



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 29, 2022 – 12:32 PM EDT

PDB ID : 6V96  
Title : Agrobacterium tumefaciens ADP-Glucose pyrophosphorylase-S72E  
Authors : Zheng, Y.; Hussien, R.; Alghamdi, M.A.; Ballicora, M.A.; Liu, D.  
Deposited on : 2019-12-13  
Resolution : 1.80 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at <http://www.wwpdb.org/validation/2017/FAQs#types>.

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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	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.29
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.29

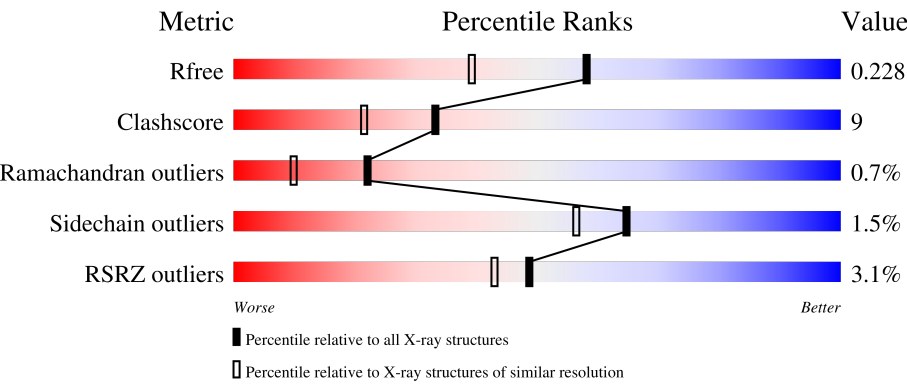
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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))
$R_{free}$	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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

Mol	Chain	Length	Quality of chain
1	A	440	<div><div>3%</div><div></div><div>78%</div><div>15%</div><div>6%</div></div>
1	B	440	<div><div>3%</div><div></div><div>77%</div><div>17%</div><div>6%</div></div>
1	C	440	<div><div>2%</div><div></div><div>78%</div><div>14%</div><div>7%</div></div>
1	D	440	<div><div>4%</div><div></div><div>79%</div><div>15%</div><div>6%</div></div>
1	E	440	<div><div>2%</div><div></div><div>80%</div><div>13%</div><div>6%</div></div>

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Mol	Chain	Length	Quality of chain
1	F	440	
1	G	440	
1	H	440	
1	I	440	
1	J	440	
1	K	440	
1	L	440	
1	M	440	
1	N	440	
1	O	440	
1	P	440	
1	Q	440	
1	R	440	
1	T	440	
1	V	440	

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
3	GOL	J	503	-	-	X	-
3	GOL	K	503	-	-	X	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 74015 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 Glucose-1-phosphate adenylyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	415	Total	C	N	O	S	0	6	0
			3291	2086	571	621	13			
1	A	415	Total	C	N	O	S	0	6	0
			3287	2084	570	620	13			
1	B	415	Total	C	N	O	S	0	6	0
			3287	2084	570	620	13			
1	C	411	Total	C	N	O	S	0	7	0
			3266	2072	564	616	14			
1	D	415	Total	C	N	O	S	0	6	0
			3281	2081	567	620	13			
1	E	415	Total	C	N	O	S	0	6	0
			3287	2084	570	620	13			
1	F	415	Total	C	N	O	S	0	6	0
			3287	2084	570	620	13			
1	G	415	Total	C	N	O	S	0	7	0
			3293	2087	571	621	14			
1	I	415	Total	C	N	O	S	0	7	0
			3293	2087	571	621	14			
1	J	408	Total	C	N	O	S	0	7	0
			3248	2062	561	611	14			
1	K	415	Total	C	N	O	S	0	6	0
			3287	2084	570	620	13			
1	L	415	Total	C	N	O	S	0	7	0
			3295	2089	571	621	14			
1	M	415	Total	C	N	O	S	0	6	0
			3287	2084	570	620	13			
1	N	415	Total	C	N	O	S	0	6	0
			3287	2084	570	620	13			
1	O	415	Total	C	N	O	S	0	8	0
			3301	2092	572	622	15			
1	P	415	Total	C	N	O	S	0	8	0
			3310	2099	575	623	13			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	415	Total	C	N	O	S	0	7	0
			3299	2093	571	622	13			
1	R	415	Total	C	N	O	S	0	7	0
			3295	2089	571	621	14			
1	V	415	Total	C	N	O	S	4	6	0
			3287	2084	570	620	13			
1	T	415	Total	C	N	O	S	12	7	0
			3295	2089	571	621	14			

There are 420 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	-19	MET	-	expression tag	UNP Q8U8L5
H	-18	GLY	-	expression tag	UNP Q8U8L5
H	-17	SER	-	expression tag	UNP Q8U8L5
H	-16	SER	-	expression tag	UNP Q8U8L5
H	-15	HIS	-	expression tag	UNP Q8U8L5
H	-14	HIS	-	expression tag	UNP Q8U8L5
H	-13	HIS	-	expression tag	UNP Q8U8L5
H	-12	HIS	-	expression tag	UNP Q8U8L5
H	-11	HIS	-	expression tag	UNP Q8U8L5
H	-10	HIS	-	expression tag	UNP Q8U8L5
H	-9	SER	-	expression tag	UNP Q8U8L5
H	-8	SER	-	expression tag	UNP Q8U8L5
H	-7	GLY	-	expression tag	UNP Q8U8L5
H	-6	LEU	-	expression tag	UNP Q8U8L5
H	-5	VAL	-	expression tag	UNP Q8U8L5
H	-4	PRO	-	expression tag	UNP Q8U8L5
H	-3	ARG	-	expression tag	UNP Q8U8L5
H	-2	GLY	-	expression tag	UNP Q8U8L5
H	-1	SER	-	expression tag	UNP Q8U8L5
H	0	HIS	-	expression tag	UNP Q8U8L5
H	72	GLU	SER	engineered mutation	UNP Q8U8L5
A	-19	MET	-	expression tag	UNP Q8U8L5
A	-18	GLY	-	expression tag	UNP Q8U8L5
A	-17	SER	-	expression tag	UNP Q8U8L5
A	-16	SER	-	expression tag	UNP Q8U8L5
A	-15	HIS	-	expression tag	UNP Q8U8L5
A	-14	HIS	-	expression tag	UNP Q8U8L5
A	-13	HIS	-	expression tag	UNP Q8U8L5
A	-12	HIS	-	expression tag	UNP Q8U8L5
A	-11	HIS	-	expression tag	UNP Q8U8L5
A	-10	HIS	-	expression tag	UNP Q8U8L5

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	SER	-	expression tag	UNP Q8U8L5
A	-8	SER	-	expression tag	UNP Q8U8L5
A	-7	GLY	-	expression tag	UNP Q8U8L5
A	-6	LEU	-	expression tag	UNP Q8U8L5
A	-5	VAL	-	expression tag	UNP Q8U8L5
A	-4	PRO	-	expression tag	UNP Q8U8L5
A	-3	ARG	-	expression tag	UNP Q8U8L5
A	-2	GLY	-	expression tag	UNP Q8U8L5
A	-1	SER	-	expression tag	UNP Q8U8L5
A	0	HIS	-	expression tag	UNP Q8U8L5
A	72	GLU	SER	engineered mutation	UNP Q8U8L5
B	-19	MET	-	expression tag	UNP Q8U8L5
B	-18	GLY	-	expression tag	UNP Q8U8L5
B	-17	SER	-	expression tag	UNP Q8U8L5
B	-16	SER	-	expression tag	UNP Q8U8L5
B	-15	HIS	-	expression tag	UNP Q8U8L5
B	-14	HIS	-	expression tag	UNP Q8U8L5
B	-13	HIS	-	expression tag	UNP Q8U8L5
B	-12	HIS	-	expression tag	UNP Q8U8L5
B	-11	HIS	-	expression tag	UNP Q8U8L5
B	-10	HIS	-	expression tag	UNP Q8U8L5
B	-9	SER	-	expression tag	UNP Q8U8L5
B	-8	SER	-	expression tag	UNP Q8U8L5
B	-7	GLY	-	expression tag	UNP Q8U8L5
B	-6	LEU	-	expression tag	UNP Q8U8L5
B	-5	VAL	-	expression tag	UNP Q8U8L5
B	-4	PRO	-	expression tag	UNP Q8U8L5
B	-3	ARG	-	expression tag	UNP Q8U8L5
B	-2	GLY	-	expression tag	UNP Q8U8L5
B	-1	SER	-	expression tag	UNP Q8U8L5
B	0	HIS	-	expression tag	UNP Q8U8L5
B	72	GLU	SER	engineered mutation	UNP Q8U8L5
C	-19	MET	-	expression tag	UNP Q8U8L5
C	-18	GLY	-	expression tag	UNP Q8U8L5
C	-17	SER	-	expression tag	UNP Q8U8L5
C	-16	SER	-	expression tag	UNP Q8U8L5
C	-15	HIS	-	expression tag	UNP Q8U8L5
C	-14	HIS	-	expression tag	UNP Q8U8L5
C	-13	HIS	-	expression tag	UNP Q8U8L5
C	-12	HIS	-	expression tag	UNP Q8U8L5
C	-11	HIS	-	expression tag	UNP Q8U8L5
C	-10	HIS	-	expression tag	UNP Q8U8L5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-9	SER	-	expression tag	UNP Q8U8L5
C	-8	SER	-	expression tag	UNP Q8U8L5
C	-7	GLY	-	expression tag	UNP Q8U8L5
C	-6	LEU	-	expression tag	UNP Q8U8L5
C	-5	VAL	-	expression tag	UNP Q8U8L5
C	-4	PRO	-	expression tag	UNP Q8U8L5
C	-3	ARG	-	expression tag	UNP Q8U8L5
C	-2	GLY	-	expression tag	UNP Q8U8L5
C	-1	SER	-	expression tag	UNP Q8U8L5
C	0	HIS	-	expression tag	UNP Q8U8L5
C	72	GLU	SER	engineered mutation	UNP Q8U8L5
D	-19	MET	-	expression tag	UNP Q8U8L5
D	-18	GLY	-	expression tag	UNP Q8U8L5
D	-17	SER	-	expression tag	UNP Q8U8L5
D	-16	SER	-	expression tag	UNP Q8U8L5
D	-15	HIS	-	expression tag	UNP Q8U8L5
D	-14	HIS	-	expression tag	UNP Q8U8L5
D	-13	HIS	-	expression tag	UNP Q8U8L5
D	-12	HIS	-	expression tag	UNP Q8U8L5
D	-11	HIS	-	expression tag	UNP Q8U8L5
D	-10	HIS	-	expression tag	UNP Q8U8L5
D	-9	SER	-	expression tag	UNP Q8U8L5
D	-8	SER	-	expression tag	UNP Q8U8L5
D	-7	GLY	-	expression tag	UNP Q8U8L5
D	-6	LEU	-	expression tag	UNP Q8U8L5
D	-5	VAL	-	expression tag	UNP Q8U8L5
D	-4	PRO	-	expression tag	UNP Q8U8L5
D	-3	ARG	-	expression tag	UNP Q8U8L5
D	-2	GLY	-	expression tag	UNP Q8U8L5
D	-1	SER	-	expression tag	UNP Q8U8L5
D	0	HIS	-	expression tag	UNP Q8U8L5
D	72	GLU	SER	engineered mutation	UNP Q8U8L5
E	-19	MET	-	expression tag	UNP Q8U8L5
E	-18	GLY	-	expression tag	UNP Q8U8L5
E	-17	SER	-	expression tag	UNP Q8U8L5
E	-16	SER	-	expression tag	UNP Q8U8L5
E	-15	HIS	-	expression tag	UNP Q8U8L5
E	-14	HIS	-	expression tag	UNP Q8U8L5
E	-13	HIS	-	expression tag	UNP Q8U8L5
E	-12	HIS	-	expression tag	UNP Q8U8L5
E	-11	HIS	-	expression tag	UNP Q8U8L5
E	-10	HIS	-	expression tag	UNP Q8U8L5

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-9	SER	-	expression tag	UNP Q8U8L5
E	-8	SER	-	expression tag	UNP Q8U8L5
E	-7	GLY	-	expression tag	UNP Q8U8L5
E	-6	LEU	-	expression tag	UNP Q8U8L5
E	-5	VAL	-	expression tag	UNP Q8U8L5
E	-4	PRO	-	expression tag	UNP Q8U8L5
E	-3	ARG	-	expression tag	UNP Q8U8L5
E	-2	GLY	-	expression tag	UNP Q8U8L5
E	-1	SER	-	expression tag	UNP Q8U8L5
E	0	HIS	-	expression tag	UNP Q8U8L5
E	72	GLU	SER	engineered mutation	UNP Q8U8L5
F	-19	MET	-	expression tag	UNP Q8U8L5
F	-18	GLY	-	expression tag	UNP Q8U8L5
F	-17	SER	-	expression tag	UNP Q8U8L5
F	-16	SER	-	expression tag	UNP Q8U8L5
F	-15	HIS	-	expression tag	UNP Q8U8L5
F	-14	HIS	-	expression tag	UNP Q8U8L5
F	-13	HIS	-	expression tag	UNP Q8U8L5
F	-12	HIS	-	expression tag	UNP Q8U8L5
F	-11	HIS	-	expression tag	UNP Q8U8L5
F	-10	HIS	-	expression tag	UNP Q8U8L5
F	-9	SER	-	expression tag	UNP Q8U8L5
F	-8	SER	-	expression tag	UNP Q8U8L5
F	-7	GLY	-	expression tag	UNP Q8U8L5
F	-6	LEU	-	expression tag	UNP Q8U8L5
F	-5	VAL	-	expression tag	UNP Q8U8L5
F	-4	PRO	-	expression tag	UNP Q8U8L5
F	-3	ARG	-	expression tag	UNP Q8U8L5
F	-2	GLY	-	expression tag	UNP Q8U8L5
F	-1	SER	-	expression tag	UNP Q8U8L5
F	0	HIS	-	expression tag	UNP Q8U8L5
F	72	GLU	SER	engineered mutation	UNP Q8U8L5
G	-19	MET	-	expression tag	UNP Q8U8L5
G	-18	GLY	-	expression tag	UNP Q8U8L5
G	-17	SER	-	expression tag	UNP Q8U8L5
G	-16	SER	-	expression tag	UNP Q8U8L5
G	-15	HIS	-	expression tag	UNP Q8U8L5
G	-14	HIS	-	expression tag	UNP Q8U8L5
G	-13	HIS	-	expression tag	UNP Q8U8L5
G	-12	HIS	-	expression tag	UNP Q8U8L5
G	-11	HIS	-	expression tag	UNP Q8U8L5
G	-10	HIS	-	expression tag	UNP Q8U8L5

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-9	SER	-	expression tag	UNP Q8U8L5
G	-8	SER	-	expression tag	UNP Q8U8L5
G	-7	GLY	-	expression tag	UNP Q8U8L5
G	-6	LEU	-	expression tag	UNP Q8U8L5
G	-5	VAL	-	expression tag	UNP Q8U8L5
G	-4	PRO	-	expression tag	UNP Q8U8L5
G	-3	ARG	-	expression tag	UNP Q8U8L5
G	-2	GLY	-	expression tag	UNP Q8U8L5
G	-1	SER	-	expression tag	UNP Q8U8L5
G	0	HIS	-	expression tag	UNP Q8U8L5
G	72	GLU	SER	engineered mutation	UNP Q8U8L5
I	-19	MET	-	expression tag	UNP Q8U8L5
I	-18	GLY	-	expression tag	UNP Q8U8L5
I	-17	SER	-	expression tag	UNP Q8U8L5
I	-16	SER	-	expression tag	UNP Q8U8L5
I	-15	HIS	-	expression tag	UNP Q8U8L5
I	-14	HIS	-	expression tag	UNP Q8U8L5
I	-13	HIS	-	expression tag	UNP Q8U8L5
I	-12	HIS	-	expression tag	UNP Q8U8L5
I	-11	HIS	-	expression tag	UNP Q8U8L5
I	-10	HIS	-	expression tag	UNP Q8U8L5
I	-9	SER	-	expression tag	UNP Q8U8L5
I	-8	SER	-	expression tag	UNP Q8U8L5
I	-7	GLY	-	expression tag	UNP Q8U8L5
I	-6	LEU	-	expression tag	UNP Q8U8L5
I	-5	VAL	-	expression tag	UNP Q8U8L5
I	-4	PRO	-	expression tag	UNP Q8U8L5
I	-3	ARG	-	expression tag	UNP Q8U8L5
I	-2	GLY	-	expression tag	UNP Q8U8L5
I	-1	SER	-	expression tag	UNP Q8U8L5
I	0	HIS	-	expression tag	UNP Q8U8L5
I	72	GLU	SER	engineered mutation	UNP Q8U8L5
J	-19	MET	-	expression tag	UNP Q8U8L5
J	-18	GLY	-	expression tag	UNP Q8U8L5
J	-17	SER	-	expression tag	UNP Q8U8L5
J	-16	SER	-	expression tag	UNP Q8U8L5
J	-15	HIS	-	expression tag	UNP Q8U8L5
J	-14	HIS	-	expression tag	UNP Q8U8L5
J	-13	HIS	-	expression tag	UNP Q8U8L5
J	-12	HIS	-	expression tag	UNP Q8U8L5
J	-11	HIS	-	expression tag	UNP Q8U8L5
J	-10	HIS	-	expression tag	UNP Q8U8L5

