Full wwPDB X-ray Structure Validation Report (i)

Oct 4, 2016 – 03:59 PM EDT

PDB ID	:	5D4C
Title	:	Crystal structure of Thermus thermophilus product complex for transcription initiation with ATP and CTP
Authors	:	Zhang, Y.; Ebright, R.H.
Deposited on	:	2015-08-07
Resolution	:	3.28 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

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

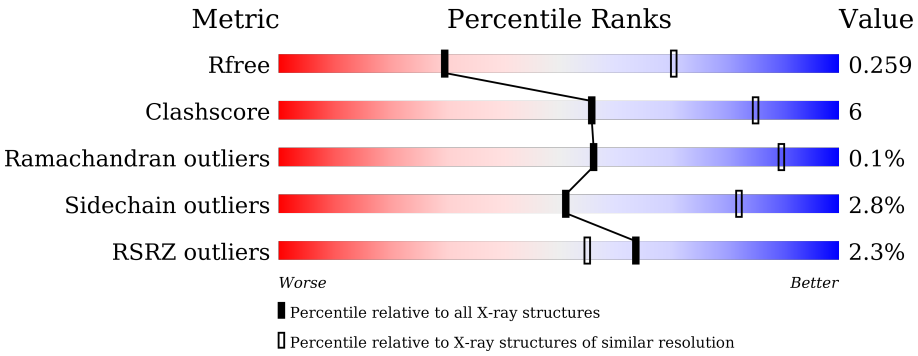
MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20027939
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20027939

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 3.28 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1756 (3.36-3.20)
Clashscore	102246	1941 (3.36-3.20)
Ramachandran outliers	100387	1905 (3.36-3.20)
Sidechain outliers	100360	1903 (3.36-3.20)
RSRZ outliers	91569	1764 (3.36-3.20)

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 >=3, 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	315	60% 13% 27%
1	B	315	59% 11% 30%
1	K	315	61% 12% 27%
1	L	315	58% 12% 30%
2	C	1119	82% 17%
2	M	1119	5% 76% 19%

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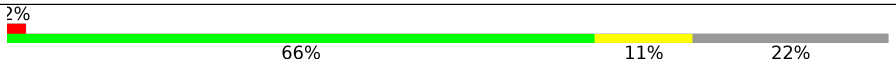
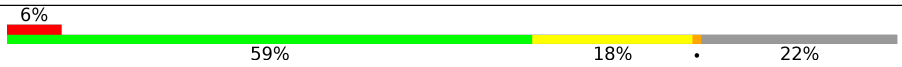
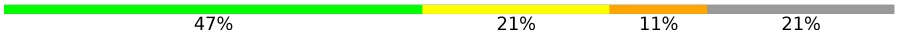
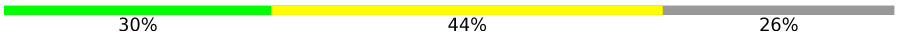
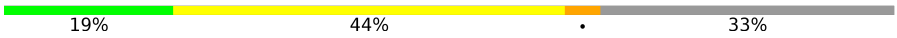
Mol	Chain	Length	Quality of chain
3	D	1524	
3	N	1524	
4	E	99	
4	O	99	
5	F	443	
5	P	443	
6	G	19	
6	R	19	
7	H	27	
7	S	27	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
11	CTP	D	2006	-	-	-	X
11	CTP	N	1606	-	-	-	X
8	MG	B	2001	-	-	-	X
8	MG	D	2007	-	-	-	X
8	MG	K	901	-	-	-	X
8	MG	N	1607	-	-	-	X

2 Entry composition

There are 13 unique types of molecules in this entry. The entry contains 56600 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 DNA-directed RNA polymerase subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	231	Total	C	N	O	S	0	0	0
			1809	1155	315	337	2			
1	B	222	Total	C	N	O	S	0	0	0
			1750	1120	303	325	2			
1	K	231	Total	C	N	O	S	0	0	0
			1809	1155	315	337	2			
1	L	222	Total	C	N	O	S	0	0	0
			1750	1120	303	325	2			

- Molecule 2 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	1112	Total	C	N	O	S	0	0	0
			8774	5550	1565	1635	24			
2	M	1080	Total	C	N	O	S	0	0	0
			8508	5375	1522	1587	24			

- Molecule 3 is a protein called DNA-directed RNA polymerase subunit beta'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	1486	Total	C	N	O	S	0	0	0
			11738	7441	2067	2195	35			
3	N	1486	Total	C	N	O	S	0	0	0
			11738	7441	2067	2195	35			

- Molecule 4 is a protein called DNA-directed RNA polymerase subunit omega.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	94	Total	C	N	O	S	0	0	0
			758	483	132	139	4			
4	O	94	Total	C	N	O	S	0	0	0
			758	483	132	139	4			

- Molecule 5 is a protein called RNA polymerase sigma factor SigA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	F	346	Total	C	N	O	S	0	0	0
			2807	1770	509	524	4			
5	P	346	Total	C	N	O	S	0	0	0
			2807	1770	509	524	4			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	-19	MET	-	initiating methionine	UNP Q5SKW1
F	-18	GLY	-	expression tag	UNP Q5SKW1
F	-17	SER	-	expression tag	UNP Q5SKW1
F	-16	SER	-	expression tag	UNP Q5SKW1
F	-15	HIS	-	expression tag	UNP Q5SKW1
F	-14	HIS	-	expression tag	UNP Q5SKW1
F	-13	HIS	-	expression tag	UNP Q5SKW1
F	-12	HIS	-	expression tag	UNP Q5SKW1
F	-11	HIS	-	expression tag	UNP Q5SKW1
F	-10	HIS	-	expression tag	UNP Q5SKW1
F	-9	SER	-	expression tag	UNP Q5SKW1
F	-8	SER	-	expression tag	UNP Q5SKW1
F	-7	GLY	-	expression tag	UNP Q5SKW1
F	-6	LEU	-	expression tag	UNP Q5SKW1
F	-5	VAL	-	expression tag	UNP Q5SKW1
F	-4	PRO	-	expression tag	UNP Q5SKW1
F	-3	ARG	-	expression tag	UNP Q5SKW1
F	-2	GLY	-	expression tag	UNP Q5SKW1
F	-1	SER	-	expression tag	UNP Q5SKW1
F	0	HIS	-	expression tag	UNP Q5SKW1
P	-19	MET	-	initiating methionine	UNP Q5SKW1
P	-18	GLY	-	expression tag	UNP Q5SKW1
P	-17	SER	-	expression tag	UNP Q5SKW1
P	-16	SER	-	expression tag	UNP Q5SKW1
P	-15	HIS	-	expression tag	UNP Q5SKW1
P	-14	HIS	-	expression tag	UNP Q5SKW1
P	-13	HIS	-	expression tag	UNP Q5SKW1
P	-12	HIS	-	expression tag	UNP Q5SKW1
P	-11	HIS	-	expression tag	UNP Q5SKW1
P	-10	HIS	-	expression tag	UNP Q5SKW1
P	-9	SER	-	expression tag	UNP Q5SKW1
P	-8	SER	-	expression tag	UNP Q5SKW1
P	-7	GLY	-	expression tag	UNP Q5SKW1
P	-6	LEU	-	expression tag	UNP Q5SKW1

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Chain	Residue	Modelled	Actual	Comment	Reference
P	-5	VAL	-	expression tag	UNP Q5SKW1
P	-4	PRO	-	expression tag	UNP Q5SKW1
P	-3	ARG	-	expression tag	UNP Q5SKW1
P	-2	GLY	-	expression tag	UNP Q5SKW1
P	-1	SER	-	expression tag	UNP Q5SKW1
P	0	HIS	-	expression tag	UNP Q5SKW1

- Molecule 6 is a DNA chain called DNA (5'-D(*CP*CP*T*GP*CP*AP*TP*CP*CP*GP*TP*GP*AP*GP*TP*AP*GP*AP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	G	15	Total	C	N	O	P	0	0	0
			308	147	60	87	14			
6	R	15	Total	C	N	O	P	0	0	0
			308	147	60	87	14			

- Molecule 7 is a DNA chain called DNA (27-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	H	20	Total	C	N	O	P	0	0	0
			414	197	82	116	19			
7	S	18	Total	C	N	O	P	0	0	0
			371	177	72	105	17			

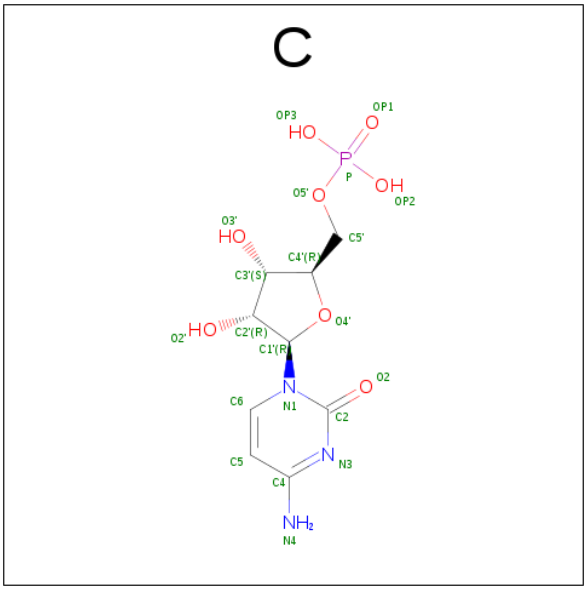
- Molecule 8 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	P	1	Total	Mg	0	0
			1	1		
8	D	3	Total	Mg	0	0
			3	3		
8	K	1	Total	Mg	0	0
			1	1		
8	B	1	Total	Mg	0	0
			1	1		
8	N	3	Total	Mg	0	0
			3	3		
8	L	1	Total	Mg	0	0
			1	1		
8	F	1	Total	Mg	0	0
			1	1		

- Molecule 9 is ZINC ION (three-letter code: ZN) (formula: Zn).

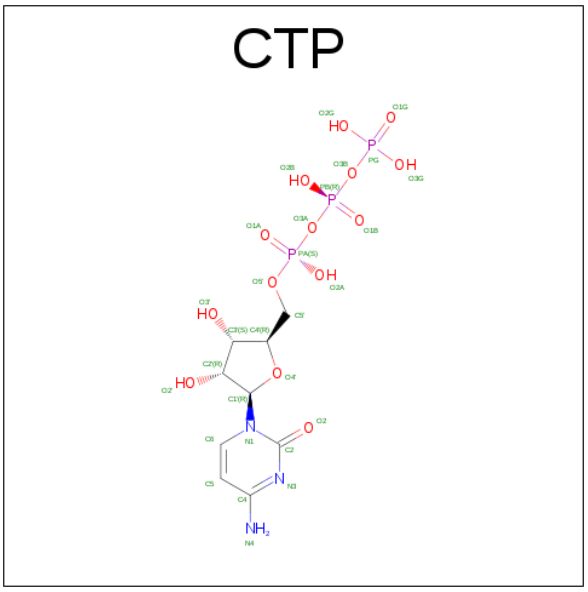
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	D	2	Total Zn 2 2	0	0
9	N	2	Total Zn 2 2	0	0

- Molecule 10 is CYTIDINE-5'-MONOPHOSPHATE (three-letter code: C) (formula: $C_9H_{14}N_3O_8P$).



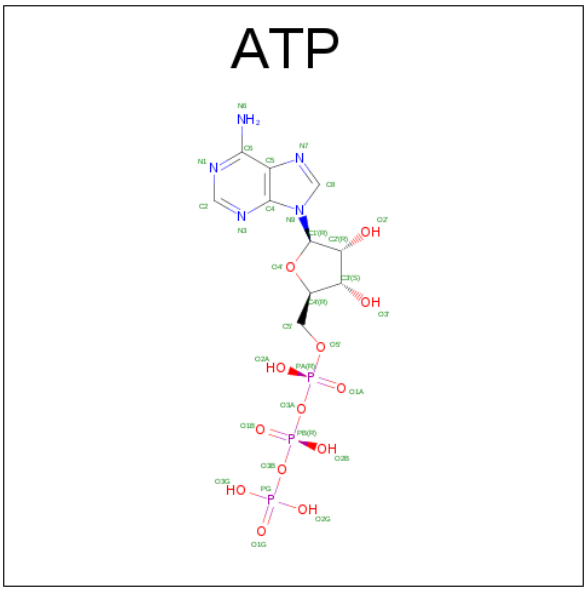
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	D	1	Total C N O P 20 9 3 7 1	0	0
10	N	1	Total C N O P 20 9 3 7 1	0	0

- Molecule 11 is CYTIDINE-5'-TRIPHOSPHATE (three-letter code: CTP) (formula: $C_9H_{16}N_3O_{14}P_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	D	1	Total	O	P	0	0
			9	7	2		
11	N	1	Total	O	P	0	0
			9	7	2		

- Molecule 12 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
12	D	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
12	N	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 13 is water.

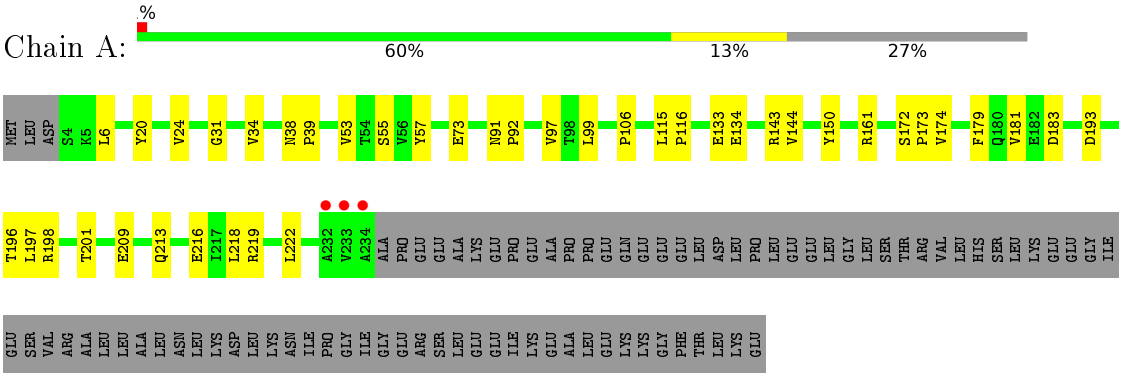
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	A	3	Total	O	0	0
			3	3		
13	B	2	Total	O	0	0
			2	2		
13	C	11	Total	O	0	0
			11	11		
13	D	18	Total	O	0	0
			18	18		
13	E	1	Total	O	0	0
			1	1		
13	G	2	Total	O	0	0
			2	2		
13	K	3	Total	O	0	0
			3	3		
13	L	1	Total	O	0	0
			1	1		
13	M	1	Total	O	0	0
			1	1		
13	N	14	Total	O	0	0
			14	14		
13	O	1	Total	O	0	0
			1	1		
13	P	1	Total	O	0	0
			1	1		

3 Residue-property plots (i)

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.

