG:HD3	7:Z:194:LEU:O	2.17	0.45
7:Z:174:CYS:SG	7:Z:176:VAL:HG13	2.56	0.45
9:g:9:ASP:O	9:g:13:ASN:ND2	2.49	0.45
11:i:15:VAL:HG23	11:i:19:TYR:HD2	1.81	0.45
7:j:244:PHE:O	7:j:248:GLN:HG2	2.16	0.45
3:G:32:ARG:HH22	4:H:32:GLU:CD	2.25	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:76:THR:HG21	5:I:57:DA:H5''	1.98	0.45
9:b:80:VAL:HG21	9:b:88:HIS:CG	2.52	0.45
4:R:89:ARG:O	4:R:93:THR:HG22	2.16	0.45
5:S:-63:DC:H2''	5:S:-62:DC:OP2	2.15	0.45
5:S:58:DG:H1	6:T:-58:DC:N4	2.12	0.45
9:b:65:GLN:O	9:b:68:SER:OG	2.34	0.45
9:g:76:CYS:SG	9:g:78:ARG:HB2	2.56	0.45
7:j:311:LYS:HD2	7:j:366:GLU:OE2	2.16	0.45
3:C:61:GLU:OE1	9:W:91:ARG:NH2	2.50	0.45
3:C:112:GLN:HG3	1:E:112:ILE:HD13	1.99	0.45
4:D:33:SER:HA	4:D:60:ASN:ND2	2.32	0.45
1:K:134:ARG:HH12	3:M:98:GLY:HA2	1.82	0.45
5:S:23:DA:H2''	5:S:24:DA:OP2	2.16	0.45
7:e:145:THR:HB	7:e:216:PHE:HZ	1.81	0.45
1:A:63:ARG:NH1	5:I:-13:DA:OP1	2.50	0.44
1:A:134:ARG:NH1	3:C:95:LYS:HA	2.32	0.44
2:L:64:ASN:HA	2:L:67:ARG:NH1	2.32	0.44
1:O:47:ALA:O	1:O:51:ILE:HG13	2.17	0.44
7:j:292:CYS:SG	7:j:338:GLU:HB2	2.57	0.44
2:F:30:THR:HB	2:F:32:PRO:HD2	1.99	0.44
7:U:21:LEU:HD11	7:U:108:LEU:HD21	2.00	0.44
7:U:131:GLU:CD	7:U:131:GLU:H	2.26	0.44
4:H:102:GLU:HG3	4:H:106:HIS:CE1	2.53	0.44
2:L:75:HIS:CE1	4:N:89:ARG:HG2	2.52	0.44
5:S:-49:DG:H2''	5:S:-48:DC:C5	2.51	0.44
7:e:46:CYS:HB3	7:e:49:CYS:SG	2.56	0.44
10:h:24:GLU:CD	10:h:53:GLY:H	2.25	0.44
5:I:-54:DA:H61	6:J:53:DC:N4	2.16	0.44
5:I:-20:DC:H42	6:J:20:DG:H1	1.66	0.44
2:P:64:ASN:HA	2:P:67:ARG:NH1	2.32	0.44
5:S:31:DT:H2''	5:S:32:DA:C8	2.53	0.44
6:T:71:DA:H1'	6:T:72:DT:H5'	1.98	0.44
7:Z:46:CYS:HB2	7:Z:68:CYS:HB3	2.00	0.44
1:A:41:TYR:OH	6:J:-66:DG:H5'	2.17	0.44
2:F:71:THR:HG22	4:H:93:THR:OG1	2.17	0.44
6:J:-63:DT:H2''	6:J:-62:DA:H8	1.81	0.44
6:J:-51:DG:H2''	6:J:-50:DT:OP2	2.17	0.44
6:J:-27:DC:H2''	6:J:-26:DC:C5	2.52	0.44
3:M:87:VAL:HG13	3:M:93:LEU:HB3	2.00	0.44
5:S:-16:DT:H4'	5:S:-15:DA:H5'	1.99	0.44
7:U:49:CYS:HB3	7:U:73:HIS:HE1	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:V:25:ILE:HD11	8:V:88:PHE:HD2	1.83	0.44
1:A:47:ALA:O	1:A:51:ILE:HG13	2.18	0.44
2:B:62:LEU:O	2:B:66:ILE:HG13	2.18	0.44
4:H:44:GLN:HE21	9:b:67:SER:CB	2.31	0.44
6:J:-31:DA:H8	6:J:-31:DA:OP2	2.00	0.44
4:R:75:SER:O	4:R:79:HIS:ND1	2.47	0.44
6:T:-15:DA:H2''	6:T:-14:DA:H8	1.83	0.44
6:T:64:DG:H1'	6:T:65:DA:C8	2.52	0.44
7:U:363:LYS:NZ	10:X:7:THR:O	2.51	0.44
7:e:235:GLN:CB	10:h:74:ARG:HB3	2.48	0.44
7:j:135:GLY:HA2	7:j:194:LEU:O	2.17	0.44
7:j:351:ILE:HG21	7:j:354:LEU:HG	1.99	0.44
6:J:54:DT:H2''	6:J:55:DC:OP2	2.18	0.44
6:J:65:DA:OP2	6:J:65:DA:H2'	2.17	0.44
4:N:104:ALA:O	4:N:108:VAL:HG23	2.17	0.44
4:R:35:ALA:HA	4:R:56:MET:SD	2.57	0.44
5:S:-62:DC:H2''	5:S:-61:DG:C8	2.52	0.44
7:Z:135:GLY:HA2	7:Z:194:LEU:O	2.17	0.44
7:j:199:ASN:HA	7:j:200:THR:HA	1.62	0.44
1:A:128:ARG:HB3	1:A:133:GLU:HB2	1.99	0.44
4:D:31:LYS:HD2	4:D:31:LYS:O	2.17	0.44
5:I:-16:DT:H4'	5:I:-15:DA:H5'	1.99	0.44
6:T:-9:DC:H2''	6:T:-8:DG:OP2	2.18	0.44
7:U:30:ILE:HD11	11:Y:36:MET:HE2	1.99	0.44
11:Y:36:MET:O	11:Y:40:LYS:HG3	2.18	0.44
11:Y:91:VAL:O	11:Y:95:GLU:HG3	2.17	0.44
7:e:290:PRO:HD2	7:e:343:GLN:HB3	2.00	0.44
7:e:311:LYS:HB3	7:e:311:LYS:HE2	1.79	0.44
7:j:228:GLN:H	7:j:228:GLN:HG3	1.65	0.44
7:j:429:ILE:HG21	7:j:441:LYS:HE3	1.99	0.44
4:R:47:PRO:CG	9:l:79:ASP:HB2	2.38	0.44
7:Z:153:GLN:HA	7:Z:156:ILE:HB	1.99	0.44
7:j:37:LYS:HD3	8:k:18:GLU:HA	1.99	0.44
10:m:37:PRO:HB2	10:m:39:ASP:OD1	2.18	0.44
1:O:51:ILE:O	1:O:55:GLN:HG3	2.18	0.43
1:O:119:ILE:HG13	2:P:50:ILE:HG13	2.00	0.43
4:R:104:ALA:O	4:R:108:VAL:HG23	2.18	0.43
6:T:-70:DC:H2''	6:T:-69:DG:OP2	2.17	0.43
7:U:429:ILE:HG21	7:U:441:LYS:HE3	2.00	0.43
11:Y:37:SER:O	11:Y:40:LYS:HB2	2.18	0.43
7:Z:411:TYR:HB3	7:Z:465:TYR:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:e:16:SER:O	7:e:20:VAL:HG23	2.18	0.43
2:F:34:ILE:HA	2:F:37:LEU:HD12	1.99	0.43
2:P:30:THR:HB	2:P:32:PRO:HD2	2.00	0.43
5:S:-53:DG:H1'	5:S:-52:DG:H5'	2.00	0.43
6:T:50:DG:H1'	6:T:51:DG:H5'	2.00	0.43
11:Y:83:LYS:HD2	11:Y:93:HIS:CE1	2.53	0.43
9:b:71:ILE:HG12	9:b:82:ALA:HB2	1.99	0.43
1:A:39:HIS:ND1	6:J:-67:DT:H5''	2.32	0.43
2:B:68:ASP:OD2	2:B:92:ARG:HD3	2.19	0.43
4:D:83:ARG:HE	5:I:-34:DG:H3'	1.83	0.43
6:T:16:DA:H1'	6:T:17:DA:C8	2.52	0.43
7:e:138:GLY:HA3	7:e:445:SER:O	2.17	0.43
7:j:54:SER:HB3	7:j:68:CYS:HB2	2.00	0.43
8:k:38:TRP:O	8:k:42:VAL:HG23	2.18	0.43
1:E:40:ARG:HH22	6:J:-8:DG:H21	1.65	0.43
11:d:90:ILE:O	11:d:94:LEU:HG	2.18	0.43
8:f:86:ARG:HE	9:g:9:ASP:HA	1.82	0.43
4:D:54:LYS:HA	4:D:54:LYS:HD3	1.67	0.43
4:N:89:ARG:O	4:N:93:THR:HG22	2.18	0.43
7:U:235:GLN:HB3	10:X:74:ARG:HB3	1.99	0.43
9:b:34:GLN:NE2	11:d:70:ASP:O	2.47	0.43
7:j:469:GLN:HG3	11:n:48:HIS:HB2	2.01	0.43
1:E:102:ALA:O	1:E:105:GLU:HB2	2.18	0.43
4:N:103:LEU:HD12	4:N:103:LEU:HA	1.84	0.43
6:T:-4:DG:H2''	6:T:-3:DA:C8	2.53	0.43
6:T:69:DT:H6	6:T:69:DT:H2'	1.63	0.43
7:U:471:ASN:CG	8:V:27:ASN:HB3	2.43	0.43
10:h:42:ARG:HD3	10:h:72:ARG:CG	2.48	0.43
7:j:16:SER:O	7:j:20:VAL:HG23	2.18	0.43
1:A:40:ARG:HH12	6:J:9:DT:H1'	1.83	0.43
1:K:119:ILE:O	2:L:47:SER:HB3	2.19	0.43
5:S:-45:DA:H2''	5:S:-44:DA:C8	2.54	0.43
5:S:54:DG:H8	5:S:54:DG:OP2	2.01	0.43
5:S:62:DT:H3	6:T:-62:DA:N6	2.16	0.43
8:a:82:LEU:HG	9:b:16:LEU:HD22	2.00	0.43
7:e:144:SER:HB3	10:h:76:GLY:HA3	2.01	0.43
9:g:38:LEU:HD23	9:g:38:LEU:HA	1.88	0.43
11:i:36:MET:O	11:i:40:LYS:HG3	2.18	0.43
3:C:66:ALA:HB2	3:C:83:LEU:HD23	2.01	0.43
5:S:31:DT:H2''	5:S:32:DA:H8	1.83	0.43
7:Z:284:GLU:HG3	7:Z:298:THR:HG23	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:ARG:NH2	5:I:-5:DA:OP1	2.48	0.43
1:A:51:ILE:O	1:A:55:GLN:HG3	2.18	0.43
4:D:54:LYS:O	4:D:58:ILE:HG13	2.19	0.43
