



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 11:56 PM GMT

PDB ID : 2C39  
Title : RNASE PH CORE OF THE ARCHAEAL EXOSOME IN COMPLEX WITH ADP  
Authors : Lorentzen, E.; Conti, E.  
Deposited on : 2005-10-05  
Resolution : 3.30 Å(reported)

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

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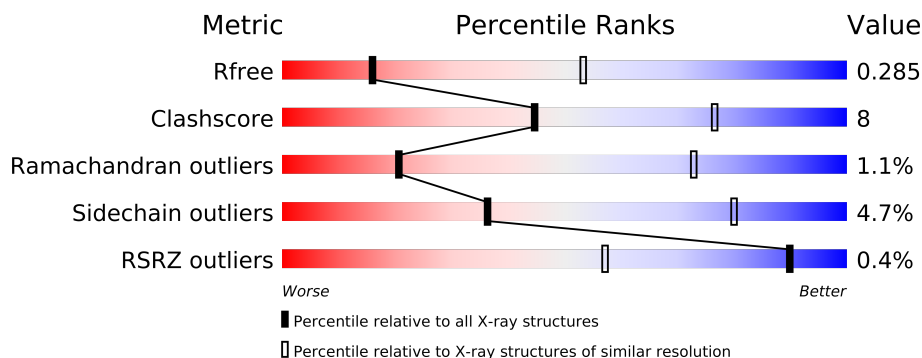
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 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 3.30 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.









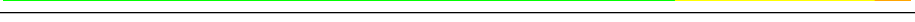

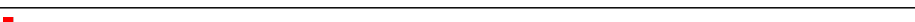

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	66092	1341 (3.40-3.20)
Clashscore	79885	1696 (3.40-3.20)
Ramachandran outliers	78287	1664 (3.40-3.20)
Sidechain outliers	78261	1662 (3.40-3.20)
RSRZ outliers	66119	1342 (3.40-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	275	
1	C	275	
1	E	275	
1	G	275	
1	I	275	
1	K	275	
1	M	275	
1	O	275	
1	Q	275	
1	S	275	
1	U	275	
1	W	275	
2	B	248	
2	D	248	

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Mol	Chain	Length	Quality of chain
2	F	248	
2	H	248	
2	J	248	
2	L	248	
2	N	248	
2	P	248	
2	R	248	
2	T	248	
2	V	248	
2	X	248	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 45814 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 PROBABLE EXOSOME COMPLEX EXONUCLEASE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	260	Total	C	N	O	S	0	0	0
			1944	1239	320	380	5			
1	C	260	Total	C	N	O	S	0	0	0
			1947	1243	321	378	5			
1	E	260	Total	C	N	O	S	0	0	0
			1968	1253	325	385	5			
1	G	260	Total	C	N	O	S	0	0	0
			1961	1250	323	383	5			
1	I	260	Total	C	N	O	S	0	0	0
			1954	1247	323	379	5			
1	K	260	Total	C	N	O	S	0	0	0
			1958	1249	323	381	5			
1	M	260	Total	C	N	O	S	0	0	0
			1960	1249	324	382	5			
1	O	260	Total	C	N	O	S	0	0	0
			1951	1246	323	377	5			
1	Q	260	Total	C	N	O	S	0	0	0
			1953	1245	322	381	5			
1	S	259	Total	C	N	O	S	0	0	0
			1955	1245	323	382	5			
1	U	255	Total	C	N	O	S	0	0	0
			1908	1219	317	367	5			
1	W	259	Total	C	N	O	S	0	0	0
			1950	1245	322	378	5			

- Molecule 2 is a protein called PROBABLE EXOSOME COMPLEX EXONUCLEASE 1.

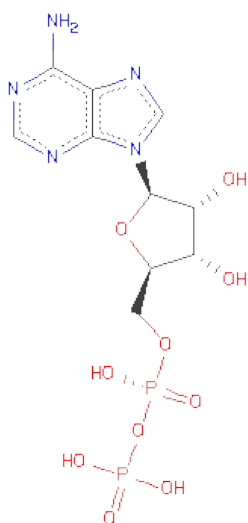
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	241	Total	C	N	O	S	0	0	0
			1812	1145	317	340	10			
2	D	248	Total	C	N	O	S	0	0	0
			1900	1198	330	360	12			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	241	Total	C	N	O	S	0	0	0
			1838	1161	317	350	10			
2	H	230	Total	C	N	O	S	0	0	0
			1732	1094	302	327	9			
2	J	241	Total	C	N	O	S	0	0	0
			1844	1165	321	348	10			
2	L	247	Total	C	N	O	S	0	0	0
			1890	1194	328	357	11			
2	N	239	Total	C	N	O	S	0	0	0
			1809	1144	313	342	10			
2	P	248	Total	C	N	O	S	0	0	0
			1904	1202	332	358	12			
2	R	248	Total	C	N	O	S	0	0	0
			1880	1188	332	348	12			
2	T	248	Total	C	N	O	S	0	0	0
			1884	1193	328	351	12			
2	V	239	Total	C	N	O	S	0	0	0
			1815	1148	318	339	10			
2	X	239	Total	C	N	O	S	0	0	0
			1821	1148	315	348	10			

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

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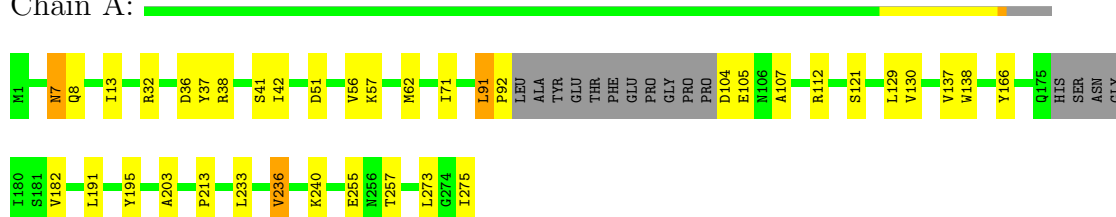
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	F	1	Total	C	O	P		0	0
			17	5	10	2			
3	H	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	J	1	Total	O	P			0	0
			9	7	2				
3	L	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	N	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	P	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	R	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	T	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	V	1	Total	C	O	P		0	0
			17	5	10	2			
3	X	1	Total	C	O	P		0	0
			17	5	10	2			

### 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.

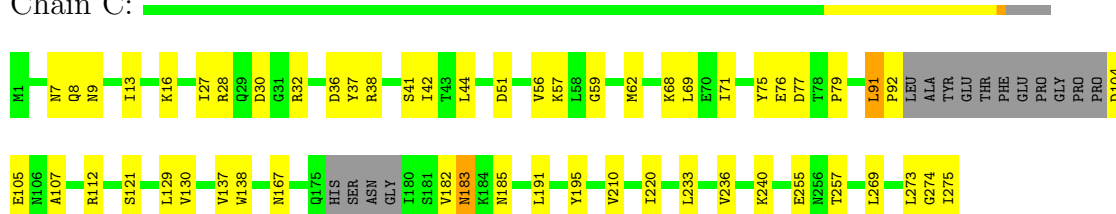
#### • Molecule 1: PROBABLE EXOSOME COMPLEX EXONUCLEASE 2

Chain A:



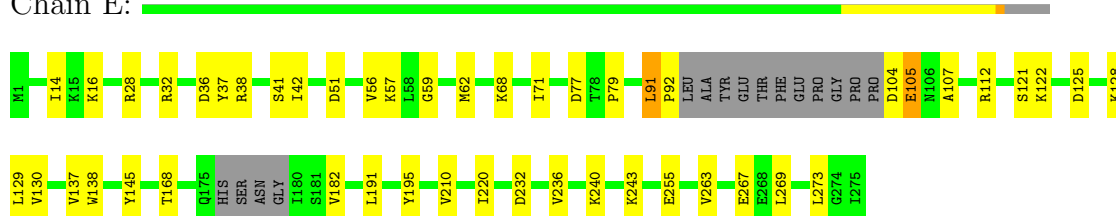
#### • Molecule 1: PROBABLE EXOSOME COMPLEX EXONUCLEASE 2

Chain C:



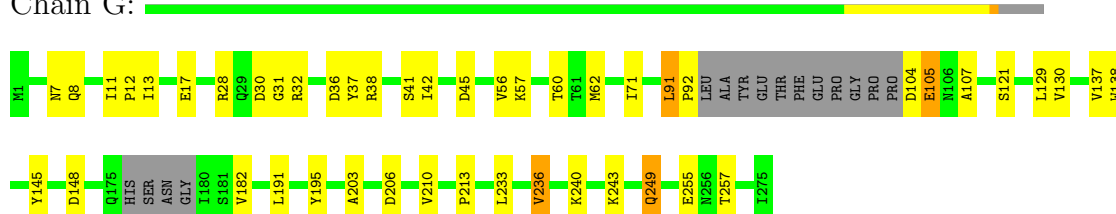
#### • Molecule 1: PROBABLE EXOSOME COMPLEX EXONUCLEASE 2

Chain E:



#### • Molecule 1: PROBABLE EXOSOME COMPLEX EXONUCLEASE 2

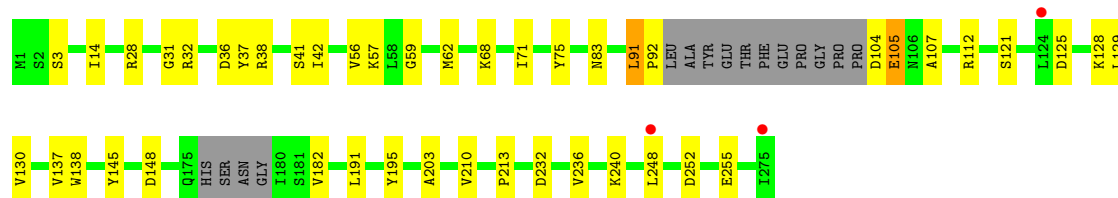
Chain G:



#### • Molecule 1: PROBABLE EXOSOME COMPLEX EXONUCLEASE 2

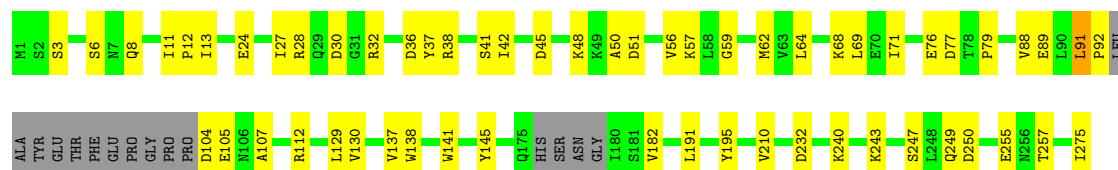
Chain I:





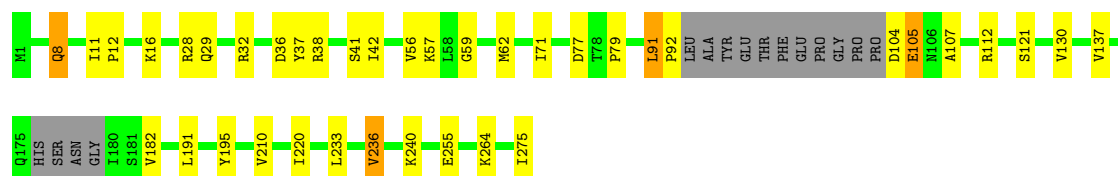
• Molecule 1: PROBABLE EXOSOME COMPLEX EXONUCLEASE 2

Chain K:



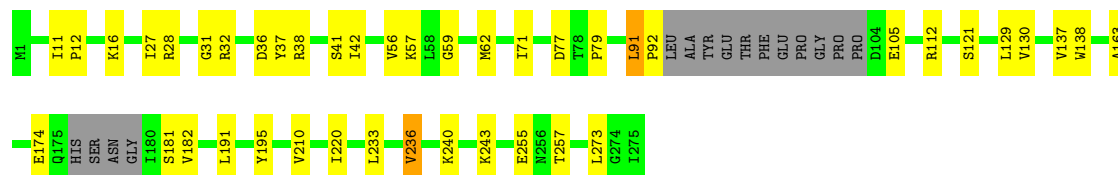
• Molecule 1: PROBABLE EXOSOME COMPLEX EXONUCLEASE 2

Chain M:



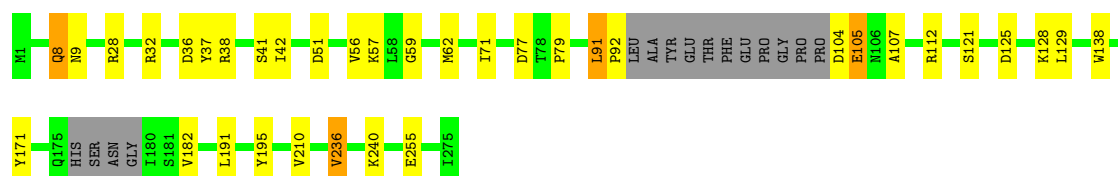
• Molecule 1: PROBABLE EXOSOME COMPLEX EXONUCLEASE 2

Chain O:



• Molecule 1: PROBABLE EXOSOME COMPLEX EXONUCLEASE 2

Chain Q:

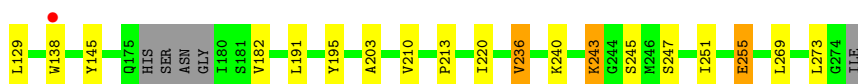


• Molecule 1: PROBABLE EXOSOME COMPLEX EXONUCLEASE 2

Chain S:

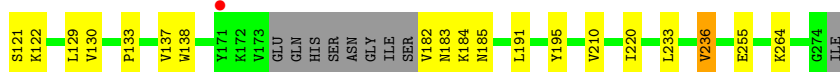






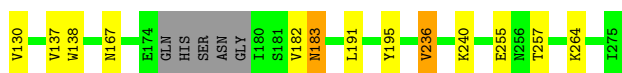
• Molecule 1: PROBABLE EXOSOME COMPLEX EXONUCLEASE 2

Chain U:



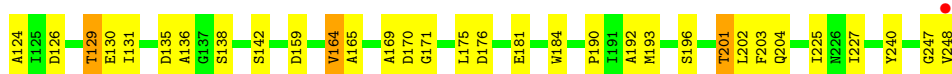
• Molecule 1: PROBABLE EXOSOME COMPLEX EXONUCLEASE 2

Chain W:



• Molecule 2: PROBABLE EXOSOME COMPLEX EXONUCLEASE 1

Chain B:



• Molecule 2: PROBABLE EXOSOME COMPLEX EXONUCLEASE 1

Chain D:



• Molecule 2: PROBABLE EXOSOME COMPLEX EXONUCLEASE 1

Chain F:



E246  
G247  
V248

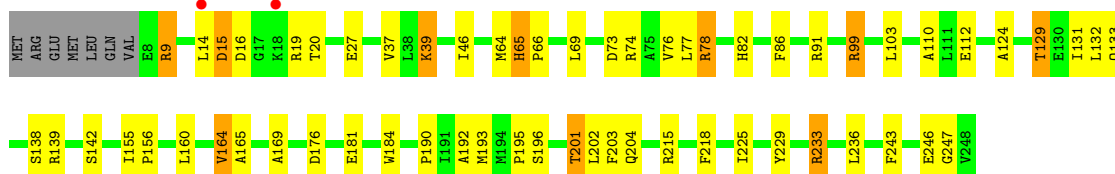
• Molecule 2: PROBABLE EXOSOME COMPLEX EXONUCLEASE 1

Chain H:



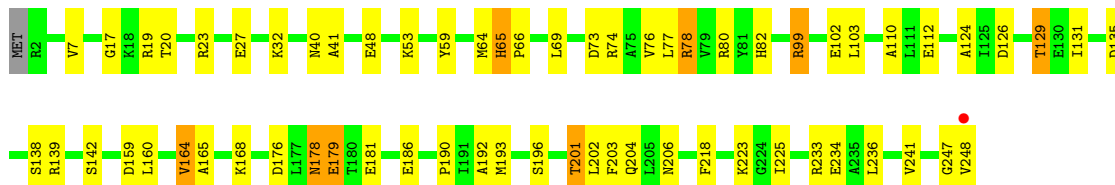
• Molecule 2: PROBABLE EXOSOME COMPLEX EXONUCLEASE 1

Chain J:



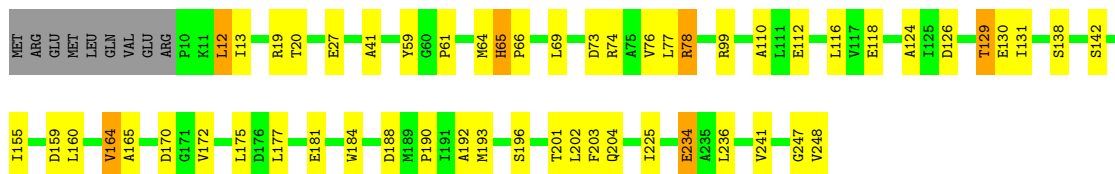
• Molecule 2: PROBABLE EXOSOME COMPLEX EXONUCLEASE 1

Chain L:



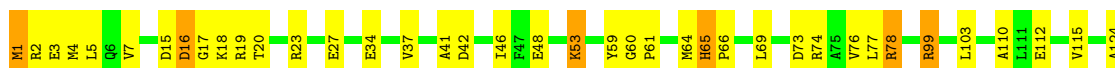
• Molecule 2: PROBABLE EXOSOME COMPLEX EXONUCLEASE 1

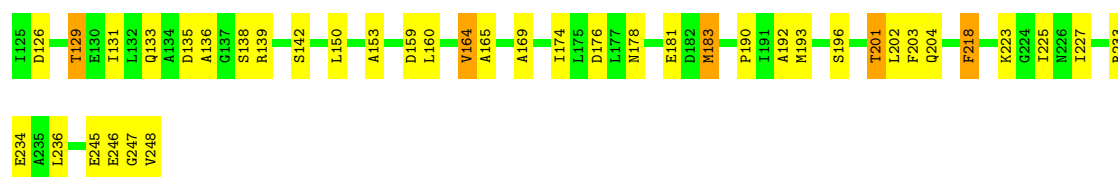
Chain N:



• Molecule 2: PROBABLE EXOSOME COMPLEX EXONUCLEASE 1

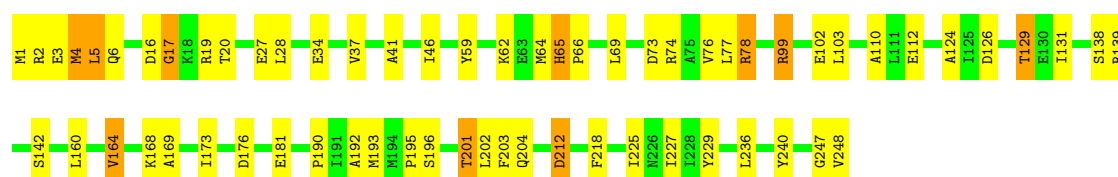
Chain P:





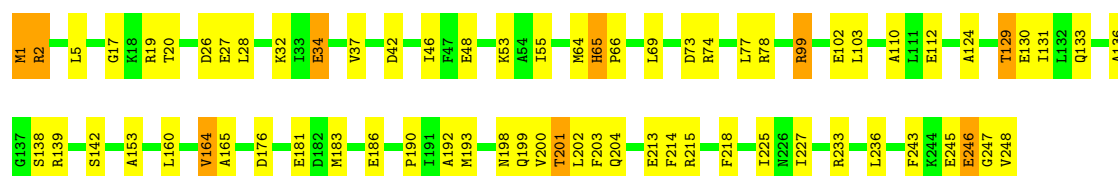
• Molecule 2: PROBABLE EXOSOME COMPLEX EXONUCLEASE 1

Chain R:



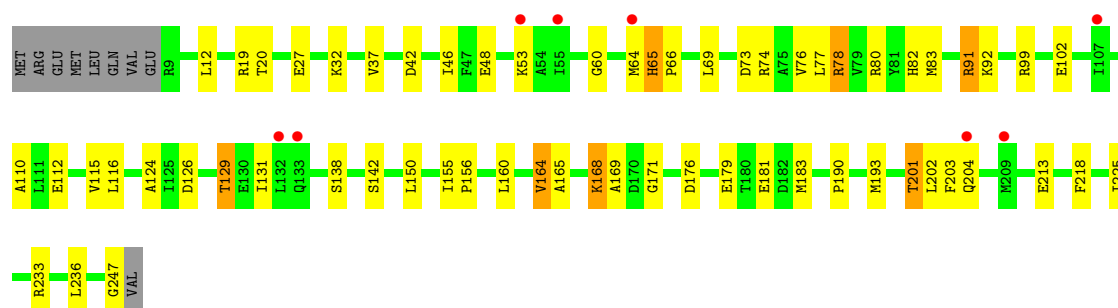
• Molecule 2: PROBABLE EXOSOME COMPLEX EXONUCLEASE 1

Chain T:



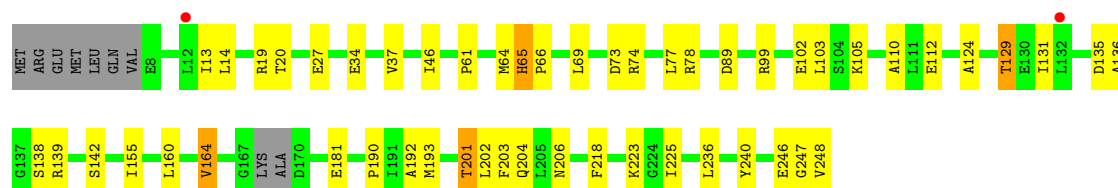
• Molecule 2: PROBABLE EXOSOME COMPLEX EXONUCLEASE 1

Chain V:



• Molecule 2: PROBABLE EXOSOME COMPLEX EXONUCLEASE 1

Chain X:



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	206.20Å 214.00Å 432.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	93.25 – 3.30 93.06 – 3.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (93.25-3.30) 93.5 (93.06-3.30)	Depositor EDS
$R_{merge}$	0.18	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.99 (at 3.33Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.274 , 0.295 0.266 , 0.285	Depositor DCC
$R_{free}$ test set	4015 reflections (3.09%)	DCC
Wilson B-factor (Å <sup>2</sup> )	86.3	Xtriage
Anisotropy	0.060	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 15.1	EDS
Estimated twinning fraction	0.048 for -k,-h,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.24$	Xtriage
Outliers	0 of 133792 reflections	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	45814	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.35% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

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

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	0.47	0/1969	0.58	0/2678
1	C	0.54	0/1972	0.59	0/2680
1	E	0.47	0/1993	0.58	0/2706
1	G	0.49	0/1986	0.57	0/2697
1	I	0.44	0/1979	0.56	0/2688
1	K	0.54	0/1983	0.60	0/2693
1	M	0.49	0/1985	0.58	0/2696
1	O	0.52	0/1976	0.60	0/2684
1	Q	0.49	0/1978	0.58	0/2688
1	S	0.47	0/1980	0.58	0/2690
1	U	0.43	0/1933	0.58	0/2628
1	W	0.47	0/1975	0.57	0/2682
2	B	0.45	0/1840	0.60	0/2494
2	D	0.51	0/1928	0.64	0/2606
2	F	0.47	0/1866	0.60	0/2524
2	H	0.45	0/1757	0.59	0/2377
2	J	0.46	0/1872	0.58	0/2530
2	L	0.51	0/1918	0.65	0/2592
2	N	0.48	0/1837	0.60	0/2487
2	P	0.51	0/1932	0.65	0/2609
2	R	0.50	0/1908	0.65	1/2578 (0.0%)
2	T	0.49	0/1912	0.63	0/2584
2	V	0.47	0/1843	0.60	0/2494
2	X	0.46	0/1848	0.60	0/2501
All	All	0.48	0/46170	0.60	1/62586 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	D	0	2
2	L	0	1
2	P	0	1
2	R	0	2
2	T	0	1
All	All	0	7

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	R	4	MET	N-CA-C	-5.96	94.91	111.00

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	1	MET	Peptide
2	D	17	GLY	Peptide
2	L	17	GLY	Peptide
2	P	17	GLY	Peptide
2	R	17	GLY	Peptide
2	R	5	LEU	Peptide
2	T	17	GLY	Peptide

## 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	1944	0	1969	20	0
1	C	1947	0	1982	30	0
1	E	1968	0	2013	35	0
1	G	1961	0	2003	33	0
1	I	1954	0	1997	33	0
1	K	1958	0	2001	35	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	M	1960	0	2003	20	0
1	O	1951	0	1995	23	0
1	Q	1953	0	1988	23	0
1	S	1955	0	1996	29	0
1	U	1908	0	1950	26	0
1	W	1950	0	1997	25	0
2	B	1812	0	1810	41	0
2	D	1900	0	1920	45	0
2	F	1838	0	1852	47	0
2	H	1732	0	1740	49	0
2	J	1844	0	1870	47	0
2	L	1890	0	1915	50	0
2	N	1809	0	1817	41	0
2	P	1904	0	1938	47	0
2	R	1880	0	1904	39	0
2	T	1884	0	1913	49	0
2	V	1815	0	1836	41	0
2	X	1821	0	1820	31	0
3	B	27	0	12	5	0
3	D	27	0	12	3	0
3	F	17	0	7	1	0
3	H	27	0	12	2	0
3	J	9	0	0	1	0
3	L	27	0	12	3	0
3	N	27	0	12	3	0
3	P	27	0	12	3	0
3	R	27	0	12	0	0
3	T	27	0	12	3	0
3	V	17	0	7	2	0
3	X	17	0	7	0	0
All	All	45814	0	46346	777	0

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 8.

All (777) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:T:1:MET:H2	2:T:2:ARG:HA	1.24	1.00
2:T:1:MET:N	2:T:2:ARG:HA	1.75	1.00
1:E:243:LYS:HG2	2:F:112:GLU:OE1	1.61	0.98
1:G:249:GLN:H	1:G:249:GLN:HE21	1.18	0.90
2:P:99:ARG:NH1	3:P:404:ADP:O2A	2.07	0.87