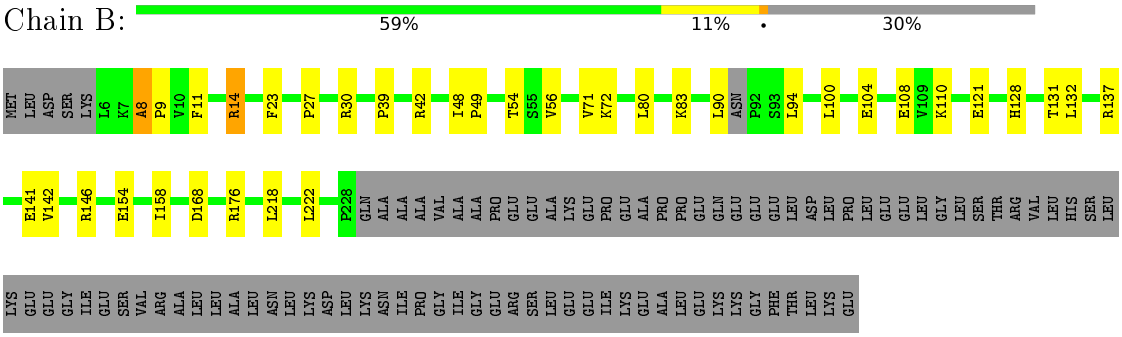
- Molecule 1: DNA-directed RNA polymerase subunit alpha

Chain A:



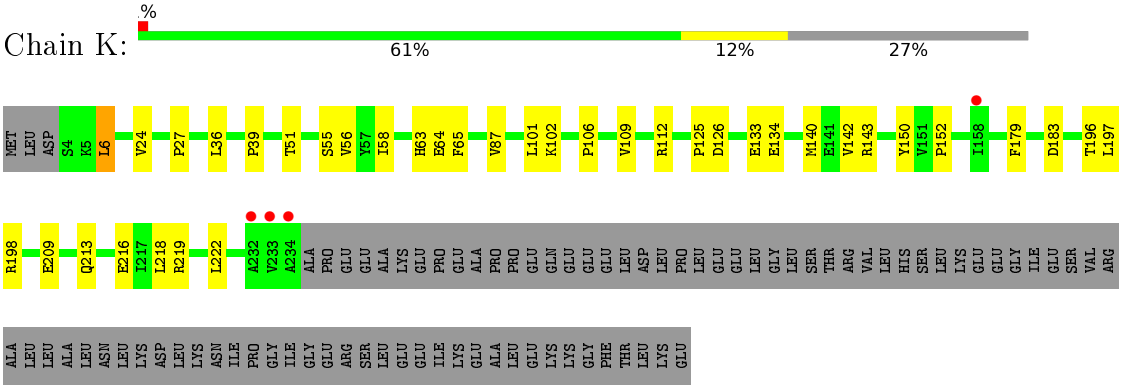
- Molecule 1: DNA-directed RNA polymerase subunit alpha

Chain B:



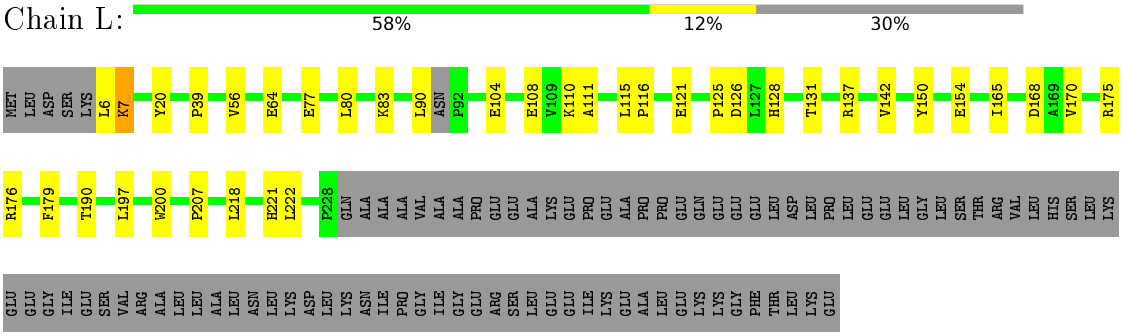
- Molecule 1: DNA-directed RNA polymerase subunit alpha

Chain K:



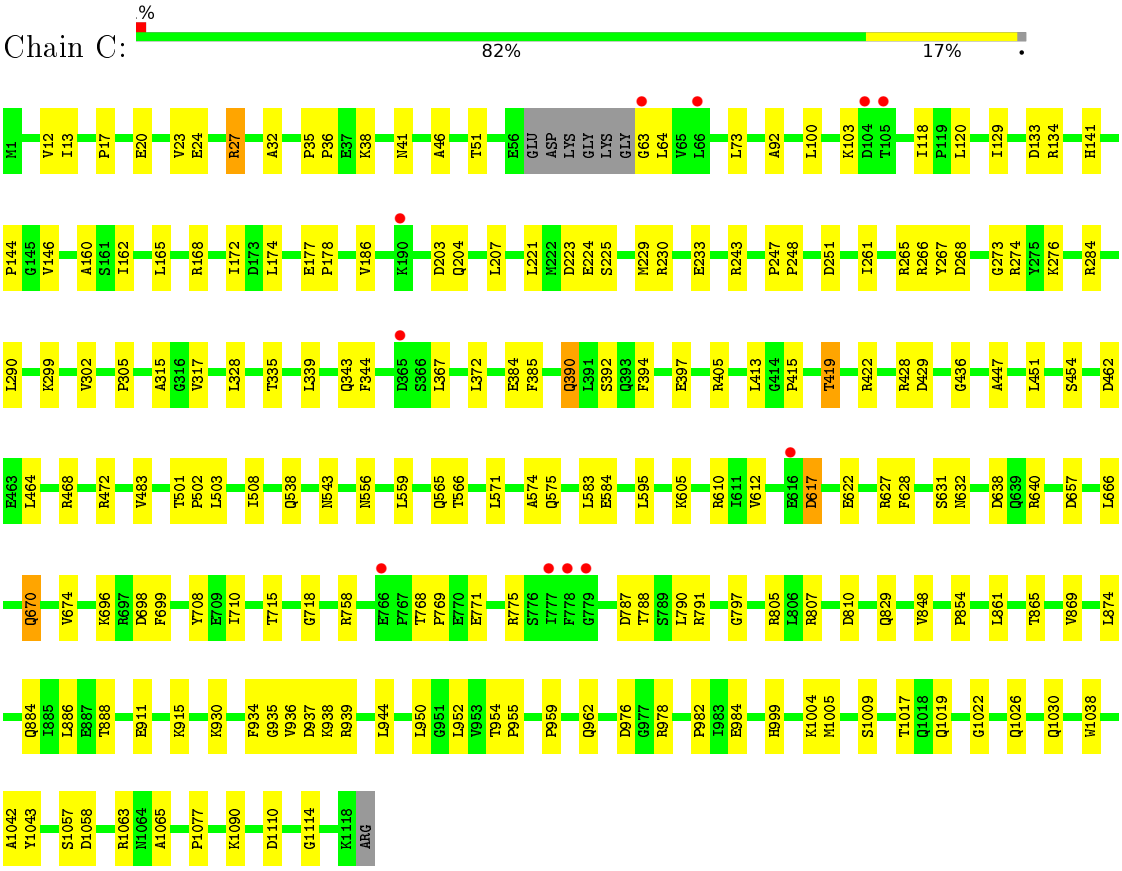
• Molecule 1: DNA-directed RNA polymerase subunit alpha

Chain L:



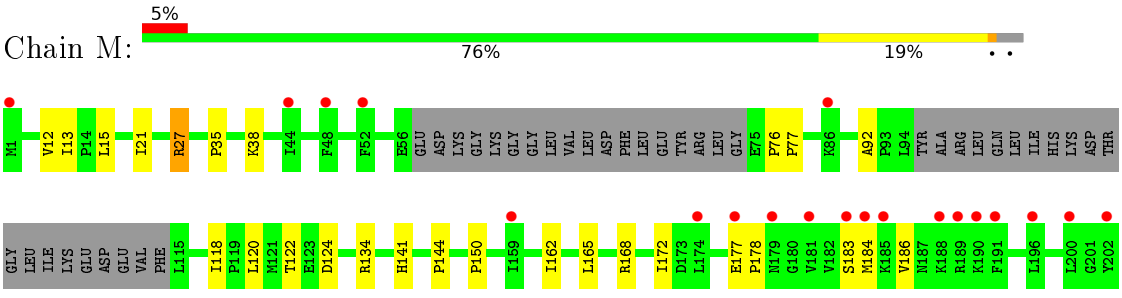
• Molecule 2: DNA-directed RNA polymerase subunit beta

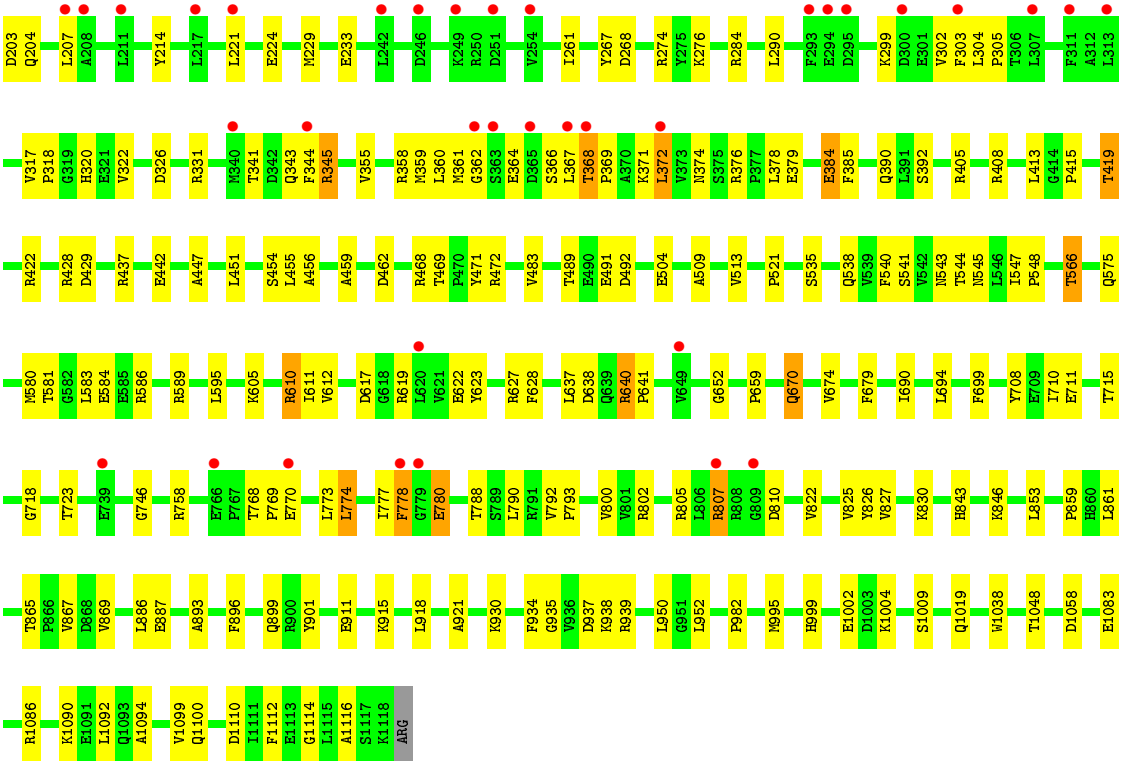
Chain C:



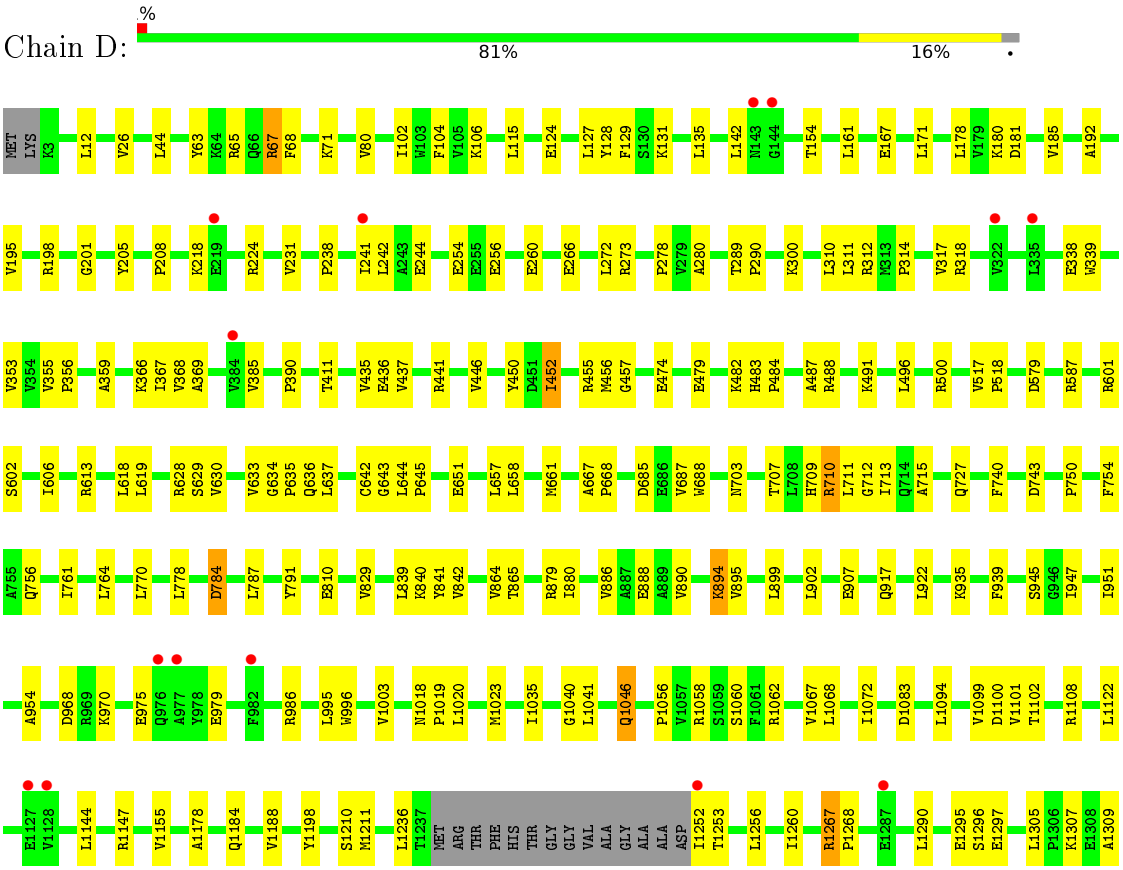
• Molecule 2: DNA-directed RNA polymerase subunit beta

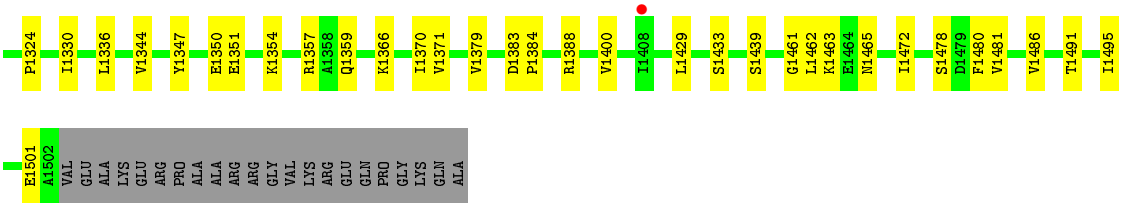
Chain M:



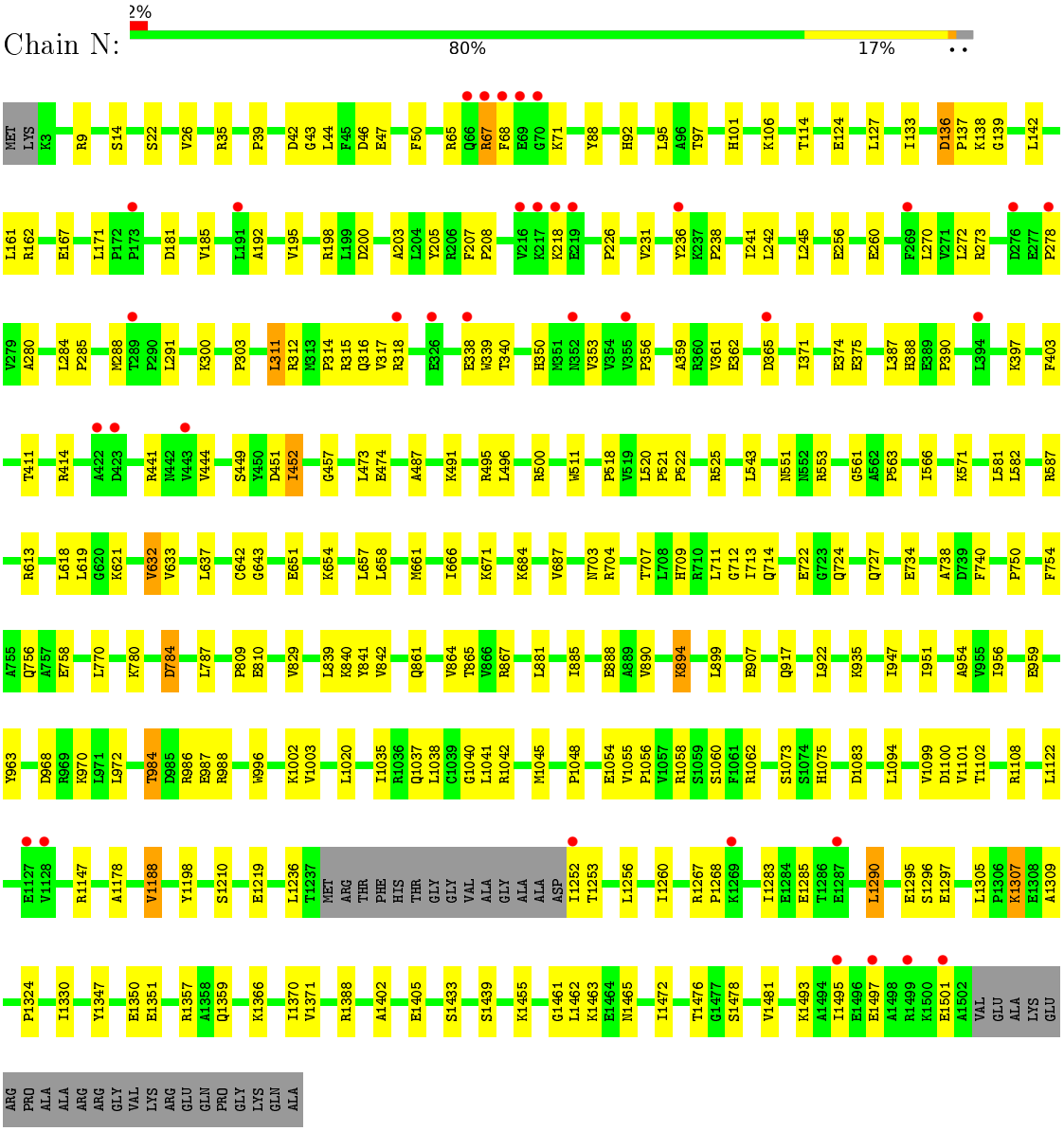


- Molecule 3: DNA-directed RNA polymerase subunit beta'

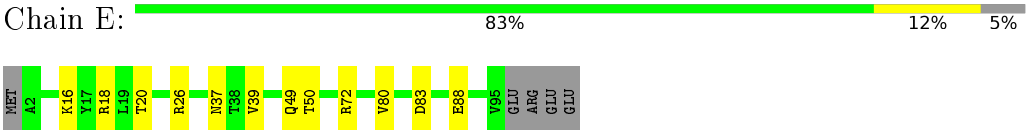




• Molecule 3: DNA-directed RNA polymerase subunit beta'

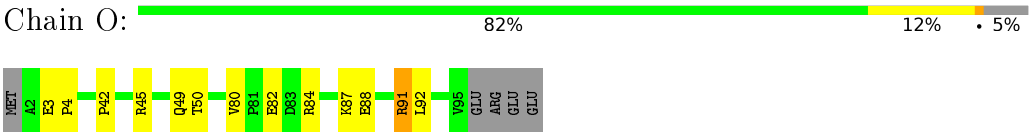


• Molecule 4: DNA-directed RNA polymerase subunit omega



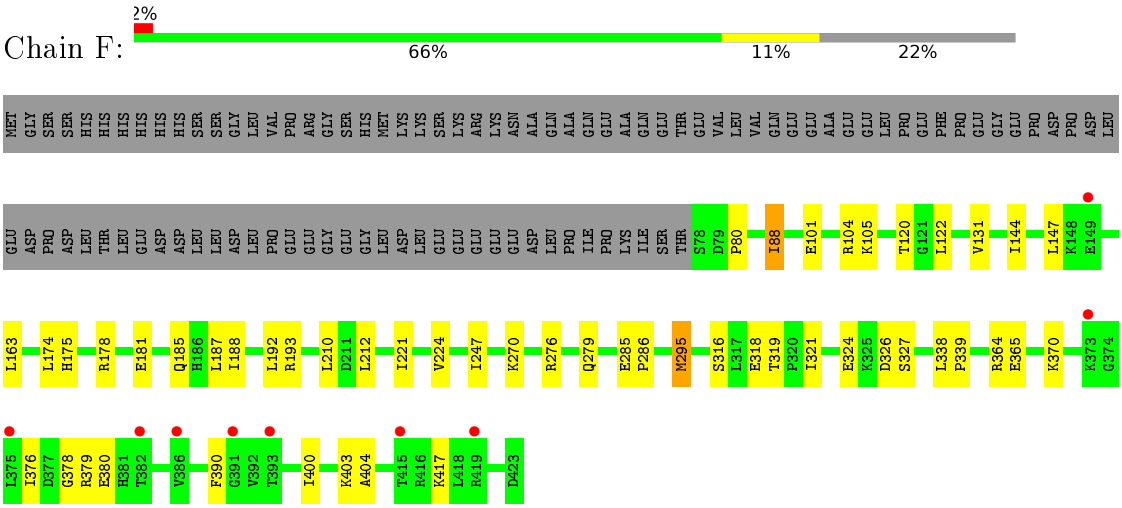
• Molecule 4: DNA-directed RNA polymerase subunit omega

Chain O:



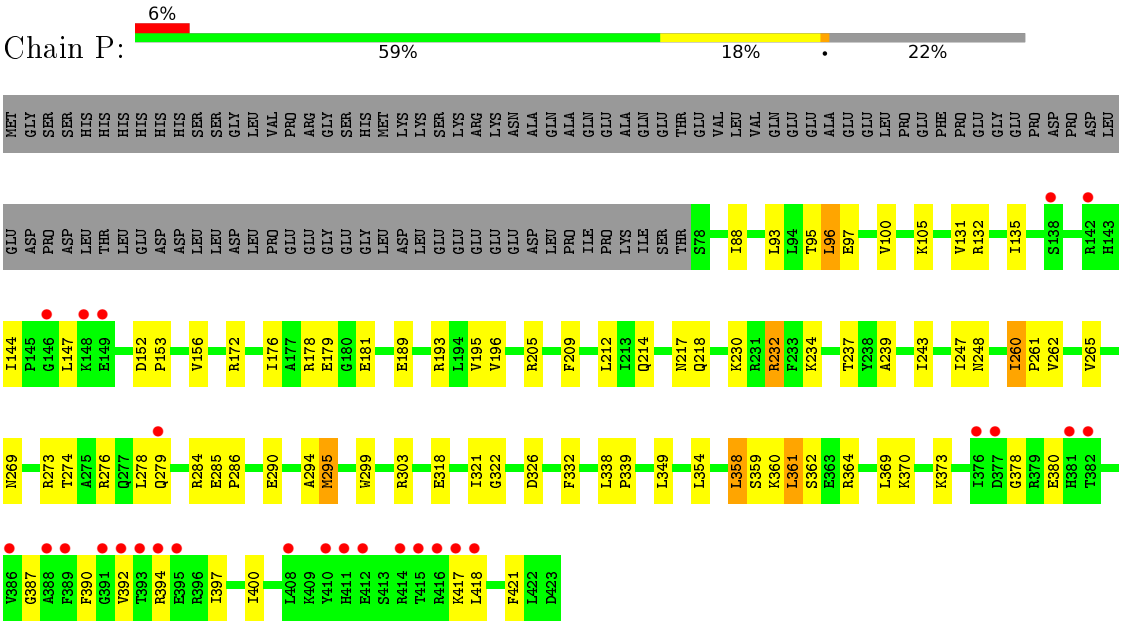
• Molecule 5: RNA polymerase sigma factor SigA

Chain F:



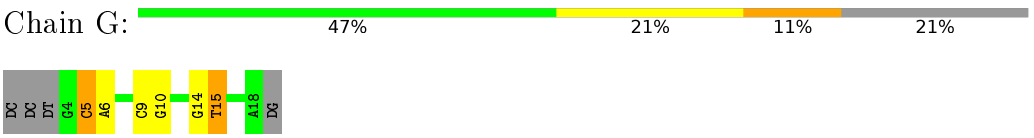
• Molecule 5: RNA polymerase sigma factor SigA

Chain P:



• Molecule 6: DNA (5'-D(*CP*CP*T*GP*CP*AP*TP*CP*CP*GP*TP*GP*AP*GP*TP*AP*GP*AP*G)-3')

Chain G:



• Molecule 6: DNA (5'-D(*CP*CP*T*GP*CP*AP*TP*CP*CP*GP*TP*GP*AP*GP*TP*AP*GP*AP*G)-3')

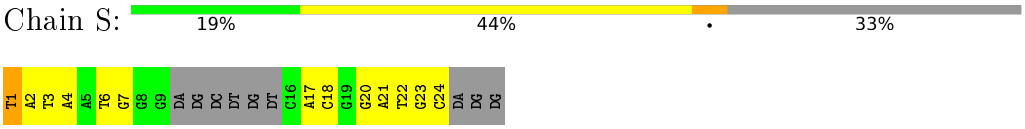
Chain R:

• Molecule 7: DNA (27-MER)

Chain H:

• Molecule 7: DNA (27-MER)

Chain S:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	185.96Å 103.64Å 297.42Å 90.00° 98.30° 90.00°	Depositor
Resolution (Å)	49.20 – 3.28 49.50 – 3.28	Depositor EDS
% Data completeness (in resolution range)	89.5 (49.20-3.28) 89.7 (49.50-3.28)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	0.18	Depositor
$< I/\sigma(I) >$ [1]	3.91 (at 3.25Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.211 , 0.257 0.212 , 0.259	Depositor DCC
R_{free} test set	7720 reflections (5.01%)	DCC
Wilson B-factor (Å^2)	57.7	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 31.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	56600	wwPDB-VP
Average B, all atoms (Å^2)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 52.73 % of the origin peak, indicating pseudo translational symmetry.The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 4.6669e-05.The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: MG, CTP, ZN, ATP

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	0.24	0/1841	0.45	0/2504
1	B	0.23	0/1781	0.47	0/2420
1	K	0.24	0/1841	0.44	0/2504
1	L	0.24	0/1781	0.45	0/2420
2	C	0.24	0/8941	0.44	0/12092
2	M	0.24	0/8669	0.45	0/11724
3	D	0.24	0/11944	0.44	0/16149
3	N	0.24	0/11944	0.44	0/16149
4	E	0.23	0/772	0.42	0/1040
4	O	0.22	0/772	0.42	0/1040
5	F	0.24	0/2852	0.40	0/3837
5	P	0.24	0/2852	0.43	0/3837
6	G	0.51	0/346	1.14	2/533 (0.4%)
6	R	0.51	0/346	1.09	1/533 (0.2%)
7	H	0.59	0/465	1.06	0/715
7	S	0.51	0/416	1.06	1/639 (0.2%)
All	All	0.25	0/57563	0.47	4/78136 (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
3	D	0	1
3	N	0	1
All	All	0	2

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	S	1	DT	O4'-C1'-N1	6.12	112.28	108.00
6	G	15	DT	O4'-C4'-C3'	-5.81	102.17	104.50
6	G	5	DC	O4'-C1'-N1	5.27	111.69	108.00
6	R	15	DT	O4'-C4'-C3'	-5.10	102.46	104.50

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	D	633	VAL	Peptide
3	N	138	LYS	Peptide

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	1809	0	1863	26	0
1	B	1750	0	1802	25	0
1	K	1809	0	1863	25	0
1	L	1750	0	1802	25	0
2	C	8774	0	8877	109	0
2	M	8508	0	8605	136	0
3	D	11738	0	11972	146	0
3	N	11738	0	11971	163	0
4	E	758	0	770	9	0
4	O	758	0	770	5	0
5	F	2807	0	2882	30	0
5	P	2807	0	2882	54	0
6	G	308	0	170	5	0
6	R	308	0	170	3	0
7	H	414	0	227	12	0
7	S	371	0	205	13	0
8	B	1	0	0	0	0
8	D	3	0	0	0	0
8	F	1	0	0	0	0
8	K	1	0	0	0	0
8	L	1	0	0	0	0
8	N	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	P	1	0	0	0	0
9	D	2	0	0	0	0
9	N	2	0	0	0	0
10	D	20	0	11	0	0
10	N	20	0	11	1	0
11	D	9	0	0	0	0
11	N	9	0	0	0	0
12	D	31	0	11	0	0
12	N	31	0	11	0	0
13	A	3	0	0	0	0
13	B	2	0	0	0	0
13	C	11	0	0	0	0
13	D	18	0	0	0	0
13	E	1	0	0	0	0
13	G	2	0	0	0	0
13	K	3	0	0	0	0
13	L	1	0	0	0	0
13	M	1	0	0	0	0
13	N	14	0	0	0	0
13	O	1	0	0	0	0
13	P	1	0	0	0	0
All	All	56600	0	56875	707	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 6.