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Chain	Residue	Modelled	Actual	Comment	Reference
J	-9	SER	-	expression tag	UNP Q8U8L5
J	-8	SER	-	expression tag	UNP Q8U8L5
J	-7	GLY	-	expression tag	UNP Q8U8L5
J	-6	LEU	-	expression tag	UNP Q8U8L5
J	-5	VAL	-	expression tag	UNP Q8U8L5
J	-4	PRO	-	expression tag	UNP Q8U8L5
J	-3	ARG	-	expression tag	UNP Q8U8L5
J	-2	GLY	-	expression tag	UNP Q8U8L5
J	-1	SER	-	expression tag	UNP Q8U8L5
J	0	HIS	-	expression tag	UNP Q8U8L5
J	72	GLU	SER	engineered mutation	UNP Q8U8L5
K	-19	MET	-	expression tag	UNP Q8U8L5
K	-18	GLY	-	expression tag	UNP Q8U8L5
K	-17	SER	-	expression tag	UNP Q8U8L5
K	-16	SER	-	expression tag	UNP Q8U8L5
K	-15	HIS	-	expression tag	UNP Q8U8L5
K	-14	HIS	-	expression tag	UNP Q8U8L5
K	-13	HIS	-	expression tag	UNP Q8U8L5
K	-12	HIS	-	expression tag	UNP Q8U8L5
K	-11	HIS	-	expression tag	UNP Q8U8L5
K	-10	HIS	-	expression tag	UNP Q8U8L5
K	-9	SER	-	expression tag	UNP Q8U8L5
K	-8	SER	-	expression tag	UNP Q8U8L5
K	-7	GLY	-	expression tag	UNP Q8U8L5
K	-6	LEU	-	expression tag	UNP Q8U8L5
K	-5	VAL	-	expression tag	UNP Q8U8L5
K	-4	PRO	-	expression tag	UNP Q8U8L5
K	-3	ARG	-	expression tag	UNP Q8U8L5
K	-2	GLY	-	expression tag	UNP Q8U8L5
K	-1	SER	-	expression tag	UNP Q8U8L5
K	0	HIS	-	expression tag	UNP Q8U8L5
K	72	GLU	SER	engineered mutation	UNP Q8U8L5
L	-19	MET	-	expression tag	UNP Q8U8L5
L	-18	GLY	-	expression tag	UNP Q8U8L5
L	-17	SER	-	expression tag	UNP Q8U8L5
L	-16	SER	-	expression tag	UNP Q8U8L5
L	-15	HIS	-	expression tag	UNP Q8U8L5
L	-14	HIS	-	expression tag	UNP Q8U8L5
L	-13	HIS	-	expression tag	UNP Q8U8L5
L	-12	HIS	-	expression tag	UNP Q8U8L5
L	-11	HIS	-	expression tag	UNP Q8U8L5
L	-10	HIS	-	expression tag	UNP Q8U8L5

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Chain	Residue	Modelled	Actual	Comment	Reference
L	-9	SER	-	expression tag	UNP Q8U8L5
L	-8	SER	-	expression tag	UNP Q8U8L5
L	-7	GLY	-	expression tag	UNP Q8U8L5
L	-6	LEU	-	expression tag	UNP Q8U8L5
L	-5	VAL	-	expression tag	UNP Q8U8L5
L	-4	PRO	-	expression tag	UNP Q8U8L5
L	-3	ARG	-	expression tag	UNP Q8U8L5
L	-2	GLY	-	expression tag	UNP Q8U8L5
L	-1	SER	-	expression tag	UNP Q8U8L5
L	0	HIS	-	expression tag	UNP Q8U8L5
L	72	GLU	SER	engineered mutation	UNP Q8U8L5
M	-19	MET	-	expression tag	UNP Q8U8L5
M	-18	GLY	-	expression tag	UNP Q8U8L5
M	-17	SER	-	expression tag	UNP Q8U8L5
M	-16	SER	-	expression tag	UNP Q8U8L5
M	-15	HIS	-	expression tag	UNP Q8U8L5
M	-14	HIS	-	expression tag	UNP Q8U8L5
M	-13	HIS	-	expression tag	UNP Q8U8L5
M	-12	HIS	-	expression tag	UNP Q8U8L5
M	-11	HIS	-	expression tag	UNP Q8U8L5
M	-10	HIS	-	expression tag	UNP Q8U8L5
M	-9	SER	-	expression tag	UNP Q8U8L5
M	-8	SER	-	expression tag	UNP Q8U8L5
M	-7	GLY	-	expression tag	UNP Q8U8L5
M	-6	LEU	-	expression tag	UNP Q8U8L5
M	-5	VAL	-	expression tag	UNP Q8U8L5
M	-4	PRO	-	expression tag	UNP Q8U8L5
M	-3	ARG	-	expression tag	UNP Q8U8L5
M	-2	GLY	-	expression tag	UNP Q8U8L5
M	-1	SER	-	expression tag	UNP Q8U8L5
M	0	HIS	-	expression tag	UNP Q8U8L5
M	72	GLU	SER	engineered mutation	UNP Q8U8L5
N	-19	MET	-	expression tag	UNP Q8U8L5
N	-18	GLY	-	expression tag	UNP Q8U8L5
N	-17	SER	-	expression tag	UNP Q8U8L5
N	-16	SER	-	expression tag	UNP Q8U8L5
N	-15	HIS	-	expression tag	UNP Q8U8L5
N	-14	HIS	-	expression tag	UNP Q8U8L5
N	-13	HIS	-	expression tag	UNP Q8U8L5
N	-12	HIS	-	expression tag	UNP Q8U8L5
N	-11	HIS	-	expression tag	UNP Q8U8L5
N	-10	HIS	-	expression tag	UNP Q8U8L5

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Chain	Residue	Modelled	Actual	Comment	Reference
N	-9	SER	-	expression tag	UNP Q8U8L5
N	-8	SER	-	expression tag	UNP Q8U8L5
N	-7	GLY	-	expression tag	UNP Q8U8L5
N	-6	LEU	-	expression tag	UNP Q8U8L5
N	-5	VAL	-	expression tag	UNP Q8U8L5
N	-4	PRO	-	expression tag	UNP Q8U8L5
N	-3	ARG	-	expression tag	UNP Q8U8L5
N	-2	GLY	-	expression tag	UNP Q8U8L5
N	-1	SER	-	expression tag	UNP Q8U8L5
N	0	HIS	-	expression tag	UNP Q8U8L5
N	72	GLU	SER	engineered mutation	UNP Q8U8L5
O	-19	MET	-	expression tag	UNP Q8U8L5
O	-18	GLY	-	expression tag	UNP Q8U8L5
O	-17	SER	-	expression tag	UNP Q8U8L5
O	-16	SER	-	expression tag	UNP Q8U8L5
O	-15	HIS	-	expression tag	UNP Q8U8L5
O	-14	HIS	-	expression tag	UNP Q8U8L5
O	-13	HIS	-	expression tag	UNP Q8U8L5
O	-12	HIS	-	expression tag	UNP Q8U8L5
O	-11	HIS	-	expression tag	UNP Q8U8L5
O	-10	HIS	-	expression tag	UNP Q8U8L5
O	-9	SER	-	expression tag	UNP Q8U8L5
O	-8	SER	-	expression tag	UNP Q8U8L5
O	-7	GLY	-	expression tag	UNP Q8U8L5
O	-6	LEU	-	expression tag	UNP Q8U8L5
O	-5	VAL	-	expression tag	UNP Q8U8L5
O	-4	PRO	-	expression tag	UNP Q8U8L5
O	-3	ARG	-	expression tag	UNP Q8U8L5
O	-2	GLY	-	expression tag	UNP Q8U8L5
O	-1	SER	-	expression tag	UNP Q8U8L5
O	0	HIS	-	expression tag	UNP Q8U8L5
O	72	GLU	SER	engineered mutation	UNP Q8U8L5
P	-19	MET	-	expression tag	UNP Q8U8L5
P	-18	GLY	-	expression tag	UNP Q8U8L5
P	-17	SER	-	expression tag	UNP Q8U8L5
P	-16	SER	-	expression tag	UNP Q8U8L5
P	-15	HIS	-	expression tag	UNP Q8U8L5
P	-14	HIS	-	expression tag	UNP Q8U8L5
P	-13	HIS	-	expression tag	UNP Q8U8L5
P	-12	HIS	-	expression tag	UNP Q8U8L5
P	-11	HIS	-	expression tag	UNP Q8U8L5
P	-10	HIS	-	expression tag	UNP Q8U8L5

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Chain	Residue	Modelled	Actual	Comment	Reference
P	-9	SER	-	expression tag	UNP Q8U8L5
P	-8	SER	-	expression tag	UNP Q8U8L5
P	-7	GLY	-	expression tag	UNP Q8U8L5
P	-6	LEU	-	expression tag	UNP Q8U8L5
P	-5	VAL	-	expression tag	UNP Q8U8L5
P	-4	PRO	-	expression tag	UNP Q8U8L5
P	-3	ARG	-	expression tag	UNP Q8U8L5
P	-2	GLY	-	expression tag	UNP Q8U8L5
P	-1	SER	-	expression tag	UNP Q8U8L5
P	0	HIS	-	expression tag	UNP Q8U8L5
P	72	GLU	SER	engineered mutation	UNP Q8U8L5
Q	-19	MET	-	expression tag	UNP Q8U8L5
Q	-18	GLY	-	expression tag	UNP Q8U8L5
Q	-17	SER	-	expression tag	UNP Q8U8L5
Q	-16	SER	-	expression tag	UNP Q8U8L5
Q	-15	HIS	-	expression tag	UNP Q8U8L5
Q	-14	HIS	-	expression tag	UNP Q8U8L5
Q	-13	HIS	-	expression tag	UNP Q8U8L5
Q	-12	HIS	-	expression tag	UNP Q8U8L5
Q	-11	HIS	-	expression tag	UNP Q8U8L5
Q	-10	HIS	-	expression tag	UNP Q8U8L5
Q	-9	SER	-	expression tag	UNP Q8U8L5
Q	-8	SER	-	expression tag	UNP Q8U8L5
Q	-7	GLY	-	expression tag	UNP Q8U8L5
Q	-6	LEU	-	expression tag	UNP Q8U8L5
Q	-5	VAL	-	expression tag	UNP Q8U8L5
Q	-4	PRO	-	expression tag	UNP Q8U8L5
Q	-3	ARG	-	expression tag	UNP Q8U8L5
Q	-2	GLY	-	expression tag	UNP Q8U8L5
Q	-1	SER	-	expression tag	UNP Q8U8L5
Q	0	HIS	-	expression tag	UNP Q8U8L5
Q	72	GLU	SER	engineered mutation	UNP Q8U8L5
R	-19	MET	-	expression tag	UNP Q8U8L5
R	-18	GLY	-	expression tag	UNP Q8U8L5
R	-17	SER	-	expression tag	UNP Q8U8L5
R	-16	SER	-	expression tag	UNP Q8U8L5
R	-15	HIS	-	expression tag	UNP Q8U8L5
R	-14	HIS	-	expression tag	UNP Q8U8L5
R	-13	HIS	-	expression tag	UNP Q8U8L5
R	-12	HIS	-	expression tag	UNP Q8U8L5
R	-11	HIS	-	expression tag	UNP Q8U8L5
R	-10	HIS	-	expression tag	UNP Q8U8L5

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Chain	Residue	Modelled	Actual	Comment	Reference
R	-9	SER	-	expression tag	UNP Q8U8L5
R	-8	SER	-	expression tag	UNP Q8U8L5
R	-7	GLY	-	expression tag	UNP Q8U8L5
R	-6	LEU	-	expression tag	UNP Q8U8L5
R	-5	VAL	-	expression tag	UNP Q8U8L5
R	-4	PRO	-	expression tag	UNP Q8U8L5
R	-3	ARG	-	expression tag	UNP Q8U8L5
R	-2	GLY	-	expression tag	UNP Q8U8L5
R	-1	SER	-	expression tag	UNP Q8U8L5
R	0	HIS	-	expression tag	UNP Q8U8L5
R	72	GLU	SER	engineered mutation	UNP Q8U8L5
V	-19	MET	-	expression tag	UNP Q8U8L5
V	-18	GLY	-	expression tag	UNP Q8U8L5
V	-17	SER	-	expression tag	UNP Q8U8L5
V	-16	SER	-	expression tag	UNP Q8U8L5
V	-15	HIS	-	expression tag	UNP Q8U8L5
V	-14	HIS	-	expression tag	UNP Q8U8L5
V	-13	HIS	-	expression tag	UNP Q8U8L5
V	-12	HIS	-	expression tag	UNP Q8U8L5
V	-11	HIS	-	expression tag	UNP Q8U8L5
V	-10	HIS	-	expression tag	UNP Q8U8L5
V	-9	SER	-	expression tag	UNP Q8U8L5
V	-8	SER	-	expression tag	UNP Q8U8L5
V	-7	GLY	-	expression tag	UNP Q8U8L5
V	-6	LEU	-	expression tag	UNP Q8U8L5
V	-5	VAL	-	expression tag	UNP Q8U8L5
V	-4	PRO	-	expression tag	UNP Q8U8L5
V	-3	ARG	-	expression tag	UNP Q8U8L5
V	-2	GLY	-	expression tag	UNP Q8U8L5
V	-1	SER	-	expression tag	UNP Q8U8L5
V	0	HIS	-	expression tag	UNP Q8U8L5
V	72	GLU	SER	engineered mutation	UNP Q8U8L5
T	-19	MET	-	expression tag	UNP Q8U8L5
T	-18	GLY	-	expression tag	UNP Q8U8L5
T	-17	SER	-	expression tag	UNP Q8U8L5
T	-16	SER	-	expression tag	UNP Q8U8L5
T	-15	HIS	-	expression tag	UNP Q8U8L5
T	-14	HIS	-	expression tag	UNP Q8U8L5
T	-13	HIS	-	expression tag	UNP Q8U8L5
T	-12	HIS	-	expression tag	UNP Q8U8L5
T	-11	HIS	-	expression tag	UNP Q8U8L5
T	-10	HIS	-	expression tag	UNP Q8U8L5