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:78:ARG:HD3	2:L:126:ASP:OD1	1.75	0.86
2:H:227:ILE:HD12	2:H:248:VAL:HG13	1.59	0.83
2:V:99:ARG:NH1	3:V:404:ADP:O2A	2.12	0.81
2:T:99:ARG:NH1	3:T:404:ADP:O2A	2.13	0.81
1:G:249:GLN:N	1:G:249:GLN:HE21	1.77	0.81
2:J:64:MET:HE2	2:J:124:ALA:HB2	1.66	0.77
2:F:99:ARG:NH1	3:F:404:ADP:O2A	2.16	0.77
2:L:64:MET:HE2	2:L:124:ALA:HB2	1.67	0.76
2:T:53:LYS:HG3	2:T:133:GLN:HB2	1.67	0.76
2:H:48:GLU:HG2	2:H:53:LYS:HG2	1.67	0.75
2:L:99:ARG:NH1	3:L:404:ADP:O2A	2.19	0.75
2:N:77:LEU:HD12	2:N:112:GLU:HG3	1.69	0.74
2:N:99:ARG:HH12	3:N:404:ADP:PA	2.09	0.74
2:R:78:ARG:HD3	2:R:126:ASP:OD1	1.88	0.74
2:X:64:MET:HE2	2:X:124:ALA:HB2	1.70	0.73
2:P:53:LYS:HG3	2:P:133:GLN:HB2	1.69	0.72
2:H:64:MET:HE2	2:H:124:ALA:HB2	1.69	0.72
2:T:243:PHE:HZ	2:T:246:GLU:HG2	1.54	0.72
2:T:64:MET:HE2	2:T:124:ALA:HB2	1.72	0.72
2:B:64:MET:HE2	2:B:124:ALA:HB2	1.70	0.72
2:L:77:LEU:HD12	2:L:112:GLU:HG3	1.72	0.72
2:H:99:ARG:NH1	3:H:404:ADP:O2A	2.22	0.72
2:J:99:ARG:HH12	3:J:404:ADP:PB	2.11	0.72
2:P:64:MET:HE2	2:P:124:ALA:HB2	1.71	0.72
2:B:175:LEU:HD22	2:B:248:VAL:HG21	1.72	0.71
1:W:51:ASP:OD2	1:W:69:LEU:HB2	1.92	0.70
2:F:77:LEU:HD12	2:F:112:GLU:HG3	1.73	0.70
2:D:64:MET:HE2	2:D:124:ALA:HB2	1.72	0.70
2:X:223:LYS:HB3	2:X:248:VAL:HG11	1.72	0.70
2:F:129:THR:HG21	2:F:142:SER:OG	1.90	0.70
2:R:77:LEU:HD12	2:R:112:GLU:HG3	1.73	0.70
2:H:78:ARG:HD3	2:H:126:ASP:OD1	1.92	0.70
2:L:48:GLU:HG2	2:L:53:LYS:HG2	1.72	0.70
2:T:66:PRO:HG2	2:T:69:LEU:HD12	1.74	0.69
2:N:99:ARG:NH1	3:N:404:ADP:O2A	2.25	0.69
2:R:64:MET:HE2	2:R:124:ALA:HB2	1.73	0.69
2:F:48:GLU:OE1	2:F:53:LYS:HE2	1.92	0.69
2:D:77:LEU:HD12	2:D:112:GLU:HG3	1.73	0.69
2:T:77:LEU:HD12	2:T:112:GLU:HG3	1.74	0.69
1:G:243:LYS:HG2	2:H:112:GLU:OE1	1.92	0.69
2:B:66:PRO:HG2	2:B:69:LEU:HD12	1.75	0.69
2:F:64:MET:HE2	2:F:124:ALA:HB2	1.75	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:65:HIS:CB	2:D:66:PRO:HD3	2.22	0.68
2:N:64:MET:HE2	2:N:124:ALA:HB2	1.74	0.68
2:X:129:THR:HG21	2:X:142:SER:OG	1.93	0.68
2:J:77:LEU:HD12	2:J:112:GLU:HG3	1.75	0.68
2:T:1:MET:N	2:T:2:ARG:CA	2.56	0.68
2:B:34:GLU:HB3	2:B:240:TYR:HD1	1.59	0.68
1:K:42:ILE:HG12	1:K:56:VAL:HG22	1.76	0.68
2:N:12:LEU:HD21	2:N:184:TRP:HB2	1.75	0.67
2:B:77:LEU:HD12	2:B:112:GLU:HG3	1.77	0.67
2:J:243:PHE:HZ	2:J:246:GLU:HG2	1.60	0.67
2:V:77:LEU:HD12	2:V:112:GLU:HG3	1.75	0.67
2:X:66:PRO:HG2	2:X:69:LEU:HD12	1.77	0.67
2:L:65:HIS:CB	2:L:66:PRO:HD3	2.23	0.67
1:I:112:ARG:NH2	2:J:99:ARG:HB2	2.10	0.67
2:V:65:HIS:CB	2:V:66:PRO:HD3	2.24	0.67
2:N:66:PRO:HG2	2:N:69:LEU:HD12	1.77	0.67
2:H:34:GLU:HB3	2:H:240:TYR:HD1	1.58	0.66
1:M:42:ILE:HG12	1:M:56:VAL:HG22	1.76	0.66
2:N:129:THR:HG21	2:N:142:SER:OG	1.95	0.66
2:P:65:HIS:CB	2:P:66:PRO:HD3	2.25	0.66
1:G:42:ILE:HG12	1:G:56:VAL:HG22	1.78	0.66
1:U:41:SER:HB2	1:U:57:LYS:HB2	1.78	0.66
2:B:136:ALA:HA	3:B:404:ADP:O3'	1.95	0.66
1:E:243:LYS:CE	2:F:108:ARG:HH21	2.08	0.66
2:L:66:PRO:HG2	2:L:69:LEU:HD12	1.76	0.66
2:H:227:ILE:HD13	2:H:248:VAL:HG22	1.78	0.66
2:F:65:HIS:CB	2:F:66:PRO:HD3	2.25	0.66
2:V:168:LYS:HD2	2:V:213:GLU:OE1	1.96	0.66
1:K:112:ARG:NH2	2:L:99:ARG:HB2	2.10	0.65
2:P:77:LEU:HD12	2:P:112:GLU:HG3	1.78	0.65
2:B:99:ARG:HH12	3:B:404:ADP:PA	2.20	0.65
2:R:66:PRO:HG2	2:R:69:LEU:HD12	1.79	0.65
2:X:77:LEU:HD12	2:X:112:GLU:HG3	1.79	0.65
2:B:78:ARG:HD3	2:B:126:ASP:OD1	1.96	0.65
2:V:64:MET:HE2	2:V:124:ALA:HB2	1.77	0.65
2:N:65:HIS:CB	2:N:66:PRO:HD3	2.27	0.65
2:H:66:PRO:HG2	2:H:69:LEU:HD12	1.79	0.65
1:Q:42:ILE:HG12	1:Q:56:VAL:HG22	1.79	0.65
2:J:65:HIS:CB	2:J:66:PRO:HD3	2.26	0.65
1:Q:51:ASP:OD1	1:Q:171:TYR:HE1	1.79	0.65
2:H:65:HIS:CB	2:H:66:PRO:HD3	2.26	0.65
2:H:129:THR:HG21	2:H:142:SER:OG	1.97	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:65:HIS:CB	2:B:66:PRO:HD3	2.26	0.64
2:X:65:HIS:CB	2:X:66:PRO:HD3	2.27	0.64
2:J:66:PRO:HG2	2:J:69:LEU:HD12	1.79	0.64
2:H:77:LEU:HD12	2:H:112:GLU:HG3	1.78	0.64
2:B:164:VAL:HG22	2:B:225:ILE:HG13	1.80	0.64
2:T:65:HIS:CB	2:T:66:PRO:HD3	2.28	0.64
2:L:19:ARG:HD2	2:L:181:GLU:OE2	1.98	0.64
2:R:164:VAL:HG22	2:R:225:ILE:HG13	1.80	0.64
2:P:66:PRO:HG2	2:P:69:LEU:HD12	1.79	0.63
2:V:78:ARG:HD3	2:V:126:ASP:OD1	1.99	0.63
2:V:129:THR:HG21	2:V:142:SER:OG	1.97	0.63
2:V:66:PRO:HG2	2:V:69:LEU:HD12	1.79	0.63
2:F:164:VAL:HG22	2:F:225:ILE:HG13	1.80	0.63
3:P:404:ADP:O2B	3:P:404:ADP:O2A	2.16	0.63
2:T:1:MET:O	2:T:1:MET:HG3	1.98	0.63
2:F:66:PRO:HG2	2:F:69:LEU:HD12	1.81	0.63
1:S:240:LYS:HB2	2:T:203:PHE:HB3	1.81	0.63
2:D:66:PRO:HG2	2:D:69:LEU:HD12	1.81	0.63
2:V:64:MET:HE1	2:V:76:VAL:HB	1.81	0.62
2:H:164:VAL:HG22	2:H:225:ILE:HG13	1.80	0.62
1:G:91:LEU:HD22	2:L:80:ARG:HH12	1.65	0.62
2:R:65:HIS:CB	2:R:66:PRO:HD3	2.29	0.62
2:B:129:THR:HG21	2:B:142:SER:OG	2.00	0.62
1:W:42:ILE:HG12	1:W:56:VAL:HG22	1.82	0.62
2:R:227:ILE:HD12	2:R:248:VAL:HG13	1.82	0.62
2:F:48:GLU:HG2	2:F:53:LYS:HG2	1.81	0.61
2:N:175:LEU:CD2	2:N:248:VAL:HG21	2.30	0.61
1:S:41:SER:HB2	1:S:57:LYS:HB2	1.81	0.61
2:J:164:VAL:HG22	2:J:225:ILE:HG13	1.82	0.61
2:T:129:THR:HG21	2:T:142:SER:OG	2.01	0.61
2:T:164:VAL:HG22	2:T:225:ILE:HG13	1.82	0.61
2:X:34:GLU:HB3	2:X:240:TYR:HD1	1.65	0.61
2:L:129:THR:HG21	2:L:142:SER:OG	2.00	0.61
2:X:73:ASP:OD2	2:X:74:ARG:HG3	2.00	0.61
1:O:41:SER:HB2	1:O:57:LYS:HB2	1.83	0.61
1:K:41:SER:HB2	1:K:57:LYS:HB2	1.83	0.61
1:M:41:SER:HB2	1:M:57:LYS:HB2	1.82	0.61
2:J:129:THR:HG21	2:J:142:SER:OG	2.01	0.61
2:B:138:SER:OG	3:B:404:ADP:O2B	2.18	0.60
2:V:179:GLU:O	2:V:183:MET:HB2	2.01	0.60
1:S:42:ILE:HG12	1:S:56:VAL:HG22	1.82	0.60
2:B:9:ARG:NH2	2:B:184:TRP:O	2.34	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:42:ILE:HG12	1:A:56:VAL:HG22	1.83	0.60
2:P:1:MET:HG3	2:P:3:GLU:H	1.67	0.60
1:C:41:SER:HB2	1:C:57:LYS:HB2	1.83	0.60
1:W:116:ARG:HD3	2:X:206:ASN:HB3	1.83	0.60
1:Q:71:ILE:HD12	1:Q:182:VAL:HG22	1.83	0.60
1:E:71:ILE:HD12	1:E:182:VAL:HG22	1.83	0.60
2:P:164:VAL:HG22	2:P:225:ILE:HG13	1.84	0.60
1:K:51:ASP:OD2	1:K:69:LEU:HB2	2.02	0.60
1:U:42:ILE:HG12	1:U:56:VAL:HG22	1.84	0.60
2:N:177:LEU:HD12	2:N:188:ASP:OD1	2.01	0.60
2:T:55:ILE:HG22	2:T:130:GLU:HB2	1.82	0.60
1:Q:41:SER:HB2	1:Q:57:LYS:HB2	1.82	0.59
1:G:148:ASP:OD1	2:L:40:ASN:HB2	2.02	0.59
1:U:112:ARG:NH2	2:V:99:ARG:HB2	2.17	0.59
1:I:42:ILE:HG12	1:I:56:VAL:HG22	1.84	0.59
2:V:164:VAL:HG22	2:V:225:ILE:HG13	1.84	0.59
1:W:41:SER:HB2	1:W:57:LYS:HB2	1.84	0.59
1:M:71:ILE:HD12	1:M:182:VAL:HG22	1.85	0.59
2:X:164:VAL:HG22	2:X:225:ILE:HG13	1.83	0.59
2:J:82:HIS:HE1	1:K:145:TYR:OH	1.86	0.59
1:O:42:ILE:HG12	1:O:56:VAL:HG22	1.85	0.59
1:I:14:ILE:HG22	2:V:116:LEU:HD22	1.85	0.59
1:C:42:ILE:HG12	1:C:56:VAL:HG22	1.85	0.59
2:L:164:VAL:HG22	2:L:225:ILE:HG13	1.85	0.59
2:N:130:GLU:HG3	1:Q:91:LEU:HD13	1.84	0.59
1:S:71:ILE:HD12	1:S:182:VAL:HG22	1.84	0.59
1:I:248:LEU:HD23	2:J:218:PHE:CZ	2.38	0.58
1:E:41:SER:HB2	1:E:57:LYS:HB2	1.86	0.58
2:F:64:MET:HE1	2:F:76:VAL:HB	1.84	0.58
2:N:164:VAL:HG22	2:N:225:ILE:HG13	1.83	0.58
1:G:41:SER:HB2	1:G:57:LYS:HB2	1.84	0.58
2:N:110:ALA:HB1	2:N:201:THR:CG2	2.34	0.58
2:L:73:ASP:OD2	2:L:74:ARG:HG3	2.04	0.58
2:L:234:GLU:HG3	2:L:241:VAL:CG2	2.33	0.58
2:H:48:GLU:HG2	2:H:53:LYS:CG	2.35	0.57
1:C:71:ILE:HD12	1:C:182:VAL:HG22	1.86	0.57
2:D:129:THR:HG21	2:D:142:SER:OG	2.03	0.57
2:P:110:ALA:HB1	2:P:201:THR:CG2	2.34	0.57
2:T:26:ASP:O	2:T:248:VAL:HB	2.03	0.57
2:R:129:THR:HG21	2:R:142:SER:OG	2.04	0.57
2:F:73:ASP:OD2	2:F:74:ARG:HG3	2.05	0.57
1:G:71:ILE:HD12	1:G:182:VAL:HG22	1.87	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:110:ALA:HB1	2:D:201:THR:CG2	2.35	0.57
1:E:14:ILE:HG21	2:N:118:GLU:OE1	2.04	0.57
2:N:129:THR:HG21	2:N:142:SER:HG	1.67	0.57
2:V:19:ARG:HD2	2:V:181:GLU:OE2	2.04	0.57
2:N:78:ARG:HD3	2:N:126:ASP:OD1	2.04	0.57
2:B:136:ALA:HA	3:B:404:ADP:HO3'	1.69	0.57
2:L:110:ALA:HB1	2:L:201:THR:CG2	2.34	0.57
2:T:19:ARG:HD2	2:T:181:GLU:OE2	2.04	0.56
2:H:89:ASP:N	2:H:89:ASP:OD2	2.27	0.56
1:E:42:ILE:HG12	1:E:56:VAL:HG22	1.86	0.56
2:R:73:ASP:OD2	2:R:74:ARG:HG3	2.06	0.56
1:A:71:ILE:HD12	1:A:182:VAL:HG22	1.87	0.56
2:T:110:ALA:HB1	2:T:201:THR:CG2	2.35	0.56
2:R:110:ALA:HB1	2:R:201:THR:CG2	2.35	0.56
1:W:71:ILE:HD12	1:W:182:VAL:HG22	1.88	0.56
2:X:110:ALA:HB1	2:X:201:THR:CG2	2.36	0.56
2:P:19:ARG:HD2	2:P:181:GLU:OE2	2.05	0.56
2:R:64:MET:HE1	2:R:76:VAL:HB	1.87	0.56
1:A:32:ARG:HB2	1:A:36:ASP:HB2	1.87	0.56
1:I:31:GLY:HA2	2:V:233:ARG:NH1	2.20	0.56
2:L:32:LYS:HB3	2:L:48:GLU:HB2	1.88	0.56
1:K:57:LYS:HG2	1:K:62:MET:HG2	1.88	0.56
1:S:255:GLU:OE2	2:T:215:ARG:NH2	2.39	0.56
2:P:129:THR:HG21	2:P:142:SER:OG	2.06	0.56
2:D:164:VAL:HG22	2:D:225:ILE:HG13	1.87	0.56
1:G:249:GLN:H	1:G:249:GLN:NE2	1.96	0.56
1:K:32:ARG:HB2	1:K:36:ASP:HB2	1.88	0.56
2:P:203:PHE:HE1	2:P:218:PHE:CE1	2.24	0.56
2:L:168:LYS:HE2	2:L:186:GLU:HB2	1.88	0.56
2:F:110:ALA:HB1	2:F:201:THR:CG2	2.36	0.56
1:O:57:LYS:HG2	1:O:62:MET:HG2	1.87	0.55
2:F:233:ARG:CZ	1:W:31:GLY:HA2	2.35	0.55
2:D:19:ARG:HD2	2:D:181:GLU:OE2	2.06	0.55
2:T:55:ILE:CG2	2:T:130:GLU:HB2	2.35	0.55
2:J:133:GLN:NE2	1:K:48:LYS:HB2	2.22	0.55
1:C:32:ARG:HB2	1:C:36:ASP:HB2	1.87	0.55
2:F:243:PHE:CZ	2:F:246:GLU:HG2	2.42	0.55
1:I:41:SER:HB2	1:I:57:LYS:HB2	1.88	0.55
2:H:227:ILE:CD1	2:H:248:VAL:HG13	2.35	0.55
1:K:6:SER:C	1:K:8:GLN:H	2.10	0.55
2:H:73:ASP:OD2	2:H:74:ARG:HG3	2.07	0.55
2:T:28:LEU:HG	2:T:248:VAL:HG21	1.89	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:71:ILE:HD12	1:I:182:VAL:HG22	1.89	0.55
2:H:135:ASP:O	2:H:136:ALA:HB3	2.07	0.55
2:P:73:ASP:OD2	2:P:74:ARG:HG3	2.07	0.55
2:H:88:THR:HG21	3:H:404:ADP:C2	2.43	0.54
2:J:86:PHE:HA	1:K:68:LYS:HD3	1.87	0.54
1:I:32:ARG:HB2	1:I:36:ASP:HB2	1.89	0.54
2:D:64:MET:HE1	2:D:76:VAL:HB	1.89	0.54
1:O:71:ILE:HD12	1:O:182:VAL:HG22	1.90	0.54
2:T:73:ASP:OD2	2:T:74:ARG:HG3	2.08	0.54
1:E:243:LYS:CE	2:F:108:ARG:NH2	2.70	0.54
1:K:71:ILE:HD12	1:K:182:VAL:HG22	1.89	0.54
1:U:32:ARG:HB2	1:U:36:ASP:HB2	1.89	0.54
2:H:110:ALA:HB1	2:H:201:THR:CG2	2.37	0.54
1:E:51:ASP:O	1:E:168:THR:HA	2.07	0.54
2:N:19:ARG:HD2	2:N:181:GLU:OE2	2.07	0.54
1:O:240:LYS:HB2	2:P:203:PHE:HB3	1.90	0.54
1:E:121:SER:O	1:E:122:LYS:HB3	2.07	0.54
1:C:240:LYS:HB2	2:D:203:PHE:HB3	1.89	0.53
2:H:80:ARG:NH1	2:H:130:GLU:OE2	2.42	0.53
2:V:73:ASP:OD2	2:V:74:ARG:HG3	2.08	0.53
1:U:71:ILE:HD12	1:U:182:VAL:HG22	1.90	0.53
2:T:243:PHE:CZ	2:T:246:GLU:HG2	2.41	0.53
2:H:227:ILE:CD1	2:H:248:VAL:HG22	2.38	0.53
1:U:57:LYS:HG2	1:U:62:MET:HG2	1.91	0.53
2:T:34:GLU:HG3	2:T:46:ILE:HB	1.90	0.53
1:O:32:ARG:HB2	1:O:36:ASP:HB2	1.91	0.53
2:P:48:GLU:HG2	2:P:53:LYS:HB3	1.91	0.53
2:J:64:MET:HE1	2:J:76:VAL:HB	1.91	0.53
1:G:32:ARG:HB2	1:G:36:ASP:HB2	1.91	0.53
2:D:223:LYS:HB3	2:D:248:VAL:CG1	2.39	0.53
2:B:110:ALA:HB1	2:B:201:THR:CG2	2.39	0.53
1:I:57:LYS:HG2	1:I:62:MET:HG2	1.91	0.53
1:E:32:ARG:HB2	1:E:36:ASP:HB2	1.90	0.53
2:P:159:ASP:OD2	2:P:196:SER:HB2	2.08	0.53
2:D:99:ARG:HH12	3:D:404:ADP:PA	2.29	0.53
2:X:223:LYS:HB3	2:X:248:VAL:CG1	2.38	0.53
2:D:233:ARG:HG3	2:D:233:ARG:NH1	2.24	0.53
2:B:51:ASN:HB2	2:B:135:ASP:OD2	2.09	0.53
2:R:212:ASP:N	2:R:212:ASP:OD1	2.39	0.53
1:E:243:LYS:HE3	2:F:108:ARG:HH21	1.73	0.52
1:M:32:ARG:HB2	1:M:36:ASP:HB2	1.90	0.52
1:G:91:LEU:H	1:G:92:PRO:CD	2.23	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:Q:57:LYS:HG2	1:Q:62:MET:HG2	1.90	0.52
2:T:42:ASP:O	2:T:153:ALA:HA	2.09	0.52
2:X:223:LYS:CB	2:X:248:VAL:HG11	2.38	0.52
2:J:91:ARG:HB2	1:K:68:LYS:HG2	1.91	0.52
2:H:131:ILE:HG12	2:H:138:SER:HB2	1.91	0.52
2:P:193:MET:HG3	2:P:225:ILE:HD13	1.92	0.52
1:E:125:ASP:HB3	1:E:128:LYS:HD2	1.91	0.52
1:W:32:ARG:HB2	1:W:36:ASP:HB2	1.91	0.52
2:H:192:ALA:HB3	2:H:202:LEU:HB3	1.92	0.52
2:F:78:ARG:HD3	2:F:126:ASP:OD1	2.10	0.52
2:H:129:THR:HG21	2:H:142:SER:HG	1.73	0.52
2:L:23:ARG:HH21	2:L:178:ASN:HD21	1.56	0.52
1:S:269:LEU:O	1:S:273:LEU:HG	2.10	0.52
1:C:112:ARG:NH1	2:D:102:GLU:OE1	2.42	0.52
2:L:193:MET:HG3	2:L:225:ILE:HD13	1.92	0.52
2:D:73:ASP:OD2	2:D:74:ARG:HG3	2.10	0.52
2:H:82:HIS:HE1	1:I:145:TYR:OH	1.92	0.52
1:E:112:ARG:NH2	2:F:99:ARG:HB2	2.25	0.52
2:L:131:ILE:HD11	2:L:142:SER:HB2	1.92	0.52
1:S:57:LYS:HG2	1:S:62:MET:HG2	1.92	0.51
2:B:193:MET:HG3	2:B:225:ILE:HD13	1.92	0.51
1:Q:32:ARG:HB2	1:Q:36:ASP:HB2	1.91	0.51
1:G:91:LEU:H	1:G:92:PRO:HD3	1.75	0.51
1:C:51:ASP:OD2	1:C:69:LEU:HB2	2.10	0.51
2:R:19:ARG:HD2	2:R:181:GLU:OE2	2.09	0.51
2:R:34:GLU:HB3	2:R:240:TYR:HD1	1.75	0.51
1:W:91:LEU:H	1:W:92:PRO:CD	2.23	0.51
1:A:57:LYS:HG2	1:A:62:MET:HG2	1.93	0.51
2:X:131:ILE:HG12	2:X:138:SER:HB2	1.93	0.51
1:S:32:ARG:HB2	1:S:36:ASP:HB2	1.91	0.51
2:J:133:GLN:OE1	1:K:48:LYS:HG2	2.10	0.51
1:S:91:LEU:H	1:S:92:PRO:CD	2.24	0.51
2:D:28:LEU:O	2:D:245:GLU:HB2	2.11	0.51
1:S:145:TYR:OH	2:V:82:HIS:HE1	1.92	0.51
2:X:135:ASP:O	2:X:136:ALA:HB3	2.11	0.51
2:P:190:PRO:HG2	2:P:204:GLN:HB2	1.92	0.51
2:P:42:ASP:O	2:P:153:ALA:HA	2.11	0.51
1:E:263:VAL:O	1:E:267:GLU:HG3	2.11	0.51
2:B:19:ARG:HD2	2:B:181:GLU:OE2	2.11	0.51
1:E:91:LEU:H	1:E:92:PRO:CD	2.23	0.51
2:T:136:ALA:N	3:T:404:ADP:O3'	2.36	0.50
2:F:131:ILE:HG12	2:F:138:SER:HB2	1.94	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:Q:240:LYS:HB2	2:R:203:PHE:HB3	1.93	0.50
2:L:135:ASP:OD1	2:L:179:GLU:HB2	2.11	0.50
1:G:30:ASP:OD1	1:G:31:GLY:N	2.44	0.50
2:B:73:ASP:OD2	2:B:74:ARG:HG3	2.10	0.50
1:A:91:LEU:H	1:A:92:PRO:HD3	1.76	0.50
2:R:168:LYS:HG3	2:R:173:ILE:HG13	1.92	0.50
2:F:19:ARG:HD2	2:F:181:GLU:OE2	2.11	0.50
1:K:91:LEU:H	1:K:92:PRO:CD	2.25	0.50
2:T:227:ILE:HD12	2:T:248:VAL:HG13	1.94	0.50
1:C:91:LEU:H	1:C:92:PRO:HD3	1.76	0.50
2:T:27:GLU:HG2	2:T:247:GLY:HA2	1.94	0.50
1:C:91:LEU:H	1:C:92:PRO:CD	2.24	0.50
2:J:110:ALA:HB1	2:J:201:THR:CG2	2.42	0.50
2:L:131:ILE:HG12	2:L:138:SER:HB2	1.92	0.50
2:P:131:ILE:HG12	2:P:138:SER:HB2	1.92	0.50
2:V:131:ILE:HD11	2:V:142:SER:HB2	1.93	0.50
1:E:91:LEU:H	1:E:92:PRO:HD3	1.76	0.50
2:V:110:ALA:HB1	2:V:201:THR:CG2	2.41	0.50
1:A:41:SER:HB2	1:A:57:LYS:HB2	1.92	0.49
2:R:2:ARG:HG2	2:R:3:GLU:H	1.77	0.49
1:E:243:LYS:HE2	2:F:108:ARG:HH21	1.74	0.49
1:C:57:LYS:HG2	1:C:62:MET:HG2	1.93	0.49
1:E:57:LYS:HG2	1:E:62:MET:HG2	1.94	0.49
1:O:91:LEU:H	1:O:92:PRO:CD	2.25	0.49
2:J:73:ASP:OD2	2:J:74:ARG:HG3	2.12	0.49
2:H:19:ARG:HD2	2:H:181:GLU:OE2	2.12	0.49
1:E:243:LYS:HE3	2:F:108:ARG:NH2	2.27	0.49
1:C:112:ARG:NH2	2:D:99:ARG:HB2	2.27	0.49
1:G:91:LEU:HD22	2:L:80:ARG:NH1	2.27	0.49
1:Q:91:LEU:H	1:Q:92:PRO:HD3	1.76	0.49
1:E:28:ARG:CZ	1:E:210:VAL:HG13	2.43	0.49
2:L:223:LYS:HB3	2:L:248:VAL:CG1	2.43	0.49
1:C:183:ASN:C	1:C:183:ASN:ND2	2.65	0.49
1:E:91:LEU:N	1:E:92:PRO:CD	2.76	0.49
2:P:192:ALA:HB3	2:P:202:LEU:HB3	1.95	0.49
2:P:1:MET:HG2	2:P:4:MET:HG3	1.95	0.49
2:J:27:GLU:HG2	2:J:247:GLY:HA2	1.95	0.49
2:T:131:ILE:HG12	2:T:138:SER:HB2	1.94	0.49
2:D:131:ILE:HD11	2:D:142:SER:HB2	1.94	0.49
1:C:68:LYS:HG2	2:F:91:ARG:HB2	1.94	0.49
2:V:27:GLU:HG2	2:V:247:GLY:HA2	1.95	0.49
1:A:91:LEU:H	1:A:92:PRO:CD	2.26	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:Q:129:LEU:HB3	1:Q:138:TRP:HB2	1.95	0.49
1:M:91:LEU:H	1:M:92:PRO:CD	2.25	0.49
2:D:131:ILE:HG12	2:D:138:SER:HB2	1.95	0.49
1:W:91:LEU:H	1:W:92:PRO:HD3	1.77	0.49
1:A:104:ASP:HB3	1:A:107:ALA:HB3	1.95	0.49
2:B:131:ILE:HG12	2:B:138:SER:HB2	1.95	0.49
1:E:38:ARG:HD2	1:E:59:GLY:HA3	1.95	0.49
1:I:240:LYS:HB2	2:J:203:PHE:HB3	1.95	0.49
1:O:121:SER:HB3	1:O:236:VAL:HG11	1.94	0.49
1:W:57:LYS:HG2	1:W:62:MET:HG2	1.94	0.48
1:M:91:LEU:H	1:M:92:PRO:HD3	1.78	0.48
2:F:168:LYS:HE2	2:F:186:GLU:HB2	1.93	0.48
2:B:169:ALA:O	2:B:171:GLY:N	2.42	0.48
2:D:233:ARG:NH2	1:G:31:GLY:HA2	2.28	0.48
1:C:27:ILE:HD11	2:T:233:ARG:HH12	1.78	0.48
1:K:112:ARG:NH1	2:L:102:GLU:OE1	2.46	0.48
2:X:129:THR:HG21	2:X:142:SER:HG	1.77	0.48
1:G:91:LEU:N	1:G:92:PRO:CD	2.76	0.48
2:H:27:GLU:HG2	2:H:247:GLY:HA2	1.96	0.48
2:N:190:PRO:HG2	2:N:204:GLN:HB2	1.94	0.48
1:S:121:SER:HB3	1:S:236:VAL:HG11	1.95	0.48
2:V:131:ILE:HG12	2:V:138:SER:HB2	1.96	0.48
1:W:91:LEU:N	1:W:92:PRO:CD	2.76	0.48
1:O:91:LEU:H	1:O:92:PRO:HD3	1.78	0.48
2:D:159:ASP:OD2	2:D:196:SER:HB2	2.13	0.48
2:N:234:GLU:HG3	2:N:241:VAL:HG21	1.94	0.48
2:X:192:ALA:HB3	2:X:202:LEU:HB3	1.96	0.48
2:N:131:ILE:HG12	2:N:138:SER:HB2	1.94	0.48
2:D:193:MET:HG3	2:D:225:ILE:HD13	1.94	0.48
2:B:175:LEU:HD22	2:B:248:VAL:CG2	2.43	0.48
2:T:190:PRO:HG2	2:T:204:GLN:HB2	1.96	0.48
2:X:103:LEU:HD21	2:X:139:ARG:CZ	2.43	0.48
1:O:163:ALA:HA	1:O:273:LEU:HD21	1.95	0.48
1:A:7:ASN:H	1:A:7:ASN:ND2	2.11	0.48
2:H:64:MET:HE1	2:H:76:VAL:HB	1.95	0.48
2:J:193:MET:HG3	2:J:225:ILE:HD13	1.95	0.48
2:R:27:GLU:HG2	2:R:247:GLY:HA2	1.95	0.48
2:T:192:ALA:HB3	2:T:202:LEU:HB3	1.96	0.48
2:F:131:ILE:HD11	2:F:142:SER:HB2	1.96	0.48
1:M:91:LEU:N	1:M:92:PRO:CD	2.77	0.48
1:U:91:LEU:H	1:U:92:PRO:HD3	1.78	0.48
1:S:68:LYS:HG2	2:V:91:ARG:HB2	1.94	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:N:73:ASP:OD2	2:N:74:ARG:HG3	2.14	0.48
2:N:12:LEU:HD22	2:N:181:GLU:HA	1.95	0.48
1:Q:91:LEU:H	1:Q:92:PRO:CD	2.26	0.48
2:D:233:ARG:HG3	2:D:233:ARG:HH11	1.79	0.48
1:A:166:TYR:HB3	1:A:273:LEU:HD22	1.96	0.48
2:P:183:MET:HE3	2:P:183:MET:HA	1.96	0.48
1:Q:112:ARG:NH2	2:R:99:ARG:HB2	2.29	0.47
1:S:251:ILE:HD12	2:T:200:VAL:HG21	1.96	0.47
2:V:37:VAL:HG11	2:V:46:ILE:HG13	1.96	0.47
2:N:202:LEU:HD23	2:N:203:PHE:N	2.29	0.47
2:N:234:GLU:HG3	2:N:241:VAL:CG2	2.44	0.47
1:W:121:SER:HB3	1:W:236:VAL:HG11	1.96	0.47