4:R:62:PHE:O	4:R:66:VAL:HG23	2.19	0.43
5:S:67:DA:H2''	5:S:68:DT:OP2	2.18	0.43
5:S:69:DC:H42	6:T:-70:DC:N4	2.16	0.43
8:k:37:GLY:O	8:k:41:LYS:HG3	2.19	0.43
3:C:35:ARG:NH2	6:J:39:DA:OP2	2.52	0.43
4:D:69:ARG:HB3	4:D:98:LEU:HD11	2.00	0.43
4:N:30:ARG:HH22	6:T:48:DG:H21	1.67	0.43
9:b:73:CYS:HB3	9:b:78:ARG:O	2.18	0.43
11:d:71:LYS:HB2	11:d:72:PRO:HD2	2.01	0.43
7:e:49:CYS:HB3	7:e:73:HIS:HE1	1.83	0.43
11:i:14:LYS:HA	11:i:17:GLU:HB2	2.01	0.43
1:A:64:LYS:HD2	1:A:90:MET:HE1	2.01	0.42
2:B:29:ILE:HG22	2:B:34:ILE:HD11	2.00	0.42
4:D:55:ALA:HA	4:D:58:ILE:HD12	2.02	0.42
4:D:92:GLN:O	4:D:95:VAL:HB	2.19	0.42
1:E:128:ARG:HH21	1:E:134:ARG:HE	1.67	0.42
6:J:4:DG:C8	6:J:5:DT:H72	2.54	0.42
3:Q:83:LEU:HD11	4:R:55:ALA:HB1	1.99	0.42
7:U:1:MET:HE3	9:W:39:LYS:NZ	2.34	0.42
7:Z:131:GLU:CD	7:Z:131:GLU:H	2.27	0.42
8:a:12:ILE:HG13	11:d:11:ILE:HD11	2.01	0.42
7:e:332:PHE:CE2	7:e:334:TYR:HB2	2.54	0.42
8:f:66:GLU:HB3	8:f:67:PRO:HD3	1.99	0.42
9:g:51:ASP:HA	9:g:52:PRO:HD2	1.86	0.42
11:i:42:THR:HG22	11:i:43:PRO:O	2.18	0.42
8:k:47:LYS:HA	8:k:50:MET:HE3	2.01	0.42
2:B:39:ARG:NH1	2:B:43:VAL:O	2.52	0.42
4:H:37:TYR:O	4:H:41:VAL:HG23	2.19	0.42
5:I:-24:DG:H1	6:J:24:DC:H42	1.65	0.42
5:I:64:DT:H3	6:J:-64:DA:N6	2.14	0.42
2:L:30:THR:HB	2:L:32:PRO:HD2	2.01	0.42
5:S:-48:DC:H42	6:T:47:DA:H62	1.67	0.42
5:S:58:DG:H2''	5:S:59:DA:C8	2.54	0.42
6:T:-65:DT:H2''	6:T:-64:DA:C8	2.54	0.42
10:c:37:PRO:HA	10:c:38:PRO:HD3	1.88	0.42
7:e:174:CYS:SG	7:e:176:VAL:HG13	2.59	0.42
7:j:428:TYR:OH	10:m:73:LEU:O	2.31	0.42
9:l:80:VAL:HB	9:l:85:LEU:HD23	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:41:VAL:HA	4:D:44:GLN:HB2	2.00	0.42
3:G:93:LEU:HD23	3:G:93:LEU:HA	1.87	0.42
5:I:-68:DG:H2''	5:I:-67:DA:H5''	2.01	0.42
6:J:39:DA:H1'	6:J:40:DC:H5'	2.01	0.42
5:S:-69:DA:H2''	5:S:-68:DG:OP2	2.18	0.42
7:j:26:ALA:HB1	11:n:36:MET:CE	2.46	0.42
1:E:48:LEU:O	1:E:52:ARG:HG3	2.18	0.42
3:G:61:GLU:O	3:G:64:GLU:HB3	2.19	0.42
5:I:-54:DA:N6	6:J:53:DC:H42	2.16	0.42
1:K:130:ILE:HG12	1:O:106:ASP:HB3	2.00	0.42
5:S:-59:DT:H3	6:T:59:DA:N6	2.14	0.42
8:V:66:GLU:HB3	8:V:67:PRO:HD3	2.02	0.42
7:j:179:PRO:HG3	7:j:189:LYS:HE2	2.00	0.42
4:D:72:GLY:O	4:D:75:SER:OG	2.27	0.42
3:G:79:ILE:HG12	3:G:82:HIS:CE1	2.54	0.42
1:K:72:ARG:HH22	5:S:-23:DC:P	2.42	0.42
3:M:54:VAL:HG13	4:N:107:ALA:HB1	2.01	0.42
11:i:36:MET:HG3	11:i:40:LYS:HE3	2.02	0.42
1:E:51:ILE:O	1:E:55:GLN:HG3	2.19	0.42
5:S:42:DC:H42	6:T:-42:DG:H1	1.67	0.42
8:V:13:GLN:O	8:V:17:VAL:HG23	2.19	0.42
9:b:75:ASN:C	9:b:77:GLY:H	2.28	0.42
3:G:17:ARG:HA	3:G:20:ARG:HD2	2.02	0.42
6:J:-70:DC:H2''	6:J:-69:DG:OP2	2.14	0.42
7:U:85:PHE:HA	7:U:95:PHE:O	2.19	0.42
10:c:63:LYS:HE3	10:c:64:GLU:OE2	2.19	0.42
11:d:93:HIS:HA	11:d:97:HIS:HB2	2.01	0.42
7:e:395:LYS:O	11:i:56:LYS:NZ	2.52	0.42
7:e:417:VAL:HB	7:e:462:LEU:HB2	2.01	0.42
8:f:67:PRO:O	8:f:71:GLU:HG3	2.19	0.42
4:H:39:TYR:HE1	4:H:43:LYS:HE3	1.82	0.42
5:I:30:DT:H2''	5:I:31:DT:OP2	2.18	0.42
5:I:67:DA:N3	6:J:-66:DG:N2	2.68	0.42
1:K:128:ARG:HB3	1:K:133:GLU:HB2	2.01	0.42
7:U:16:SER:O	7:U:20:VAL:HG23	2.20	0.42
7:e:271:CYS:HB3	7:e:276:HIS:ND1	2.34	0.42
11:i:12:LYS:O	11:i:15:VAL:HG12	2.20	0.42
11:i:13:PRO:O	11:i:17:GLU:HB2	2.20	0.42
1:E:63:ARG:HB3	5:I:17:DA:O3'	2.19	0.42
5:I:54:DG:H8	5:I:54:DG:OP2	2.03	0.42
4:N:62:PHE:O	4:N:66:VAL:HG23	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:T:54:DT:H2"	6:T:55:DC:OP2	2.18	0.42
7:U:28:ARG:NH2	7:U:109:ILE:O	2.42	0.42
7:Z:247:ASN:O	7:Z:251:GLN:HG2	2.19	0.42
10:c:6:LYS:NZ	10:c:10:GLY:O	2.48	0.42
7:e:41:LEU:HD21	8:f:17:VAL:HB	2.02	0.42
7:e:58:PHE:HA	7:e:87:ILE:O	2.20	0.42
9:g:49:TYR:HB3	9:g:57:ASP:O	2.20	0.42
11:i:37:SER:O	11:i:40:LYS:HB2	2.20	0.42
4:D:56:MET:HE3	4:D:56:MET:O	2.20	0.42
3:G:16:THR:CG2	5:I:47:DG:H5"	2.50	0.42
4:R:77:LEU:O	4:R:81:ASN:ND2	2.52	0.42
5:S:7:DC:N4	6:T:-7:DG:H1	2.17	0.42
7:Z:85:PHE:HA	7:Z:95:PHE:O	2.20	0.42
7:Z:195:TYR:HB3	11:d:75:TYR:CZ	2.55	0.42
4:D:60:ASN:O	4:D:64:ASN:ND2	2.53	0.41
1:O:42:ARG:NH2	6:T:-5:DG:OP1	2.46	0.41
8:f:58:PHE:CZ	8:f:62:LEU:HD22	2.54	0.41
4:H:103:LEU:HA	4:H:103:LEU:HD12	1.86	0.41
5:I:-2:DC:H1'	5:I:-1:DG:C8	2.55	0.41
3:M:43:VAL:N	6:T:39:DA:OP1	2.50	0.41
3:M:61:GLU:O	3:M:64:GLU:HB3	2.20	0.41
7:Z:161:PHE:CZ	7:Z:187:LEU:HD13	2.54	0.41
10:c:11:LYS:HE2	10:c:11:LYS:HB3	1.94	0.41
7:e:159:PRO:HG3	11:i:68:PRO:HB2	2.02	0.41
8:k:66:GLU:HB3	8:k:67:PRO:HD3	2.00	0.41
9:l:72:HIS:HD2	9:l:79:ASP:OD2	2.03	0.41
1:E:121:PRO:HB3	2:F:53:GLU:HG3	2.02	0.41
6:J:11:DC:H2"	6:J:12:DG:C8	2.55	0.41
4:N:92:GLN:O	4:N:95:VAL:HB	2.20	0.41
6:T:55:DC:H2"	6:T:56:DG:C8	2.55	0.41
7:U:81:ILE:HG22	7:U:83:HIS:CE1	2.55	0.41
7:e:191:VAL:HG12	11:i:84:PRO:HG2	2.03	0.41
7:j:271:CYS:HB3	7:j:276:HIS:ND1	2.34	0.41
9:l:46:ARG:NH2	9:l:58:ILE:O	2.52	0.41
9:l:66:GLU:HA	9:l:68:SER:H	1.86	0.41
4:D:103:LEU:HD12	4:D:103:LEU:HA	1.91	0.41
5:I:67:DA:H2"	5:I:68:DT:OP2	2.19	0.41
5:I:71:DA:H61	6:J:-71:DT:H3	1.66	0.41
2:L:34:ILE:H	2:L:34:ILE:HG13	1.66	0.41
2:P:34:ILE:H	2:P:34:ILE:HG13	1.65	0.41
10:X:42:ARG:HD3	10:X:72:ARG:CG	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:Z:368:LEU:HD12	7:Z:372:SER:OG	2.21	0.41
2:B:102:GLY:OXT	4:H:64:ASN:HB2	2.20	0.41
4:D:75:SER:O	4:D:79:HIS:ND1	2.52	0.41
2:F:95:ARG:HB3	2:F:95:ARG:NH1	2.36	0.41
3:G:43:VAL:HG12	3:G:44:GLY:O	2.21	0.41
4:N:89:ARG:O	4:N:92:GLN:HB3	2.21	0.41
1:O:54:TYR:HD2	2:P:39:ARG:HB2	1.85	0.41
7:e:50:HIS:ND1	8:f:10:SER:HB2	2.36	0.41
7:j:243:GLN:HE22	10:m:44:ILE:HG23	1.86	0.41
8:k:33:LEU:HB3	8:k:38:TRP:CE3	2.55	0.41
1:E:99:TYR:OH	1:E:133:GLU:OE1	2.34	0.41
3:G:108:LEU:HA	3:G:109:PRO:HD3	1.91	0.41
5:I:-69:DA:H2''	5:I:-68:DG:OP2	2.21	0.41
6:J:-37:DG:OP2	6:J:-37:DG:H2'	2.20	0.41