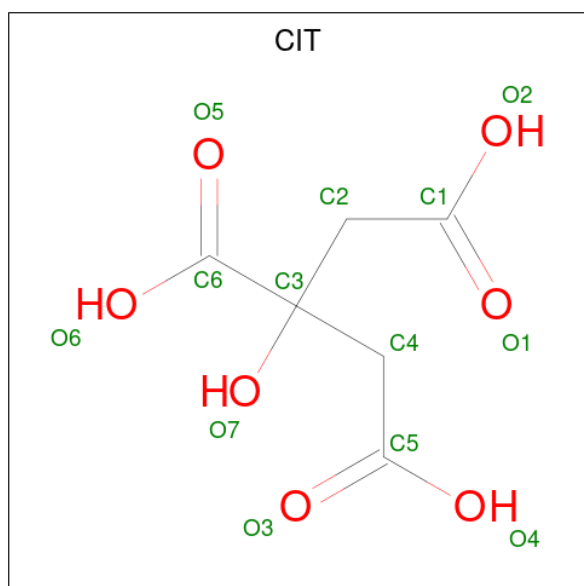
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Chain	Residue	Modelled	Actual	Comment	Reference
T	-9	SER	-	expression tag	UNP Q8U8L5
T	-8	SER	-	expression tag	UNP Q8U8L5
T	-7	GLY	-	expression tag	UNP Q8U8L5
T	-6	LEU	-	expression tag	UNP Q8U8L5
T	-5	VAL	-	expression tag	UNP Q8U8L5
T	-4	PRO	-	expression tag	UNP Q8U8L5
T	-3	ARG	-	expression tag	UNP Q8U8L5
T	-2	GLY	-	expression tag	UNP Q8U8L5
T	-1	SER	-	expression tag	UNP Q8U8L5
T	0	HIS	-	expression tag	UNP Q8U8L5
T	72	GLU	SER	engineered mutation	UNP Q8U8L5

- Molecule 2 is CITRIC ACID (three-letter code: CIT) (formula:  $C_6H_8O_7$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	H	1	Total C O 13 6 7	0	0
2	A	1	Total C O 13 6 7	0	0
2	B	1	Total C O 13 6 7	0	0
2	C	1	Total C O 13 6 7	0	0
2	D	1	Total C O 13 6 7	0	0
2	E	1	Total C O 13 6 7	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	F	1	Total	C	O	0	0
			13	6	7		
2	G	1	Total	C	O	0	0
			13	6	7		
2	I	1	Total	C	O	0	0
			13	6	7		
2	J	1	Total	C	O	0	0
			13	6	7		
2	K	1	Total	C	O	0	0
			13	6	7		
2	L	1	Total	C	O	0	0
			13	6	7		
2	M	1	Total	C	O	0	0
			13	6	7		
2	N	1	Total	C	O	0	0
			13	6	7		
2	O	1	Total	C	O	0	0
			13	6	7		
2	P	1	Total	C	O	0	0
			13	6	7		
2	Q	1	Total	C	O	0	0
			13	6	7		
2	R	1	Total	C	O	0	0
			13	6	7		
2	V	1	Total	C	O	0	0
			13	6	7		
2	T	1	Total	C	O	0	0
			13	6	7		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	H	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	F	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	I	1	Total	C	O	0	0
			6	3	3		
3	J	1	Total	C	O	0	0
			6	3	3		
3	J	1	Total	C	O	0	0
			6	3	3		
3	K	1	Total	C	O	0	0
			6	3	3		
3	K	1	Total	C	O	0	0
			6	3	3		
3	L	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	M	1	Total	C	O	0	0
			6	3	3		
3	N	1	Total	C	O	0	0
			6	3	3		
3	O	1	Total	C	O	0	0
			6	3	3		
3	P	1	Total	C	O	0	0
			6	3	3		
3	Q	1	Total	C	O	0	0
			6	3	3		
3	R	1	Total	C	O	0	0
			6	3	3		
3	V	1	Total	C	O	0	0
			6	3	3		
3	T	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	394	Total	O	0	0
			394	394		
4	A	364	Total	O	0	0
			364	364		
4	B	346	Total	O	0	0
			346	346		
4	C	434	Total	O	0	0
			434	434		
4	D	393	Total	O	0	0
			393	393		
4	E	385	Total	O	0	0
			385	385		
4	F	378	Total	O	0	0
			378	378		
4	G	392	Total	O	0	0
			392	392		
4	I	412	Total	O	0	0
			412	412		
4	J	412	Total	O	0	0
			412	412		
4	K	404	Total	O	0	0
			404	404		

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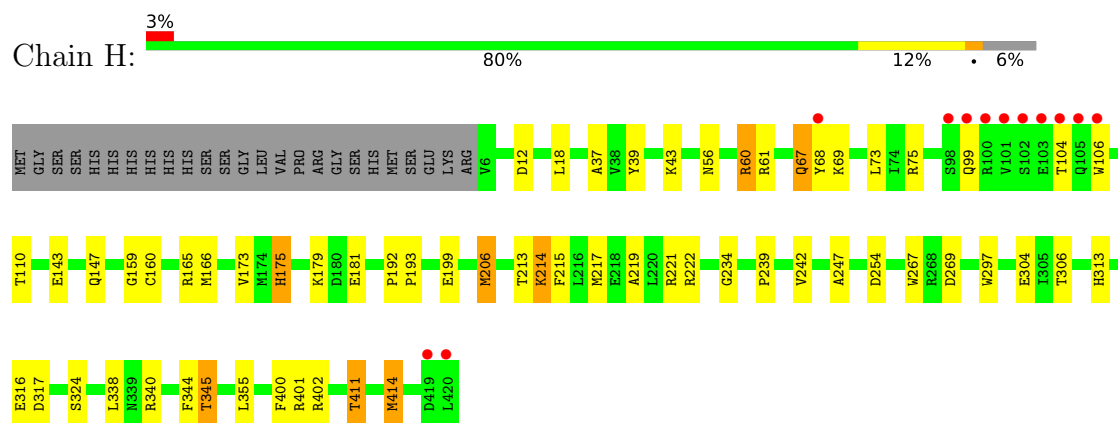
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	372	Total 372	O 372	0	0
4	M	361	Total 361	O 361	0	0
4	N	384	Total 384	O 384	0	0
4	O	348	Total 348	O 348	0	0
4	P	400	Total 400	O 400	0	0
4	Q	446	Total 446	O 446	0	0
4	R	470	Total 470	O 470	0	0
4	V	361	Total 361	O 361	0	0
4	T	404	Total 404	O 404	0	0

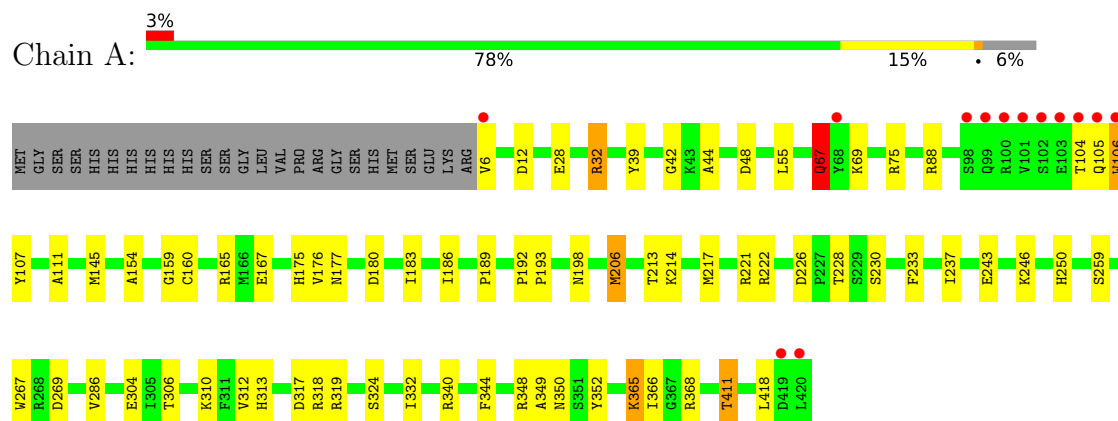
### 3 Residue-property plots [i](#)

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

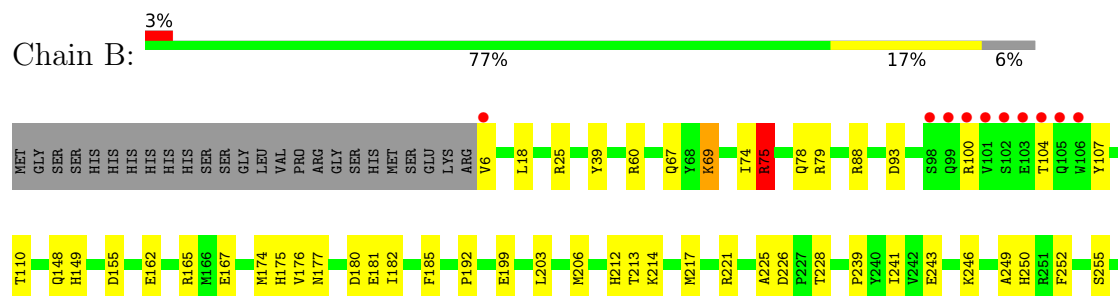
- Molecule 1: Glucose-1-phosphate adenylyltransferase



- Molecule 1: Glucose-1-phosphate adenylyltransferase

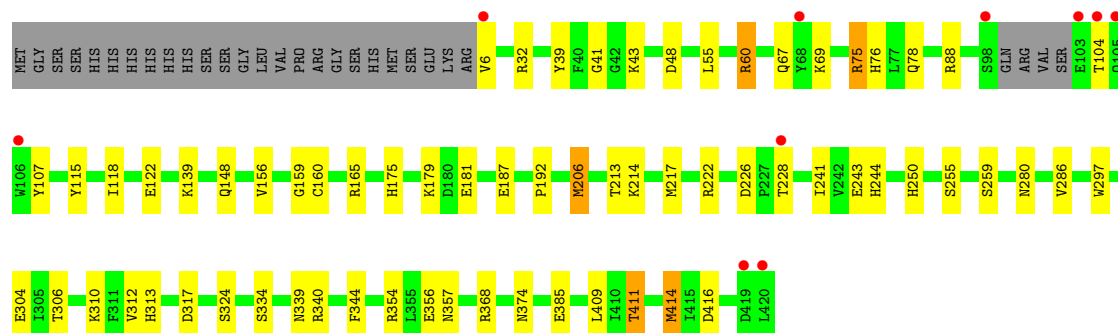
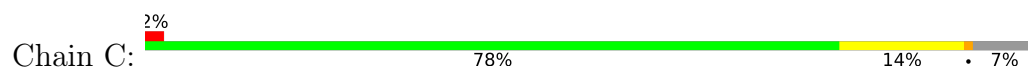


- Molecule 1: Glucose-1-phosphate adenylyltransferase

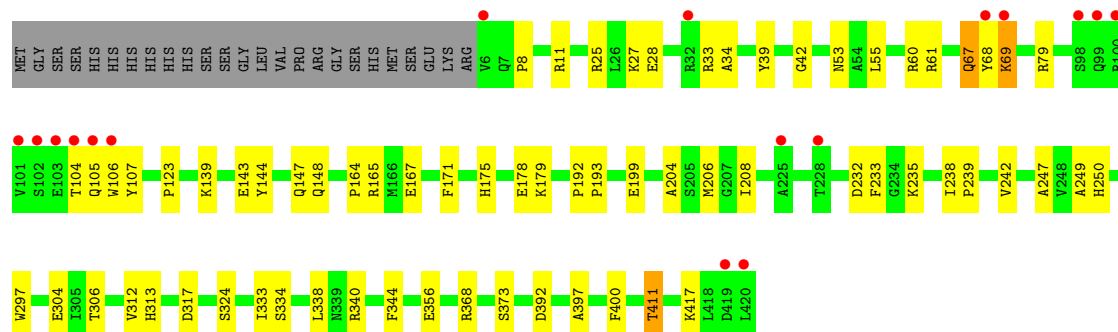
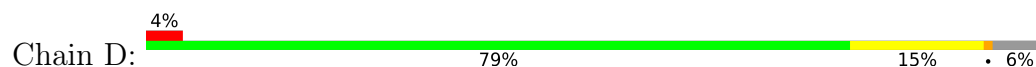




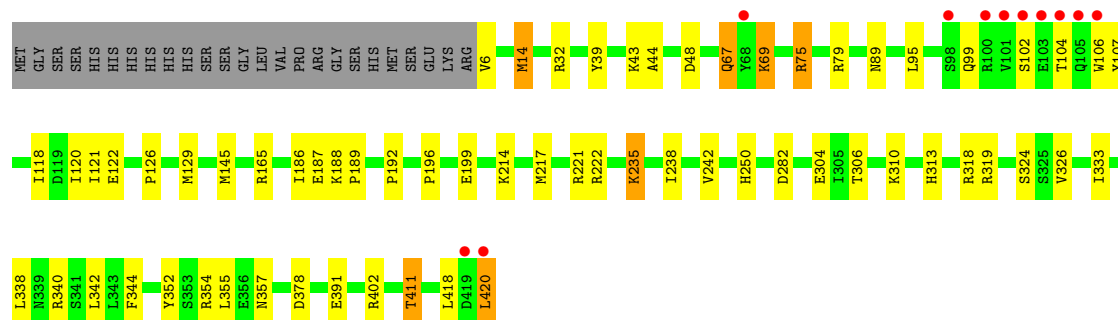
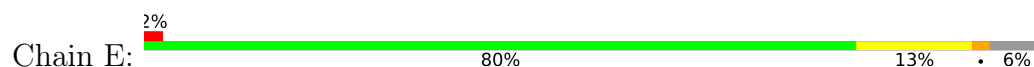
- Molecule 1: Glucose-1-phosphate adenylyltransferase



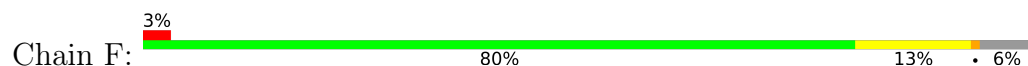
- Molecule 1: Glucose-1-phosphate adenylyltransferase

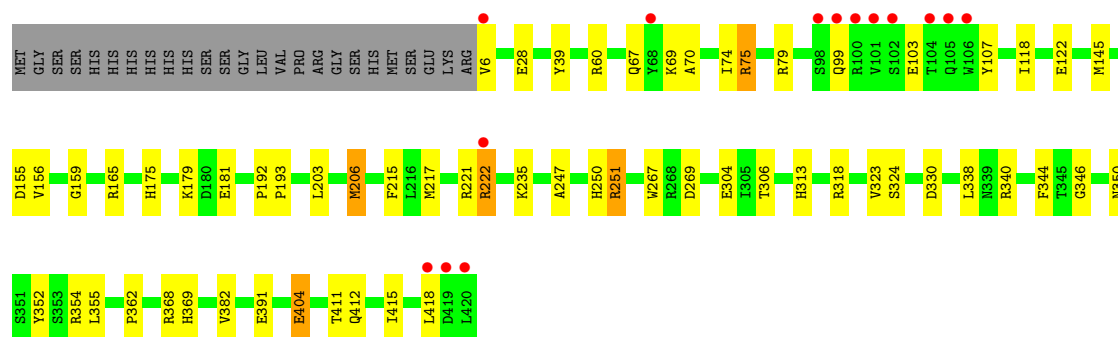


- Molecule 1: Glucose-1-phosphate adenylyltransferase

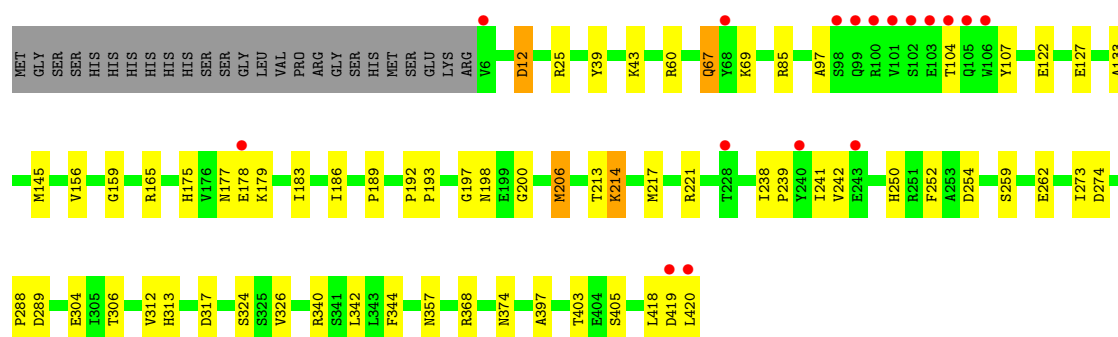
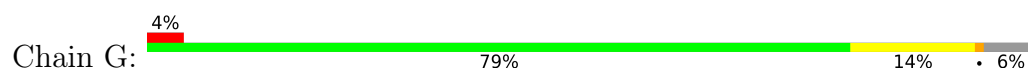