2:J:14:LEU:O	2:J:16:ASP:N	2.44	0.47
2:R:2:ARG:HG2	2:R:3:GLU:N	2.29	0.47
1:U:121:SER:O	1:U:122:LYS:HB2	2.14	0.47
1:U:104:ASP:HB3	1:U:107:ALA:HB3	1.97	0.47
1:I:104:ASP:HB3	1:I:107:ALA:HB3	1.96	0.47
2:V:99:ARG:HH12	3:V:404:ADP:PA	2.37	0.47
2:N:131:ILE:HD11	2:N:142:SER:HB2	1.95	0.47
2:T:28:LEU:O	2:T:245:GLU:HB3	2.15	0.47
2:P:135:ASP:O	2:P:136:ALA:HB3	2.14	0.47
2:V:83:MET:HB2	2:V:92:LYS:HD3	1.97	0.47
2:B:27:GLU:HG2	2:B:247:GLY:HA2	1.97	0.47
1:E:129:LEU:HB3	1:E:138:TRP:HB2	1.96	0.47
1:S:129:LEU:HB3	1:S:138:TRP:HB2	1.96	0.47
2:D:13:ILE:HD13	2:D:19:ARG:HG2	1.95	0.47
2:H:86:PHE:HA	1:I:68:LYS:HD3	1.97	0.47
1:A:121:SER:HB3	1:A:236:VAL:HG11	1.95	0.47
1:I:38:ARG:HD2	1:I:59:GLY:HA3	1.97	0.47
1:S:91:LEU:H	1:S:92:PRO:HD3	1.79	0.47
1:A:91:LEU:N	1:A:92:PRO:CD	2.77	0.47
2:F:27:GLU:HG2	2:F:247:GLY:HA2	1.95	0.47
2:J:190:PRO:HG2	2:J:204:GLN:HB2	1.96	0.47
1:C:38:ARG:HD2	1:C:59:GLY:HA3	1.97	0.47
2:V:32:LYS:HB3	2:V:48:GLU:HB2	1.97	0.47
1:W:183:ASN:C	1:W:183:ASN:ND2	2.66	0.47
2:F:26:ASP:HB2	2:F:248:VAL:HG12	1.95	0.47
2:H:131:ILE:HD11	2:H:142:SER:HB2	1.97	0.47
2:H:193:MET:HG3	2:H:225:ILE:HD13	1.97	0.47
1:I:121:SER:HB3	1:I:236:VAL:HG11	1.96	0.47
1:M:104:ASP:HB3	1:M:107:ALA:HB3	1.97	0.47
1:I:28:ARG:CZ	1:I:210:VAL:HG13	2.45	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:179:GLU:HG3	3:L:404:ADP:O2'	2.15	0.46
2:P:41:ALA:HA	2:P:59:TYR:CE2	2.50	0.46
1:Q:91:LEU:N	1:Q:92:PRO:CD	2.78	0.46
1:U:91:LEU:H	1:U:92:PRO:CD	2.28	0.46
1:O:77:ASP:O	1:O:79:PRO:HD3	2.15	0.46
2:N:27:GLU:HG2	2:N:247:GLY:HA2	1.98	0.46
2:B:34:GLU:HB3	2:B:240:TYR:CD1	2.46	0.46
1:S:91:LEU:N	1:S:92:PRO:CD	2.78	0.46
2:J:19:ARG:HD2	2:J:181:GLU:OE2	2.15	0.46
2:L:234:GLU:HG3	2:L:241:VAL:HG21	1.98	0.46
1:Q:38:ARG:HD2	1:Q:59:GLY:HA3	1.97	0.46
1:K:91:LEU:N	1:K:92:PRO:CD	2.79	0.46
2:P:227:ILE:HD12	2:P:248:VAL:HG22	1.97	0.46
1:G:91:LEU:CD2	2:L:80:ARG:HH12	2.28	0.46
2:V:193:MET:HG3	2:V:225:ILE:HD13	1.97	0.46
1:C:91:LEU:N	1:C:92:PRO:CD	2.78	0.46
2:J:192:ALA:HB3	2:J:202:LEU:HB3	1.96	0.46
1:W:129:LEU:HB3	1:W:138:TRP:HB2	1.97	0.46
2:J:196:SER:HB3	1:O:27:ILE:HG21	1.96	0.46
2:P:23:ARG:HH21	2:P:178:ASN:ND2	2.14	0.46
2:X:193:MET:HG3	2:X:225:ILE:HD13	1.98	0.46
2:R:131:ILE:HD11	2:R:142:SER:HB2	1.98	0.46
1:U:38:ARG:HD2	1:U:59:GLY:HA3	1.98	0.46
2:P:27:GLU:HG2	2:P:247:GLY:HA2	1.98	0.46
2:H:28:LEU:HG	2:H:248:VAL:HG21	1.98	0.46
2:D:202:LEU:HD23	2:D:203:PHE:N	2.31	0.46
1:K:91:LEU:H	1:K:92:PRO:HD3	1.80	0.46
2:R:192:ALA:HB3	2:R:202:LEU:HB3	1.97	0.46
1:E:104:ASP:HB3	1:E:107:ALA:HB3	1.98	0.46
2:L:64:MET:HE1	2:L:76:VAL:HB	1.98	0.46
1:E:32:ARG:CZ	1:E:38:ARG:HG3	2.46	0.46
2:H:202:LEU:HD23	2:H:203:PHE:N	2.31	0.46
2:R:190:PRO:HG2	2:R:204:GLN:HB2	1.97	0.46
2:J:160:LEU:HD11	2:J:236:LEU:HD22	1.98	0.46
1:C:28:ARG:CZ	1:C:210:VAL:HG13	2.46	0.46
2:P:16:ASP:C	2:P:18:LYS:H	2.19	0.46
1:M:57:LYS:HG2	1:M:62:MET:HG2	1.97	0.45
1:G:57:LYS:HG2	1:G:62:MET:HG2	1.98	0.45
2:F:192:ALA:HB3	2:F:202:LEU:HB3	1.98	0.45
2:P:64:MET:HE1	2:P:76:VAL:HB	1.98	0.45
2:D:65:HIS:CB	2:D:66:PRO:CD	2.94	0.45
1:O:91:LEU:N	1:O:92:PRO:CD	2.79	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:X:27:GLU:HG2	2:X:247:GLY:HA2	1.97	0.45
1:W:240:LYS:HB2	2:X:203:PHE:HB3	1.98	0.45
1:W:77:ASP:O	1:W:79:PRO:HD3	2.16	0.45
2:P:37:VAL:HG11	2:P:46:ILE:HG13	1.98	0.45
2:V:168:LYS:HE3	2:V:171:GLY:HA2	1.97	0.45
1:K:243:LYS:HD3	2:L:74:ARG:HH12	1.81	0.45
1:I:125:ASP:OD2	1:I:128:LYS:HE3	2.16	0.45
1:C:269:LEU:O	1:C:273:LEU:HG	2.17	0.45
1:K:3:SER:HB3	2:L:78:ARG:HG3	1.97	0.45
2:N:64:MET:HE1	2:N:76:VAL:HB	1.97	0.45
1:W:38:ARG:HD2	1:W:59:GLY:HA3	1.99	0.45
2:V:48:GLU:OE1	2:V:53:LYS:HE2	2.17	0.45
2:V:160:LEU:HD11	2:V:236:LEU:HD22	1.99	0.45
2:J:233:ARG:NH1	1:O:31:GLY:HA2	2.31	0.45
1:Q:8:GLN:HG2	1:Q:8:GLN:O	2.16	0.45
2:L:23:ARG:HH12	2:L:176:ASP:HB3	1.82	0.45
1:I:91:LEU:H	1:I:92:PRO:HD3	1.82	0.45
1:Q:121:SER:HB3	1:Q:236:VAL:HG11	1.97	0.45
1:C:75:TYR:CD1	1:C:75:TYR:N	2.85	0.45
1:A:129:LEU:HB3	1:A:138:TRP:HB2	1.99	0.45
2:J:131:ILE:HG12	2:J:138:SER:HB2	1.97	0.45
1:I:75:TYR:CE2	1:I:83:ASN:OD1	2.70	0.45
1:K:77:ASP:O	1:K:79:PRO:HD3	2.16	0.45
2:B:192:ALA:HB3	2:B:202:LEU:HB3	1.99	0.45
2:T:213:GLU:O	2:T:214:PHE:C	2.54	0.45
3:D:404:ADP:H5'2	3:D:404:ADP:O3B	2.16	0.45
2:R:131:ILE:HG12	2:R:138:SER:HB2	1.99	0.45
1:A:32:ARG:CZ	1:A:38:ARG:HG3	2.47	0.45
2:D:192:ALA:HB3	2:D:202:LEU:HB3	1.99	0.45
1:I:91:LEU:H	1:I:92:PRO:CD	2.30	0.45
2:R:195:PRO:HG2	2:R:229:TYR:CD1	2.52	0.45
2:X:61:PRO:HD3	2:X:155:ILE:HD11	1.99	0.45
2:F:210:THR:HG23	2:F:213:GLU:OE1	2.16	0.45
2:B:131:ILE:HD11	2:B:142:SER:HB2	1.99	0.45
2:J:131:ILE:HD11	2:J:142:SER:HB2	1.99	0.45
1:E:14:ILE:HG22	2:N:116:LEU:HD22	1.99	0.45
2:X:19:ARG:HD2	2:X:181:GLU:OE2	2.16	0.45
2:T:32:LYS:HB3	2:T:48:GLU:HB2	1.98	0.44
2:P:15:ASP:C	2:P:16:ASP:O	2.55	0.44
1:C:129:LEU:HB3	1:C:138:TRP:HB2	1.99	0.44
1:Q:104:ASP:CG	1:Q:105:GLU:N	2.71	0.44
2:J:37:VAL:HG11	2:J:46:ILE:HG13	1.99	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:91:ARG:HB2	1:I:68:LYS:HG2	1.99	0.44
1:I:130:VAL:HA	1:I:137:VAL:HG12	1.98	0.44
1:G:104:ASP:HB3	1:G:107:ALA:HB3	1.99	0.44
2:V:190:PRO:HG2	2:V:204:GLN:HB2	1.99	0.44
2:J:86:PHE:HB2	1:K:50:ALA:HB2	1.98	0.44
1:E:121:SER:HB3	1:E:236:VAL:HG11	1.98	0.44
2:T:37:VAL:HG11	2:T:46:ILE:HG13	1.99	0.44
1:U:91:LEU:N	1:U:92:PRO:CD	2.79	0.44
2:B:91:ARG:HB2	1:E:68:LYS:HG2	2.00	0.44
1:O:130:VAL:HA	1:O:137:VAL:HG12	1.99	0.44
1:W:104:ASP:HB3	1:W:107:ALA:HB3	1.99	0.44
1:K:130:VAL:HA	1:K:137:VAL:HG12	1.99	0.44
2:L:27:GLU:HG2	2:L:247:GLY:HA2	2.00	0.44
2:R:62:LYS:NZ	2:R:126:ASP:OD2	2.48	0.44
2:B:99:ARG:NH1	3:B:404:ADP:O2A	2.50	0.44
2:L:41:ALA:HA	2:L:59:TYR:CE2	2.53	0.44
1:S:16:LYS:HG3	1:S:220:ILE:HB	2.00	0.44
1:M:130:VAL:HA	1:M:137:VAL:HG12	1.99	0.44
2:H:48:GLU:CG	2:H:53:LYS:HG2	2.41	0.44
1:O:32:ARG:CZ	1:O:38:ARG:HG3	2.47	0.44
1:W:32:ARG:CZ	1:W:38:ARG:HG3	2.48	0.44
2:B:190:PRO:HG2	2:B:204:GLN:HB2	1.99	0.44
1:S:28:ARG:CZ	1:S:210:VAL:HG13	2.48	0.44
1:W:264:LYS:HG3	1:W:264:LYS:H	1.66	0.44
2:J:39:LYS:HA	2:J:39:LYS:HD2	1.85	0.44
2:X:131:ILE:HD11	2:X:142:SER:HB2	2.00	0.44
1:E:104:ASP:CG	1:E:105:GLU:N	2.71	0.44
2:L:192:ALA:HB3	2:L:202:LEU:HB3	1.99	0.44
2:F:159:ASP:OD2	2:F:196:SER:HB2	2.17	0.44
1:O:112:ARG:NH2	2:P:99:ARG:HB2	2.33	0.44
2:P:1:MET:HG3	2:P:3:GLU:N	2.30	0.44
1:U:104:ASP:CG	1:U:105:GLU:N	2.71	0.44
2:D:160:LEU:HD11	2:D:236:LEU:HD22	2.00	0.44
1:K:28:ARG:CZ	1:K:210:VAL:HG13	2.48	0.44
1:G:240:LYS:HB2	2:H:203:PHE:HB3	2.00	0.44
1:S:38:ARG:HD2	1:S:59:GLY:HA3	1.99	0.44
1:U:129:LEU:HB3	1:U:138:TRP:HB2	2.00	0.44
2:H:227:ILE:HD12	2:H:248:VAL:CG1	2.41	0.44
2:T:64:MET:HB2	2:T:64:MET:HE3	1.82	0.44
2:N:118:GLU:CD	2:N:118:GLU:H	2.20	0.44
1:A:240:LYS:HB2	2:B:203:PHE:HB3	2.00	0.44
2:P:169:ALA:HB2	2:P:174:ILE:HD13	2.00	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:104:ASP:CG	1:I:105:GLU:N	2.71	0.43
1:E:130:VAL:HA	1:E:137:VAL:HG12	2.00	0.43
2:F:135:ASP:O	2:F:136:ALA:HB3	2.18	0.43
2:T:165:ALA:HA	2:T:190:PRO:HA	2.00	0.43
2:L:165:ALA:HA	2:L:190:PRO:HA	1.99	0.43
2:N:13:ILE:CD1	2:N:172:VAL:HB	2.48	0.43
1:I:3:SER:HB3	2:J:78:ARG:CG	2.47	0.43
1:K:27:ILE:HG21	2:R:196:SER:HB3	2.00	0.43
2:V:12:LEU:HD22	2:V:181:GLU:HG3	2.00	0.43
1:C:121:SER:HB3	1:C:236:VAL:HG11	2.00	0.43
2:D:214:PHE:O	2:D:218:PHE:HB2	2.18	0.43
2:P:160:LEU:HD11	2:P:236:LEU:HD22	2.00	0.43
1:M:121:SER:HB3	1:M:236:VAL:HG11	2.01	0.43
2:B:41:ALA:HA	2:B:59:TYR:CE2	2.54	0.43
1:G:203:ALA:HB2	1:G:213:PRO:HG3	1.99	0.43
1:U:13:ILE:H	1:U:13:ILE:HG13	1.61	0.43
1:O:38:ARG:HD2	1:O:59:GLY:HA3	2.00	0.43
1:I:91:LEU:N	1:I:92:PRO:CD	2.81	0.43
1:U:80:ASN:OD1	1:U:133:PRO:HB3	2.18	0.43
2:J:103:LEU:HD21	2:J:139:ARG:CZ	2.48	0.43
2:F:41:ALA:HA	2:F:59:TYR:CE2	2.54	0.43
1:S:112:ARG:NH1	2:T:102:GLU:OE1	2.52	0.43
1:A:112:ARG:NH1	2:B:102:GLU:OE1	2.51	0.43
1:K:38:ARG:HD2	1:K:59:GLY:HA3	2.00	0.43
1:Q:104:ASP:HB3	1:Q:107:ALA:HB3	2.00	0.43
2:L:190:PRO:HG2	2:L:204:GLN:HB2	2.00	0.43
1:S:11:ILE:HA	1:S:12:PRO:HD3	1.93	0.43
1:Q:77:ASP:O	1:Q:79:PRO:HD3	2.18	0.43
2:X:190:PRO:HG2	2:X:204:GLN:HB2	1.99	0.43
2:B:202:LEU:HD23	2:B:203:PHE:N	2.33	0.43
1:G:121:SER:HB3	1:G:236:VAL:HG11	1.99	0.43
1:U:77:ASP:O	1:U:79:PRO:HD3	2.18	0.43
2:D:190:PRO:HG2	2:D:204:GLN:HB2	1.99	0.43
2:N:61:PRO:HD3	2:N:155:ILE:HD11	2.01	0.43
1:Q:125:ASP:CG	1:Q:128:LYS:HG3	2.39	0.43
1:A:130:VAL:HA	1:A:137:VAL:HG12	2.00	0.43
2:T:131:ILE:HD11	2:T:142:SER:HB2	2.01	0.43
2:T:160:LEU:HD11	2:T:236:LEU:HD22	2.00	0.43
1:G:145:TYR:OH	2:L:82:HIS:HE1	2.01	0.43
2:D:27:GLU:HG2	2:D:247:GLY:HA2	2.01	0.43
2:N:12:LEU:HD12	2:N:170:ASP:OD1	2.17	0.43
1:S:32:ARG:CZ	1:S:38:ARG:HG3	2.49	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:S:91:LEU:HD22	2:V:80:ARG:HH12	1.83	0.43
2:P:165:ALA:HA	2:P:190:PRO:HA	2.00	0.43
1:W:105:GLU:OE1	2:X:105:LYS:HG3	2.19	0.43
2:F:65:HIS:CB	2:F:66:PRO:CD	2.97	0.43
1:M:32:ARG:CZ	1:M:38:ARG:HG3	2.48	0.43
2:L:23:ARG:HH21	2:L:178:ASN:ND2	2.15	0.43
1:E:240:LYS:HB2	2:F:203:PHE:HB3	2.00	0.43
2:H:61:PRO:HD3	2:H:155:ILE:HD11	2.01	0.43
2:J:155:ILE:HA	2:J:156:PRO:HD3	1.88	0.43
1:G:130:VAL:HA	1:G:137:VAL:HG12	2.00	0.43
2:N:193:MET:HG3	2:N:225:ILE:HD13	2.00	0.43
2:P:223:LYS:HB3	2:P:248:VAL:CG1	2.48	0.43
2:V:155:ILE:HA	2:V:156:PRO:HD3	1.88	0.43
1:C:104:ASP:HB3	1:C:107:ALA:HB3	2.01	0.43
1:U:112:ARG:NH1	2:V:102:GLU:OE1	2.49	0.42
2:V:165:ALA:HA	2:V:190:PRO:HA	2.01	0.42
1:C:130:VAL:HA	1:C:137:VAL:HG12	2.01	0.42
1:S:104:ASP:HB3	1:S:107:ALA:HB3	2.01	0.42
2:L:65:HIS:CB	2:L:66:PRO:CD	2.95	0.42
1:I:32:ARG:CZ	1:I:38:ARG:HG3	2.48	0.42
2:H:160:LEU:HD11	2:H:236:LEU:HD22	2.00	0.42
1:U:11:ILE:HA	1:U:12:PRO:HD3	1.94	0.42
1:S:77:ASP:O	1:S:79:PRO:HD3	2.19	0.42
2:D:135:ASP:O	2:D:136:ALA:HB3	2.20	0.42
2:X:34:GLU:HB3	2:X:240:TYR:CD1	2.50	0.42
1:U:32:ARG:CZ	1:U:38:ARG:HG3	2.49	0.42
1:U:121:SER:HB3	1:U:236:VAL:HG11	2.00	0.42
2:B:82:HIS:HB3	2:B:130:GLU:OE2	2.19	0.42
2:V:115:VAL:CG2	2:V:150:LEU:HD13	2.49	0.42
2:R:41:ALA:HA	2:R:59:TYR:CE2	2.54	0.42
2:H:165:ALA:HA	2:H:190:PRO:HA	2.02	0.42
2:B:159:ASP:OD2	2:B:196:SER:HB2	2.19	0.42
2:X:160:LEU:HD11	2:X:236:LEU:HD22	2.02	0.42
1:M:28:ARG:CZ	1:M:210:VAL:HG13	2.49	0.42
1:A:13:ILE:HG13	1:A:13:ILE:H	1.69	0.42
1:G:60:THR:HB	2:L:40:ASN:ND2	2.35	0.42
2:P:131:ILE:HD11	2:P:142:SER:HB2	2.01	0.42
2:V:202:LEU:HD23	2:V:203:PHE:N	2.34	0.42
1:M:11:ILE:HA	1:M:12:PRO:HD3	1.92	0.42
1:E:16:LYS:HG3	1:E:220:ILE:HB	2.01	0.42
2:N:159:ASP:OD2	2:N:196:SER:HB2	2.19	0.42
2:D:136:ALA:HA	3:D:404:ADP:H5'2	2.02	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:224:GLY:HA2	2:H:248:VAL:HG11	2.01	0.42
2:F:129:THR:HG21	2:F:142:SER:HG	1.80	0.42
2:F:19:ARG:HD3	2:F:174:ILE:HG21	2.01	0.42
2:N:192:ALA:HB3	2:N:202:LEU:HB3	2.01	0.42
1:C:77:ASP:O	1:C:79:PRO:HD3	2.19	0.42
1:O:174:GLU:O	1:O:181:SER:N	2.47	0.42
1:K:13:ILE:H	1:K:13:ILE:HG13	1.67	0.42
3:N:404:ADP:O3B	3:N:404:ADP:O3'	2.30	0.42
1:M:112:ARG:NH2	2:N:99:ARG:HB2	2.34	0.42
1:G:60:THR:HB	2:L:40:ASN:HD21	1.83	0.42
2:P:23:ARG:HH12	2:P:176:ASP:HB3	1.85	0.42
2:B:82:HIS:HE1	1:E:145:TYR:OH	2.02	0.42
2:F:190:PRO:HG2	2:F:204:GLN:HB2	2.02	0.42
1:U:130:VAL:HA	1:U:137:VAL:HG12	2.00	0.42
1:G:13:ILE:O	1:G:17:GLU:HG3	2.20	0.42
2:N:41:ALA:HA	2:N:59:TYR:CE2	2.55	0.42
2:L:159:ASP:OD1	2:L:160:LEU:N	2.53	0.42
1:Q:112:ARG:NH1	2:R:102:GLU:OE1	2.53	0.42
2:L:99:ARG:HH12	3:L:404:ADP:PA	2.40	0.42
2:R:193:MET:HG3	2:R:225:ILE:HD13	2.00	0.42
2:D:233:ARG:CG	2:D:233:ARG:HH11	2.33	0.42
1:G:32:ARG:CZ	1:G:38:ARG:HG3	2.50	0.42
2:F:160:LEU:HD11	2:F:236:LEU:HD22	2.01	0.42
1:U:28:ARG:CZ	1:U:210:VAL:HG13	2.50	0.42
2:B:227:ILE:HD12	2:B:248:VAL:HG13	2.01	0.42
1:G:30:ASP:OD1	1:G:32:ARG:HG2	2.20	0.42
2:D:227:ILE:HD12	2:D:248:VAL:HG22	2.01	0.42
2:R:2:ARG:CD	2:R:4:MET:HG2	2.50	0.42
2:N:165:ALA:HA	2:N:190:PRO:HA	2.01	0.42
1:Q:28:ARG:CZ	1:Q:210:VAL:HG13	2.50	0.42
1:E:77:ASP:O	1:E:79:PRO:HD3	2.20	0.42
1:K:88:VAL:HG12	1:K:89:GLU:N	2.35	0.42
1:S:247:SER:HA	2:T:198:ASN:O	2.20	0.42
2:D:9:ARG:HG3	2:D:184:TRP:HB3	2.01	0.42
1:U:3:SER:HB3	2:V:78:ARG:CG	2.50	0.42
2:F:165:ALA:HA	2:F:190:PRO:HA	2.02	0.42
2:P:78:ARG:HD3	2:P:126:ASP:OD1	2.20	0.42
2:D:83:MET:HE2	2:D:99:ARG:HH21	1.84	0.41
2:R:129:THR:O	2:R:129:THR:HG22	2.20	0.41
1:M:38:ARG:HD2	1:M:59:GLY:HA3	2.02	0.41
2:R:3:GLU:CB	2:R:5:LEU:N	2.83	0.41
1:A:166:TYR:HB2	1:A:273:LEU:HD21	2.01	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:75:TYR:HD1	1:C:75:TYR:N	2.17	0.41
2:L:159:ASP:OD2	2:L:196:SER:HB2	2.19	0.41
1:I:129:LEU:HB3	1:I:138:TRP:HB2	2.02	0.41
2:J:132:LEU:HD22	1:K:64:LEU:CD2	2.50	0.41
2:R:103:LEU:HD21	2:R:139:ARG:CZ	2.50	0.41
1:W:130:VAL:HA	1:W:137:VAL:HG12	2.02	0.41
2:J:243:PHE:CZ	2:J:246:GLU:HG2	2.47	0.41
2:T:193:MET:HG3	2:T:225:ILE:HD13	2.01	0.41
2:D:129:THR:O	2:D:129:THR:HG22	2.20	0.41
2:R:16:ASP:HB3	2:R:17:GLY:H	1.76	0.41
2:D:42:ASP:O	2:D:153:ALA:HA	2.19	0.41
1:O:16:LYS:HG3	1:O:220:ILE:HB	2.02	0.41
1:O:11:ILE:HA	1:O:12:PRO:HD3	1.92	0.41
1:C:13:ILE:H	1:C:13:ILE:HG13	1.69	0.41
2:D:37:VAL:HG11	2:D:46:ILE:HG13	2.02	0.41
2:P:99:ARG:NH2	3:P:404:ADP:O2B	2.53	0.41
2:F:48:GLU:CG	2:F:53:LYS:HG2	2.48	0.41
1:O:28:ARG:CZ	1:O:210:VAL:HG13	2.50	0.41
2:R:37:VAL:HG11	2:R:46:ILE:HG13	2.02	0.41
2:B:37:VAL:HG11	2:B:46:ILE:HG13	2.02	0.41
2:R:160:LEU:HD11	2:R:236:LEU:HD22	2.03	0.41
1:M:104:ASP:CG	1:M:105:GLU:N	2.73	0.41
2:H:59:TYR:CE1	1:I:92:PRO:HA	2.55	0.41
2:L:202:LEU:HD23	2:L:203:PHE:N	2.35	0.41
1:I:203:ALA:HB2	1:I:213:PRO:HG3	2.03	0.41
2:X:37:VAL:HG11	2:X:46:ILE:HG13	2.02	0.41
1:U:16:LYS:HG3	1:U:220:ILE:HB	2.03	0.41
2:J:65:HIS:CB	2:J:66:PRO:CD	2.97	0.41
2:H:59:TYR:CZ	1:I:92:PRO:HA	2.55	0.41
1:G:104:ASP:CG	1:G:105:GLU:N	2.73	0.41
2:P:103:LEU:HD21	2:P:139:ARG:CZ	2.51	0.41
1:O:129:LEU:HB3	1:O:138:TRP:HB2	2.02	0.41
2:F:37:VAL:HG11	2:F:46:ILE:HG13	2.01	0.41
1:K:11:ILE:HA	1:K:12:PRO:HD3	1.94	0.41
2:B:165:ALA:HA	2:B:190:PRO:HA	2.03	0.41
1:C:16:LYS:HG3	1:C:220:ILE:HB	2.03	0.41
2:D:155:ILE:HA	2:D:156:PRO:HD3	1.88	0.41
2:F:155:ILE:HA	2:F:156:PRO:HD3	1.89	0.41
2:R:28:LEU:HG	2:R:248:VAL:HG21	2.03	0.41
2:T:32:LYS:HE2	2:T:34:GLU:OE2	2.20	0.41
2:H:195:PRO:HG2	2:H:229:TYR:CD1	2.55	0.41
1:G:11:ILE:HA	1:G:12:PRO:HD3	1.93	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:K:104:ASP:HB3	1:K:107:ALA:HB3	2.03	0.41
2:J:195:PRO:HG2	2:J:229:TYR:CD1	2.56	0.41
1:I:252:ASP:OD1	2:J:215:ARG:NH1	2.54	0.41
1:S:245:SER:HB2	2:T:199:GLN:HB3	2.02	0.41
1:M:77:ASP:O	1:M:79:PRO:HD3	2.21	0.41
1:C:183:ASN:ND2	1:C:185:ASN:OD1	2.38	0.41
1:M:240:LYS:HB2	2:N:203:PHE:HB3	2.03	0.41
2:P:169:ALA:HB2	2:P:174:ILE:CD1	2.51	0.41
2:N:159:ASP:OD1	2:N:160:LEU:N	2.54	0.41
2:D:1:MET:C	2:D:3:GLU:N	2.73	0.41
2:B:26:ASP:O	2:B:248:VAL:HG23	2.21	0.41
2:X:65:HIS:CB	2:X:66:PRO:CD	2.98	0.41
1:I:248:LEU:CD2	2:J:218:PHE:CZ	3.04	0.41
1:Q:32:ARG:CZ	1:Q:38:ARG:HG3	2.50	0.41
2:H:190:PRO:HG2	2:H:204:GLN:HB2	2.02	0.41
2:J:9:ARG:HG2	2:J:184:TRP:CE3	2.56	0.41
1:M:16:LYS:HG3	1:M:220:ILE:HB	2.03	0.41
1:S:203:ALA:HB2	1:S:213:PRO:HG3	2.02	0.41
2:D:30:SER:N	2:D:245:GLU:HB3	2.36	0.41
2:J:202:LEU:HD23	2:J:203:PHE:N	2.36	0.41
1:K:240:LYS:HB2	2:L:203:PHE:HB3	2.02	0.41
2:D:165:ALA:HA	2:D:190:PRO:HA	2.03	0.41
2:F:13:ILE:HD11	2:F:169:ALA:HB3	2.03	0.41
2:P:60:GLY:O	2:P:61:PRO:C	2.60	0.41
1:K:129:LEU:HB3	1:K:138:TRP:HB2	2.03	0.41
1:C:274:GLY:O	1:C:275:ILE:HB	2.21	0.41
2:H:116:LEU:HD12	2:H:156:PRO:HB2	2.03	0.41
1:E:269:LEU:O	1:E:273:LEU:HG	2.20	0.41
2:T:138:SER:OG	3:T:404:ADP:O3B	2.29	0.40
2:B:64:MET:HE1	2:B:76:VAL:HB	2.03	0.40
2:J:129:THR:HG21	2:J:142:SER:HG	1.87	0.40
2:L:160:LEU:HD11	2:L:236:LEU:HD22	2.03	0.40
2:T:103:LEU:HD21	2:T:139:ARG:CZ	2.51	0.40
1:W:75:TYR:HD2	1:W:75:TYR:N	2.20	0.40
2:V:42:ASP:CG	2:V:60:GLY:H	2.24	0.40
1:S:243:LYS:HG2	2:T:112:GLU:OE1	2.21	0.40
2:R:2:ARG:HD2	2:R:4:MET:HG2	2.02	0.40
2:J:165:ALA:HA	2:J:190:PRO:HA	2.03	0.40
2:D:41:ALA:HA	2:D:59:TYR:CE2	2.56	0.40
1:A:203:ALA:HB2	1:A:213:PRO:HG3	2.04	0.40
2:L:103:LEU:HD21	2:L:139:ARG:CZ	2.51	0.40
2:P:245:GLU:O	2:P:246:GLU:HB3	2.22	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:V:129:THR:HG21	2:V:142:SER:HG	1.83	0.40
1:K:32:ARG:CZ	1:K:38:ARG:HG3	2.51	0.40
2:N:160:LEU:HD11	2:N:236:LEU:HD22	2.04	0.40
2:P:115:VAL:CG2	2:P:150:LEU:HD13	2.51	0.40
1:U:183:ASN:O	1:U:185:ASN:N	2.52	0.40
2:F:193:MET:HG3	2:F:225:ILE:HD13	2.03	0.40
2:J:91:ARG:HD2	1:K:141:TRP:CG	2.57	0.40
1:K:247:SER:O	1:K:250:ASP:HB2	2.21	0.40
1:C:44:LEU:HD22	1:C:167:ASN:HB2	2.03	0.40
1:W:112:ARG:NH1	2:X:102:GLU:OE1	2.53	0.40
2:H:39:LYS:HG2	1:I:148:ASP:OD2	2.22	0.40
1:G:129:LEU:HB3	1:G:138:TRP:HB2	2.02	0.40
1:G:28:ARG:CZ	1:G:210:VAL:HG13	2.51	0.40
2:H:135:ASP:O	2:H:136:ALA:CB	2.70	0.40
2:F:202:LEU:HD23	2:F:203:PHE:N	2.37	0.40
1:W:44:LEU:HD22	1:W:167:ASN:HB2	2.03	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	254/275 (92%)	241 (95%)	10 (4%)	3 (1%)	19	71
1	C	254/275 (92%)	240 (94%)	8 (3%)	6 (2%)	9	53
1	E	254/275 (92%)	241 (95%)	11 (4%)	2 (1%)	27	78
1	G	254/275 (92%)	239 (94%)	11 (4%)	4 (2%)	14	64
1	I	254/275 (92%)	242 (95%)	10 (4%)	2 (1%)	27	78
1	K	254/275 (92%)	240 (94%)	11 (4%)	3 (1%)	19	71
1	M	254/275 (92%)	240 (94%)	11 (4%)	3 (1%)	19	71
1	O	254/275 (92%)	245 (96%)	7 (3%)	2 (1%)	27	78
1	Q	254/275 (92%)	238 (94%)	13 (5%)	3 (1%)	19	71