6:J:-31:DA:H2''	6:J:-30:DA:C8	2.56	0.41
5:S:5:DC:H2''	5:S:6:DC:C5	2.56	0.41
6:T:-27:DC:H4'	6:T:-26:DC:H5'	2.02	0.41
7:U:365:PHE:HA	7:U:374:ARG:O	2.21	0.41
11:Y:16:ILE:HA	11:Y:20:SER:HA	2.02	0.41
7:Z:235:GLN:NE2	7:Z:425:GLU:OE2	2.54	0.41
10:h:43:LEU:O	10:h:50:LEU:HG	2.21	0.41
7:j:131:GLU:CD	7:j:131:GLU:H	2.28	0.41
11:n:86:ALA:O	11:n:90:ILE:HG13	2.21	0.41
3:C:50:TYR:O	3:C:53:ALA:HB3	2.21	0.41
2:F:45:ARG:HH22	5:I:6:DC:H1'	1.83	0.41
5:I:-71:DT:H3	6:J:71:DA:N6	2.18	0.41
5:I:-25:DA:H1'	5:I:-24:DG:C8	2.56	0.41
6:J:42:DA:H2''	6:J:43:DA:OP2	2.20	0.41
6:T:34:DC:H2''	6:T:35:DT:OP2	2.20	0.41
6:T:42:DA:H2''	6:T:43:DA:OP2	2.21	0.41
9:W:30:ARG:O	9:W:34:GLN:HG3	2.21	0.41
10:X:63:LYS:HE3	10:X:64:GLU:OE2	2.20	0.41
7:Z:27:ALA:O	7:Z:31:LEU:HG	2.21	0.41
7:Z:397:LYS:NZ	7:Z:397:LYS:HA	2.35	0.41
9:b:38:LEU:HD23	9:b:38:LEU:HA	1.89	0.41
7:e:367:HIS:NE2	10:h:35:GLY:O	2.54	0.41
11:n:71:LYS:HB2	11:n:72:PRO:HD2	2.03	0.41
2:L:77:LYS:HG3	4:N:89:ARG:HH12	1.86	0.41
7:U:277:THR:O	11:Y:62:ALA:HA	2.19	0.41
7:U:368:LEU:HD11	7:U:374:ARG:NE	2.36	0.41
11:d:12:LYS:O	11:d:15:VAL:HG12	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:e:28:ARG:HA	7:e:31:LEU:HD12	2.03	0.41
7:e:199:ASN:HA	7:e:200:THR:HA	1.58	0.41
7:e:397:LYS:HA	7:e:397:LYS:NZ	2.35	0.41
7:j:149:SER:O	7:j:153:GLN:HG2	2.20	0.41
11:n:93:HIS:HA	11:n:97:HIS:HB2	2.03	0.41
1:K:134:ARG:NH1	3:M:98:GLY:HA2	2.36	0.41
3:M:63:LEU:HD22	4:N:42:LEU:HD13	2.03	0.41
3:Q:18:SER:HB3	3:Q:23:LEU:O	2.21	0.41
3:Q:42:ARG:O	4:R:85:THR:HA	2.21	0.41
7:U:81:ILE:HD12	7:U:81:ILE:N	2.35	0.41
7:U:107:ASP:OD1	9:W:41:ARG:NH2	2.36	0.41
7:U:174:CYS:SG	7:U:176:VAL:HG13	2.60	0.41
7:U:332:PHE:CD2	7:U:334:TYR:HB2	2.56	0.41
8:V:38:TRP:O	8:V:42:VAL:HG23	2.21	0.41
11:Y:75:TYR:HB2	11:Y:85:LEU:O	2.21	0.41
7:Z:151:ILE:HD11	7:Z:238:ALA:HB1	2.02	0.41
10:c:1:MET:N	10:c:16:GLU:OE2	2.50	0.41
7:e:147:PHE:HD1	10:h:76:GLY:HA2	1.86	0.41
7:j:47:GLY:HA3	7:j:65:PHE:CE1	2.56	0.41
9:l:51:ASP:HA	9:l:52:PRO:HD2	1.95	0.41
1:A:73:GLU:O	1:A:76:GLN:N	2.54	0.41
5:I:10:DC:H2"	5:I:11:DG:C8	2.56	0.41
2:L:62:LEU:O	2:L:66:ILE:HG13	2.21	0.41
1:O:52:ARG:HB3	1:O:56:LYS:NZ	2.36	0.41
5:S:-68:DG:N2	6:T:69:DT:O2	2.54	0.41
7:U:52:ILE:HG12	8:V:13:GLN:OE1	2.21	0.41
7:Z:21:LEU:HD13	7:Z:115:ALA:HB1	2.02	0.41
8:a:38:TRP:O	8:a:42:VAL:HG23	2.21	0.41
7:e:131:GLU:H	7:e:131:GLU:CD	2.29	0.41
8:f:38:TRP:O	8:f:42:VAL:HG23	2.21	0.41
10:h:42:ARG:HD3	10:h:72:ARG:HG2	2.02	0.41
7:j:158:ASN:HD21	7:j:357:VAL:HG21	1.84	0.41
4:D:62:PHE:O	4:D:66:VAL:HG23	2.21	0.40
5:I:-59:DT:H3	6:J:59:DA:H61	1.68	0.40
1:K:73:GLU:O	1:K:76:GLN:N	2.54	0.40
3:M:57:TYR:O	3:M:60:ALA:HB3	2.21	0.40
5:S:-36:DT:H3	6:T:36:DA:N6	2.17	0.40
7:U:145:THR:HG22	7:U:232:GLY:HA3	2.02	0.40
7:Z:449:SER:HB2	9:b:50:PHE:HA	2.03	0.40
7:e:7:ILE:HG21	7:e:103:ILE:HD13	2.03	0.40
7:e:81:ILE:HD12	7:e:81:ILE:N	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:j:150:SER:HB2	7:j:430:ALA:HB2	2.03	0.40
7:j:285:SER:O	7:j:298:THR:HA	2.21	0.40
2:B:91:LYS:HE3	4:H:68:GLU:OE1	2.21	0.40
3:C:47:ALA:N	3:C:48:PRO:HD2	2.37	0.40
3:C:77:ARG:HD2	5:I:-54:DA:H4'	2.02	0.40
4:D:100:PRO:O	4:D:102:GLU:N	2.53	0.40
5:I:-35:DA:H2''	5:I:-34:DG:OP2	2.21	0.40
5:S:-35:DA:H2''	5:S:-34:DG:OP2	2.22	0.40
6:T:-9:DC:H6	6:T:-9:DC:H2'	1.70	0.40
7:U:153:GLN:HG3	7:U:442:PHE:HB3	2.02	0.40
7:Z:326:LYS:HG2	7:Z:348:GLN:HB2	2.03	0.40
9:b:7:THR:N	9:b:10:SER:HG	2.20	0.40
7:e:162:ILE:HG23	7:e:166:MET:HE3	2.04	0.40
7:j:406:VAL:HG13	11:n:57:TYR:CZ	2.57	0.40
7:j:422:THR:N	7:j:425:GLU:O	2.46	0.40
2:F:90:LEU:HB3	2:F:95:ARG:O	2.20	0.40
5:I:10:DC:H2''	5:I:11:DG:H8	1.86	0.40
1:O:42:ARG:HA	1:O:43:PRO:HD3	1.92	0.40
5:S:-38:DC:H6	5:S:-38:DC:H2'	1.73	0.40
5:S:10:DC:H2''	5:S:11:DG:H8	1.86	0.40
5:S:50:DA:N6	6:T:-50:DT:H3	2.20	0.40
7:U:413:LEU:HB2	7:U:465:TYR:CE1	2.56	0.40
8:V:61:ILE:O	8:V:65:VAL:HG23	2.21	0.40
7:Z:250:HIS:CD2	7:Z:250:HIS:C	2.98	0.40
7:j:142:MET:SD	9:l:82:ALA:HB1	2.61	0.40
3:C:88:ARG:HA	3:C:88:ARG:HD3	1.93	0.40
4:H:92:GLN:HE21	4:H:96:ARG:NH1	2.20	0.40
4:N:106:HIS:O	4:N:109:SER:HB3	2.21	0.40
3:Q:90:ASP:OD1	3:Q:93:LEU:N	2.47	0.40
5:S:-49:DG:N2	6:T:49:DC:N3	2.67	0.40
7:U:292:CYS:SG	7:U:338:GLU:HB2	2.62	0.40
7:Z:81:ILE:HD12	7:Z:81:ILE:N	2.36	0.40
9:b:49:TYR:CE2	9:b:56:LEU:HD12	2.57	0.40
7:j:81:ILE:N	7:j:81:ILE:HD12	2.37	0.40
6:J:-65:DT:H2''	6:J:-64:DA:C8	2.55	0.40
1:O:102:ALA:HB1	1:O:131:ARG:HH22	1.86	0.40
3:Q:31:HIS:CG	3:Q:35:ARG:HH21	2.40	0.40
7:U:124:LYS:NZ	11:Y:94:LEU:HD22	2.36	0.40
11:Y:42:THR:HG22	11:Y:43:PRO:O	2.22	0.40
7:Z:52:ILE:HG13	8:a:17:VAL:CG2	2.51	0.40
10:c:45:PHE:HB2	10:c:67:LEU:HD22	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:j:432:CYS:HB2	7:j:442:PHE:HE1	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	95/136 (70%)	84 (88%)	8 (8%)	3 (3%)	3	25
1	E	95/136 (70%)	86 (90%)	9 (10%)	0	100	100
1	K	95/136 (70%)	84 (88%)	8 (8%)	3 (3%)	3	25
1	O	95/136 (70%)	84 (88%)	8 (8%)	3 (3%)	3	25
2	B	81/103 (79%)	71 (88%)	8 (10%)	2 (2%)	4	28
2	F	76/103 (74%)	69 (91%)	7 (9%)	0	100	100
2	L	81/103 (79%)	71 (88%)	7 (9%)	3 (4%)	2	21
2	P	76/103 (74%)	65 (86%)	8 (10%)	3 (4%)	2	21
3	C	101/130 (78%)	89 (88%)	9 (9%)	3 (3%)	3	26
3	G	103/130 (79%)	91 (88%)	12 (12%)	0	100	100
3	M	101/130 (78%)	90 (89%)	8 (8%)	3 (3%)	3	26
3	Q	103/130 (79%)	91 (88%)	9 (9%)	3 (3%)	3	26
4	D	93/123 (76%)	84 (90%)	7 (8%)	2 (2%)	5	30
4	H	91/123 (74%)	81 (89%)	9 (10%)	1 (1%)	11	42
4	N	93/123 (76%)	83 (89%)	8 (9%)	2 (2%)	5	30
4	R	91/123 (74%)	82 (90%)	7 (8%)	2 (2%)	5	30
7	U	439/472 (93%)	430 (98%)	9 (2%)	0	100	100
7	Z	443/472 (94%)	434 (98%)	9 (2%)	0	100	100
7	e	445/472 (94%)	435 (98%)	9 (2%)	1 (0%)	43	73