- Molecule 1: Glucose-1-phosphate adenylyltransferase

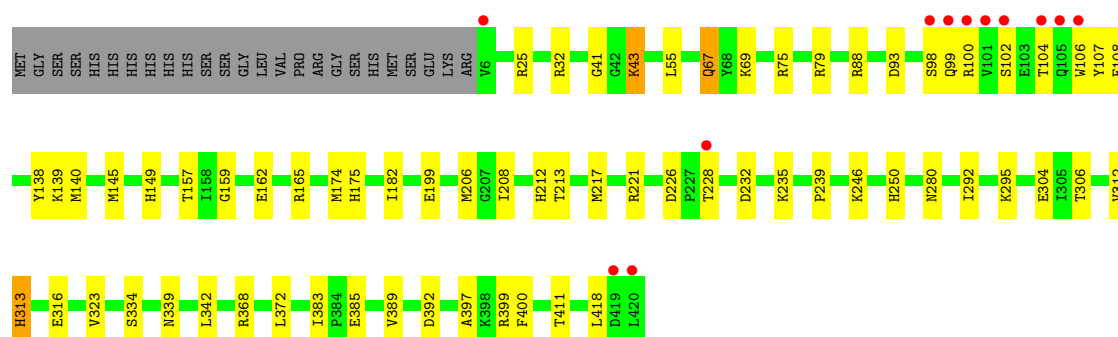
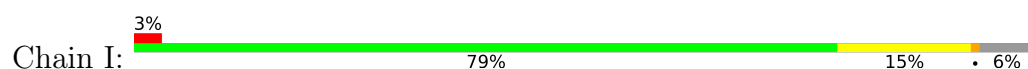




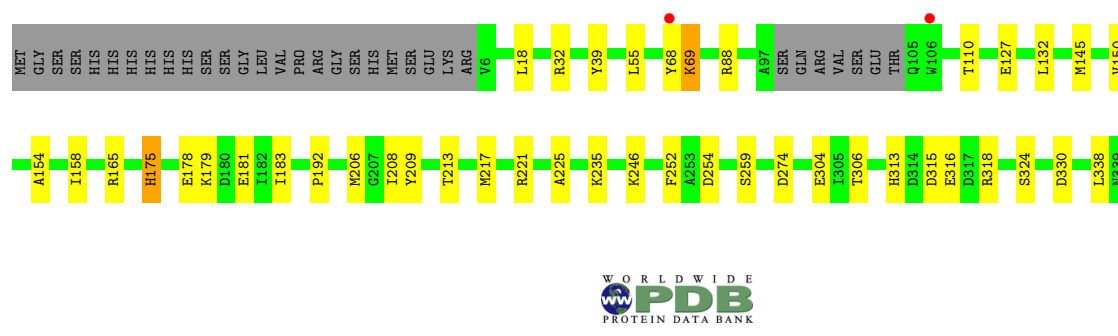
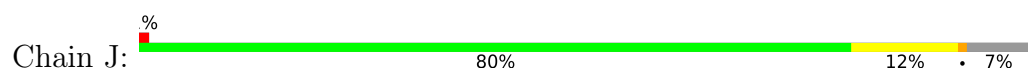
• Molecule 1: Glucose-1-phosphate adenylyltransferase



• Molecule 1: Glucose-1-phosphate adenylyltransferase

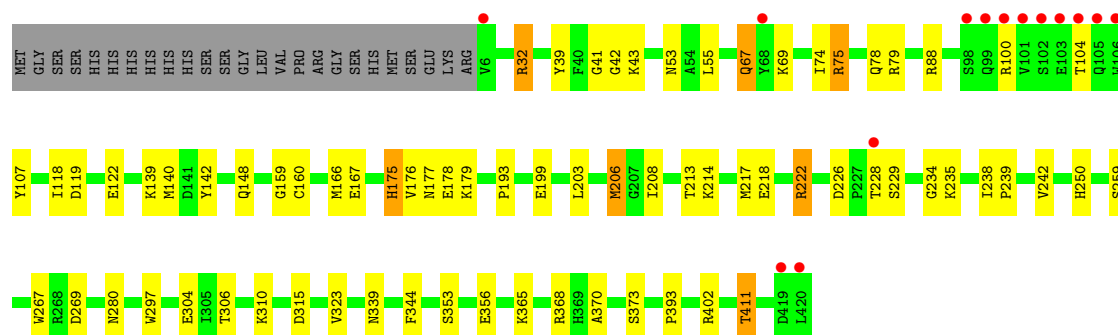
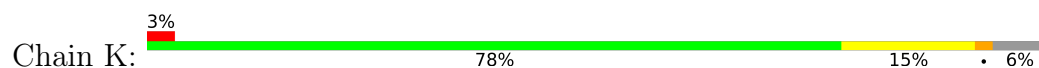


• Molecule 1: Glucose-1-phosphate adenylyltransferase

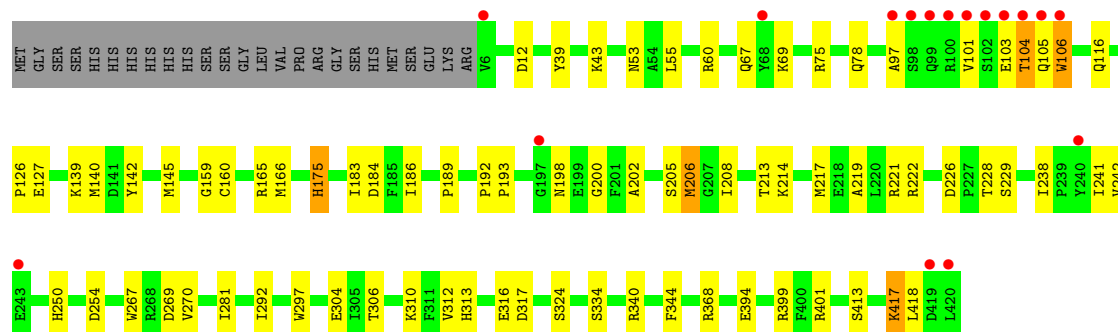
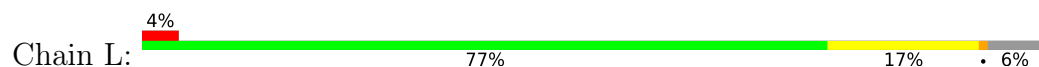




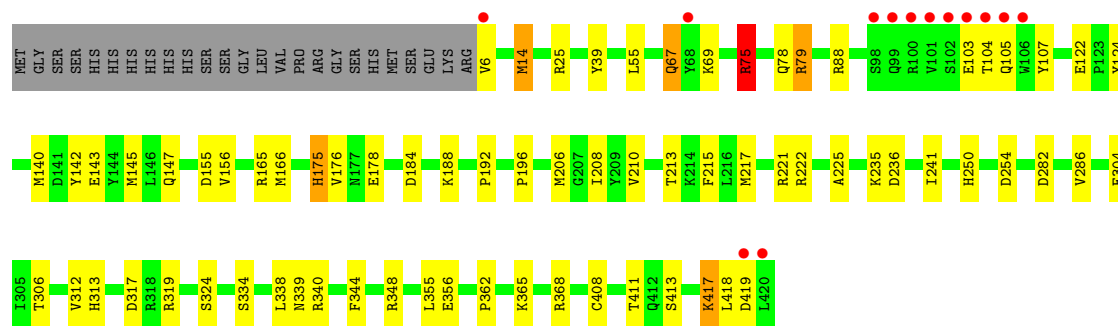
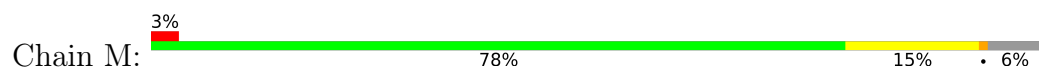
- Molecule 1: Glucose-1-phosphate adenylyltransferase



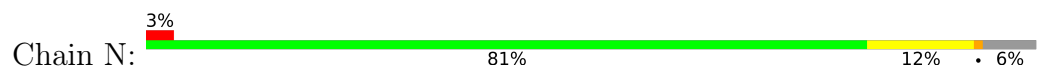
- Molecule 1: Glucose-1-phosphate adenylyltransferase

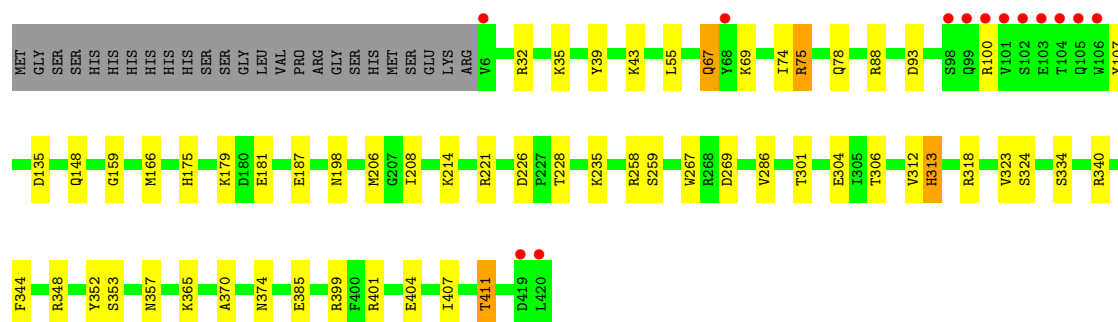


- Molecule 1: Glucose-1-phosphate adenylyltransferase

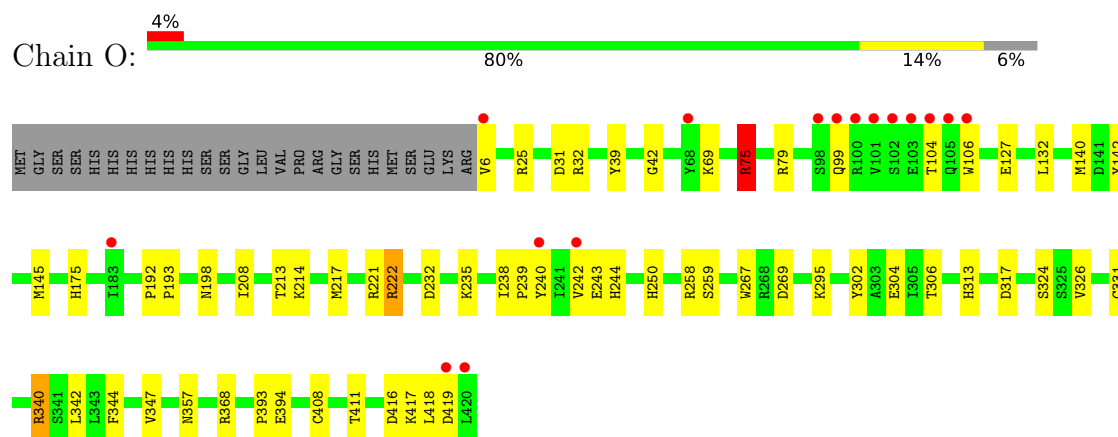


- Molecule 1: Glucose-1-phosphate adenylyltransferase

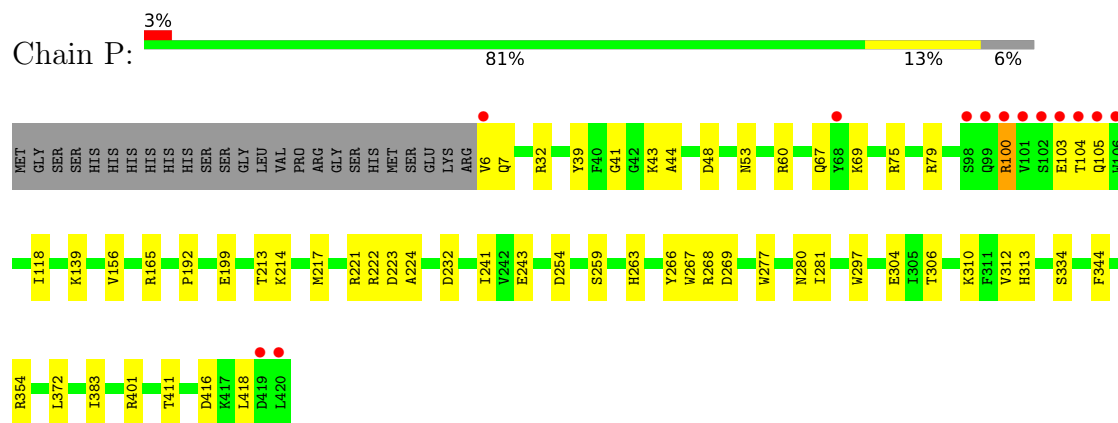




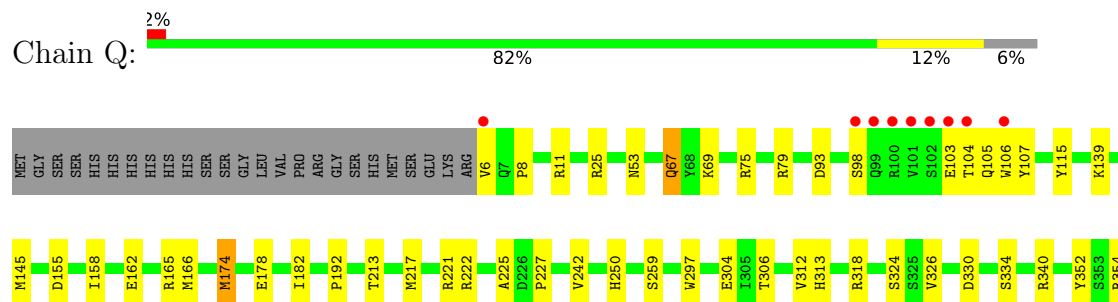
- Molecule 1: Glucose-1-phosphate adenylyltransferase



- Molecule 1: Glucose-1-phosphate adenylyltransferase



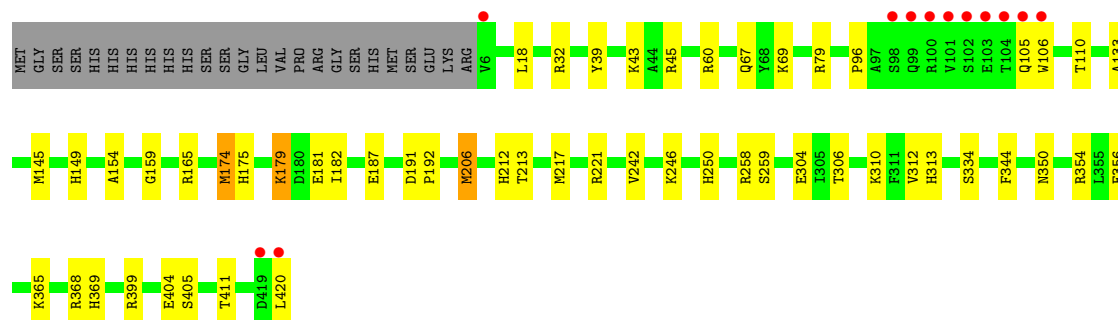
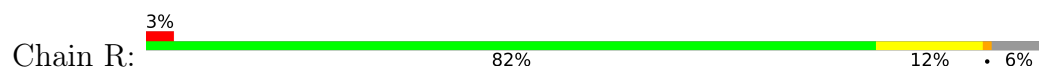
- Molecule 1: Glucose-1-phosphate adenylyltransferase



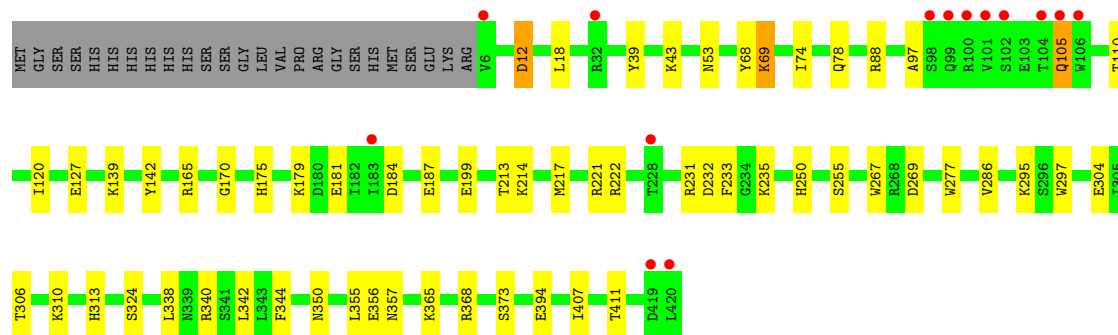
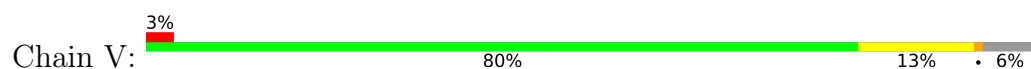