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	S	253/275 (92%)	242 (96%)	9 (4%)	2 (1%)	27	78
1	U	249/275 (90%)	236 (95%)	11 (4%)	2 (1%)	27	78
1	W	253/275 (92%)	241 (95%)	9 (4%)	3 (1%)	19	71
2	B	239/248 (96%)	221 (92%)	15 (6%)	3 (1%)	18	69
2	D	246/248 (99%)	231 (94%)	10 (4%)	5 (2%)	11	58
2	F	239/248 (96%)	226 (95%)	10 (4%)	3 (1%)	18	69
2	H	226/248 (91%)	216 (96%)	8 (4%)	2 (1%)	25	76
2	J	239/248 (96%)	225 (94%)	10 (4%)	4 (2%)	14	62
2	L	245/248 (99%)	231 (94%)	13 (5%)	1 (0%)	43	89
2	N	237/248 (96%)	224 (94%)	12 (5%)	1 (0%)	43	89
2	P	246/248 (99%)	230 (94%)	13 (5%)	3 (1%)	19	71
2	R	246/248 (99%)	230 (94%)	13 (5%)	3 (1%)	19	71
2	T	246/248 (99%)	233 (95%)	11 (4%)	2 (1%)	27	78
2	V	237/248 (96%)	218 (92%)	15 (6%)	4 (2%)	14	62
2	X	235/248 (95%)	223 (95%)	11 (5%)	1 (0%)	43	89
All	All	5922/6276 (94%)	5593 (94%)	262 (4%)	67 (1%)	21	73