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
7	j	437/472 (93%)	426 (98%)	11 (2%)	0	100	100
8	V	87/96 (91%)	87 (100%)	0	0	100	100
8	a	86/96 (90%)	86 (100%)	0	0	100	100
8	f	81/96 (84%)	81 (100%)	0	0	100	100
8	k	84/96 (88%)	84 (100%)	0	0	100	100
9	W	88/99 (89%)	86 (98%)	2 (2%)	0	100	100
9	b	87/99 (88%)	84 (97%)	2 (2%)	1 (1%)	11	42
9	g	88/99 (89%)	86 (98%)	2 (2%)	0	100	100
9	l	91/99 (92%)	89 (98%)	2 (2%)	0	100	100
10	X	74/76 (97%)	73 (99%)	1 (1%)	0	100	100
10	c	74/76 (97%)	73 (99%)	1 (1%)	0	100	100
10	h	74/76 (97%)	72 (97%)	2 (3%)	0	100	100
10	m	74/76 (97%)	73 (99%)	1 (1%)	0	100	100
11	Y	78/104 (75%)	75 (96%)	3 (4%)	0	100	100
11	d	85/104 (82%)	80 (94%)	4 (5%)	1 (1%)	10	40
11	i	85/104 (82%)	82 (96%)	3 (4%)	0	100	100
11	n	82/104 (79%)	79 (96%)	3 (4%)	0	100	100
All	All	4552/5356 (85%)	4320 (95%)	196 (4%)	36 (1%)	16	49