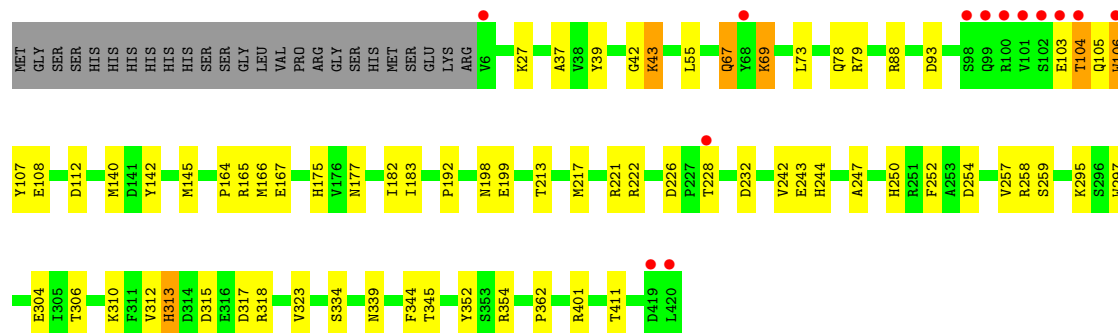
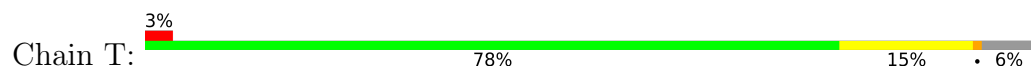
• Molecule 1: Glucose-1-phosphate adenylyltransferase



• Molecule 1: Glucose-1-phosphate adenylyltransferase



• Molecule 1: Glucose-1-phosphate adenylyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	93.61Å 141.19Å 229.43Å 107.93° 101.72° 89.97°	Depositor
Resolution (Å)	37.39 – 1.80 37.39 – 1.80	Depositor EDS
% Data completeness (in resolution range)	76.7 (37.39-1.80) 70.5 (37.39-1.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.92 (at 1.79Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, $R_{free}$	0.193 , 0.228 0.193 , 0.228	Depositor DCC
$R_{free}$ test set	38499 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	17.6	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 47.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.59$ , $\langle L^2 \rangle = 0.44$	Xtriage
Estimated twinning fraction	0.477 for h,-k,-h-l 0.477 for -h,k,-k-l 0.477 for -h,-k,h+k+l	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	74015	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	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 72.17 % 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 2.3109e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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

## 5 Model quality

### 5.1 Standard geometry

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

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	0.52	0/3367	0.67	2/4575 (0.0%)
1	B	0.50	0/3367	0.65	1/4575 (0.0%)
1	C	0.53	0/3345	0.69	3/4543 (0.1%)
1	D	0.53	0/3361	0.66	1/4568 (0.0%)
1	E	0.51	0/3367	0.63	0/4575
1	F	0.55	3/3367 (0.1%)	0.84	5/4575 (0.1%)
1	G	0.50	0/3373	0.67	2/4583 (0.0%)
1	H	0.53	0/3371	0.68	4/4580 (0.1%)
1	I	0.52	0/3373	0.67	2/4583 (0.0%)
1	J	0.54	0/3327	0.68	0/4518
1	K	0.55	2/3367 (0.1%)	0.73	5/4575 (0.1%)
1	L	0.54	0/3375	0.67	1/4585 (0.0%)
1	M	0.57	2/3367 (0.1%)	0.70	3/4575 (0.1%)
1	N	0.49	0/3367	0.63	0/4575
1	O	0.56	2/3381 (0.1%)	0.66	1/4593 (0.0%)
1	P	0.56	2/3391 (0.1%)	0.69	3/4607 (0.1%)
1	Q	0.57	1/3380 (0.0%)	0.70	3/4593 (0.1%)
1	R	0.56	0/3375	0.68	2/4585 (0.0%)
1	T	0.55	0/3375	0.66	2/4585 (0.0%)
1	V	0.52	0/3367	0.65	0/4575
All	All	0.54	12/67363 (0.0%)	0.68	40/91523 (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
1	M	0	1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	P	222	ARG	CZ-NH2	10.30	1.46	1.33
1	M	79	ARG	CZ-NH1	9.73	1.45	1.33
1	P	222	ARG	CD-NE	-8.80	1.31	1.46
1	Q	354	ARG	CD-NE	8.69	1.61	1.46
1	K	222	ARG	CZ-NH2	8.58	1.44	1.33
1	O	75	ARG	CZ-NH2	8.45	1.44	1.33
1	F	222	ARG	CD-NE	-8.39	1.32	1.46
1	M	75	ARG	CD-NE	7.65	1.59	1.46
1	F	75	ARG	CD-NE	6.35	1.57	1.46
1	F	222	ARG	CZ-NH2	5.78	1.40	1.33
1	K	222	ARG	CD-NE	-5.55	1.37	1.46
1	O	408	CYS	CB-SG	-5.30	1.73	1.81

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	222	ARG	NE-CZ-NH1	-27.94	106.33	120.30
1	F	222	ARG	NE-CZ-NH2	20.36	130.48	120.30
1	K	222	ARG	NE-CZ-NH1	18.11	129.35	120.30
1	P	222	ARG	NE-CZ-NH2	-12.28	114.16	120.30
1	A	32	ARG	NE-CZ-NH2	9.94	125.27	120.30
1	Q	354	ARG	NE-CZ-NH2	-9.64	115.48	120.30
1	K	222	ARG	NE-CZ-NH2	-9.24	115.68	120.30
1	M	75	ARG	NE-CZ-NH1	-8.47	116.07	120.30
1	M	79	ARG	NE-CZ-NH2	8.34	124.47	120.30
1	A	206	MET	CG-SD-CE	8.10	113.16	100.20
1	M	14	MET	CG-SD-CE	7.78	112.65	100.20
1	H	206	MET	CG-SD-CE	7.64	112.42	100.20
1	F	206	MET	CG-SD-CE	7.64	112.42	100.20
1	T	222	ARG	NE-CZ-NH1	7.27	123.93	120.30
1	L	206	MET	CG-SD-CE	7.23	111.77	100.20
1	G	206	MET	CG-SD-CE	6.87	111.19	100.20
1	C	206	MET	CG-SD-CE	6.75	110.99	100.20
1	K	206	MET	CG-SD-CE	6.60	110.76	100.20
1	Q	174	MET	CG-SD-CE	6.33	110.32	100.20
1	F	75	ARG	NE-CZ-NH2	6.27	123.43	120.30
1	D	392	ASP	CB-CG-OD1	5.76	123.48	118.30
1	O	222	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	C	60	ARG	NE-CZ-NH2	-5.68	117.46	120.30
1	P	222	ARG	NE-CZ-NH1	5.67	123.14	120.30
1	I	392	ASP	CB-CG-OD1	5.66	123.40	118.30
1	I	392	ASP	CB-CG-OD2	-5.64	113.22	118.30
1	R	206	MET	CG-SD-CE	5.63	109.21	100.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Q	354	ARG	NE-CZ-NH1	5.58	123.09	120.30
1	H	61	ARG	NE-CZ-NH1	-5.51	117.54	120.30
1	R	174	MET	CG-SD-CE	5.50	109.00	100.20
1	H	414	MET	CB-CG-SD	-5.49	95.92	112.40
1	H	414	MET	CA-CB-CG	5.41	122.50	113.30
1	K	365	LYS	CG-CD-CE	-5.29	96.03	111.90
1	B	75	ARG	NE-CZ-NH1	-5.24	117.68	120.30
1	K	365	LYS	CB-CG-CD	5.21	125.14	111.60
1	T	222	ARG	CG-CD-NE	-5.19	100.89	111.80
1	F	222	ARG	CG-CD-NE	5.18	122.69	111.80
1	C	60	ARG	NE-CZ-NH1	5.13	122.87	120.30
1	P	222	ARG	CD-NE-CZ	5.09	130.72	123.60
1	G	85	ARG	NE-CZ-NH1	5.02	122.81	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	M	75	ARG	Sidechain

## 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3287	0	3192	60	0
1	B	3287	0	3192	70	0
1	C	3266	0	3170	62	0
1	D	3281	0	3181	60	0
1	E	3287	0	3192	52	0
1	F	3287	0	3192	44	1
1	G	3293	0	3196	53	0
1	H	3291	0	3198	55	0
1	I	3293	0	3196	61	0
1	J	3248	0	3156	53	0
1	K	3287	0	3192	64	0
1	L	3295	0	3200	79	0
1	M	3287	0	3192	61	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	N	3287	0	3192	55	0
1	O	3301	0	3204	51	0
1	P	3310	0	3212	45	0
1	Q	3299	0	3200	52	1
1	R	3295	0	3200	49	0
1	T	3295	0	3200	69	0
1	V	3287	0	3192	65	0
2	A	13	0	5	3	0
2	B	13	0	5	2	0
2	C	13	0	5	2	0
2	D	13	0	5	2	0
2	E	13	0	5	2	0
2	F	13	0	5	3	0
2	G	13	0	5	2	0
2	H	13	0	5	1	0
2	I	13	0	5	1	0
2	J	13	0	5	4	0
2	K	13	0	5	1	0
2	L	13	0	5	4	0
2	M	13	0	5	2	0
2	N	13	0	5	1	0
2	O	13	0	5	2	0
2	P	13	0	5	1	0
2	Q	13	0	5	2	0
2	R	13	0	5	3	0
2	T	13	0	5	2	0
2	V	13	0	5	2	0
3	A	6	0	8	0	0
3	B	6	0	8	0	0
3	C	6	0	8	0	0
3	D	6	0	8	0	0
3	E	6	0	8	0	0
3	F	6	0	8	0	0
3	G	6	0	8	1	0
3	H	6	0	8	0	0
3	I	6	0	8	0	0
3	J	12	0	16	4	0
3	K	12	0	16	5	0
3	L	6	0	8	0	0
3	M	6	0	8	0	0
3	N	6	0	8	0	0
3	O	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	P	6	0	8	0	0
3	Q	6	0	8	0	0
3	R	6	0	8	0	0
3	T	6	0	8	0	0
3	V	6	0	8	0	0
4	A	364	0	0	13	0
4	B	346	0	0	27	0
4	C	434	0	0	20	1
4	D	393	0	0	23	4
4	E	385	0	0	14	2
4	F	378	0	0	8	6
4	G	392	0	0	18	2
4	H	394	0	0	15	2
4	I	412	0	0	22	1
4	J	412	0	0	21	0
4	K	404	0	0	16	0
4	L	372	0	0	31	0
4	M	361	0	0	19	0
4	N	384	0	0	16	0
4	O	348	0	0	17	2
4	P	400	0	0	19	6
4	Q	446	0	0	14	5
4	R	470	0	0	21	4
4	T	404	0	0	28	1
4	V	361	0	0	21	0
All	All	74015	0	64125	1119	19

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 9.