All (707) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:243:ARG:NH1	7:H:9:DG:O6	1.99	0.93
2:M:165:LEU:HB2	2:M:168:ARG:HG3	1.59	0.84
2:M:758:ARG:HH21	2:M:788:THR:HB	1.45	0.80
3:D:1495:ILE:HG12	4:E:88:GLU:HG3	1.67	0.77
2:M:770:GLU:HB3	5:P:354:LEU:HG	1.65	0.77
2:C:63:GLY:HA3	2:C:100:LEU:HD21	1.68	0.76
2:C:12:VAL:HG21	2:C:472:ARG:HD3	1.66	0.75
2:C:628:PHE:H	2:C:638:ASP:HB3	1.49	0.75
2:M:802:ARG:HB2	2:M:826:TYR:HB2	1.71	0.71
3:N:127:LEU:HA	3:N:457:GLY:HA2	1.71	0.71
7:H:21:DA:H2"	7:H:22:DT:H5"	1.73	0.71
7:S:21:DA:H2"	7:S:22:DT:H5"	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:39:VAL:O	4:E:72:ARG:NH1	2.24	0.69
3:D:260:GLU:OE1	3:D:273:ARG:NH1	2.26	0.69
5:P:274:THR:HG21	5:P:295:MET:HG2	1.74	0.69
2:M:92:ALA:HB2	2:M:120:LEU:HD11	1.74	0.68
3:N:65:ARG:NH1	5:P:378:GLY:O	2.26	0.68
3:D:1461:GLY:O	3:D:1465:ASN:ND2	2.26	0.68
2:M:950:LEU:HB3	2:M:952:LEU:HD13	1.75	0.68
2:M:628:PHE:H	2:M:638:ASP:HB3	1.58	0.67
2:C:1019:GLN:HG2	2:C:1058:ASP:HB3	1.75	0.67
2:C:261:ILE:HG23	2:C:290:LEU:HB2	1.74	0.67
1:A:24:VAL:HG22	1:A:196:THR:HG23	1.75	0.67
3:D:65:ARG:NH1	5:F:378:GLY:O	2.28	0.67
5:P:273:ARG:HG2	5:P:276:ARG:HH12	1.59	0.67
2:M:939:ARG:HG2	2:M:982:PRO:HD3	1.77	0.66
2:C:343:GLN:HG3	2:C:385:PHE:HB2	1.76	0.66
2:M:408:ARG:NH1	2:M:456:ALA:O	2.29	0.66
1:L:7:LYS:NZ	1:L:7:LYS:HB3	2.10	0.66
1:B:176:ARG:NH2	3:D:888:GLU:OE1	2.28	0.65
3:D:241:ILE:HA	3:D:312:ARG:HG2	1.78	0.65
2:M:674:VAL:HG12	2:M:869:VAL:HB	1.79	0.65
3:N:956:ILE:HD11	3:N:1062:ARG:HG2	1.78	0.65
2:C:939:ARG:HG2	2:C:982:PRO:HD3	1.78	0.65
3:D:356:PRO:HG2	3:D:359:ALA:HB2	1.79	0.65
3:N:356:PRO:HG2	3:N:359:ALA:HB2	1.79	0.65
2:C:950:LEU:HB3	2:C:952:LEU:HD13	1.79	0.64
3:D:1495:ILE:HD13	4:E:80:VAL:HG21	1.78	0.64
3:N:563:PRO:HD2	3:N:566:ILE:HD12	1.79	0.64
2:C:428:ARG:NH2	2:C:447:ALA:O	2.31	0.63
3:D:366:LYS:HD3	3:D:369:ALA:HB2	1.80	0.63
2:C:165:LEU:HB2	2:C:168:ARG:HG3	1.81	0.63
2:M:168:ARG:HD3	2:M:268:ASP:HB3	1.81	0.63
3:N:273:ARG:HB3	3:N:278:PRO:HA	1.79	0.63
3:N:1495:ILE:HD13	4:O:80:VAL:HG21	1.80	0.63
1:K:55:SER:HB3	1:K:143:ARG:HB3	1.81	0.63
2:M:846:LYS:NZ	10:N:1605:C:OP1	2.31	0.63
2:C:807:ARG:NH1	2:C:810:ASP:OD2	2.33	0.62
3:D:256:GLU:HG3	3:D:300:LYS:HG3	1.80	0.62
2:M:341:THR:HG22	2:M:345:ARG:HH12	1.64	0.62
3:D:643:GLY:HA3	3:D:727:GLN:HB2	1.81	0.62
2:C:674:VAL:HG12	2:C:869:VAL:HB	1.82	0.61
2:M:1019:GLN:HG2	2:M:1058:ASP:HB3	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:N:643:GLY:HA3	3:N:727:GLN:HB2	1.81	0.61
3:N:242:LEU:HB3	3:N:311:LEU:HD12	1.80	0.61
2:M:807:ARG:NH1	2:M:810:ASP:OD2	2.33	0.61
3:N:124:GLU:OE2	3:N:587:ARG:NH2	2.33	0.61
3:N:241:ILE:HA	3:N:312:ARG:HG2	1.83	0.61
1:K:24:VAL:HG22	1:K:196:THR:HG23	1.82	0.61
5:P:322:GLY:HA3	5:P:326:ASP:HB2	1.83	0.61
2:C:715:THR:OG1	2:C:718:GLY:O	2.17	0.61
3:D:266:GLU:HG3	3:D:314:PRO:HB3	1.82	0.61
1:B:108:GLU:HG2	1:B:131:THR:HG22	1.83	0.61
5:P:397:ILE:HD12	5:P:400:ILE:HD12	1.83	0.60
5:P:131:VAL:HG13	5:P:178:ARG:HD3	1.84	0.60
3:D:1480:PHE:O	4:E:18:ARG:NH2	2.34	0.60
2:M:15:LEU:O	2:M:586:ARG:NH1	2.33	0.60
1:A:216:GLU:OE2	1:A:219:ARG:NH2	2.34	0.60
5:P:260:ILE:HD11	5:P:265:VAL:HG22	1.83	0.60
3:D:711:LEU:HD22	3:D:778:LEU:HD23	1.84	0.60
2:M:637:LEU:HG	2:M:659:PRO:HG3	1.84	0.60
3:D:894:LYS:HD3	3:D:894:LYS:H	1.66	0.60
3:N:954:ALA:O	3:N:1062:ARG:NH2	2.35	0.59
3:N:707:THR:HG23	3:N:712:GLY:HA3	1.83	0.59
2:C:930:LYS:HE3	2:C:935:GLY:HA2	1.84	0.59
3:N:161:LEU:HB3	3:N:452:ILE:HD11	1.84	0.59
3:N:894:LYS:H	3:N:894:LYS:HD3	1.67	0.59
1:B:94:LEU:O	1:B:146:ARG:NH2	2.36	0.59
2:C:92:ALA:HB2	2:C:120:LEU:HD11	1.84	0.59
3:D:1003:VAL:HG21	3:D:1041:LEU:HG	1.83	0.59
3:D:171:LEU:HD12	3:D:390:PRO:HG2	1.85	0.59
5:F:316:SER:HB3	5:F:319:THR:HG23	1.85	0.58
5:P:195:VAL:HG12	5:P:243:ILE:HD12	1.86	0.58
3:N:1461:GLY:O	3:N:1465:ASN:ND2	2.35	0.58
2:C:134:ARG:NH1	2:C:392:SER:OG	2.36	0.58
1:L:104:GLU:OE2	1:L:137:ARG:NH1	2.36	0.58
2:M:12:VAL:HG12	2:M:13:ILE:HG23	1.84	0.58
2:M:21:ILE:HD12	2:M:455:LEU:HD22	1.85	0.58
2:C:266:ARG:NH1	7:H:11:DG:O6	2.37	0.57
2:M:711:GLU:O	2:M:758:ARG:NH1	2.37	0.57
1:L:56:VAL:HG22	1:L:142:VAL:HG12	1.86	0.57
3:D:208:PRO:HA	3:D:390:PRO:HA	1.85	0.57
1:L:108:GLU:HG2	1:L:131:THR:HG22	1.86	0.57
1:B:8:ALA:HB1	1:B:9:PRO:HA	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:172:ILE:HG12	2:C:186:VAL:HG22	1.86	0.57
2:M:690:ILE:HG22	2:M:869:VAL:HG22	1.87	0.57
3:N:543:LEU:HD13	3:N:581:LEU:HA	1.87	0.57
2:C:146:VAL:HG22	2:C:162:ILE:HG12	1.87	0.57
3:N:1040:GLY:O	3:N:1060:SER:HB3	2.04	0.57
1:A:218:LEU:HD23	1:B:222:LEU:HD21	1.86	0.56
2:C:769:PRO:HG3	3:D:65:ARG:HH12	1.70	0.56
2:M:35:PRO:HG2	2:M:38:LYS:HB2	1.87	0.56
1:B:56:VAL:HG22	1:B:142:VAL:HG12	1.86	0.56
5:F:270:LYS:HG2	5:F:295:MET:HE1	1.86	0.56
3:N:1324:PRO:HG3	3:N:1330:ILE:HD11	1.88	0.56
2:M:462:ASP:HB3	2:M:468:ARG:HD2	1.87	0.56
3:N:561:GLY:HA3	5:P:132:ARG:HD3	1.88	0.56
7:S:22:DT:H2"	7:S:23:DG:C8	2.41	0.56
2:M:715:THR:OG1	2:M:718:GLY:O	2.24	0.56
7:H:23:DG:H2"	7:H:24:DC:H5"	1.87	0.56
5:F:188:ILE:HD13	5:F:221:ILE:HG12	1.87	0.56
3:D:273:ARG:HB3	3:D:278:PRO:HA	1.88	0.56
3:D:124:GLU:OE2	3:D:587:ARG:NH2	2.39	0.56
2:C:207:LEU:HD13	2:C:221:LEU:HD21	1.88	0.56
2:C:605:LYS:HB2	2:C:612:VAL:HB	1.88	0.56
3:D:1100:ASP:OD2	3:D:1463:LYS:NZ	2.38	0.56
2:M:361:MET:SD	2:M:362:GLY:N	2.79	0.55
3:D:657:LEU:HG	3:D:661:MET:HE2	1.88	0.55
3:N:42:ASP:N	3:N:46:ASP:OD2	2.35	0.55
3:D:1478:SER:HB3	3:D:1481:VAL:HG22	1.88	0.55
5:P:321:ILE:HG21	5:P:332:PHE:HE2	1.71	0.55
2:C:1038:TRP:CE2	3:D:1099:VAL:HG11	2.42	0.55
2:M:1092:LEU:HD13	2:M:1099:VAL:HG21	1.88	0.55
2:M:326:ASP:HA	2:M:331:ARG:HD2	1.88	0.55
3:N:968:ASP:OD2	3:N:1058:ARG:NH2	2.40	0.55
5:P:153:PRO:HA	5:P:156:VAL:HG22	1.89	0.55
2:C:367:LEU:HD13	2:C:372:LEU:HD21	1.88	0.55
2:M:390:GLN:HB3	2:M:415:PRO:HD3	1.89	0.55
2:M:428:ARG:NH2	2:M:447:ALA:O	2.34	0.55
2:M:711:GLU:HG2	2:M:822:VAL:HG22	1.89	0.55
5:F:131:VAL:HG13	5:F:178:ARG:HD3	1.88	0.55
2:M:207:LEU:HD13	2:M:221:LEU:HD21	1.89	0.55
3:N:1147:ARG:HD3	3:N:1188:VAL:HG11	1.88	0.55
1:A:198:ARG:HD3	2:C:934:PHE:CZ	2.42	0.55
2:M:541:SER:O	2:M:545:ASN:ND2	2.31	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:N:1100:ASP:OD2	3:N:1463:LYS:NZ	2.38	0.55
2:C:168:ARG:HD3	2:C:268:ASP:HB3	1.89	0.54
1:K:218:LEU:HD23	1:L:222:LEU:HD21	1.88	0.54
3:N:97:THR:HG21	3:N:571:LYS:HG2	1.88	0.54
2:C:937:ASP:OD1	2:C:939:ARG:HD3	2.06	0.54
3:N:996:TRP:CE2	3:N:1056:PRO:HG3	2.41	0.54
5:P:96:LEU:HD22	5:P:234:LYS:HD3	1.89	0.54
3:D:899:LEU:HD22	3:D:917:GLN:HB3	1.88	0.54
2:M:12:VAL:HG21	2:M:472:ARG:HD3	1.88	0.54
2:M:504:GLU:HG2	2:M:509:ALA:HB2	1.89	0.54
5:P:237:THR:HG21	7:S:3:DT:H5"	1.88	0.54
3:D:317:VAL:HG23	3:D:339:TRP:HB3	1.88	0.54
3:D:1236:LEU:HA	3:D:1359:GLN:HG3	1.88	0.54
3:D:1102:THR:HG21	3:D:1371:VAL:HG22	1.88	0.54
3:D:127:LEU:HA	3:D:457:GLY:HA2	1.90	0.54
2:M:177:GLU:HG3	2:M:178:PRO:HD2	1.90	0.54
3:N:474:GLU:HG3	3:N:496:LEU:HD11	1.90	0.54
1:K:222:LEU:HD21	1:L:218:LEU:HD23	1.90	0.54
3:D:1020:LEU:HB3	3:D:1035:ILE:HD12	1.90	0.54
5:F:80:PRO:HB2	5:F:210:LEU:HD11	1.90	0.54
3:D:181:ASP:HB2	3:D:205:TYR:CD2	2.43	0.53
3:D:474:GLU:HG3	3:D:496:LEU:HD11	1.91	0.53
7:H:10:DA:H2"	7:H:11:DG:OP2	2.08	0.53
2:C:462:ASP:HB3	2:C:468:ARG:HD2	1.89	0.53
3:D:1462:LEU:HD22	3:D:1472:ILE:HB	1.90	0.53
3:D:238:PRO:HD3	3:D:318:ARG:HG3	1.90	0.53
5:F:187:LEU:HD23	5:F:224:VAL:HG13	1.91	0.53
1:L:80:LEU:HD21	3:N:842:VAL:HG12	1.90	0.53
3:N:1003:VAL:HG21	3:N:1041:LEU:HG	1.90	0.53
3:D:63:TYR:HB2	3:D:80:VAL:HG21	1.91	0.53
2:M:1100:GLN:HG3	3:N:9:ARG:HH21	1.72	0.53
3:N:703:ASN:HB2	3:N:713:ILE:HG12	1.90	0.53
5:P:358:LEU:HD12	5:P:370:LYS:HE3	1.89	0.53
3:N:1305:LEU:HD13	3:N:1309:ALA:HB3	1.91	0.53
3:D:970:LYS:HD3	3:D:995:LEU:HD13	1.90	0.53
3:N:657:LEU:HG	3:N:661:MET:HE2	1.90	0.53
5:F:212:LEU:HD22	5:F:247:ILE:HG23	1.91	0.53
1:K:51:THR:OG1	1:K:87:VAL:O	2.21	0.53
2:M:456:ALA:HB3	2:M:459:ALA:HB2	1.91	0.53
2:C:768:THR:OG1	2:C:771:GLU:OE1	2.26	0.53
3:D:890:VAL:HB	3:D:922:LEU:HD13	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:F:365:GLU:HB2	5:F:404:ALA:HB2	1.91	0.53
3:N:520:LEU:HD12	3:N:521:PRO:HD2	1.90	0.53
2:C:41:ASN:O	2:C:46:ALA:HB2	2.07	0.53
2:M:930:LYS:HE3	2:M:935:GLY:HA2	1.90	0.53
3:N:285:PRO:HD2	3:N:288:MET:SD	2.49	0.53
1:A:222:LEU:HD21	1:B:218:LEU:HD23	1.91	0.53
3:D:1040:GLY:O	3:D:1060:SER:HB3	2.09	0.52
1:L:176:ARG:NH2	3:N:888:GLU:OE1	2.41	0.52
3:D:128:TYR:OH	3:D:579:ASP:OD2	2.25	0.52
1:K:216:GLU:OE2	1:K:219:ARG:NH2	2.42	0.52
2:M:358:ARG:HG2	2:M:372:LEU:HA	1.91	0.52
3:D:314:PRO:HB2	3:D:317:VAL:HG12	1.92	0.52
2:M:589:ARG:NH2	2:M:652:GLY:O	2.38	0.52
5:P:96:LEU:O	5:P:100:VAL:HG23	2.09	0.52
3:D:367:ILE:HG22	3:D:368:VAL:HG23	1.90	0.52
2:M:229:MET:HB2	2:M:233:GLU:HB2	1.91	0.52
2:M:710:ILE:HD12	2:M:790:LEU:HB2	1.91	0.52
2:M:937:ASP:OD1	2:M:939:ARG:HD3	2.09	0.52
1:K:183:ASP:HA	2:M:938:LYS:HE3	1.92	0.52
2:M:612:VAL:HG22	2:M:622:GLU:HG3	1.90	0.52
3:N:899:LEU:HD22	3:N:917:GLN:HB3	1.90	0.52
5:P:239:ALA:O	5:P:243:ILE:HG12	2.09	0.52
3:D:167:GLU:OE2	3:D:198:ARG:NH1	2.42	0.52
4:E:37:ASN:N	4:E:37:ASN:OD1	2.41	0.52
2:M:899:GLN:NE2	2:M:901:TYR:OH	2.42	0.52
2:C:617:ASP:OD1	2:C:617:ASP:N	2.43	0.52
2:C:886:LEU:HD21	3:D:951:ILE:HG12	1.92	0.52
2:C:24:GLU:OE2	2:C:27:ARG:NH2	2.42	0.52
3:D:242:LEU:HB3	3:D:311:LEU:HD12	1.92	0.52
1:B:90:LEU:HD21	1:B:121:GLU:HB2	1.92	0.51
3:N:192:ALA:HB3	3:N:195:VAL:HB	1.93	0.51
3:D:954:ALA:O	3:D:1062:ARG:NH2	2.44	0.51
5:P:417:LYS:HB3	5:P:418:LEU:HD12	1.92	0.51
2:M:343:GLN:NE2	2:M:384:GLU:OE2	2.44	0.51
3:N:658:LEU:HA	3:N:661:MET:HE3	1.91	0.51
2:C:1110:ASP:OD2	2:C:1114:GLY:N	2.34	0.51
1:K:6:LEU:HD11	1:K:27:PRO:HG2	1.92	0.51
3:N:95:LEU:HA	3:N:551:ASN:HD21	1.76	0.51
2:M:172:ILE:HG12	2:M:186:VAL:HG22	1.93	0.51
2:C:848:VAL:HG22	3:D:740:PHE:O	2.11	0.51
2:M:605:LYS:HB2	2:M:612:VAL:HB	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:N:1020:LEU:HB3	3:N:1035:ILE:HD12	1.92	0.51
1:A:106:PRO:HG3	1:A:134:GLU:HG2	1.93	0.51
7:S:23:DG:H2"	7:S:24:DC:H5"	1.93	0.51
2:C:670:GLN:HG2	2:C:699:PHE:CD2	2.46	0.51
2:M:605:LYS:HB3	2:M:610:ARG:HH11	1.76	0.51
2:M:843:HIS:NE2	2:M:887:GLU:OE2	2.44	0.51
3:D:1366:LYS:O	3:D:1370:ILE:HG12	2.10	0.50
3:N:181:ASP:HB2	3:N:205:TYR:CD2	2.46	0.50
3:D:787:LEU:HD21	3:D:947:ILE:HG21	1.93	0.50
3:D:975:GLU:O	3:D:979:GLU:HG2	2.10	0.50
3:N:988:ARG:NH2	3:N:1054:GLU:OE2	2.44	0.50
3:D:483:HIS:CE1	3:D:488:ARG:HD3	2.46	0.50
1:B:80:LEU:HD21	3:D:842:VAL:HG12	1.94	0.50
3:N:22:SER:HB2	3:N:92:HIS:HB3	1.94	0.50
2:M:168:ARG:O	2:M:267:TYR:HA	2.11	0.50
3:N:411:THR:O	5:P:178:ARG:NH1	2.38	0.50
3:N:758:GLU:OE1	3:N:1476:THR:OG1	2.23	0.50
5:P:279:GLN:HB3	5:P:286:PRO:HD3	1.94	0.50
3:D:644:LEU:HD12	3:D:645:PRO:HD2	1.92	0.49
2:C:612:VAL:HG22	2:C:622:GLU:HG3	1.94	0.49
3:D:1101:VAL:HG13	3:D:1102:THR:HG23	1.94	0.49