All (36) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	73	GLU
2	B	29	ILE
4	D	101	GLY
1	K	73	GLU
2	L	29	ILE
4	N	101	GLY
4	R	100	PRO
9	b	76	CYS
11	d	47	ASP
4	H	101	GLY
1	K	77	ASP
4	N	109	SER
1	O	73	GLU
1	O	77	ASP

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Mol	Chain	Res	Type
2	P	29	ILE
1	A	77	ASP
4	D	109	SER
4	R	109	SER
7	e	292	CYS
2	B	26	ILE
3	C	64	GLU
3	M	64	GLU
3	Q	64	GLU
1	A	74	ILE
1	K	74	ILE
2	L	26	ILE
1	O	74	ILE
3	C	113	SER
3	M	113	SER
3	Q	113	SER
2	P	26	ILE
3	Q	114	VAL
3	C	114	VAL
2	P	34	ILE
2	L	34	ILE
3	M	114	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	85/111 (77%)	78 (92%)	7 (8%)	10	34
1	E	84/111 (76%)	83 (99%)	1 (1%)	63	71
1	K	85/111 (77%)	79 (93%)	6 (7%)	13	39
1	O	84/111 (76%)	77 (92%)	7 (8%)	10	34
2	B	68/79 (86%)	62 (91%)	6 (9%)	9	33
2	F	63/79 (80%)	60 (95%)	3 (5%)	23	47
2	L	68/79 (86%)	62 (91%)	6 (9%)	9	33