All (1119) 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)
1:F:60:ARG:NH2	4:F:601:HOH:O	1.73	1.17
1:H:206:MET:SD	4:H:919:HOH:O	2.09	1.10
1:H:60:ARG:NH2	4:H:601:HOH:O	1.85	1.09
1:Q:174:MET:HE2	1:Q:182:ILE:HD13	1.35	1.08
1:L:60:ARG:NH2	4:L:601:HOH:O	1.86	1.05
1:D:206:MET:HE3	1:D:208:ILE:HD12	1.31	1.04
1:O:6:VAL:N	4:O:601:HOH:O	1.88	1.03
1:L:222:ARG:NH2	4:L:604:HOH:O	1.93	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:206:MET:SD	4:A:875:HOH:O	2.20	1.00
1:G:206:MET:SD	4:G:898:HOH:O	2.18	1.00
1:J:206:MET:SD	4:J:940:HOH:O	2.22	0.96
1:K:199:GLU:OE2	4:K:601:HOH:O	1.84	0.96
1:K:226:ASP:OD2	1:K:228:THR:HG22	1.66	0.96
1:R:133:ALA:O	4:R:601:HOH:O	1.82	0.96
1:F:122:GLU:OE1	4:F:602:HOH:O	1.82	0.95
1:A:418:LEU:O	4:A:601:HOH:O	1.83	0.95
1:J:330:ASP:OD2	4:J:601:HOH:O	1.83	0.94
1:L:75:ARG:NH1	4:L:606:HOH:O	1.98	0.94
1:G:133:ALA:O	4:G:601:HOH:O	1.85	0.94
1:V:175:HIS:ND1	4:V:601:HOH:O	1.93	0.93
1:E:79:ARG:NH1	4:E:603:HOH:O	2.01	0.93
1:H:401:ARG:H	1:H:414:MET:HE1	1.34	0.92
1:K:78:GLN:HE21	1:V:78:GLN:HE21	0.95	0.92
1:L:206:MET:SD	4:L:888:HOH:O	2.28	0.91
1:J:158:ILE:HD13	1:J:209:TYR:HD2	1.36	0.91
1:L:78:GLN:HE21	1:T:78:GLN:HE21	0.93	0.91
1:D:206:MET:SD	4:D:893:HOH:O	2.28	0.91
1:F:206:MET:SD	4:F:878:HOH:O	2.28	0.90
1:K:78:GLN:HE21	1:V:78:GLN:NE2	1.68	0.90
1:N:206:MET:SD	4:N:902:HOH:O	2.28	0.89
1:T:354:ARG:NH1	4:T:605:HOH:O	2.04	0.89
1:L:229:SER:O	4:L:602:HOH:O	1.90	0.88
1:A:167:GLU:O	4:A:602:HOH:O	1.91	0.88
1:K:234:GLY:O	4:K:602:HOH:O	1.91	0.88
1:K:78:GLN:NE2	1:V:78:GLN:HE21	1.71	0.87
1:R:242:VAL:O	4:R:602:HOH:O	1.90	0.87
1:B:162:GLU:O	4:B:602:HOH:O	1.92	0.87
1:C:317:ASP:OD1	4:C:601:HOH:O	1.92	0.87
1:E:235:LYS:NZ	4:E:604:HOH:O	2.09	0.86
1:R:96:PRO:O	4:R:603:HOH:O	1.93	0.86
1:Q:67:GLN:NE2	1:Q:107:TYR:H	1.73	0.85
1:D:60:ARG:NH2	4:D:604:HOH:O	2.08	0.85
1:C:159:GLY:C	1:C:206:MET:HE3	1.97	0.85
1:D:242:VAL:O	4:D:602:HOH:O	1.95	0.84
1:R:174:MET:SD	4:R:949:HOH:O	2.36	0.84
1:I:221:ARG:NH1	4:I:601:HOH:O	2.01	0.84
1:D:144:TYR:O	4:D:601:HOH:O	1.95	0.84
1:R:174:MET:HE2	1:R:182:ILE:HD13	1.58	0.84
1:N:206:MET:HE3	1:N:208:ILE:HD12	1.58	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:243:GLU:O	4:A:603:HOH:O	1.96	0.83
1:D:206:MET:CE	1:D:208:ILE:HD12	2.07	0.83
1:J:158:ILE:HD13	1:J:209:TYR:CD2	2.14	0.83
1:I:75:ARG:NE	4:I:606:HOH:O	2.11	0.83
1:L:254:ASP:OD2	4:L:605:HOH:O	1.96	0.83
1:V:110:THR:HG22	1:V:232:ASP:HA	1.61	0.82
1:E:89:ASN:OD1	4:E:601:HOH:O	1.97	0.82
1:T:67:GLN:NE2	4:T:609:HOH:O	2.12	0.82
1:M:6:VAL:N	4:M:606:HOH:O	2.13	0.82
1:B:60:ARG:HD2	4:B:656:HOH:O	1.80	0.82
1:V:110:THR:CG2	1:V:233:PHE:H	1.92	0.82
1:T:315:ASP:OD2	4:T:602:HOH:O	1.97	0.81
1:R:354:ARG:NH2	4:R:607:HOH:O	2.12	0.80
1:J:221:ARG:NH1	4:J:603:HOH:O	2.01	0.80
1:N:75:ARG:NH2	4:N:603:HOH:O	2.08	0.80
1:L:417:LYS:NZ	4:L:611:HOH:O	2.15	0.80
1:K:75:ARG:NH1	4:K:607:HOH:O	2.15	0.80
1:L:78:GLN:HE21	1:T:78:GLN:NE2	1.77	0.80
4:Q:688:HOH:O	1:R:310:LYS:HE2	1.82	0.80
1:I:140:MET:HE3	1:I:208:ILE:HD13	1.64	0.79
1:T:258:ARG:NH2	4:T:612:HOH:O	2.15	0.79
1:C:222:ARG:NH2	4:C:604:HOH:O	2.13	0.79
1:P:354:ARG:NH1	4:P:605:HOH:O	2.14	0.79
1:A:177:ASN:HB3	1:A:183:ILE:HD11	1.64	0.79
1:J:150:VAL:O	4:J:602:HOH:O	2.00	0.79
1:Q:67:GLN:HE22	1:Q:107:TYR:H	1.26	0.79
1:D:27:LYS:NZ	4:D:603:HOH:O	2.00	0.78
1:J:206:MET:CE	1:J:252:PHE:HB2	2.13	0.78
1:V:175:HIS:O	4:V:603:HOH:O	2.01	0.78
1:T:199:GLU:OE2	4:T:603:HOH:O	2.00	0.78
1:D:417:LYS:NZ	4:D:610:HOH:O	2.16	0.77
1:I:108:GLU:OE2	4:I:602:HOH:O	2.02	0.77
1:B:167:GLU:O	4:B:603:HOH:O	2.00	0.77
1:B:249:ALA:O	4:B:604:HOH:O	2.02	0.77
1:K:175:HIS:HA	3:K:503:GOL:H12	1.64	0.77
1:C:60:ARG:NH2	4:C:606:HOH:O	2.17	0.77
2:T:501:CIT:O4	4:T:604:HOH:O	2.03	0.77
1:O:317:ASP:OD2	4:O:602:HOH:O	2.02	0.77
1:E:391:GLU:OE1	4:E:602:HOH:O	2.01	0.77
1:L:75:ARG:NH1	1:T:93:ASP:OD2	2.18	0.77
1:V:313:HIS:NE2	2:V:501:CIT:O4	2.14	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:326:VAL:HB	1:E:342:LEU:HD23	1.67	0.77
1:L:368:ARG:HH21	2:L:501:CIT:H41	1.48	0.76
1:I:138:TYR:HE2	1:I:140:MET:HE2	1.46	0.76
1:L:78:GLN:NE2	1:T:78:GLN:HE21	1.78	0.76
1:D:313:HIS:HE1	2:D:501:CIT:O3	1.68	0.76
1:B:6:VAL:N	4:B:610:HOH:O	2.17	0.76
1:Q:400:PHE:HD1	1:Q:414:MET:HE3	1.50	0.76
1:F:159:GLY:HA3	1:F:206:MET:HE3	1.68	0.75
1:P:53:ASN:HD22	1:P:297:TRP:HE1	1.32	0.75
1:B:78:GLN:NE2	4:B:613:HOH:O	2.20	0.75
1:B:313:HIS:NE2	2:B:501:CIT:O4	2.17	0.75
1:G:122:GLU:O	4:G:602:HOH:O	2.04	0.75
1:I:165:ARG:HD2	1:I:199:GLU:O	1.86	0.75
1:N:181:GLU:OE2	4:N:601:HOH:O	2.02	0.75
1:G:403:THR:HG22	1:G:405:SER:H	1.52	0.75
2:J:501:CIT:O6	4:J:604:HOH:O	2.03	0.75
1:Q:155:ASP:OD1	4:Q:602:HOH:O	2.05	0.75
1:L:413:SER:O	1:L:417:LYS:HE2	1.87	0.74
1:L:221:ARG:O	4:L:608:HOH:O	2.06	0.74
1:H:316:GLU:OE2	4:H:604:HOH:O	2.06	0.74
1:I:397:ALA:O	4:I:603:HOH:O	2.04	0.74
1:K:167:GLU:O	4:K:603:HOH:O	2.05	0.74
1:L:394:GLU:OE2	4:L:607:HOH:O	2.04	0.73
1:M:175:HIS:O	4:M:602:HOH:O	2.05	0.73
1:J:404:GLU:CD	1:J:404:GLU:H	1.89	0.73
1:R:221:ARG:NH1	4:R:605:HOH:O	2.10	0.73
1:P:79[B]:ARG:NH2	4:P:604:HOH:O	2.14	0.73
1:H:304:GLU:O	1:H:306:THR:HG23	1.89	0.73
1:Q:330:ASP:OD2	4:Q:601:HOH:O	2.03	0.73
1:K:53:ASN:HD21	1:K:139:LYS:HA	1.52	0.73
1:V:184:ASP:O	4:V:603:HOH:O	2.06	0.73
1:Q:115:TYR:OH	4:Q:603:HOH:O	2.05	0.73
1:G:60:ARG:NH2	4:G:608:HOH:O	2.22	0.72
1:K:42[B]:GLY:O	4:K:604:HOH:O	2.06	0.72
1:T:43[B]:LYS:HG3	4:T:668:HOH:O	1.89	0.72
1:T:254:ASP:OD1	4:T:606:HOH:O	2.06	0.72
1:I:400:PHE:O	4:I:603:HOH:O	2.07	0.72
1:J:354:ARG:NH1	4:J:611:HOH:O	2.19	0.72
1:H:179:LYS:HB2	1:H:181:GLU:HG3	1.71	0.72
1:T:67:GLN:OE1	1:T:107:TYR:HB2	1.88	0.72
1:E:32:ARG:NH1	1:E:378:ASP:OD2	2.22	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:401:ARG:NH1	4:L:610:HOH:O	2.10	0.72
1:J:179:LYS:HG2	1:J:181:GLU:HG3	1.71	0.72
1:J:313:HIS:NE2	2:J:501:CIT:O3	2.15	0.72
1:P:263:HIS:HD2	4:P:904:HOH:O	1.72	0.72
1:Q:313:HIS:HD2	4:Q:950:HOH:O	1.72	0.72
1:B:155:ASP:OD1	4:B:605:HOH:O	2.06	0.71
1:R:356:GLU:OE1	4:R:604:HOH:O	2.06	0.71
1:P:100:ARG:HB2	4:P:618:HOH:O	1.90	0.71
1:B:67:GLN:OE1	1:B:107:TYR:N	2.22	0.71
1:L:313:HIS:HE1	2:L:501:CIT:O3	1.72	0.71
1:M:221:ARG:NH1	4:M:608:HOH:O	2.18	0.71
1:A:6:VAL:N	4:A:608:HOH:O	2.23	0.71
1:J:183:ILE:O	4:J:605:HOH:O	2.08	0.71
1:V:187:GLU:OE1	4:V:604:HOH:O	2.07	0.71
1:I:316:GLU:OE1	4:I:604:HOH:O	2.07	0.71
1:M:78:GLN:NE2	4:M:610:HOH:O	2.22	0.71
1:P:354:ARG:NH1	4:P:610:HOH:O	2.22	0.70
1:I:67:GLN:OE1	1:I:107:TYR:N	2.19	0.70
1:K:32:ARG:NH2	4:K:612:HOH:O	2.24	0.70
1:T:317:ASP:OD2	4:T:607:HOH:O	2.08	0.70
1:I:232:ASP:OD1	4:I:605:HOH:O	2.08	0.70
1:P:60:ARG:NH2	4:P:611:HOH:O	2.24	0.70
1:T:313:HIS:HD2	4:T:919:HOH:O	1.73	0.70
1:P:232:ASP:OD2	4:P:601:HOH:O	2.09	0.70
1:B:182:ILE:HD11	1:B:249:ALA:HB2	1.72	0.70
1:J:315:ASP:OD2	4:J:606:HOH:O	2.08	0.70
1:T:232:ASP:OD1	4:T:608:HOH:O	2.10	0.70
1:D:317:ASP:OD2	4:D:605:HOH:O	2.09	0.70
1:V:170:GLY:O	4:V:608:HOH:O	2.10	0.70
1:O:243:GLU:OE2	4:O:603:HOH:O	2.10	0.70
1:D:312:VAL:HG13	1:D:334:SER:HA	1.73	0.70
1:B:225:ALA:O	4:B:606:HOH:O	2.08	0.69
1:R:174:MET:CE	1:R:182:ILE:HG21	2.22	0.69
1:K:140:MET:HE2	1:K:142:TYR:HE1	1.56	0.69
1:K:226:ASP:CG	1:K:228:THR:HG22	2.12	0.69
1:D:232:ASP:OD2	4:D:606:HOH:O	2.10	0.69
1:O:416:ASP:OD1	4:O:605:HOH:O	2.10	0.69
1:P:53:ASN:HD21	1:P:139:LYS:HA	1.56	0.69
1:D:53:ASN:HD21	1:D:139:LYS:HA	1.56	0.69
1:L:200:GLY:O	4:L:609:HOH:O	2.09	0.69
1:B:75:ARG:HE	1:B:79:ARG:HD2	1.58	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:222:ARG:NH1	4:C:604:HOH:O	2.26	0.69
1:D:397:ALA:O	4:D:607:HOH:O	2.11	0.69
1:K:53:ASN:HD22	1:K:297:TRP:HE1	1.38	0.69
1:F:156:VAL:CG1	1:F:247:ALA:HA	2.22	0.69
1:J:225:ALA:O	4:J:607:HOH:O	2.10	0.68
1:K:41:GLY:H	1:K:280:ASN:HD21	1.39	0.68
1:L:53:ASN:HD22	1:L:297:TRP:HE1	1.41	0.68
1:O:75:ARG:HE	1:O:79:ARG:HD2	1.58	0.68
1:V:295:LYS:NZ	4:V:611:HOH:O	2.22	0.68
1:K:78:GLN:OE1	4:K:606:HOH:O	2.12	0.68
1:H:254:ASP:OD2	4:H:605:HOH:O	2.11	0.68
1:O:127:GLU:OE1	4:O:604:HOH:O	2.10	0.68
1:I:25:ARG:HD3	4:I:612:HOH:O	1.93	0.68
1:M:14:MET:HE3	1:M:124:TYR:HD2	1.59	0.68
1:F:382:VAL:H	1:F:412:GLN:HE22	1.41	0.68
1:M:184:ASP:N	4:M:602:HOH:O	2.08	0.68
1:M:225:ALA:O	4:M:605:HOH:O	2.11	0.68
1:E:122:GLU:OE2	4:E:605:HOH:O	2.11	0.67
1:M:122:GLU:OE1	4:M:604:HOH:O	2.11	0.67
1:R:365:LYS:NZ	4:R:609:HOH:O	2.14	0.67
1:A:42[B]:GLY:O	4:A:604:HOH:O	2.11	0.67
1:G:127:GLU:O	1:G:214:LYS:NZ	2.27	0.67
1:G:317:ASP:HB2	4:G:755:HOH:O	1.95	0.67
1:K:178:GLU:OE1	4:K:605:HOH:O	2.12	0.67
1:B:243:GLU:OE1	4:B:607:HOH:O	2.11	0.67
1:C:286:VAL:HG22	1:D:312:VAL:HG23	1.77	0.67
1:C:115:TYR:OH	4:C:603:HOH:O	2.10	0.67
1:G:254:ASP:OD2	4:G:604:HOH:O	2.12	0.67
1:N:206:MET:CE	1:N:208:ILE:HD12	2.24	0.67
1:Q:222:ARG:NH1	4:Q:616:HOH:O	2.27	0.67
2:R:501:CIT:O6	4:R:606:HOH:O	2.12	0.67
1:C:313:HIS:HD2	4:C:947:HOH:O	1.77	0.66
1:J:235:LYS:NZ	4:J:615:HOH:O	2.25	0.66
1:T:317:ASP:HB2	4:T:822:HOH:O	1.95	0.66
1:H:234:GLY:O	4:H:606:HOH:O	2.14	0.66
1:D:53:ASN:HD22	1:D:297:TRP:HE1	1.41	0.66
1:I:138:TYR:CE2	1:I:140:MET:HE2	2.29	0.66
1:G:67:GLN:NE2	4:G:613:HOH:O	2.27	0.66
1:E:14:MET:HE2	1:E:126:PRO:HG3	1.77	0.66
1:K:43[B]:LYS:NZ	1:L:306:THR:O	2.24	0.66
1:G:403:THR:HG22	1:G:405:SER:N	2.11	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:223:ASP:OD1	4:P:603:HOH:O	2.14	0.66