All (67) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	8	GLN
2	J	15	ASP
2	P	16	ASP
1	Q	8	GLN
2	V	91	ARG
2	V	169	ALA
1	A	8	GLN
1	A	105	GLU
1	C	8	GLN
1	C	9	ASN
1	C	105	GLU
2	D	7	VAL
1	E	105	GLU
1	G	7	ASN
1	G	105	GLU
1	I	105	GLU
1	K	105	GLU

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Mol	Chain	Res	Type
1	M	105	GLU
1	O	105	GLU
1	Q	105	GLU
2	R	169	ALA
1	S	105	GLU
1	U	105	GLU
1	U	184	LYS
1	W	105	GLU
1	C	7	ASN
1	C	76	GLU
2	F	169	ALA
1	W	9	ASN
2	B	170	ASP
2	D	176	ASP
2	H	176	ASP
2	V	176	ASP
2	B	176	ASP
2	D	2	ARG
2	D	16	ASP
2	F	176	ASP
2	J	65	HIS
2	J	169	ALA
1	K	76	GLU
2	P	7	VAL
2	R	176	ASP
2	T	176	ASP
2	V	65	HIS
2	B	65	HIS
2	D	65	HIS
2	F	65	HIS
2	H	65	HIS
2	J	176	ASP
2	L	65	HIS
1	M	8	GLN
2	N	65	HIS
2	P	65	HIS
2	R	65	HIS
2	X	65	HIS
2	T	65	HIS
1	C	91	LEU
1	K	91	LEU
1	O	91	LEU