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	P	63/79 (80%)	56 (89%)	7 (11%)	6	24
3	C	82/102 (80%)	76 (93%)	6 (7%)	13	38
3	G	83/102 (81%)	79 (95%)	4 (5%)	23	47
3	M	82/102 (80%)	78 (95%)	4 (5%)	22	47
3	Q	83/102 (81%)	80 (96%)	3 (4%)	31	54
4	D	81/103 (79%)	72 (89%)	9 (11%)	6	24
4	H	79/103 (77%)	72 (91%)	7 (9%)	9	32
4	N	81/103 (79%)	72 (89%)	9 (11%)	6	24
4	R	79/103 (77%)	70 (89%)	9 (11%)	5	23
7	U	406/428 (95%)	393 (97%)	13 (3%)	34	56
7	Z	409/428 (96%)	396 (97%)	13 (3%)	34	56
7	e	411/428 (96%)	401 (98%)	10 (2%)	43	62
7	j	405/428 (95%)	394 (97%)	11 (3%)	39	59
8	V	84/91 (92%)	81 (96%)	3 (4%)	31	54
8	a	86/91 (94%)	84 (98%)	2 (2%)	44	63
8	f	81/91 (89%)	79 (98%)	2 (2%)	42	61
8	k	81/91 (89%)	79 (98%)	2 (2%)	42	61
9	W	82/89 (92%)	79 (96%)	3 (4%)	30	53
9	b	81/89 (91%)	78 (96%)	3 (4%)	30	53
9	g	82/89 (92%)	80 (98%)	2 (2%)	43	62
9	l	84/89 (94%)	80 (95%)	4 (5%)	23	47
10	X	68/68 (100%)	67 (98%)	1 (2%)	57	69
10	c	68/68 (100%)	67 (98%)	1 (2%)	57	69
10	h	68/68 (100%)	67 (98%)	1 (2%)	57	69
10	m	68/68 (100%)	68 (100%)	0	100	100
11	Y	77/90 (86%)	72 (94%)	5 (6%)	15	42
11	d	81/90 (90%)	77 (95%)	4 (5%)	22	47
11	i	81/90 (90%)	76 (94%)	5 (6%)	16	42
11	n	79/90 (88%)	77 (98%)	2 (2%)	42	61
All	All	4132/4644 (89%)	3951 (96%)	181 (4%)	25	49