1:V:110:THR:HG21	4:V:860:HOH:O	1.95	0.66
1:B:176:VAL:HG13	1:B:180:ASP:HA	1.76	0.66
1:K:41:GLY:H	1:K:280:ASN:ND2	1.93	0.66
1:P:254:ASP:OD2	4:P:602:HOH:O	2.13	0.66
1:Q:25:ARG:NH1	4:Q:615:HOH:O	2.27	0.66
1:J:127:GLU:OE1	4:J:608:HOH:O	2.13	0.66
1:N:313:HIS:NE2	2:N:501:CIT:O4	2.26	0.66
1:O:214:LYS:NZ	4:O:612:HOH:O	2.28	0.65
1:R:60:ARG:NH2	4:R:617:HOH:O	2.28	0.65
1:P:41:GLY:H	1:P:280:ASN:ND2	1.94	0.65
1:T:198:ASN:OD1	4:T:610:HOH:O	2.14	0.65
1:D:25:ARG:HD3	4:D:635:HOH:O	1.96	0.65
1:D:235:LYS:N	4:D:608:HOH:O	2.20	0.65
1:M:67:GLN:OE1	1:M:107:TYR:N	2.28	0.65
1:H:219:ALA:HA	1:H:222:ARG:HD2	1.79	0.65
1:C:411[A]:THR:H	1:C:414[A]:MET:HE2	1.61	0.65
1:H:401:ARG:N	1:H:414:MET:HE1	2.09	0.65
1:K:229:SER:O	4:K:609:HOH:O	2.15	0.65
1:B:382:VAL:H	1:B:412:GLN:HE22	1.42	0.64
1:F:156:VAL:HG13	1:F:247:ALA:HA	1.79	0.64
1:R:405:SER:N	4:R:613:HOH:O	2.23	0.64
1:A:177:ASN:HB3	1:A:183:ILE:CD1	2.28	0.64
1:G:67:GLN:NE2	4:G:614:HOH:O	2.30	0.64
1:I:138:TYR:CE2	1:I:140:MET:CE	2.81	0.64
1:K:218:GLU:O	1:K:222:ARG:HG2	1.96	0.64
1:Q:53:ASN:HD22	1:Q:297:TRP:HE1	1.45	0.64
1:I:41:GLY:H	1:I:280:ASN:HD21	1.45	0.64
1:L:140:MET:HE2	1:L:142:TYR:CE1	2.32	0.64
1:L:401:ARG:HD2	4:L:610:HOH:O	1.97	0.64
1:M:319:ARG:HD2	1:N:323:VAL:HG11	1.80	0.64
1:V:53:ASN:HD21	1:V:139:LYS:HA	1.61	0.64
1:K:315:ASP:OD2	4:K:608:HOH:O	2.15	0.64
1:H:12:ASP:OD2	4:H:607:HOH:O	2.14	0.64
1:H:67:GLN:NE2	4:H:602:HOH:O	2.01	0.64
1:D:232:ASP:OD1	4:D:608:HOH:O	2.15	0.64
1:O:221:ARG:NH1	4:O:607:HOH:O	2.18	0.64
1:Q:313:HIS:HE1	2:Q:501:CIT:O3	1.81	0.63
1:O:222:ARG:NH2	1:O:240:TYR:CE1	2.66	0.63
1:T:42[B]:GLY:O	4:T:611:HOH:O	2.14	0.63
1:Q:306:THR:O	1:R:43[B]:LYS:NZ	2.25	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:159:GLY:CA	1:C:206:MET:HE3	2.29	0.63
1:M:408:CYS:SG	4:M:857:HOH:O	2.37	0.63
1:K:67:GLN:OE1	1:K:107:TYR:N	2.30	0.63
1:T:313:HIS:HE1	2:T:501:CIT:O3	1.80	0.63
1:K:140:MET:CE	1:K:208:ILE:HD13	2.28	0.63
1:D:233:PHE:O	1:D:238:ILE:HG12	1.99	0.62
1:K:304:GLU:O	1:K:306:THR:HG23	1.99	0.62
1:V:350:ASN:HD21	1:V:365:LYS:HE3	1.64	0.62
1:B:78:GLN:OE1	1:C:78:GLN:NE2	2.24	0.62
1:Q:165:ARG:NH1	1:Q:192:PRO:O	2.27	0.62
1:T:177:ASN:HB3	1:T:183:ILE:HD11	1.80	0.62
1:L:175:HIS:CD2	1:L:183:ILE:HD11	2.33	0.62
1:R:174:MET:HE3	1:R:182:ILE:HG21	1.82	0.62
1:B:221:ARG:NH1	4:B:617:HOH:O	2.31	0.62
1:O:75:ARG:HD3	1:O:75:ARG:C	2.20	0.62
1:V:110:THR:HG23	1:V:233:PHE:HD1	1.64	0.62
1:V:139:LYS:NZ	4:V:609:HOH:O	2.14	0.62
1:A:226:ASP:N	4:A:613:HOH:O	2.33	0.62
1:B:175:HIS:HA	4:B:624:HOH:O	1.98	0.62
1:E:14:MET:CE	1:E:126:PRO:HG3	2.29	0.62
1:A:175:HIS:HA	4:A:619:HOH:O	2.00	0.61
1:P:41:GLY:H	1:P:280:ASN:HD21	1.46	0.61
1:Q:400:PHE:HD1	1:Q:414:MET:CE	2.12	0.61
1:A:105:GLN:O	1:A:106:TRP:HE3	1.83	0.61
1:J:316:GLU:OE2	4:J:609:HOH:O	2.16	0.61
1:O:198:ASN:ND2	4:O:616:HOH:O	2.34	0.61
1:P:53:ASN:ND2	1:P:297:TRP:HE1	1.97	0.61
1:Q:53:ASN:HD21	1:Q:139:LYS:HA	1.65	0.61
1:G:368:ARG:HH21	2:G:501:CIT:H41	1.65	0.61
1:L:53:ASN:ND2	1:L:297:TRP:HE1	1.97	0.61
1:V:97:ALA:HB1	1:V:105:GLN:O	2.00	0.61
1:E:14:MET:HE1	1:E:121:ILE:HG23	1.82	0.61
1:F:6:VAL:N	4:F:609:HOH:O	2.32	0.61
1:A:348:ARG:HG3	1:A:365:LYS:HE3	1.81	0.61
1:C:222:ARG:CZ	4:C:604:HOH:O	2.49	0.61
1:C:385:GLU:HB3	1:Q:166:MET:HE1	1.82	0.61
4:O:720:HOH:O	1:P:310:LYS:HE2	2.01	0.61
1:D:69:LYS:O	4:D:609:HOH:O	2.16	0.61
1:G:288:PRO:HD2	3:G:502:GOL:H11	1.82	0.61
1:V:18:LEU:HD13	1:V:110:THR:OG1	2.01	0.61
1:K:140:MET:HE2	1:K:142:TYR:CE1	2.34	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:14:MET:CE	1:M:124:TYR:HD2	2.13	0.61
1:T:257:VAL:HG21	1:T:297:TRP:CG	2.36	0.61
1:C:385:GLU:CB	1:Q:166:MET:HE1	2.31	0.61
1:A:32:ARG:HG2	4:A:778:HOH:O	2.01	0.60
3:J:503:GOL:H11	4:J:838:HOH:O	2.01	0.60
1:H:43[B]:LYS:NZ	4:H:608:HOH:O	2.22	0.60
1:D:368:ARG:HH21	2:D:501:CIT:H41	1.67	0.60
1:M:254:ASP:OD2	4:M:607:HOH:O	2.16	0.60
1:H:214:LYS:HD3	1:H:215:PHE:H	1.66	0.60
1:J:68:TYR:CZ	1:J:69:LYS:HG2	2.36	0.60
1:L:241:ILE:HG13	4:L:625:HOH:O	2.00	0.60
1:C:67:GLN:OE1	1:C:107:TYR:N	2.28	0.60
1:G:97:ALA:O	4:G:605:HOH:O	2.15	0.60
2:L:501:CIT:O4	4:L:612:HOH:O	2.17	0.60
1:E:214:LYS:HA	1:E:214:LYS:HE3	1.82	0.60
1:Q:6:VAL:N	4:Q:624:HOH:O	2.35	0.60
1:R:365:LYS:HD3	4:R:609:HOH:O	2.01	0.60
1:T:140:MET:CE	1:T:252:PHE:HD1	2.14	0.60
1:G:397:ALA:O	4:G:606:HOH:O	2.16	0.60
1:I:67:GLN:NE2	4:I:613:HOH:O	2.27	0.60
1:T:401:ARG:NH1	4:T:601:HOH:O	1.86	0.60
1:E:282:ASP:OD1	4:E:607:HOH:O	2.15	0.60
1:M:75:ARG:NH2	1:M:79:ARG:HD3	2.16	0.60
1:H:401:ARG:H	1:H:414:MET:CE	2.11	0.60
3:J:503:GOL:O1	4:J:610:HOH:O	2.17	0.60
1:Q:98:SER:HB3	4:Q:636:HOH:O	2.01	0.60
1:A:317:ASP:HB2	4:A:703:HOH:O	2.01	0.59
1:G:357:ASN:HD22	1:G:374:ASN:HD21	1.50	0.59
1:I:41:GLY:H	1:I:280:ASN:ND2	2.00	0.59
1:L:140:MET:HE1	1:L:208:ILE:HD13	1.83	0.59
1:V:235:LYS:NZ	4:V:602:HOH:O	1.97	0.59
1:R:313:HIS:NE2	2:R:501:CIT:O4	2.28	0.59
1:T:295:LYS:NZ	4:T:616:HOH:O	2.35	0.59
1:H:313:HIS:HE1	2:H:501:CIT:O3	1.85	0.59
1:N:67:GLN:OE1	1:N:107:TYR:N	2.35	0.59
1:N:304:GLU:O	1:N:306:THR:HG23	2.02	0.59
1:H:214:LYS:HD3	1:H:215:PHE:N	2.17	0.59
1:F:382:VAL:H	1:F:412:GLN:NE2	1.99	0.59
1:K:140:MET:CE	1:K:142:TYR:HE1	2.14	0.59
1:B:174:MET:O	4:B:609:HOH:O	2.16	0.59
1:R:145[A]:MET:HG3	1:R:250:HIS:CD2	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:6:VAL:O	4:E:608:HOH:O	2.17	0.59
1:N:312:VAL:HG13	1:N:334:SER:HA	1.85	0.59
1:O:340:ARG:HD3	1:O:357:ASN:ND2	2.17	0.59
1:D:42[B]:GLY:O	4:D:611:HOH:O	2.17	0.58
1:V:53:ASN:HD22	1:V:297:TRP:HE1	1.51	0.58
1:K:368:ARG:HH21	2:K:501:CIT:H41	1.68	0.58
1:B:199:GLU:OE1	4:B:608:HOH:O	2.16	0.58
1:V:68:TYR:CZ	1:V:69:LYS:HG3	2.38	0.58
1:T:140:MET:HE3	1:T:252:PHE:CD1	2.39	0.58
1:A:159:GLY:HA3	1:A:206:MET:HE3	1.85	0.58
1:A:286:VAL:HG22	1:B:312:VAL:HG23	1.84	0.58
1:J:55:LEU:HD21	1:N:88:ARG:HD2	1.86	0.58
1:P:165:ARG:HD3	1:P:199:GLU:O	2.03	0.58
1:R:79:ARG:NH1	4:R:619:HOH:O	2.31	0.58
1:M:286:VAL:HG22	1:N:312:VAL:HG23	1.85	0.58
1:R:174:MET:HE2	1:R:182:ILE:HG21	1.85	0.58
1:C:250:HIS:HE1	1:C:255:SER:OG	1.85	0.58
1:O:42[B]:GLY:O	4:O:606:HOH:O	2.17	0.58
1:I:174:MET:CE	1:I:182:ILE:HG21	2.34	0.58
1:E:67:GLN:OE1	1:E:107:TYR:N	2.37	0.57
1:G:177:ASN:HB3	1:G:183:ILE:HD11	1.85	0.57
1:J:206:MET:HE3	1:J:252:PHE:HB2	1.84	0.57
1:L:214:LYS:HE3	1:L:214:LYS:HA	1.86	0.57
1:C:139:LYS:HG2	1:C:297:TRP:HD1	1.69	0.57
1:J:400:PHE:O	3:J:503:GOL:H32	2.04	0.57
1:T:43[A]:LYS:NZ	4:T:618:HOH:O	2.37	0.57
1:A:310:LYS:HE2	4:B:692:HOH:O	2.05	0.57
1:T:226:ASP:OD1	1:T:228:THR:HG22	2.04	0.57
1:F:67:GLN:OE1	1:F:107:TYR:N	2.37	0.57
1:B:386:GLY:O	4:B:611:HOH:O	2.17	0.57
1:C:222:ARG:NH1	4:C:626:HOH:O	2.38	0.57
1:L:140:MET:HE2	1:L:142:TYR:HE1	1.70	0.57
1:T:69:LYS:HD2	4:T:747:HOH:O	2.03	0.57
1:C:156:VAL:HG12	1:C:241:ILE:HG21	1.86	0.57
1:Q:312:VAL:HG23	1:Q:334:SER:HA	1.86	0.57
1:V:250:HIS:HE1	1:V:255:SER:OG	1.87	0.57
1:K:118:ILE:O	1:K:122:GLU:HG3	2.05	0.57
1:Q:174:MET:HE3	1:Q:182:ILE:HG21	1.85	0.57
1:C:304:GLU:O	1:C:306:THR:HG23	2.05	0.57
1:F:304:GLU:O	1:F:306:THR:HG23	2.04	0.57
1:I:100:ARG:O	1:N:69:LYS:NZ	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:140:MET:HE1	1:K:208:ILE:HD13	1.85	0.57
1:F:338:LEU:HD22	1:F:355:LEU:HD12	1.86	0.56
1:K:53:ASN:ND2	1:K:297:TRP:HE1	2.02	0.56
1:L:205:SER:OG	4:L:603:HOH:O	1.90	0.56
1:M:317:ASP:HB2	4:M:790:HOH:O	2.05	0.56
1:P:221:ARG:NH1	4:P:612:HOH:O	2.24	0.56
1:Q:242:VAL:O	4:Q:605:HOH:O	2.17	0.56
1:Q:410:ILE:HA	1:Q:414:MET:CE	2.35	0.56
1:B:312:VAL:HG13	1:B:334:SER:HA	1.87	0.56
1:E:165:ARG:O	1:E:192:PRO:HG2	2.05	0.56
1:G:67:GLN:OE1	1:G:107:TYR:N	2.38	0.56
1:B:368:ARG:HH21	2:B:501:CIT:H41	1.70	0.56
1:R:145[B]:MET:HG3	1:R:250:HIS:CG	2.40	0.56
1:T:140:MET:HE3	1:T:252:PHE:HD1	1.70	0.56
1:C:357:ASN:HD22	1:C:374:ASN:HD21	1.51	0.56
1:R:350:ASN:ND2	1:R:365:LYS:HE3	2.21	0.56
1:F:75:ARG:HH22	1:F:79:ARG:HD3	1.70	0.56
1:G:197:GLY:O	4:G:607:HOH:O	2.17	0.56
1:A:304:GLU:O	1:A:306:THR:HG23	2.05	0.56
1:C:43[B]:LYS:NZ	4:C:602:HOH:O	2.06	0.56
1:F:75:ARG:NH2	1:F:79:ARG:HD3	2.21	0.56
1:N:401:ARG:NH1	4:N:613:HOH:O	2.28	0.56
1:O:258:ARG:NH2	4:O:619:HOH:O	2.39	0.56
1:B:340:ARG:HD3	1:B:357:ASN:HD21	1.71	0.56
1:A:111:ALA:HB2	1:A:237:ILE:HD11	1.87	0.56
1:C:165:ARG:O	1:C:192:PRO:HG2	2.06	0.56
1:D:239:PRO:HA	1:D:242:VAL:HG22	1.87	0.56
1:G:419:ASP:CG	1:G:420:LEU:H	2.09	0.56
1:N:258:ARG:NH2	4:N:622:HOH:O	2.38	0.56
1:R:221:ARG:HD2	4:R:605:HOH:O	2.06	0.56
1:J:411[A]:THR:H	1:J:414[A]:MET:HE2	1.71	0.56
1:V:232:ASP:HB3	1:V:235:LYS:HB2	1.87	0.56
1:H:175:HIS:HA	4:H:624:HOH:O	2.07	0.55
1:G:165:ARG:HD2	1:G:192:PRO:O	2.06	0.55
1:G:326:VAL:HB	1:G:342:LEU:HD12	1.87	0.55
1:D:249:ALA:O	4:D:612:HOH:O	2.18	0.55
1:K:226:ASP:OD2	1:K:228:THR:N	2.40	0.55
1:O:75:ARG:NE	1:O:79:ARG:HD2	2.21	0.55
1:P:416:ASP:OD2	4:P:606:HOH:O	2.17	0.55
1:M:312:VAL:HG13	1:M:334:SER:HA	1.89	0.55
1:I:79:ARG:CZ	4:I:638:HOH:O	2.54	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:75:ARG:HD2	4:O:769:HOH:O	2.06	0.55
1:J:354:ARG:NE	4:J:619:HOH:O	2.37	0.55
1:M:103:GLU:O	1:M:105:GLN:N	2.40	0.55
1:O:198:ASN:CG	4:O:616:HOH:O	2.45	0.55
1:Q:162:GLU:OE1	4:Q:606:HOH:O	2.17	0.55
1:A:226:ASP:OD2	1:A:228:THR:OG1	2.23	0.55
1:D:148:GLN:NE2	4:D:629:HOH:O	2.39	0.55
1:Q:221:ARG:NH1	4:Q:613:HOH:O	2.26	0.55
1:I:174:MET:HE3	1:I:182:ILE:HG21	1.89	0.55
1:J:206:MET:HE1	1:J:252:PHE:HB2	1.87	0.55