1:K:112:ARG:HG3	1:K:125:PRO:HB2	1.94	0.49
2:M:1094:ALA:HA	3:N:518:PRO:HB2	1.93	0.49
3:N:208:PRO:HG2	3:N:353:VAL:HG21	1.94	0.49
2:M:611:ILE:HD11	2:M:641:PRO:HB3	1.94	0.49
5:P:135:ILE:HG13	5:P:181:GLU:HB2	1.93	0.49
3:D:106:LYS:HE3	3:D:587:ARG:HG3	1.95	0.49
2:M:540:PHE:HB3	2:M:544:THR:HB	1.94	0.49
2:M:896:PHE:HB2	2:M:921:ALA:HB1	1.94	0.49
3:N:1094:LEU:HD22	3:N:1260:ILE:HG12	1.93	0.49
2:M:773:LEU:HD13	5:P:373:LYS:HG3	1.93	0.49
2:C:13:ILE:HD13	2:C:483:VAL:HG11	1.93	0.49
2:M:343:GLN:HG3	2:M:385:PHE:HB2	1.93	0.49
5:P:284:ARG:NH2	5:P:290:GLU:OE2	2.44	0.49
3:D:1147:ARG:HD3	3:D:1188:VAL:HG11	1.94	0.49
1:K:179:PHE:HB3	1:K:197:LEU:HD23	1.93	0.49
1:K:209:GLU:O	1:K:213:GLN:HG2	2.12	0.49
2:M:419:THR:HG23	2:M:422:ARG:HG3	1.95	0.49
3:N:959:GLU:HB3	3:N:963:TYR:CE2	2.47	0.49
3:N:996:TRP:CD2	3:N:1056:PRO:HG3	2.47	0.49
2:C:51:THR:O	2:C:265:ARG:NH2	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:864:VAL:HG22	3:D:865:THR:H	1.77	0.49
3:D:945:SER:OG	3:D:947:ILE:HG12	2.13	0.49
1:L:90:LEU:HD21	1:L:121:GLU:HB2	1.94	0.49
3:N:1037:GLN:HG2	3:N:1042:ARG:HD2	1.95	0.49
2:C:571:LEU:HB2	2:C:574:ALA:HB2	1.95	0.49
3:D:996:TRP:CD2	3:D:1056:PRO:HG3	2.47	0.49
1:K:106:PRO:HG3	1:K:134:GLU:HG2	1.94	0.49
3:N:162:ARG:NH1	3:N:451:ASP:OD1	2.45	0.49
3:N:238:PRO:HB3	3:N:315:ARG:HA	1.95	0.49
1:B:83:LYS:HE2	1:B:168:ASP:HB2	1.94	0.48
2:C:805:ARG:HE	2:C:807:ARG:HE	1.61	0.48
3:N:114:THR:HG23	3:N:495:ARG:HG2	1.95	0.48
3:N:350:HIS:HE1	5:P:232:ARG:HG3	1.76	0.48
3:N:840:LYS:HE3	3:N:841:TYR:CZ	2.48	0.48
3:N:787:LEU:HD21	3:N:947:ILE:HG21	1.94	0.48
1:A:209:GLU:O	1:A:213:GLN:HG2	2.13	0.48
2:C:787:ASP:OD2	2:C:791:ARG:NH2	2.44	0.48
2:M:617:ASP:HB2	2:M:619:ARG:HG2	1.94	0.48
2:C:1017:THR:HB	2:C:1019:GLN:HG3	1.95	0.48
2:C:229:MET:HB2	2:C:233:GLU:HB2	1.95	0.48
1:L:111:ALA:HB3	1:L:125:PRO:HA	1.95	0.48
2:C:670:GLN:HG2	2:C:699:PHE:CG	2.49	0.48
3:D:178:LEU:HG	3:D:192:ALA:HA	1.95	0.48
2:M:437:ARG:NH1	2:M:491:GLU:OE2	2.41	0.48
2:M:368:THR:HG22	2:M:369:PRO:HD2	1.95	0.48
2:M:584:GLU:N	2:M:584:GLU:OE2	2.43	0.48
3:N:361:VAL:HG23	3:N:365:ASP:HB2	1.95	0.48
3:D:840:LYS:HE3	3:D:841:TYR:CZ	2.49	0.48
3:N:666:ILE:HG21	3:N:687:VAL:HG12	1.96	0.48
2:M:203:ASP:OD1	2:M:204:GLN:N	2.46	0.48
2:M:800:VAL:HG22	2:M:827:VAL:HG22	1.96	0.48
3:N:487:ALA:O	3:N:491:LYS:HG2	2.13	0.48
3:D:1305:LEU:HD13	3:D:1309:ALA:HB3	1.95	0.48
2:M:405:ARG:HD2	2:M:442:GLU:OE2	2.13	0.48
3:N:260:GLU:OE1	3:N:273:ARG:NH1	2.47	0.48
2:C:627:ARG:HD3	2:C:638:ASP:HB2	1.96	0.48
3:D:1350:GLU:OE2	3:D:1357:ARG:NH1	2.44	0.48
3:N:473:LEU:HD21	3:N:495:ARG:HH21	1.79	0.48
1:A:55:SER:HB3	1:A:143:ARG:HB3	1.95	0.47
2:C:32:ALA:HB2	2:C:73:LEU:HD12	1.96	0.47
3:N:39:PRO:HG2	3:N:47:GLU:HG3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:P:152:ASP:N	5:P:152:ASP:OD1	2.47	0.47
5:P:105:LYS:HD3	5:P:179:GLU:HG2	1.96	0.47
2:C:1057:SER:O	2:C:1063:ARG:NH1	2.45	0.47
3:D:1379:VAL:HG21	3:D:1400:VAL:HG11	1.96	0.47
3:D:185:VAL:N	3:D:201:GLY:O	2.48	0.47
3:D:707:THR:HG23	3:D:712:GLY:HA3	1.95	0.47
2:M:1009:SER:HB3	3:N:651:GLU:O	2.15	0.47
2:C:1030:GLN:OE1	3:D:628:ARG:NH1	2.41	0.47
2:C:503:LEU:HD23	2:C:508:ILE:HA	1.95	0.47
3:D:1324:PRO:HG3	3:D:1330:ILE:HD11	1.96	0.47
2:M:134:ARG:NH1	2:M:392:SER:OG	2.47	0.47
6:R:4:DG:H1	7:S:24:DC:H42	1.63	0.47
2:C:168:ARG:HA	2:C:168:ARG:HE	1.78	0.47
2:C:413:LEU:HD21	2:C:451:LEU:HD13	1.97	0.47
3:D:895:VAL:HG11	3:D:922:LEU:HD21	1.97	0.47
6:R:14:DG:H2'	6:R:15:DT:C6	2.50	0.47
2:C:129:ILE:HB	2:C:134:ARG:HD2	1.97	0.47
2:C:758:ARG:HH21	2:C:788:THR:HB	1.79	0.47
1:K:198:ARG:HD3	2:M:934:PHE:CZ	2.49	0.47
3:N:1366:LYS:O	3:N:1370:ILE:HG12	2.14	0.47
3:N:226:PRO:HG2	3:N:245:LEU:HD11	1.97	0.47
5:P:172:ARG:O	5:P:176:ILE:HG12	2.15	0.47
6:R:6:DA:H5"	6:R:6:DA:H8	1.80	0.47
1:B:71:VAL:HG22	1:B:132:LEU:HG	1.97	0.47
3:N:1350:GLU:OE2	3:N:1357:ARG:NH1	2.43	0.47
3:N:67:ARG:HD3	3:N:68:PHE:CE2	2.50	0.47
7:S:17:DA:H2'	7:S:18:DC:C6	2.50	0.47
7:S:20:DG:H4'	7:S:21:DA:OP1	2.14	0.47
1:A:34:VAL:HB	1:B:42:ARG:NH2	2.30	0.47
3:D:1046:GLN:HB3	3:D:1046:GLN:HE21	1.56	0.47
2:M:774:LEU:HD12	5:P:418:LEU:HD23	1.97	0.47
3:N:26:VAL:HG11	3:N:44:LEU:HD23	1.96	0.47
3:D:435:VAL:HG22	3:D:446:VAL:HG22	1.97	0.46
6:G:5:DC:H42	7:H:23:DG:H1	1.61	0.46
3:D:135:LEU:HD22	3:D:455:ARG:HE	1.80	0.46
3:D:355:VAL:HG11	3:D:385:VAL:HG21	1.97	0.46
3:D:637:LEU:HD13	3:D:642:CYS:HA	1.97	0.46
4:E:83:ASP:OD1	4:E:83:ASP:N	2.48	0.46
5:F:101:GLU:HG3	5:F:104:ARG:HH21	1.80	0.46
5:F:88:ILE:HD11	5:F:192:LEU:HD13	1.97	0.46
1:K:39:PRO:HG3	1:L:39:PRO:HG3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:N:403:PHE:CD1	3:N:444:VAL:HG23	2.50	0.46
6:G:5:DC:H2"	6:G:6:DA:C8	2.50	0.46
3:N:272:LEU:HB2	3:N:280:ALA:HB3	1.97	0.46
1:L:83:LYS:NZ	3:N:842:VAL:O	2.49	0.46
5:P:193:ARG:HB2	7:S:6:DT:H1'	1.97	0.46
3:D:129:PHE:CD2	3:D:456:MET:HB3	2.51	0.46
3:D:770:LEU:HB2	3:D:1210:SER:HA	1.97	0.46
5:F:400:ILE:HA	5:F:403:LYS:HG2	1.98	0.46
3:N:14:SER:HB3	3:N:511:TRP:CE2	2.51	0.46
1:A:179:PHE:HB3	1:A:197:LEU:HD23	1.97	0.46
2:C:160:ALA:HB3	2:C:174:LEU:HB2	1.96	0.46
3:D:761:ILE:HD12	4:E:20:THR:HA	1.97	0.46
1:K:64:GLU:OE2	2:M:830:LYS:NZ	2.40	0.46
3:N:231:VAL:O	3:N:236:TYR:OH	2.33	0.46
2:C:144:PRO:HB2	2:C:273:GLY:HA3	1.97	0.46
3:D:131:LYS:NZ	3:D:154:THR:HG22	2.30	0.46
3:D:224:ARG:NE	3:D:254:GLU:OE2	2.33	0.46
2:M:144:PRO:HG2	2:M:165:LEU:HD23	1.98	0.46
1:A:133:GLU:HG2	1:A:134:GLU:H	1.81	0.46
1:A:73:GLU:OE1	1:A:73:GLU:N	2.49	0.46
1:L:175:ARG:N	1:L:200:TRP:O	2.44	0.46
2:M:627:ARG:HH22	2:M:640:ARG:HG3	1.81	0.46
2:M:770:GLU:CD	2:M:770:GLU:H	2.19	0.46
2:C:397:GLU:HG3	2:C:631:SER:HB2	1.97	0.46
3:D:1068:LEU:O	3:D:1072:ILE:HG12	2.15	0.46
1:L:64:GLU:HA	1:L:165:ILE:HD13	1.98	0.46
1:A:53:VAL:HG22	1:A:144:VAL:HG22	1.98	0.46
1:A:181:VAL:HG12	2:C:938:LYS:HD2	1.96	0.46
3:D:500:ARG:NH1	3:D:1388:ARG:O	2.45	0.46
3:N:350:HIS:CE1	5:P:232:ARG:HG3	2.50	0.46
2:C:911:GLU:O	2:C:915:LYS:HG2	2.16	0.46
3:D:1336:LEU:HB2	3:D:1344:VAL:HG21	1.97	0.46
3:D:1347:TYR:CZ	3:D:1351:GLU:HG3	2.50	0.46
2:M:368:THR:H	2:M:371:LYS:HD3	1.81	0.46
3:N:1462:LEU:HD22	3:N:1472:ILE:HB	1.98	0.46
3:N:316:GLN:NE2	3:N:340:THR:OG1	2.46	0.46
3:N:520:LEU:O	3:N:525:ARG:NE	2.41	0.46
3:N:750:PRO:HG2	3:N:756:GLN:NE2	2.31	0.46
3:N:784:ASP:N	3:N:784:ASP:OD1	2.49	0.46
2:M:184:MET:HE3	2:M:303:PHE:HE1	1.81	0.45
2:M:670:GLN:HG2	2:M:699:PHE:CD2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:N:1101:VAL:HG13	3:N:1102:THR:HG23	1.99	0.45
3:N:959:GLU:HB3	3:N:963:TYR:HE2	1.80	0.45
1:A:115:LEU:HA	1:A:116:PRO:HD3	1.83	0.45
2:C:861:LEU:HD12	2:C:865:THR:HB	1.98	0.45
3:D:356:PRO:HB3	3:D:441:ARG:HA	1.97	0.45
3:D:619:LEU:HD11	3:D:1439:SER:HB2	1.98	0.45
2:C:1009:SER:HB3	3:D:651:GLU:O	2.17	0.45
3:D:658:LEU:HA	3:D:661:MET:HE3	1.98	0.45
5:F:120:THR:HG22	5:F:122:LEU:HD13	1.98	0.45
3:N:101:HIS:CE1	3:N:582:LEU:HD13	2.51	0.45
5:P:265:VAL:HG12	5:P:269:ASN:HD21	1.81	0.45
7:S:1:DT:H2"	7:S:2:DA:H8	1.80	0.45
1:B:128:HIS:CE1	1:B:131:THR:HG23	2.52	0.45
1:B:8:ALA:HB1	1:B:27:PRO:HD2	1.97	0.45
2:C:64:LEU:HD12	2:C:103:LYS:HB2	1.99	0.45
2:C:936:VAL:HG11	2:C:959:PRO:HB2	1.98	0.45
3:D:1296:SER:OG	3:D:1297:GLU:N	2.50	0.45
3:D:634:GLY:HA2	3:D:635:PRO:HD3	1.54	0.45
5:P:144:ILE:HB	5:P:147:LEU:HD13	1.98	0.45
2:M:778:PHE:O	2:M:780:GLU:HB2	2.16	0.45
3:N:43:GLY:H	3:N:46:ASP:HB2	1.80	0.45
5:P:195:VAL:HG21	5:P:217:ASN:OD1	2.16	0.45
2:C:328:LEU:HA	2:C:328:LEU:HD23	1.81	0.45
2:C:708:TYR:HB3	2:C:790:LEU:HD21	1.98	0.45
3:D:703:ASN:HB2	3:D:713:ILE:HG12	1.98	0.45
2:M:214:TYR:HE1	2:M:317:VAL:HG21	1.81	0.45
3:N:1478:SER:HB3	3:N:1481:VAL:HG22	1.99	0.45
3:N:317:VAL:HB	3:N:339:TRP:HB3	1.98	0.45
3:N:637:LEU:HD13	3:N:642:CYS:HA	1.99	0.45
1:A:39:PRO:HG3	1:B:39:PRO:HG3	1.99	0.45
2:C:884:GLN:O	2:C:888:THR:OG1	2.29	0.45
3:D:613:ARG:HG3	3:D:618:LEU:HD23	1.98	0.45
5:F:181:GLU:O	5:F:185:GLN:HG2	2.17	0.45
2:M:214:TYR:CE1	2:M:317:VAL:HG21	2.52	0.45
3:N:1236:LEU:HA	3:N:1359:GLN:HG3	1.98	0.45
3:D:1144:LEU:HA	3:D:1144:LEU:HD23	1.76	0.45
1:L:150:TYR:CE2	1:L:170:VAL:HG22	2.52	0.45
3:N:1102:THR:HG21	3:N:1371:VAL:HG22	1.98	0.45
3:N:218:LYS:HG2	3:N:338:GLU:HG2	1.99	0.45
3:N:829:VAL:HG21	3:N:839:LEU:HD11	1.99	0.45
1:A:20:TYR:OH	1:A:198:ARG:HD2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:685:ASP:HA	3:D:688:TRP:HD1	1.82	0.45
7:H:17:DA:H2'	7:H:18:DC:C6	2.52	0.45
1:K:56:VAL:HG22	1:K:142:VAL:HG12	1.98	0.45
2:C:17:PRO:HB2	2:C:20:GLU:HB3	1.98	0.44
7:H:20:DG:H4'	7:H:21:DA:OP1	2.16	0.44
3:N:1108:ARG:NH2	3:N:1198:TYR:O	2.50	0.44
3:N:1256:LEU:O	3:N:1260:ILE:HG13	2.17	0.44
2:C:177:GLU:HG3	2:C:178:PRO:HD2	1.98	0.44
2:C:299:LYS:HE3	2:C:299:LYS:HB2	1.84	0.44
2:C:419:THR:HG23	2:C:422:ARG:HG3	1.98	0.44
2:C:954:THR:HA	2:C:955:PRO:HD3	1.87	0.44
1:L:110:LYS:HD3	1:L:128:HIS:HA	1.98	0.44
3:N:1048:PRO:HD3	3:N:1075:HIS:CG	2.52	0.44
3:N:356:PRO:HB3	3:N:441:ARG:HA	1.98	0.44
3:N:704:ARG:HD2	3:N:738:ALA:HB2	2.00	0.44
5:P:196:VAL:HG11	7:S:7:DG:H4'	1.98	0.44
1:L:83:LYS:HE2	1:L:168:ASP:HB2	1.99	0.44
2:M:1090:LYS:HE2	2:M:1112:PHE:CZ	2.53	0.44
3:N:256:GLU:OE2	3:N:300:LYS:NZ	2.37	0.44
3:N:270:LEU:HD12	3:N:284:LEU:HD11	1.99	0.44
1:A:150:TYR:CD1	2:C:696:LYS:HG2	2.53	0.44
1:B:14:ARG:HE	1:B:14:ARG:HB2	1.53	0.44
5:F:326:ASP:OD1	5:F:327:SER:N	2.50	0.44
1:K:102:LYS:HE3	1:K:102:LYS:HB2	1.86	0.44
2:M:670:GLN:HG2	2:M:699:PHE:CG	2.53	0.44
3:N:619:LEU:HD11	3:N:1439:SER:HB2	1.99	0.44
2:C:944:LEU:HD22	2:C:962:GLN:HB3	1.99	0.44
3:D:601:ARG:HD3	5:F:318:GLU:HG2	1.99	0.44
1:K:36:LEU:HD11	1:L:221:HIS:HB3	1.98	0.44
2:M:261:ILE:HG23	2:M:290:LEU:HB2	1.99	0.44
2:M:413:LEU:HD21	2:M:451:LEU:HD13	2.00	0.44
3:N:500:ARG:NH1	3:N:1388:ARG:O	2.45	0.44
3:N:864:VAL:HG22	3:N:865:THR:H	1.83	0.44
3:D:968:ASP:OD1	3:D:1058:ARG:NH2	2.51	0.44
1:L:7:LYS:HZ2	1:L:7:LYS:HB3	1.81	0.44
2:M:15:LEU:HB2	2:M:586:ARG:HH11	1.82	0.44
7:S:3:DT:H2'	7:S:4:DA:C8	2.53	0.44
2:M:1116:ALA:HB2	3:N:88:TYR:HB3	2.00	0.44
2:M:224:GLU:CD	2:M:224:GLU:H	2.21	0.44
2:M:462:ASP:OD2	2:M:468:ARG:NH1	2.41	0.44
3:N:1285:GLU:HG3	3:N:1290:LEU:HG	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:N:162:ARG:O	3:N:449:SER:HB2	2.17	0.44
3:N:200:ASP:O	3:N:397:LYS:HG2	2.18	0.44
3:N:671:LYS:HZ2	5:P:421:PHE:HA	1.82	0.44
5:P:212:LEU:HD22	5:P:247:ILE:HG23	1.99	0.44
5:P:285:GLU:HA	5:P:286:PRO:HD3	1.79	0.44
2:C:1038:TRP:NE1	3:D:1099:VAL:HG11	2.33	0.44
2:M:861:LEU:HD12	2:M:865:THR:HB	2.00	0.44
5:P:338:LEU:HA	5:P:339:PRO:HD3	1.86	0.44
1:L:77:GLU:OE1	3:N:867:ARG:NH2	2.46	0.44
4:O:87:LYS:HA	4:O:87:LYS:HD3	1.75	0.44
2:C:168:ARG:O	2:C:267:TYR:HA	2.18	0.43
2:M:437:ARG:HH22	2:M:491:GLU:HG3	1.83	0.43
3:N:314:PRO:HB2	3:N:317:VAL:HG12	1.99	0.43
4:O:88:GLU:OE2	4:O:91:ARG:NH1	2.51	0.43
1:A:183:ASP:HA	2:C:938:LYS:HE3	1.99	0.43
3:D:629:SER:OG	3:D:630:VAL:N	2.50	0.43
3:D:935:LYS:HE2	3:D:935:LYS:HB3	1.79	0.43