All (181) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	VAL
1	A	48	LEU
1	A	59	GLU
1	A	63	ARG
1	A	65	LEU
1	A	105	GLU
1	A	129	ARG
2	B	47	SER
2	B	73	THR
2	B	80	THR
2	B	91	LYS
2	B	92	ARG
2	B	95	ARG
3	C	29	ARG
3	C	59	THR
3	C	73	ASN
3	C	81	ARG
3	C	114	VAL
3	C	118	LYS
4	D	39	TYR
4	D	49	THR
4	D	77	LEU
4	D	84	SER
4	D	85	THR
4	D	98	LEU
4	D	109	SER
4	D	116	THR
4	D	119	THR
1	E	59	GLU
2	F	26	ILE
2	F	73	THR
2	F	92	ARG
3	G	72	ASP
3	G	81	ARG
3	G	88	ARG
3	G	114	VAL
4	H	49	THR
4	H	88	SER
4	H	93	THR
4	H	95	VAL
4	H	99	LEU
4	H	102	GLU
4	H	103	LEU

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Mol	Chain	Res	Type
1	K	46	VAL
1	K	59	GLU
1	K	63	ARG
1	K	65	LEU
1	K	105	GLU
1	K	129	ARG
2	L	47	SER
2	L	73	THR
2	L	80	THR
2	L	91	LYS
2	L	92	ARG
2	L	95	ARG
3	M	29	ARG
3	M	81	ARG
3	M	91	GLU
3	M	114	VAL
4	N	49	THR
4	N	60	ASN
4	N	77	LEU
4	N	84	SER
4	N	85	THR
4	N	93	THR
4	N	98	LEU
4	N	116	THR
4	N	119	THR
1	O	46	VAL
1	O	59	GLU
1	O	63	ARG
1	O	65	LEU
1	O	90	MET
1	O	105	GLU
1	O	134	ARG
2	P	26	ILE
2	P	47	SER
2	P	73	THR
2	P	80	THR
2	P	85	ASP
2	P	92	ARG
2	P	95	ARG
3	Q	81	ARG
3	Q	114	VAL
3	Q	118	LYS

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Mol	Chain	Res	Type
4	R	39	TYR
4	R	49	THR
4	R	77	LEU
4	R	84	SER
4	R	93	THR
4	R	97	LEU
4	R	109	SER
4	R	116	THR
4	R	119	THR
7	U	1	MET
7	U	98	LYS
7	U	124	LYS
7	U	142	MET
7	U	156	ILE
7	U	176	VAL
7	U	210	THR
7	U	228	GLN
7	U	269	LYS
7	U	296	SER
7	U	365	PHE
7	U	378	ASP
7	U	425	GLU
8	V	28	GLU
8	V	68	LYS
8	V	91	GLU
9	W	65	GLN
9	W	67	SER
9	W	85	LEU
10	X	39	ASP
11	Y	11	ILE
11	Y	18	GLU
11	Y	29	ASN
11	Y	87	LEU
11	Y	96	ASN
7	Z	1	MET
7	Z	98	LYS
7	Z	116	LYS
7	Z	142	MET
7	Z	156	ILE
7	Z	176	VAL
7	Z	184	SER
7	Z	210	THR

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Mol	Chain	Res	Type
7	Z	228	GLN
7	Z	269	LYS
7	Z	365	PHE
7	Z	378	ASP
7	Z	397	LYS
8	a	28	GLU
8	a	68	LYS
9	b	67	SER
9	b	75	ASN
9	b	85	LEU
10	c	39	ASP
11	d	18	GLU
11	d	29	ASN
11	d	87	LEU
11	d	96	ASN
7	e	98	LYS
7	e	142	MET
7	e	156	ILE
7	e	176	VAL
7	e	184	SER
7	e	210	THR
7	e	228	GLN
7	e	269	LYS
7	e	365	PHE
7	e	397	LYS
8	f	28	GLU
8	f	68	LYS
9	g	65	GLN
9	g	85	LEU
10	h	39	ASP
11	i	18	GLU
11	i	29	ASN
11	i	87	LEU
11	i	92	ASP
11	i	96	ASN
7	j	98	LYS
7	j	142	MET
7	j	156	ILE
7	j	176	VAL
7	j	210	THR
7	j	228	GLN
7	j	269	LYS

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Mol	Chain	Res	Type
7	j	296	SER
7	j	365	PHE
7	j	378	ASP
7	j	425	GLU
8	k	28	GLU
8	k	68	LYS
9	l	65	GLN
9	l	67	SER
9	l	85	LEU
9	l	92	CYS
11	n	87	LEU
11	n	96	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (51) such sidechains are listed below:

Mol	Chain	Res	Type
3	C	38	ASN
1	E	39	HIS
3	G	31	HIS
4	H	92	GLN
1	K	55	GLN
1	K	113	HIS
2	L	75	HIS
4	N	60	ASN
1	O	39	HIS
3	Q	82	HIS
3	Q	112	GLN
4	R	64	ASN
7	U	42	ASN
7	U	62	GLN
7	U	83	HIS
7	U	88	ASN
7	U	90	ASN
7	U	168	GLN
7	U	248	GLN
7	U	352	HIS
7	U	373	ASN
7	U	459	GLN
8	V	57	ASN
11	Y	97	HIS
7	Z	53	ASN
7	Z	229	ASN