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Mol	Chain	Res	Type
1	S	91	LEU
1	W	91	LEU
1	E	91	LEU
1	G	91	LEU
1	Q	91	LEU
1	A	91	LEU
1	I	91	LEU
1	M	91	LEU

### 5.3.2 Protein sidechains ⓘ

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	212/242 (88%)	202 (95%)	10 (5%)	36	80
1	C	212/242 (88%)	204 (96%)	8 (4%)	44	84
1	E	218/242 (90%)	213 (98%)	5 (2%)	63	91
1	G	216/242 (89%)	206 (95%)	10 (5%)	37	80
1	I	214/242 (88%)	209 (98%)	5 (2%)	63	91
1	K	215/242 (89%)	204 (95%)	11 (5%)	33	77
1	M	216/242 (89%)	206 (95%)	10 (5%)	37	80
1	O	213/242 (88%)	205 (96%)	8 (4%)	44	84
1	Q	214/242 (88%)	208 (97%)	6 (3%)	56	89
1	S	216/242 (89%)	208 (96%)	8 (4%)	45	85
1	U	208/242 (86%)	199 (96%)	9 (4%)	40	82
1	W	214/242 (88%)	207 (97%)	7 (3%)	50	87
2	B	186/208 (89%)	176 (95%)	10 (5%)	31	76
2	D	200/208 (96%)	186 (93%)	14 (7%)	21	66
2	F	193/208 (93%)	182 (94%)	11 (6%)	29	74
2	H	179/208 (86%)	169 (94%)	10 (6%)	30	75
2	J	194/208 (93%)	184 (95%)	10 (5%)	32	76
2	L	199/208 (96%)	187 (94%)	12 (6%)	27	72

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	N	188/208 (90%)	182 (97%)	6 (3%)	51	88
2	P	201/208 (97%)	186 (92%)	15 (8%)	19	62
2	R	194/208 (93%)	184 (95%)	10 (5%)	32	76
2	T	196/208 (94%)	182 (93%)	14 (7%)	21	65
2	V	189/208 (91%)	182 (96%)	7 (4%)	45	85
2	X	191/208 (92%)	180 (94%)	11 (6%)	28	73
All	All	4878/5400 (90%)	4651 (95%)	227 (5%)	36	80

All (227) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	37	TYR
1	A	51	ASP
1	A	191	LEU
1	A	195	TYR
1	A	233	LEU
1	A	236	VAL
1	A	255	GLU
1	A	257	THR
1	A	275	ILE
2	B	12	LEU
2	B	13	ILE
2	B	20	THR
2	B	26	ASP
2	B	78	ARG
2	B	89	ASP
2	B	99	ARG
2	B	129	THR
2	B	164	VAL
2	B	201	THR
1	C	30	ASP
1	C	37	TYR
1	C	183	ASN
1	C	191	LEU
1	C	195	TYR
1	C	233	LEU
1	C	255	GLU
1	C	257	THR
2	D	1	MET

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Mol	Chain	Res	Type
2	D	2	ARG
2	D	7	VAL
2	D	8	GLU
2	D	20	THR
2	D	34	GLU
2	D	78	ARG
2	D	99	ARG
2	D	129	THR
2	D	164	VAL
2	D	201	THR
2	D	212	ASP
2	D	233	ARG
2	D	245	GLU
1	E	37	TYR
1	E	191	LEU
1	E	195	TYR
1	E	232	ASP
1	E	255	GLU
2	F	15	ASP
2	F	18	LYS
2	F	20	THR
2	F	34	GLU
2	F	39	LYS
2	F	78	ARG
2	F	99	ARG
2	F	129	THR
2	F	164	VAL
2	F	201	THR
2	F	218	PHE
1	G	37	TYR
1	G	45	ASP
1	G	191	LEU
1	G	195	TYR
1	G	206	ASP
1	G	233	LEU
1	G	236	VAL
1	G	249	GLN
1	G	255	GLU
1	G	257	THR
2	H	20	THR
2	H	39	LYS
2	H	78	ARG

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Mol	Chain	Res	Type
2	H	89	ASP
2	H	99	ARG
2	H	129	THR
2	H	164	VAL
2	H	178	ASN
2	H	201	THR
2	H	212	ASP
1	I	37	TYR
1	I	191	LEU
1	I	195	TYR
1	I	232	ASP
1	I	255	GLU
2	J	9	ARG
2	J	15	ASP
2	J	20	THR
2	J	39	LYS
2	J	78	ARG
2	J	99	ARG
2	J	129	THR
2	J	164	VAL
2	J	201	THR
2	J	233	ARG
1	K	24	GLU
1	K	30	ASP
1	K	37	TYR
1	K	45	ASP
1	K	191	LEU
1	K	195	TYR
1	K	232	ASP
1	K	249	GLN
1	K	255	GLU
1	K	257	THR
1	K	275	ILE
2	L	7	VAL
2	L	20	THR
2	L	78	ARG
2	L	99	ARG
2	L	129	THR
2	L	164	VAL
2	L	178	ASN
2	L	179	GLU
2	L	201	THR

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Mol	Chain	Res	Type
2	L	206	ASN
2	L	218	PHE
2	L	233	ARG
1	M	8	GLN
1	M	29	GLN
1	M	37	TYR
1	M	191	LEU
1	M	195	TYR
1	M	233	LEU
1	M	236	VAL
1	M	255	GLU
1	M	264	LYS
1	M	275	ILE
2	N	12	LEU
2	N	20	THR
2	N	78	ARG
2	N	129	THR
2	N	164	VAL
2	N	234	GLU
1	O	37	TYR
1	O	191	LEU
1	O	195	TYR
1	O	233	LEU
1	O	236	VAL
1	O	243	LYS
1	O	255	GLU
1	O	257	THR
2	P	1	MET
2	P	2	ARG
2	P	5	LEU
2	P	20	THR
2	P	34	GLU
2	P	53	LYS
2	P	78	ARG
2	P	99	ARG
2	P	129	THR
2	P	164	VAL
2	P	183	MET
2	P	201	THR
2	P	218	PHE
2	P	233	ARG
2	P	234	GLU

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Mol	Chain	Res	Type
1	Q	9	ASN
1	Q	37	TYR
1	Q	191	LEU
1	Q	195	TYR
1	Q	236	VAL
1	Q	255	GLU
2	R	1	MET
2	R	6	GLN
2	R	20	THR
2	R	78	ARG
2	R	99	ARG
2	R	129	THR
2	R	164	VAL
2	R	201	THR
2	R	212	ASP
2	R	218	PHE
1	S	18	SER
1	S	30	ASP
1	S	37	TYR
1	S	191	LEU
1	S	195	TYR
1	S	236	VAL
1	S	243	LYS
1	S	255	GLU
2	T	1	MET
2	T	2	ARG
2	T	5	LEU
2	T	20	THR
2	T	34	GLU
2	T	78	ARG
2	T	99	ARG
2	T	129	THR
2	T	164	VAL
2	T	183	MET
2	T	186	GLU
2	T	201	THR
2	T	218	PHE
2	T	246	GLU
1	U	8	GLN
1	U	29	GLN
1	U	37	TYR
1	U	191	LEU