5:F:193:ARG:NH1	7:H:7:DG:N7	2.67	0.43
2:M:317:VAL:HA	2:M:318:PRO:HD3	1.88	0.43
2:C:35:PRO:HA	2:C:36:PRO:HD3	1.90	0.43
2:C:35:PRO:HG2	2:C:38:LYS:HD2	2.00	0.43
2:C:394:PHE:CE2	2:C:632:ASN:HB3	2.52	0.43
2:C:710:ILE:HD12	2:C:790:LEU:HB2	2.00	0.43
3:D:602:SER:O	3:D:606:ILE:HG13	2.18	0.43
2:M:150:PRO:HG3	2:M:322:VAL:HG11	2.00	0.43
2:M:792:VAL:HA	2:M:793:PRO:HD3	1.88	0.43
3:N:1347:TYR:CZ	3:N:1351:GLU:HG3	2.53	0.43
3:N:142:LEU:HD13	3:N:161:LEU:HD11	1.99	0.43
3:N:185:VAL:HG21	3:N:203:ALA:HB2	2.00	0.43
3:N:50:PHE:CD1	3:N:522:PRO:HD3	2.53	0.43
5:P:387:GLY:O	5:P:392:VAL:N	2.44	0.43
1:B:11:PHE:CE1	1:B:23:PHE:HB3	2.53	0.43
2:C:224:GLU:CD	2:C:224:GLU:H	2.22	0.43
2:C:27:ARG:HH21	2:C:27:ARG:HB3	1.83	0.43
2:C:999:HIS:HB3	2:C:1004:LYS:NZ	2.33	0.43
3:D:26:VAL:HG11	3:D:44:LEU:HD23	1.99	0.43
5:F:88:ILE:HG23	5:F:193:ARG:HG2	2.00	0.43
1:L:7:LYS:HZ3	1:L:7:LYS:HB3	1.82	0.43
3:N:1296:SER:OG	3:N:1297:GLU:N	2.51	0.43
2:M:893:ALA:HB2	2:M:918:LEU:HD23	2.00	0.43
3:N:136:ASP:HA	3:N:137:PRO:HD3	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:N:563:PRO:HB3	5:P:189:GLU:HG3	2.01	0.43
2:C:405:ARG:HD3	2:C:566:THR:HG21	2.00	0.43
1:B:30:ARG:NH2	2:C:854:PRO:HB3	2.33	0.43
3:D:450:TYR:HB2	3:D:452:ILE:HG13	2.01	0.43
1:K:63:HIS:HD2	2:M:746:GLY:HA2	1.82	0.43
3:N:684:LYS:HE2	3:N:684:LYS:HB3	1.81	0.43
1:A:91:ASN:HA	1:A:92:PRO:HD3	1.89	0.43
1:K:150:TYR:CE2	1:K:152:PRO:HG3	2.53	0.43
2:M:367:LEU:HA	2:M:371:LYS:HD3	2.00	0.43
3:N:633:VAL:HB	3:N:740:PHE:CE2	2.54	0.43
5:P:261:PRO:O	5:P:265:VAL:HG23	2.19	0.43
5:P:364:ARG:HG3	5:P:390:PHE:CE1	2.54	0.43
2:M:595:LEU:HD11	2:M:623:TYR:HB3	2.01	0.43
2:M:825:VAL:HG12	2:M:827:VAL:HG23	2.01	0.43
3:N:1455:LYS:HE3	3:N:1455:LYS:HB2	1.74	0.43
3:N:935:LYS:HE2	3:N:935:LYS:HB3	1.83	0.43
2:M:886:LEU:HD21	3:N:951:ILE:HG12	2.01	0.43
1:A:57:TYR:CD1	1:A:161:ARG:HD2	2.54	0.43
3:D:208:PRO:HG2	3:D:353:VAL:HG21	2.00	0.43
3:N:106:LYS:HD3	3:N:106:LYS:HA	1.88	0.43
2:M:1038:TRP:CE2	3:N:1099:VAL:HG11	2.54	0.43
3:N:658:LEU:HD23	3:N:661:MET:HE1	2.01	0.43
3:N:711:LEU:HB3	3:N:714:GLN:HE21	1.83	0.43
1:A:31:GLY:N	1:A:193:ASP:OD2	2.52	0.43
2:C:1022:GLY:O	2:C:1026:GLN:NE2	2.51	0.43
2:C:390:GLN:HB3	2:C:415:PRO:HD3	2.01	0.43
2:C:797:GLY:O	2:C:829:GLN:NE2	2.51	0.43
3:D:1290:LEU:HD22	3:D:1305:LEU:HD11	2.01	0.43
3:D:142:LEU:HB2	3:D:161:LEU:HD11	2.01	0.43
3:D:778:LEU:HA	3:D:778:LEU:HD12	1.85	0.43
3:D:886:VAL:O	3:D:890:VAL:HG22	2.18	0.43
2:M:122:THR:OG1	2:M:124:ASP:OD1	2.32	0.43
3:D:1463:LYS:HB3	3:D:1463:LYS:HE2	1.85	0.42
3:D:1486:VAL:HG11	4:E:26:ARG:HB2	1.99	0.42
3:D:829:VAL:HG21	3:D:839:LEU:HD11	2.01	0.42
3:N:734:GLU:OE2	3:N:780:LYS:NZ	2.52	0.42
1:A:97:VAL:HG12	1:A:99:LEU:HD12	2.00	0.42
1:B:110:LYS:HD3	1:B:128:HIS:HA	2.02	0.42
2:C:302:VAL:O	2:C:305:PRO:HD2	2.20	0.42
5:F:364:ARG:HG3	5:F:390:PHE:CE1	2.55	0.42
2:M:355:VAL:HG12	2:M:359:MET:HE2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:N:1122:LEU:HD13	3:N:1178:ALA:HB2	2.01	0.42
5:P:360:LYS:HE3	5:P:360:LYS:HB2	1.71	0.42
1:A:172:SER:HA	1:A:173:PRO:HD2	1.89	0.42
3:D:1094:LEU:HD22	3:D:1260:ILE:HG12	2.01	0.42
3:D:411:THR:HG23	3:D:436:GLU:HA	2.00	0.42
2:M:1083:GLU:OE1	2:M:1086:ARG:NH1	2.50	0.42
2:M:1002:GLU:HA	3:N:724:GLN:HE22	1.85	0.42
2:C:584:GLU:HB3	2:C:666:LEU:H	1.84	0.42
5:F:370:LYS:HB3	5:F:376:ILE:HG12	2.01	0.42
6:G:14:DG:H2'	6:G:15:DT:C6	2.54	0.42
2:M:27:ARG:HB3	2:M:27:ARG:HH21	1.85	0.42
2:M:768:THR:N	2:M:769:PRO:HD3	2.35	0.42
3:N:167:GLU:OE2	3:N:198:ARG:NH1	2.53	0.42
3:N:291:LEU:HD12	3:N:303:PRO:HB2	2.00	0.42
1:B:72:LYS:HB3	1:B:131:THR:OG1	2.19	0.42
3:D:12:LEU:HD21	3:D:104:PHE:CZ	2.54	0.42
5:F:120:THR:HG21	5:F:122:LEU:HD22	2.02	0.42
3:N:171:LEU:HD12	3:N:390:PRO:HG2	2.01	0.42
3:N:890:VAL:HB	3:N:922:LEU:HD13	2.01	0.42
5:P:299:TRP:CE3	5:P:303:ARG:HD3	2.55	0.42
2:C:1090:LYS:HA	2:C:1090:LYS:HD3	1.78	0.42
3:D:479:GLU:HA	3:D:482:LYS:HE2	2.01	0.42
3:D:880:ILE:HD13	3:D:880:ILE:HA	1.94	0.42
7:H:9:DG:H2"	7:H:10:DA:OP2	2.19	0.42
2:M:299:LYS:HB2	2:M:299:LYS:HE3	1.79	0.42
5:P:93:LEU:HD21	5:P:193:ARG:HD2	2.01	0.42
2:C:874:LEU:HD23	3:D:1023:MET:SD	2.59	0.42
3:D:1491:THR:O	3:D:1495:ILE:HG13	2.19	0.42
3:D:487:ALA:O	3:D:491:LYS:HG2	2.20	0.42
5:F:193:ARG:HB2	7:H:6:DT:H1'	2.00	0.42
1:K:101:LEU:HD21	1:K:109:VAL:HG11	2.01	0.42
1:K:58:ILE:HG12	1:K:140:MET:HG2	2.01	0.42
1:K:63:HIS:CD2	2:M:746:GLY:HA2	2.55	0.42
2:M:374:ASN:OD1	5:P:276:ARG:HD2	2.18	0.42
3:N:553:ARG:HD3	5:P:214:GLN:HB3	2.01	0.42
2:C:1065:ALA:HB1	2:C:1077:PRO:HG3	2.01	0.42
3:D:1256:LEU:O	3:D:1260:ILE:HG13	2.19	0.42
3:D:171:LEU:HD23	3:D:171:LEU:HA	1.86	0.42
3:D:218:LYS:HG2	3:D:338:GLU:HG2	2.01	0.42
3:D:667:ALA:HA	3:D:668:PRO:HD3	1.93	0.42
2:C:1042:ALA:HB3	3:D:710:ARG:HB3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:715:ALA:HB3	3:D:764:LEU:HA	2.01	0.42
1:B:11:PHE:HE1	1:B:23:PHE:HB3	1.84	0.42
2:C:976:ASP:OD1	2:C:978:ARG:HD3	2.20	0.42
3:D:289:THR:HA	3:D:290:PRO:HD2	1.96	0.42
5:F:188:ILE:HG12	5:F:224:VAL:HG21	2.00	0.42
2:M:774:LEU:HA	2:M:777:ILE:HG12	2.02	0.42
2:C:315:ALA:HB3	2:C:317:VAL:HG23	2.02	0.42
2:C:501:THR:HA	2:C:502:PRO:HD3	1.91	0.42
2:M:999:HIS:HB3	2:M:1004:LYS:NZ	2.34	0.42
3:N:1045:MET:HG2	3:N:1073:SER:HA	2.02	0.42
3:N:1402:ALA:HA	3:N:1405:GLU:HG2	2.02	0.42
2:C:118:ILE:HD11	2:C:344:PHE:CE1	2.54	0.41
2:C:556:ASN:O	2:C:559:LEU:HB3	2.20	0.41
3:D:272:LEU:HB2	3:D:280:ALA:HB3	2.02	0.41
1:L:115:LEU:HA	1:L:116:PRO:HD3	1.84	0.41
2:M:543:ASN:HD21	2:M:566:THR:HG22	1.84	0.41
3:N:1307:LYS:HG3	3:N:1307:LYS:H	1.61	0.41
3:N:1493:LYS:O	3:N:1497:GLU:HG2	2.20	0.41
3:N:632:VAL:O	3:N:727:GLN:HA	2.20	0.41
3:N:881:LEU:O	3:N:885:ILE:HG13	2.20	0.41
1:B:104:GLU:OE2	1:B:137:ARG:NH1	2.53	0.41
3:D:115:LEU:HD23	3:D:115:LEU:HA	1.85	0.41
2:M:344:PHE:CD1	2:M:378:LEU:HD11	2.55	0.41
2:M:859:PRO:O	2:M:867:VAL:HG22	2.20	0.41
3:N:963:TYR:CE1	3:N:1002:LYS:HD3	2.56	0.41
3:N:1038:LEU:O	3:N:1060:SER:HB2	2.20	0.41
2:M:1019:GLN:NE2	3:N:621:LYS:HG2	2.35	0.41
3:N:654:LYS:O	3:N:658:LEU:HG	2.20	0.41
3:D:244:GLU:HG3	3:D:310:LEU:HG	2.01	0.41
3:D:483:HIS:CG	3:D:484:PRO:HD2	2.54	0.41
1:K:133:GLU:HG2	1:K:134:GLU:H	1.85	0.41
2:M:805:ARG:O	2:M:807:ARG:NH2	2.39	0.41
5:P:214:GLN:O	5:P:217:ASN:HB2	2.19	0.41
5:P:276:ARG:O	5:P:279:GLN:HG3	2.20	0.41
1:B:54:THR:OG1	1:B:158:ILE:HD11	2.20	0.41
1:A:38:ASN:ND2	2:C:978:ARG:O	2.54	0.41
3:D:750:PRO:HG2	3:D:756:GLN:NE2	2.35	0.41
2:M:580:MET:HB3	2:M:584:GLU:CD	2.40	0.41
1:K:196:THR:HG21	2:M:934:PHE:HE1	1.86	0.41
1:L:190:THR:HG21	3:N:722:GLU:CD	2.40	0.41
2:C:223:ASP:OD2	2:C:225:SER:OG	2.32	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:984:GLU:OE2	3:D:791:TYR:OH	2.35	0.41
5:F:101:GLU:HG2	5:F:105:LYS:HE2	2.02	0.41
3:D:437:VAL:HG11	5:F:175:HIS:CD2	2.55	0.41
5:F:338:LEU:HA	5:F:339:PRO:HD3	1.84	0.41
2:M:366:SER:O	2:M:371:LYS:NZ	2.36	0.41
2:M:376:ARG:NH1	2:M:379:GLU:OE1	2.53	0.41
3:N:236:TYR:CE1	3:N:242:LEU:HD12	2.56	0.41
3:N:984:THR:HG23	3:N:987:GLU:HB2	2.01	0.41
3:D:1018:ASN:HA	3:D:1019:PRO:HD3	1.91	0.41
3:D:1122:LEU:HD13	3:D:1178:ALA:HB2	2.02	0.41
3:D:784:ASP:HB2	3:D:939:PHE:CE1	2.55	0.41
1:L:179:PHE:HB3	1:L:197:LEU:HD13	2.02	0.41
2:M:302:VAL:O	2:M:305:PRO:HD2	2.20	0.41
3:N:397:LYS:HE2	3:N:397:LYS:HB3	1.82	0.41
5:P:193:ARG:HB3	7:S:7:DG:H5'	2.02	0.41
1:B:100:LEU:HG	1:B:141:GLU:HG2	2.03	0.41
2:M:76:PRO:HA	2:M:77:PRO:HD2	1.97	0.41
2:M:708:TYR:HB3	2:M:790:LEU:HD21	2.02	0.41
5:P:358:LEU:HD12	5:P:370:LYS:HG3	2.02	0.41
3:D:1383:ASP:HA	3:D:1384:PRO:HD3	1.78	0.41
3:D:1429:LEU:HD23	3:D:1429:LEU:HA	1.95	0.41
2:M:177:GLU:OE2	2:M:183:SER:OG	2.35	0.41
3:N:1267:ARG:HA	3:N:1268:PRO:HD3	1.88	0.41
3:N:238:PRO:HD3	3:N:318:ARG:HG3	2.03	0.41
3:N:809:PRO:HG3	3:N:829:VAL:HG11	2.02	0.41
5:P:394:ARG:O	5:P:397:ILE:HG22	2.20	0.41
3:D:1267:ARG:HA	3:D:1268:PRO:HD3	1.87	0.41
6:G:9:DC:H2"	6:G:10:DG:C8	2.56	0.41
1:L:20:TYR:C	1:L:207:PRO:HG2	2.40	0.41
2:M:1048:THR:HG22	3:N:758:GLU:OE2	2.20	0.41
2:M:317:VAL:O	2:M:320:HIS:ND1	2.49	0.41
2:M:679:PHE:CE2	2:M:853:LEU:HD11	2.56	0.41
5:F:276:ARG:O	5:F:279:GLN:HG3	2.21	0.41
5:F:285:GLU:HA	5:F:286:PRO:HD3	1.81	0.41
2:M:304:LEU:HB3	2:M:305:PRO:HD3	2.02	0.41
3:N:770:LEU:HB2	3:N:1210:SER:HA	2.03	0.41
5:P:278:LEU:HD11	5:P:294:ALA:HB3	2.03	0.41
7:S:1:DT:H2"	7:S:2:DA:O5'	2.21	0.41
1:A:174:VAL:HA	1:A:201:THR:HG22	2.03	0.41
3:D:1211:MET:HE1	4:E:16:LYS:HD2	2.01	0.41
3:D:192:ALA:HB3	3:D:195:VAL:HB	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:13:ILE:HD13	2:M:483:VAL:HG11	2.02	0.41
2:M:535:SER:O	2:M:538:GLN:HG2	2.20	0.41
3:N:256:GLU:HG3	3:N:300:LYS:HG3	2.02	0.41
1:B:48:ILE:HA	1:B:49:PRO:HD3	1.83	0.40
2:C:247:PRO:HA	2:C:248:PRO:HD3	1.67	0.40
2:C:335:THR:O	2:C:339:LEU:HG	2.21	0.40
3:D:67:ARG:HD3	3:D:68:PHE:CE2	2.56	0.40
3:D:879:ARG:HD3	3:D:902:LEU:O	2.21	0.40
5:F:144:ILE:HB	5:F:147:LEU:HD13	2.02	0.40
3:N:414:ARG:HD3	3:N:451:ASP:HB2	2.03	0.40
3:N:613:ARG:HG3	3:N:618:LEU:HD23	2.02	0.40
3:N:861:GLN:OE1	3:N:861:GLN:N	2.54	0.40
3:N:956:ILE:HG12	3:N:956:ILE:H	1.72	0.40
3:N:970:LYS:HD3	3:N:970:LYS:O	2.21	0.40
5:P:373:LYS:HA	5:P:373:LYS:HD3	1.90	0.40
2:C:436:GLY:HA2	2:C:538:GLN:O	2.22	0.40
3:D:180:LYS:HE3	3:D:180:LYS:HB3	1.97	0.40
3:D:318:ARG:NH1	3:D:338:GLU:OE1	2.54	0.40
5:F:163:LEU:HD13	5:F:174:LEU:HD13	2.03	0.40
2:M:690:ILE:HB	2:M:694:LEU:HD12	2.02	0.40
2:M:911:GLU:O	2:M:915:LYS:HG2	2.20	0.40
2:M:521:PRO:HB2	3:N:1055:VAL:HG11	2.03	0.40
3:N:972:LEU:HD23	3:N:972:LEU:HA	1.92	0.40
3:D:1108:ARG:NH2	3:D:1198:TYR:O	2.54	0.40
2:M:1110:ASP:OD2	2:M:1114:GLY:N	2.34	0.40
3:N:137:PRO:HB2	3:N:139:GLY:HA2	2.02	0.40
4:O:3:GLU:HA	4:O:4:PRO:HD3	1.95	0.40
2:C:1005:MET:HB3	2:C:1005:MET:HE2	1.96	0.40
2:C:203:ASP:OD1	2:C:204:GLN:N	2.51	0.40
2:C:543:ASN:HD21	2:C:566:THR:HG22	1.86	0.40
3:D:517:VAL:HA	3:D:518:PRO:HD3	1.96	0.40
7:H:22:DT:H2"	7:H:23:DG:C8	2.56	0.40
2:M:469:THR:HG23	2:M:471:TYR:CE2	2.56	0.40
2:M:547:ILE:HA	2:M:548:PRO:HD2	1.96	0.40
4:O:42:PRO:HA	4:O:45:ARG:HD2	2.04	0.40
3:D:1350:GLU:O	3:D:1354:LYS:HG3	2.21	0.40
6:G:6:DA:H5"	6:G:6:DA:H8	1.86	0.40
3:N:1283:ILE:HG13	3:N:1283:ILE:H	1.70	0.40
3:N:841:TYR:HB2	3:N:864:VAL:HG23	2.03	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	229/315 (73%)	226 (99%)	3 (1%)	0	100	100
1	B	218/315 (69%)	212 (97%)	5 (2%)	1 (0%)	34	74
1	K	229/315 (73%)	226 (99%)	3 (1%)	0	100	100
1	L	218/315 (69%)	212 (97%)	6 (3%)	0	100	100
2	C	1108/1119 (99%)	1085 (98%)	22 (2%)	1 (0%)	56	90
2	M	1074/1119 (96%)	1049 (98%)	25 (2%)	0	100	100
3	D	1482/1524 (97%)	1451 (98%)	31 (2%)	0	100	100
3	N	1482/1524 (97%)	1446 (98%)	35 (2%)	1 (0%)	56	90
4	E	92/99 (93%)	89 (97%)	3 (3%)	0	100	100
4	O	92/99 (93%)	89 (97%)	3 (3%)	0	100	100
5	F	344/443 (78%)	340 (99%)	4 (1%)	0	100	100
5	P	344/443 (78%)	338 (98%)	5 (2%)	1 (0%)	46	82
All	All	6912/7630 (91%)	6763 (98%)	145 (2%)	4 (0%)	56	90