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Mol	Chain	Res	Type
7	Z	247	ASN
7	Z	335	HIS
7	Z	459	GLN
7	Z	469	GLN
8	a	57	ASN
9	b	75	ASN
11	d	97	HIS
7	e	168	GLN
7	e	328	GLN
7	e	348	GLN
7	e	427	HIS
8	f	57	ASN
9	g	17	ASN
11	i	61	ASN
7	j	8	GLN
7	j	42	ASN
7	j	157	HIS
7	j	229	ASN
7	j	469	GLN
8	k	27	ASN
8	k	57	ASN
9	l	72	HIS
9	l	83	ASN
9	l	88	HIS
11	n	29	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 32 ligands modelled in this entry, 32 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	97/136 (71%)	-0.72	0 100 100	134, 166, 198, 212	0
1	E	97/136 (71%)	-0.61	0 100 100	139, 172, 213, 249	0
1	K	97/136 (71%)	-0.53	0 100 100	146, 175, 209, 216	0
1	O	97/136 (71%)	-0.80	0 100 100	141, 177, 219, 249	0
2	B	83/103 (80%)	-0.73	0 100 100	124, 155, 185, 198	0
2	F	78/103 (75%)	-0.73	0 100 100	124, 166, 189, 201	0
2	L	83/103 (80%)	-0.85	0 100 100	131, 152, 192, 204	0
2	P	78/103 (75%)	-0.49	0 100 100	153, 183, 219, 230	0
3	C	103/130 (79%)	-0.65	0 100 100	128, 162, 187, 197	0
3	G	105/130 (80%)	-0.50	0 100 100	148, 177, 198, 207	0
3	M	103/130 (79%)	-0.56	0 100 100	143, 168, 195, 216	0
3	Q	105/130 (80%)	-0.65	0 100 100	141, 179, 205, 222	0
4	D	95/123 (77%)	-0.70	1 (1%) 78 57	125, 155, 201, 243	0
4	H	93/123 (75%)	-0.62	0 100 100	145, 172, 197, 209	0
4	N	95/123 (77%)	-0.65	0 100 100	137, 165, 190, 239	0
4	R	93/123 (75%)	-0.57	0 100 100	149, 172, 196, 207	0
5	I	145/145 (100%)	-0.71	0 100 100	176, 210, 241, 255	0
5	S	145/145 (100%)	-0.77	1 (0%) 84 66	177, 208, 241, 257	0
6	J	145/145 (100%)	-0.81	0 100 100	169, 211, 243, 269	0
6	T	145/145 (100%)	-0.80	0 100 100	177, 211, 243, 264	0
7	U	447/472 (94%)	-0.66	1 (0%) 91 82	133, 183, 235, 285	0
7	Z	451/472 (95%)	-0.51	0 100 100	161, 205, 261, 302	0
7	e	453/472 (95%)	-0.67	0 100 100	148, 192, 244, 276	0
7	j	447/472 (94%)	-0.73	0 100 100	146, 189, 232, 286	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
8	V	89/96 (92%)	-0.50	0	100	100	160, 196, 260, 275	0
8	a	90/96 (93%)	-0.21	0	100	100	181, 238, 267, 283	0
8	f	85/96 (88%)	-0.47	0	100	100	196, 227, 267, 298	0
8	k	86/96 (89%)	-0.68	0	100	100	154, 217, 249, 277	1 (1%)
9	W	90/99 (90%)	-0.76	0	100	100	163, 196, 217, 231	0
9	b	89/99 (89%)	-0.41	1 (1%)	78	57	173, 219, 252, 260	0
9	g	90/99 (90%)	-0.66	0	100	100	170, 211, 264, 282	1 (1%)
9	l	93/99 (93%)	-0.75	1 (1%)	78	57	168, 201, 263, 275	0
10	X	76/76 (100%)	-0.79	0	100	100	157, 193, 221, 237	0
10	c	76/76 (100%)	-0.75	0	100	100	159, 192, 218, 252	0
10	h	76/76 (100%)	-0.78	0	100	100	170, 195, 214, 239	0
10	m	76/76 (100%)	-0.78	0	100	100	159, 188, 218, 233	0
11	Y	82/104 (78%)	-0.51	0	100	100	159, 193, 257, 301	0
11	d	89/104 (85%)	-0.39	2 (2%)	62	44	174, 212, 279, 315	0
11	i	89/104 (85%)	-0.53	0	100	100	162, 200, 266, 297	0
11	n	86/104 (82%)	-0.58	0	100	100	167, 197, 281, 349	0
All	All	5242/5936 (88%)	-0.64	7 (0%)	92	87	124, 192, 248, 349	2 (0%)

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
11	d	37	SER	4.9
9	l	9	ASP	4.0
11	d	43	PRO	2.3
9	b	65	GLN	2.2
7	U	38	GLU	2.2
4	D	29	THR	2.2
5	S	60	DT	2.1

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
12	ZN	Z	503	1/1	0.96	0.07	237,237,237,237	0
12	ZN	Z	502	1/1	0.98	0.08	233,233,233,233	0
12	ZN	U	502	1/1	0.98	0.04	195,195,195,195	0
12	ZN	e	501	1/1	0.98	0.04	251,251,251,251	0
12	ZN	U	506	1/1	0.99	0.03	252,252,252,252	0
12	ZN	Z	505	1/1	0.99	0.02	272,272,272,272	0
12	ZN	Z	506	1/1	0.99	0.03	290,290,290,290	0
12	ZN	b	101	1/1	0.99	0.03	218,218,218,218	0
12	ZN	U	501	1/1	0.99	0.03	228,228,228,228	0
12	ZN	e	502	1/1	0.99	0.05	279,279,279,279	0
12	ZN	e	503	1/1	0.99	0.03	229,229,229,229	0
12	ZN	e	506	1/1	0.99	0.03	251,251,251,251	0
12	ZN	g	101	1/1	0.99	0.03	206,206,206,206	0
12	ZN	j	501	1/1	0.99	0.04	286,286,286,286	0
12	ZN	j	504	1/1	0.99	0.03	198,198,198,198	0
12	ZN	j	505	1/1	0.99	0.02	212,212,212,212	0
12	ZN	n	201	1/1	0.99	0.03	216,216,216,216	0
12	ZN	U	504	1/1	1.00	0.02	201,201,201,201	0
12	ZN	U	505	1/1	1.00	0.03	160,160,160,160	0
12	ZN	e	504	1/1	1.00	0.01	188,188,188,188	0
12	ZN	e	505	1/1	1.00	0.02	152,152,152,152	0
12	ZN	Z	504	1/1	1.00	0.02	215,215,215,215	0
12	ZN	U	503	1/1	1.00	0.02	209,209,209,209	0
12	ZN	i	201	1/1	1.00	0.01	195,195,195,195	0
12	ZN	W	101	1/1	1.00	0.02	199,199,199,199	0
12	ZN	j	502	1/1	1.00	0.04	160,160,160,160	0
12	ZN	j	503	1/1	1.00	0.02	179,179,179,179	0
12	ZN	Y	201	1/1	1.00	0.03	184,184,184,184	0
12	ZN	d	201	1/1	1.00	0.03	226,226,226,226	0
12	ZN	j	506	1/1	1.00	0.03	192,192,192,192	0
12	ZN	l	101	1/1	1.00	0.01	173,173,173,173	0
12	ZN	Z	501	1/1	1.00	0.02	234,234,234,234	0

6.5 Other polymers [i](#)

There are no such residues in this entry.