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Mol	Chain	Res	Type
1	U	195	TYR
1	U	233	LEU
1	U	236	VAL
1	U	255	GLU
1	U	264	LYS
2	V	20	THR
2	V	78	ARG
2	V	129	THR
2	V	164	VAL
2	V	168	LYS
2	V	201	THR
2	V	218	PHE
1	W	37	TYR
1	W	183	ASN
1	W	191	LEU
1	W	195	TYR
1	W	236	VAL
1	W	255	GLU
1	W	257	THR
2	X	13	ILE
2	X	14	LEU
2	X	20	THR
2	X	78	ARG
2	X	89	ASP
2	X	99	ARG
2	X	129	THR
2	X	164	VAL
2	X	201	THR
2	X	218	PHE
2	X	246	GLU

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

Mol	Chain	Res	Type
1	A	7	ASN
2	B	178	ASN
2	D	178	ASN
2	F	82	HIS
1	G	249	GLN
2	H	178	ASN
1	I	29	GLN
2	J	82	HIS

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Mol	Chain	Res	Type
2	L	40	ASN
2	L	82	HIS
2	L	178	ASN
1	O	8	GLN
2	P	178	ASN
2	R	6	GLN
2	R	82	HIS
1	S	29	GLN
2	T	82	HIS
2	V	82	HIS
2	X	82	HIS

### 5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

12 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
3	ADP	B	404	-	29,29,29	1.21	3 (10%)	45,45,45	1.74	6 (13%)
3	ADP	D	404	2	29,29,29	1.16	2 (6%)	45,45,45	2.17	8 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	ADP	F	404	-	15,17,29	0.91	0	22,26,45	1.28	1 (4%)
3	ADP	H	404	-	29,29,29	1.05	1 (3%)	45,45,45	2.03	9 (20%)
3	ADP	J	404	-	8,8,29	1.06	0	13,13,45	0.99	1 (7%)
3	ADP	L	404	-	29,29,29	1.13	2 (6%)	45,45,45	2.06	9 (20%)
3	ADP	N	404	-	29,29,29	1.20	2 (6%)	45,45,45	1.82	7 (15%)
3	ADP	P	404	-	29,29,29	1.12	2 (6%)	45,45,45	2.14	10 (22%)
3	ADP	R	404	2	29,29,29	1.24	3 (10%)	45,45,45	1.83	9 (20%)
3	ADP	T	404	-	29,29,29	1.13	2 (6%)	45,45,45	1.87	10 (22%)
3	ADP	V	404	-	15,17,29	0.91	0	22,26,45	1.25	2 (9%)
3	ADP	X	404	-	15,17,29	0.89	0	22,26,45	1.25	2 (9%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ADP	B	404	-	-	0/16/32/32	0/1/3/3
3	ADP	D	404	2	-	0/16/32/32	0/1/3/3
3	ADP	F	404	-	-	0/12/25/32	0/1/1/3
3	ADP	H	404	-	-	0/16/32/32	0/1/3/3
3	ADP	J	404	-	-	0/6/6/32	0/0/0/3
3	ADP	L	404	-	-	0/16/32/32	0/1/3/3
3	ADP	N	404	-	-	0/16/32/32	0/1/3/3
3	ADP	P	404	-	-	0/16/32/32	0/1/3/3
3	ADP	R	404	2	-	0/16/32/32	0/1/3/3
3	ADP	T	404	-	-	0/16/32/32	0/1/3/3
3	ADP	V	404	-	-	0/12/25/32	0/1/1/3
3	ADP	X	404	-	-	0/12/25/32	0/1/1/3

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	404	ADP	C5-C4	3.56	1.48	1.40
3	R	404	ADP	C5-C4	3.51	1.48	1.40
3	N	404	ADP	C5-C4	3.43	1.48	1.40
3	L	404	ADP	C5-C4	3.41	1.48	1.40
3	T	404	ADP	C4-N9	-3.39	1.32	1.37
3	H	404	ADP	C5-C4	3.29	1.47	1.40
3	D	404	ADP	C5-C4	3.13	1.47	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	404	ADP	C5-C4	2.99	1.47	1.40
3	T	404	ADP	C5-C4	2.75	1.46	1.40
3	L	404	ADP	C4-N9	-2.72	1.33	1.37
3	D	404	ADP	C4-N9	-2.61	1.33	1.37
3	P	404	ADP	C4-N9	-2.49	1.34	1.37
3	R	404	ADP	C4-N9	-2.45	1.34	1.37
3	B	404	ADP	PB-O3A	2.12	1.63	1.60
3	N	404	ADP	O4'-C1'	2.07	1.44	1.41
3	R	404	ADP	O4'-C1'	2.03	1.44	1.41
3	B	404	ADP	C4-N9	-2.02	1.34	1.37

All (74) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	404	ADP	N3-C2-N1	-7.63	122.33	128.71
3	D	404	ADP	N3-C2-N1	-7.07	122.80	128.71
3	D	404	ADP	O4'-C1'-N9	7.03	114.98	108.44
3	N	404	ADP	N3-C2-N1	-6.77	123.05	128.71
3	H	404	ADP	N3-C2-N1	-6.62	123.17	128.71
3	B	404	ADP	N3-C2-N1	-6.46	123.31	128.71
3	T	404	ADP	N3-C2-N1	-6.09	123.62	128.71
3	L	404	ADP	O4'-C1'-N9	5.88	113.91	108.44
3	L	404	ADP	N3-C2-N1	-5.76	123.90	128.71
3	H	404	ADP	O4'-C1'-N9	5.72	113.77	108.44
3	P	404	ADP	N3-C4-N9	5.59	135.53	125.43
3	N	404	ADP	N3-C4-N9	5.58	135.51	125.43
3	R	404	ADP	N3-C2-N1	-5.58	124.04	128.71
3	B	404	ADP	N3-C4-N9	5.54	135.43	125.43
3	H	404	ADP	N3-C4-N9	5.51	135.37	125.43
3	T	404	ADP	N3-C4-N9	5.39	135.16	125.43
3	R	404	ADP	N3-C4-N9	5.00	134.46	125.43
3	R	404	ADP	O4'-C1'-N9	4.95	113.04	108.44
3	D	404	ADP	N3-C4-N9	4.88	134.25	125.43
3	L	404	ADP	N3-C4-N9	4.84	134.16	125.43
3	L	404	ADP	C4-C5-N7	-4.60	105.58	109.52
3	P	404	ADP	PA-O3A-PB	-4.14	119.53	131.68
3	F	404	ADP	C4'-C3'-C2'	3.99	106.61	101.77
3	N	404	ADP	O4'-C1'-N9	3.91	112.08	108.44
3	L	404	ADP	PA-O3A-PB	-3.88	120.29	131.68
3	D	404	ADP	C4-C5-N7	-3.59	106.45	109.52
3	P	404	ADP	C4-C5-N7	-3.50	106.53	109.52
3	T	404	ADP	C2'-C1'-N9	-3.50	104.29	113.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	404	ADP	C8-N9-C4	3.49	109.56	106.90
3	T	404	ADP	C8-N9-C4	3.44	109.53	106.90
3	B	404	ADP	C5-C4-N3	-3.42	118.25	125.70
3	P	404	ADP	C2'-C1'-N9	-3.35	104.66	113.27
3	H	404	ADP	C5-C4-N3	-3.34	118.42	125.70
3	V	404	ADP	C4'-C3'-C2'	3.32	105.80	101.77
3	N	404	ADP	C5-C4-N3	-3.31	118.49	125.70
3	H	404	ADP	C2'-C1'-N9	-3.26	104.89	113.27
3	B	404	ADP	C4-C5-N7	-3.23	106.75	109.52
3	R	404	ADP	C4-C5-N7	-3.20	106.78	109.52
3	L	404	ADP	C5-C4-N3	-3.18	118.77	125.70
3	R	404	ADP	C5-C4-N3	-3.13	118.89	125.70
3	P	404	ADP	C5-C4-N3	-3.12	118.90	125.70
3	T	404	ADP	PA-O3A-PB	-3.12	122.54	131.68
3	D	404	ADP	PA-O3A-PB	-3.12	122.54	131.68
3	D	404	ADP	C5-C4-N3	-3.02	119.12	125.70
3	X	404	ADP	PA-O3A-PB	-2.93	123.10	131.68
3	H	404	ADP	PA-O3A-PB	-2.88	123.24	131.68
3	T	404	ADP	C5-C4-N3	-2.83	119.54	125.70
3	H	404	ADP	C4-C5-N7	-2.76	107.16	109.52
3	B	404	ADP	O4'-C1'-N9	2.74	110.99	108.44
3	X	404	ADP	C4'-C3'-C2'	2.72	105.07	101.77
3	H	404	ADP	C2-N3-C4	2.70	121.69	114.01
3	N	404	ADP	C2-N3-C4	2.68	121.65	114.01
3	P	404	ADP	C2-N3-C4	2.66	121.58	114.01
3	R	404	ADP	O2'-C2'-C1'	-2.58	103.41	111.23
3	B	404	ADP	C2-N3-C4	2.55	121.26	114.01
3	J	404	ADP	O3B-PB-O1B	2.52	118.67	110.44
3	N	404	ADP	C4-C5-N7	-2.49	107.39	109.52
3	D	404	ADP	C2-N3-C4	2.48	121.06	114.01
3	L	404	ADP	O3B-PB-O2B	2.46	117.19	107.61
3	L	404	ADP	C2-N3-C4	2.43	120.92	114.01
3	D	404	ADP	C2'-C1'-N9	-2.36	107.21	113.27
3	R	404	ADP	C2-N3-C4	2.33	120.63	114.01
3	T	404	ADP	C4'-O4'-C1'	-2.29	107.26	109.75
3	T	404	ADP	O4'-C1'-N9	2.29	110.57	108.44
3	V	404	ADP	PA-O3A-PB	-2.26	125.04	131.68
3	P	404	ADP	O4'-C1'-N9	2.26	110.55	108.44
3	T	404	ADP	C2-N3-C4	2.16	120.17	114.01
3	T	404	ADP	C4-C5-N7	-2.15	107.68	109.52
3	H	404	ADP	C3'-C2'-C1'	2.14	104.26	100.91
3	N	404	ADP	PA-O3A-PB	-2.14	125.41	131.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	R	404	ADP	O3B-PB-O2B	2.11	115.84	107.61
3	L	404	ADP	C2'-C1'-N9	-2.09	107.89	113.27
3	P	404	ADP	O3B-PB-O2B	2.06	115.65	107.61
3	R	404	ADP	C4'-O4'-C1'	2.06	111.99	109.75

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	260/275 (94%)	0.29	0 100 100	46, 53, 67, 74	0
1	C	260/275 (94%)	0.36	0 100 100	46, 53, 67, 75	0
1	E	260/275 (94%)	0.30	0 100 100	46, 53, 67, 74	0
1	G	260/275 (94%)	0.34	0 100 100	46, 53, 67, 74	0
1	I	260/275 (94%)	0.35	3 (1%) 75 29	46, 53, 67, 73	0
1	K	260/275 (94%)	0.39	0 100 100	46, 53, 67, 75	0
1	M	260/275 (94%)	0.32	0 100 100	46, 53, 67, 75	0
1	O	260/275 (94%)	0.36	0 100 100	46, 53, 67, 75	0
1	Q	260/275 (94%)	0.34	0 100 100	46, 53, 67, 74	0
1	S	259/275 (94%)	0.38	2 (0%) 83 39	46, 53, 67, 73	0
1	U	255/275 (92%)	0.39	3 (1%) 75 29	46, 53, 67, 73	0
1	W	259/275 (94%)	0.38	0 100 100	46, 53, 67, 75	0
2	B	241/248 (97%)	0.40	2 (0%) 83 39	46, 52, 68, 80	0
2	D	248/248 (100%)	0.36	0 100 100	45, 52, 68, 80	0
2	F	241/248 (97%)	0.34	0 100 100	46, 52, 68, 80	0
2	H	230/248 (92%)	0.47	3 (1%) 74 27	46, 52, 69, 80	0
2	J	241/248 (97%)	0.43	2 (0%) 83 39	46, 52, 68, 80	0
2	L	247/248 (99%)	0.35	1 (0%) 90 57	45, 52, 68, 80	0
2	N	239/248 (96%)	0.34	0 100 100	46, 52, 69, 79	0
2	P	248/248 (100%)	0.38	0 100 100	45, 52, 68, 80	0
2	R	248/248 (100%)	0.34	0 100 100	45, 52, 69, 80	0
2	T	248/248 (100%)	0.35	0 100 100	45, 52, 69, 79	0
2	V	239/248 (96%)	0.51	8 (3%) 44 10	46, 52, 68, 79	0
2	X	239/248 (96%)	0.36	2 (0%) 83 39	46, 52, 69, 79	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
All	All	6022/6276 (95%)	0.37	26 (0%) 90 57	45, 52, 68, 80	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	U	171	TYR	3.3
2	V	132	LEU	3.0
1	S	90	LEU	2.8
1	U	1	MET	2.7
2	J	18	LYS	2.5
2	V	55	ILE	2.4
2	L	248	VAL	2.3
1	I	124	LEU	2.3
2	H	55	ILE	2.3
1	I	275	ILE	2.3
1	I	248	LEU	2.2
2	X	12	LEU	2.2
2	V	64	MET	2.2
2	V	209	MET	2.2
2	V	53	LYS	2.2
1	S	138	TRP	2.2
2	V	204	GLN	2.1
2	B	248	VAL	2.1
2	V	107	ILE	2.1
2	J	14	LEU	2.1
2	H	80	ARG	2.1
1	U	44	LEU	2.1
2	H	248	VAL	2.1
2	V	133	GLN	2.0
2	B	15	ASP	2.0
2	X	132	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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	ADP	V	404	17/27	0.28	-0.35	90,93,95,95	0
3	ADP	T	404	27/27	0.24	-0.48	55,60,63,63	0
3	ADP	D	404	27/27	0.25	-0.57	48,50,51,51	0
3	ADP	H	404	27/27	0.23	-0.62	116,117,119,119	0
3	ADP	R	404	27/27	0.22	-0.69	47,55,58,58	0
3	ADP	N	404	27/27	0.25	-0.80	59,62,64,64	0
3	ADP	L	404	27/27	0.24	-0.81	52,54,56,58	0
3	ADP	P	404	27/27	0.23	-0.94	33,39,41,42	0
3	ADP	B	404	27/27	0.20	-1.42	79,90,95,95	0
3	ADP	F	404	17/27	0.22	-1.46	91,94,95,95	0
3	ADP	X	404	17/27	0.20	-1.46	86,87,89,89	0
3	ADP	J	404	9/27	0.14	-2.35	100,100,101,101	0

## 6.5 Other polymers

There are no such residues in this entry.