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
5	P	361	LEU
3	N	207	PHE
1	B	8	ALA
2	C	23	VAL

5.3.2 Protein sidechains (i)

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	200/273 (73%)	199 (100%)	1 (0%)	92 96
1	B	195/273 (71%)	193 (99%)	2 (1%)	82 93
1	K	200/273 (73%)	197 (98%)	3 (2%)	72 90
1	L	195/273 (71%)	191 (98%)	4 (2%)	61 86
2	C	936/941 (100%)	910 (97%)	26 (3%)	51 82
2	M	908/941 (96%)	876 (96%)	32 (4%)	43 78
3	D	1253/1279 (98%)	1225 (98%)	28 (2%)	60 86
3	N	1253/1279 (98%)	1221 (97%)	32 (3%)	54 83
4	E	82/88 (93%)	80 (98%)	2 (2%)	57 84
4	O	82/88 (93%)	76 (93%)	6 (7%)	17 55
5	F	301/388 (78%)	294 (98%)	7 (2%)	58 85
5	P	301/388 (78%)	280 (93%)	21 (7%)	19 57
All	All	5906/6484 (91%)	5742 (97%)	164 (3%)	51 82

All (164) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
1	B	14	ARG
1	B	154	GLU
2	C	27	ARG
2	C	133	ASP
2	C	141	HIS
2	C	230	ARG
2	C	251	ASP
2	C	274	ARG
2	C	276	LYS
2	C	284	ARG
2	C	384	GLU
2	C	390	GLN
2	C	419	THR
2	C	429	ASP
2	C	454	SER
2	C	464	LEU
2	C	565	GLN
2	C	575	GLN
2	C	583	LEU

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Mol	Chain	Res	Type
2	C	595	LEU
2	C	610	ARG
2	C	617	ASP
2	C	640	ARG
2	C	657	ASP
2	C	670	GLN
2	C	698	ASP
2	C	775	ARG
2	C	1043	TYR
3	D	67	ARG
3	D	71	LYS
3	D	102	ILE
3	D	231	VAL
3	D	452	ILE
3	D	636	GLN
3	D	687	VAL
3	D	709	HIS
3	D	710	ARG
3	D	743	ASP
3	D	754	PHE
3	D	784	ASP
3	D	810	GLU
3	D	894	LYS
3	D	907	GLU
3	D	986	ARG
3	D	1046	GLN
3	D	1067	VAL
3	D	1083	ASP
3	D	1155	VAL
3	D	1184	GLN
3	D	1252	ILE
3	D	1253	THR
3	D	1267	ARG
3	D	1295	GLU
3	D	1307	LYS
3	D	1433	SER
3	D	1501	GLU
4	E	49	GLN
4	E	50	THR
5	F	88	ILE
5	F	295	MET
5	F	321	ILE

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Mol	Chain	Res	Type
5	F	324	GLU
5	F	379	ARG
5	F	380	GLU
5	F	417	LYS
1	K	6	LEU
1	K	65	PHE
1	K	126	ASP
1	L	6	LEU
1	L	7	LYS
1	L	126	ASP
1	L	154	GLU
2	M	27	ARG
2	M	118	ILE
2	M	141	HIS
2	M	162	ILE
2	M	274	ARG
2	M	276	LYS
2	M	284	ARG
2	M	345	ARG
2	M	360	LEU
2	M	364	GLU
2	M	368	THR
2	M	372	LEU
2	M	384	GLU
2	M	419	THR
2	M	429	ASP
2	M	454	SER
2	M	489	THR
2	M	492	ASP
2	M	513	VAL
2	M	566	THR
2	M	575	GLN
2	M	581	THR
2	M	583	LEU
2	M	610	ARG
2	M	640	ARG
2	M	670	GLN
2	M	723	THR
2	M	774	LEU
2	M	778	PHE
2	M	780	GLU
2	M	807	ARG

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Mol	Chain	Res	Type
2	M	995	MET
3	N	35	ARG
3	N	67	ARG
3	N	71	LYS
3	N	133	ILE
3	N	136	ASP
3	N	311	LEU
3	N	362	GLU
3	N	371	ILE
3	N	374	GLU
3	N	375	GLU
3	N	387	LEU
3	N	388	HIS
3	N	452	ILE
3	N	632	VAL
3	N	709	HIS
3	N	754	PHE
3	N	784	ASP
3	N	810	GLU
3	N	894	LYS
3	N	907	GLU
3	N	984	THR
3	N	986	ARG
3	N	1083	ASP
3	N	1188	VAL
3	N	1219	GLU
3	N	1252	ILE
3	N	1253	THR
3	N	1290	LEU
3	N	1295	GLU
3	N	1307	LYS
3	N	1433	SER
3	N	1501	GLU
4	O	49	GLN
4	O	50	THR
4	O	82	GLU
4	O	84	ARG
4	O	91	ARG
4	O	92	LEU
5	P	88	ILE
5	P	95	THR
5	P	96	LEU

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Mol	Chain	Res	Type
5	P	97	GLU
5	P	205	ARG
5	P	209	PHE
5	P	218	GLN
5	P	230	LYS
5	P	232	ARG
5	P	248	ASN
5	P	260	ILE
5	P	262	VAL
5	P	295	MET
5	P	318	GLU
5	P	349	LEU
5	P	358	LEU
5	P	359	SER
5	P	361	LEU
5	P	362	SER
5	P	369	LEU
5	P	380	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	212	ASN
2	C	219	GLN
2	C	1026	GLN
2	C	1047	HIS
3	D	66	GLN
3	D	669	ASN
3	D	696	HIS
3	D	724	GLN
3	D	994	GLN
3	D	1046	GLN
3	D	1184	GLN
5	F	83	GLN
1	K	63	HIS
1	L	212	ASN
2	M	219	GLN
2	M	343	GLN
2	M	538	GLN
2	M	899	GLN
2	M	1047	HIS
3	N	66	GLN

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Mol	Chain	Res	Type
3	N	350	HIS
3	N	611	GLN
3	N	636	GLN
3	N	669	ASN
3	N	714	GLN
3	N	724	GLN
3	N	976	GLN
3	N	1046	GLN
3	N	1172	HIS
5	P	83	GLN
5	P	218	GLN
5	P	248	ASN
5	P	269	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 carbohydrates in this entry.

5.6 Ligand geometry

Of 21 ligands modelled in this entry, 15 are monoatomic - leaving 6 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
10	C	D	2005	8,12	14,21,22	0.70	0	18,30,33	0.60	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
11	CTP	D	2006	8	6,8,30	1.61	1 (16%)	6,13,47	0.61	0
12	ATP	D	2008	10	26,33,33	1.57	5 (19%)	26,52,52	2.13	1 (3%)
10	C	N	1605	8,12	14,21,22	0.75	0	18,30,33	0.56	0
11	CTP	N	1606	8	6,8,30	1.06	1 (16%)	6,13,47	0.26	0
12	ATP	N	1608	10	26,33,33	1.60	6 (23%)	26,52,52	2.18	2 (7%)

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
10	C	D	2005	8,12	-	0/3/25/26	0/2/2/2
11	CTP	D	2006	8	-	0/6/6/38	0/0/0/2
12	ATP	D	2008	10	-	0/18/38/38	0/3/3/3
10	C	N	1605	8,12	-	0/3/25/26	0/2/2/2
11	CTP	N	1606	8	-	0/6/6/38	0/0/0/2
12	ATP	N	1608	10	-	0/18/38/38	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	N	1608	ATP	C2'-C1'	-3.20	1.48	1.53
12	D	2008	ATP	C2'-C1'	-3.18	1.48	1.53
12	N	1608	ATP	O3'-C3'	-2.21	1.37	1.43
12	D	2008	ATP	O3'-C3'	-2.11	1.38	1.43
12	N	1608	ATP	C2'-C3'	-2.01	1.48	1.53
11	N	1606	CTP	PA-O5'	2.02	1.61	1.54
12	N	1608	ATP	C2-N1	2.26	1.38	1.33
12	D	2008	ATP	C2-N1	2.27	1.38	1.33
12	D	2008	ATP	C6-N6	2.30	1.43	1.34
12	N	1608	ATP	C6-N6	2.32	1.43	1.34
11	D	2006	CTP	PA-O1A	3.45	1.61	1.50
12	N	1608	ATP	C2-N3	4.30	1.39	1.32
12	D	2008	ATP	C2-N3	4.43	1.40	1.32

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
12	N	1608	ATP	N3-C2-N1	-10.20	120.86	128.87
12	D	2008	ATP	N3-C2-N1	-9.86	121.12	128.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
12	N	1608	ATP	C2-N1-C6	2.03	122.39	118.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	N	1605	C	1	0

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^{th} 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	231/315 (73%)	-0.32	3 (1%) 79 72	13, 36, 63, 128	1 (0%)
1	B	222/315 (70%)	-0.25	0 100 100	16, 42, 80, 94	0
1	K	231/315 (73%)	-0.14	4 (1%) 73 65	18, 47, 73, 123	1 (0%)
1	L	222/315 (70%)	-0.27	0 100 100	17, 46, 87, 114	0
2	C	1112/1119 (99%)	-0.28	11 (0%) 84 79	4, 31, 89, 122	3 (0%)
2	M	1080/1119 (96%)	0.18	55 (5%) 32 24	5, 58, 121, 135	2 (0%)
3	D	1486/1524 (97%)	-0.23	15 (1%) 84 79	3, 32, 90, 122	4 (0%)
3	N	1486/1524 (97%)	-0.08	35 (2%) 62 53	3, 40, 103, 143	4 (0%)
4	E	94/99 (94%)	-0.43	0 100 100	11, 29, 67, 94	0
4	O	94/99 (94%)	-0.40	0 100 100	13, 40, 81, 99	0
5	F	346/443 (78%)	-0.10	9 (2%) 59 51	13, 47, 108, 124	0
5	P	346/443 (78%)	0.30	27 (7%) 16 12	29, 74, 142, 167	0
6	G	15/19 (78%)	-0.33	0 100 100	15, 34, 132, 135	0
6	R	15/19 (78%)	-0.41	0 100 100	24, 43, 131, 133	0
7	H	20/27 (74%)	-0.30	0 100 100	36, 66, 117, 125	0
7	S	18/27 (66%)	-0.22	0 100 100	63, 80, 125, 140	0
All	All	7018/7722 (90%)	-0.12	159 (2%) 64 55	3, 42, 105, 167	15 (0%)

All (159) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	P	377	ASP	7.9
5	P	391	GLY	7.4
5	P	392	VAL	5.9
5	P	381	HIS	5.9
3	N	70	GLY	5.2

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Mol	Chain	Res	Type	RSRZ
2	M	200	LEU	4.6
5	P	414	ARG	4.4
3	N	1127	GLU	4.2
2	M	368	THR	4.1
2	M	293	PHE	4.0
2	M	367	LEU	4.0
1	K	233	VAL	4.0
2	M	196	LEU	4.0
1	A	234	ALA	3.9
1	A	233	VAL	3.9
2	C	365	ASP	3.9
2	M	191	PHE	3.8
2	M	251	ASP	3.7
3	N	67	ARG	3.7
5	P	386	VAL	3.7
2	M	188	LYS	3.7
5	P	417	LYS	3.7
2	C	63	GLY	3.6
2	M	362	GLY	3.6
2	M	778	PHE	3.6
2	M	344	PHE	3.6
3	N	1497	GLU	3.6
2	C	778	PHE	3.5
1	K	232	ALA	3.5
2	M	809	GLY	3.5
2	M	48	PHE	3.5
2	M	221	LEU	3.5
2	M	211	LEU	3.4
5	P	394	ARG	3.4
2	M	217	LEU	3.4
1	A	232	ALA	3.3
2	M	311	PHE	3.3
3	N	365	ASP	3.3
3	N	219	GLU	3.3
5	P	138	SER	3.3
2	M	207	LEU	3.3
2	M	300	ASP	3.2
2	M	190	LYS	3.2
3	D	1252	ILE	3.2
2	M	242	LEU	3.1
2	M	181	VAL	3.1
2	C	779	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
2	M	246	ASP	3.0
1	K	234	ALA	3.0
5	P	279	GLN	3.0
5	P	382	THR	3.0
5	P	389	PHE	3.0
5	P	393	THR	2.9
3	N	1252	ILE	2.9
5	P	412	GLU	2.9
3	D	322	VAL	2.9
2	M	52	PHE	2.9
2	M	307	LEU	2.9
5	P	408	LEU	2.8
3	N	1128	VAL	2.8
3	N	1495	ILE	2.8
3	D	1128	VAL	2.8
2	C	777	ILE	2.8
2	M	159	ILE	2.7
2	M	365	ASP	2.7
5	P	146	GLY	2.7
2	M	766	GLU	2.7
2	M	185	LYS	2.7
3	N	173	PRO	2.7
3	D	977	ALA	2.7
2	M	249	LYS	2.7
2	M	86	LYS	2.6
2	M	202	TYR	2.6
2	M	184	MET	2.6
3	N	1501	GLU	2.6
3	N	278	PRO	2.6
3	N	68	PHE	2.6
2	M	1	MET	2.6
2	M	649	VAL	2.6
2	M	739	GLU	2.6
3	D	144	GLY	2.6
3	N	276	ASP	2.6
3	D	982	PHE	2.6
3	D	1127	GLU	2.6
3	D	143	ASN	2.6
3	N	191	LEU	2.5
3	N	326	GLU	2.5
2	C	104	ASP	2.5
3	D	241	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
3	N	218	LYS	2.5
3	N	394	LEU	2.5
5	P	149	GLU	2.5
5	F	415	THR	2.5
3	N	66	GLN	2.5
2	M	340	MET	2.5
2	M	189	ARG	2.5
3	N	338	GLU	2.5
5	F	373	LYS	2.4
3	D	1287	GLU	2.4
5	P	415	THR	2.4
3	D	335	LEU	2.4
3	N	69	GLU	2.4
5	P	416	ARG	2.4
2	M	770	GLU	2.4
2	C	190	LYS	2.4
5	F	382	THR	2.4
2	M	254	VAL	2.3
3	N	355	VAL	2.3
3	D	1408	ILE	2.3
3	D	976	GLN	2.3
2	M	363	SER	2.3
2	C	766	GLU	2.3
3	N	443	VAL	2.3
5	P	142	ARG	2.3
5	P	395	GLU	2.3
3	N	269	PHE	2.3
5	P	148	LYS	2.3
5	P	388	ALA	2.3
3	D	219	GLU	2.2
2	M	807	ARG	2.2
5	F	393	THR	2.2
5	F	386	VAL	2.2
3	N	289	THR	2.2
3	N	352	ASN	2.2
5	F	419	ARG	2.2
3	N	318	ARG	2.2
2	M	174	LEU	2.2
2	M	303	PHE	2.2
2	M	294	GLU	2.2
5	P	410	TYR	2.2
5	P	376	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
2	M	295	ASP	2.2
5	F	149	GLU	2.2
2	M	372	LEU	2.2
3	N	1287	GLU	2.2
2	M	183	SER	2.2
2	C	616	GLU	2.1
3	N	1499	ARG	2.1
3	N	423	ASP	2.1
3	N	217	LYS	2.1
2	M	177	GLU	2.1
2	M	179	ASN	2.1
3	N	1269	LYS	2.1
2	M	208	ALA	2.1
2	M	44	ILE	2.1
1	K	158	ILE	2.1
5	F	375	LEU	2.1
5	P	411	HIS	2.1
3	N	216	VAL	2.1
2	C	66	LEU	2.1
5	F	391	GLY	2.0
3	D	384	VAL	2.0
3	N	422	ALA	2.0
3	N	236	TYR	2.0
2	M	620	LEU	2.0
2	M	779	GLY	2.0
2	M	313	LEU	2.0
5	P	418	LEU	2.0
2	C	105	THR	2.0

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

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

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
8	MG	D	2007	1/1	0.95	0.34	16.04	17,17,17,17	0
11	CTP	N	1606	9/29	0.97	0.32	14.11	34,40,54,72	0
8	MG	N	1607	1/1	0.95	0.30	11.60	30,30,30,30	0
11	CTP	D	2006	9/29	0.94	0.30	8.15	27,37,51,69	0
8	MG	B	2001	1/1	0.37	0.49	6.10	70,70,70,70	0
8	MG	K	901	1/1	0.96	0.26	2.45	36,36,36,36	0
10	C	D	2005	20/21	0.94	0.17	0.20	16,22,28,34	0
10	C	N	1605	20/21	0.93	0.17	-0.12	23,34,41,41	0
12	ATP	N	1608	31/31	0.92	0.17	-0.44	28,39,97,99	0
8	MG	D	2004	1/1	0.95	0.18	-0.58	29,29,29,29	0
12	ATP	D	2008	31/31	0.95	0.15	-0.83	18,28,68,77	0
9	ZN	N	1602	1/1	0.99	0.14	-0.96	8,8,8,8	0
8	MG	N	1601	1/1	0.99	0.16	-1.04	25,25,25,25	0
9	ZN	D	2001	1/1	0.99	0.12	-1.37	10,10,10,10	0
9	ZN	N	1603	1/1	0.91	0.06	-1.88	128,128,128,128	0
8	MG	P	2001	1/1	0.98	0.06	-2.33	77,77,77,77	0
9	ZN	D	2002	1/1	0.98	0.03	-5.75	92,92,92,92	0
8	MG	D	2003	1/1	0.99	0.26	-	6,6,6,6	0
8	MG	N	1604	1/1	0.99	0.21	-	12,12,12,12	0
8	MG	F	2001	1/1	0.95	0.05	-	31,31,31,31	0
8	MG	L	2001	1/1	0.69	0.29	-	72,72,72,72	0

6.5 Other polymers (i)

There are no such residues in this entry.