1:J:304:GLU:O	1:J:306:THR:HG23	2.06	0.55
1:N:226:ASP:OD2	1:N:228:THR:HG23	2.07	0.55
1:R:179:LYS:HD3	1:R:181:GLU:CD	2.27	0.55
1:C:214:LYS:HG2	4:C:820:HOH:O	2.06	0.55
1:K:193:PRO:O	3:K:503:GOL:H31	2.07	0.55
1:D:53:ASN:ND2	1:D:297:TRP:HE1	2.04	0.54
1:I:139:LYS:NZ	4:I:624:HOH:O	2.39	0.54
1:M:313:HIS:NE2	2:M:501:CIT:O3	2.27	0.54
1:V:231:ARG:HD3	4:V:612:HOH:O	2.07	0.54
1:B:185:PHE:CD2	1:B:239:PRO:HG3	2.42	0.54
1:E:118:ILE:HG12	1:E:217:MET:SD	2.47	0.54
1:G:25:ARG:NH1	4:G:621:HOH:O	2.40	0.54
1:R:32:ARG:NH2	4:R:621:HOH:O	2.39	0.54
1:B:382:VAL:H	1:B:412:GLN:NE2	2.05	0.54
1:L:313:HIS:CE1	2:L:501:CIT:O3	2.58	0.54
4:V:736:HOH:O	1:T:310:LYS:HE2	2.07	0.54
1:B:176:VAL:HG21	1:B:203:LEU:HD12	1.89	0.54
1:L:401:ARG:CD	4:L:610:HOH:O	2.54	0.54
1:M:206:MET:SD	1:M:208:ILE:HD12	2.48	0.54
1:C:317:ASP:OD2	4:C:605:HOH:O	2.17	0.54
1:C:368:ARG:HH21	2:C:501:CIT:H41	1.73	0.54
1:F:99:GLN:HA	1:F:103:GLU:HA	1.89	0.54
1:I:246:LYS:NZ	4:I:626:HOH:O	2.40	0.54
1:I:368:ARG:HH21	2:I:501:CIT:H41	1.72	0.54
1:C:122:GLU:O	4:C:608:HOH:O	2.18	0.54
1:Q:174:MET:CE	1:Q:182:ILE:HG21	2.38	0.54
1:T:242:VAL:HG22	1:T:247:ALA:HB3	1.89	0.54
1:B:93:ASP:OD2	1:C:75:ARG:NH2	2.40	0.54
1:C:48:ASP:OD1	1:C:76:HIS:HE1	1.91	0.54
1:P:165:ARG:O	1:P:192:PRO:HG2	2.08	0.54
1:I:25:ARG:NH2	4:I:623:HOH:O	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:53:ASN:HD21	1:L:139:LYS:HA	1.73	0.54
1:O:340:ARG:HD3	1:O:357:ASN:HD21	1.72	0.54
1:D:179:LYS:NZ	1:D:179:LYS:HB3	2.23	0.53
1:N:159:GLY:HA3	1:N:206:MET:HE1	1.90	0.53
1:J:69:LYS:H	1:J:69:LYS:HD2	1.74	0.53
1:L:43[A]:LYS:HE2	1:L:292:ILE:HG22	1.90	0.53
1:V:39:TYR:HB3	1:V:344:PHE:CE2	2.43	0.53
1:A:67:GLN:OE1	1:A:106:TRP:HA	2.08	0.53
1:B:69:LYS:O	4:B:612:HOH:O	2.18	0.53
1:L:160:CYS:N	1:L:206:MET:HE2	2.22	0.53
1:M:140:MET:HE2	1:M:142:TYR:HE1	1.72	0.53
1:T:304:GLU:O	1:T:306:THR:HG23	2.08	0.53
1:A:222:ARG:NH2	4:A:623:HOH:O	2.41	0.53
1:F:203:LEU:HD21	1:F:251:ARG:HD3	1.90	0.53
1:V:179:LYS:HB2	1:V:181:GLU:OE2	2.08	0.53
1:D:206:MET:CG	4:D:893:HOH:O	2.56	0.53
1:G:274:ASP:OD2	1:G:405:SER:OG	2.26	0.53
1:J:39:TYR:HB3	1:J:344:PHE:CE1	2.42	0.53
1:E:221:ARG:HD2	4:E:616:HOH:O	2.08	0.53
1:O:304:GLU:O	1:O:306:THR:HG23	2.09	0.53
1:V:68:TYR:CE1	1:V:69:LYS:HG3	2.43	0.53
1:F:145:MET:HG3	1:F:250:HIS:CD2	2.43	0.53
1:Q:53:ASN:ND2	1:Q:297:TRP:HE1	2.07	0.53
1:V:187:GLU:HG2	4:V:602:HOH:O	2.08	0.53
1:B:241:ILE:HG13	4:B:723:HOH:O	2.09	0.53
1:C:187:GLU:OE2	4:C:607:HOH:O	2.18	0.53
1:L:304:GLU:O	1:L:306:THR:HG23	2.09	0.53
1:P:304:GLU:O	1:P:306:THR:HG23	2.08	0.53
1:O:140:MET:HE2	1:O:142:TYR:HE1	1.74	0.52
1:B:214:LYS:HE3	4:B:721:HOH:O	2.07	0.52
1:D:8:PRO:HG2	1:D:11:ARG:HG3	1.91	0.52
1:T:103:GLU:O	1:T:105:GLN:N	2.43	0.52
1:B:250:HIS:HE1	1:B:255:SER:OG	1.92	0.52
1:M:176:VAL:HG23	1:M:196:PRO:HD2	1.91	0.52
1:R:258:ARG:NH2	4:R:625:HOH:O	2.42	0.52
1:Q:400:PHE:CD1	1:Q:414:MET:HE3	2.39	0.52
1:V:214:LYS:HD3	4:V:633:HOH:O	2.08	0.52
1:E:338:LEU:HD22	1:E:355:LEU:HD12	1.92	0.52
1:L:140:MET:CE	1:L:142:TYR:HE1	2.22	0.52
1:R:213:THR:O	1:R:217:MET:HG2	2.09	0.52
1:T:243:GLU:HG2	1:T:244:HIS:CE1	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:165:ARG:HD3	1:H:199:GLU:O	2.10	0.52
1:H:401:ARG:N	1:H:414:MET:CE	2.71	0.52
1:B:100:ARG:NH1	4:B:627:HOH:O	2.43	0.52
1:C:139:LYS:HG2	1:C:297:TRP:CD1	2.44	0.52
1:K:79:ARG:NH2	4:K:619:HOH:O	2.32	0.52
1:Q:103:GLU:O	1:Q:105:GLN:N	2.43	0.52
1:R:187:GLU:OE1	4:R:611:HOH:O	2.19	0.52
1:B:213:THR:O	1:B:217:MET:HG2	2.10	0.51
1:E:75:ARG:CZ	1:E:79:ARG:HH12	2.24	0.51
1:I:75:ARG:CZ	4:I:606:HOH:O	2.51	0.51
1:K:323:VAL:HG23	1:K:339:ASN:HD22	1.74	0.51
1:N:32:ARG:NH2	4:N:624:HOH:O	2.42	0.51
1:N:159:GLY:HA3	1:N:206:MET:CE	2.40	0.51
1:O:140:MET:HE1	1:O:208:ILE:CD1	2.40	0.51
1:T:140:MET:CE	1:T:252:PHE:CD1	2.93	0.51
1:A:233:PHE:HA	1:A:237:ILE:HD13	1.93	0.51
1:I:221:ARG:NH1	4:I:611:HOH:O	2.25	0.51
1:R:304:GLU:O	1:R:306:THR:HG23	2.09	0.51
1:C:39:TYR:HB3	1:C:344:PHE:CE2	2.45	0.51
1:E:75:ARG:NH2	1:E:79:ARG:HH12	2.09	0.51
1:V:350:ASN:ND2	1:V:365:LYS:HE3	2.25	0.51
1:I:93:ASP:OD2	1:N:75:ARG:NH1	2.44	0.51
1:I:149:HIS:HE1	1:I:212:HIS:CD2	2.29	0.51
1:J:145:MET:HB2	4:J:704:HOH:O	2.10	0.51
1:M:213:THR:O	1:M:217:MET:HG2	2.10	0.51
1:Q:8:PRO:HG2	1:Q:11:ARG:HG3	1.92	0.51
1:I:43[B]:LYS:HG2	4:I:649:HOH:O	2.10	0.51
1:O:324:SER:O	1:O:340:ARG:HA	2.10	0.51
1:V:110:THR:HG23	1:V:233:PHE:CD1	2.45	0.51
1:V:368:ARG:HH21	2:V:501:CIT:H41	1.76	0.51
1:H:159:GLY:CA	1:H:206:MET:HE3	2.41	0.51
1:L:238:ILE:O	1:L:242:VAL:HG23	2.10	0.51
1:V:338:LEU:HD22	1:V:355:LEU:HD12	1.93	0.51
1:K:119:ASP:OD1	4:K:610:HOH:O	2.19	0.51
1:M:145:MET:HG3	4:M:712:HOH:O	2.11	0.51
1:V:221:ARG:NH1	4:V:619:HOH:O	2.32	0.51
1:C:286:VAL:HG22	1:D:312:VAL:CG2	2.40	0.51
1:F:179:LYS:HB2	1:F:181:GLU:HG3	1.93	0.51
1:L:221:ARG:NH1	4:L:616:HOH:O	2.28	0.51
1:L:324:SER:O	1:L:340:ARG:HA	2.10	0.51
1:H:39:TYR:HB3	1:H:344:PHE:CE2	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:354:ARG:NH1	4:C:633:HOH:O	2.44	0.51
1:A:267:TRP:CH2	1:A:269:ASP:HB3	2.46	0.51
1:G:214:LYS:CE	1:G:214:LYS:H	2.23	0.51
1:L:140:MET:CE	1:L:208:ILE:HD13	2.40	0.50
1:P:32:ARG:NH2	4:P:627:HOH:O	2.44	0.50
1:C:6:VAL:N	4:C:635:HOH:O	2.44	0.50
1:M:39:TYR:HB3	1:M:344:PHE:CE2	2.46	0.50
1:D:178:GLU:H	1:D:178:GLU:CD	2.13	0.50
1:E:196:PRO:HG3	4:E:672:HOH:O	2.11	0.50
1:M:75:ARG:NH1	4:M:622:HOH:O	2.43	0.50
1:R:159:GLY:HA3	1:R:206:MET:HE3	1.94	0.50
1:B:148:GLN:NE2	1:B:250:HIS:HD2	2.09	0.50
1:C:41:GLY:H	1:C:280:ASN:ND2	2.10	0.50
1:D:304:GLU:O	1:D:306:THR:HG23	2.10	0.50
1:I:304:GLU:O	1:I:306:THR:HG23	2.11	0.50
1:J:409:LEU:O	1:J:414[A]:MET:HE1	2.11	0.50
1:M:312:VAL:HG23	1:N:286:VAL:HG22	1.92	0.50
1:T:27:LYS:NZ	4:T:620:HOH:O	2.39	0.50
1:G:324:SER:O	1:G:340:ARG:HA	2.12	0.50
1:L:219:ALA:HA	1:L:222:ARG:HG2	1.94	0.50
1:E:411[A]:THR:HG22	4:E:628:HOH:O	2.12	0.50
1:F:118:ILE:HG12	1:F:217:MET:SD	2.52	0.50
1:P:165:ARG:NH1	1:P:192:PRO:O	2.42	0.50
1:K:39:TYR:HB3	1:K:344:PHE:CE1	2.47	0.50
1:H:160:CYS:N	1:H:206:MET:CE	2.75	0.50
1:G:403:THR:CG2	1:G:405:SER:H	2.24	0.50
1:N:148:GLN:NE2	4:N:627:HOH:O	2.44	0.50
1:V:127:GLU:OE2	4:V:610:HOH:O	2.18	0.50
1:A:67:GLN:HE22	1:A:107:TYR:N	2.10	0.50
1:B:317:ASP:HB2	4:B:714:HOH:O	2.11	0.50
1:J:179:LYS:HE3	1:J:181:GLU:CD	2.32	0.50
1:J:254:ASP:OD2	4:J:612:HOH:O	2.19	0.49
1:L:103:GLU:O	1:L:105:GLN:N	2.45	0.49
1:L:214:LYS:HA	1:L:214:LYS:CE	2.42	0.49
1:A:176:VAL:N	4:A:619:HOH:O	2.38	0.49
1:C:41:GLY:H	1:C:280:ASN:HD21	1.58	0.49
1:F:156:VAL:HG13	1:F:156:VAL:O	2.11	0.49
1:T:312:VAL:HG13	1:T:334:SER:HA	1.94	0.49
1:H:402:ARG:NH1	4:H:603:HOH:O	2.05	0.49
1:D:123:PRO:HG2	4:D:804:HOH:O	2.13	0.49
1:K:74:ILE:O	1:K:78:GLN:HG3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:14:MET:CE	1:M:124:TYR:CD2	2.95	0.49
1:M:178:GLU:CD	1:M:178:GLU:H	2.15	0.49
1:M:222:ARG:NH1	4:M:618:HOH:O	2.38	0.49
1:H:160:CYS:C	1:H:206:MET:HE2	2.32	0.49
1:C:312:VAL:HG23	1:C:334:SER:HA	1.93	0.49
1:O:145:MET:HG3	1:O:250:HIS:CD2	2.47	0.49
1:V:74:ILE:O	1:V:78:GLN:HG3	2.13	0.49
1:B:174:MET:HB3	1:B:182:ILE:HD12	1.94	0.49
1:I:43[A]:LYS:HE2	1:I:292:ILE:HG22	1.95	0.49
1:J:165:ARG:O	1:J:192:PRO:HG2	2.12	0.49
1:O:140:MET:CE	1:O:142:TYR:HE1	2.25	0.49
1:P:39[B]:TYR:HB3	1:P:344:PHE:CE1	2.47	0.49
1:Q:178:GLU:OE2	4:Q:607:HOH:O	2.20	0.49
1:V:340:ARG:HH11	1:V:357:ASN:ND2	2.11	0.49
1:T:165:ARG:O	1:T:192:PRO:HG2	2.13	0.49
1:B:165:ARG:O	1:B:192:PRO:HG2	2.12	0.49
1:C:226:ASP:OD1	1:C:228:THR:HG22	2.12	0.49
1:E:32:ARG:NH1	4:E:628:HOH:O	2.46	0.49
1:E:420:LEU:O	4:E:609:HOH:O	2.18	0.49
1:M:368:ARG:HH21	2:M:501:CIT:H41	1.76	0.49
1:N:411[A]:THR:HG22	4:N:668:HOH:O	2.13	0.49
1:O:75:ARG:NH1	1:O:302:TYR:CE2	2.81	0.49
1:P:213:THR:O	1:P:217:MET:HG2	2.11	0.49
1:Q:67:GLN:HE22	1:Q:107:TYR:N	2.04	0.49
1:C:76:HIS:HD2	4:C:888:HOH:O	1.96	0.49
1:C:160:CYS:N	1:C:206:MET:HE3	2.27	0.49
1:A:176:VAL:HG22	1:A:180:ASP:HA	1.94	0.49
1:D:28:GLU:HB3	1:D:411[B]:THR:HG21	1.95	0.49
1:K:140:MET:HE3	1:K:208:ILE:HD13	1.95	0.49
1:K:175:HIS:HA	3:K:503:GOL:C1	2.38	0.49
1:O:419:ASP:N	1:O:419:ASP:OD1	2.45	0.49
1:C:213:THR:O	1:C:217:MET:HG2	2.12	0.49
1:D:165:ARG:HD2	1:D:199:GLU:O	2.12	0.49
1:E:310:LYS:HE2	4:F:711:HOH:O	2.13	0.49
1:V:53:ASN:ND2	1:V:297:TRP:HE1	2.10	0.49
1:B:394:GLU:CD	1:B:394:GLU:H	2.15	0.48
1:I:75:ARG:NH1	1:N:93:ASP:OD2	2.46	0.48
1:Q:225:ALA:O	1:Q:227:PRO:HD3	2.13	0.48
1:I:69:LYS:HE3	1:N:100:ARG:HA	1.94	0.48
1:I:145:MET:HG3	1:I:250:HIS:CD2	2.48	0.48
1:L:267:TRP:CH2	1:L:269:ASP:HB3	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:7:GLN:NE2	4:P:614:HOH:O	2.27	0.48
1:R:312:VAL:HG23	1:R:334:SER:HA	1.94	0.48
1:B:88:ARG:HD2	1:D:55:LEU:HD21	1.94	0.48
1:D:206:MET:HG2	4:D:893:HOH:O	2.13	0.48
1:K:148:GLN:NE2	4:K:628:HOH:O	2.46	0.48
1:N:35:LYS:HD3	1:N:135:ASP:HB2	1.95	0.48
1:B:318:ARG:HB3	1:B:352:TYR:CE2	2.48	0.48
1:L:189:PRO:HG2	1:L:192:PRO:HG3	1.95	0.48
1:M:25:ARG:NH1	4:M:609:HOH:O	2.20	0.48
1:E:402:ARG:NH1	4:E:606:HOH:O	2.14	0.48
1:N:221:ARG:HD2	4:N:604:HOH:O	2.12	0.48
1:Q:410:ILE:HA	1:Q:414:MET:HE2	1.95	0.48
1:B:25:ARG:NH1	4:B:632:HOH:O	2.46	0.48
1:C:409:LEU:O	1:C:414[A]:MET:HE1	2.13	0.48
1:M:145:MET:HE1	1:M:210:VAL:HG23	1.96	0.48
1:V:267:TRP:CH2	1:V:269:ASP:HB3	2.49	0.48
1:H:159:GLY:C	1:H:206:MET:HE3	2.34	0.48
1:I:399:ARG:HD2	1:I:418:LEU:HD22	1.96	0.48
1:N:324:SER:O	1:N:340:ARG:HA	2.13	0.48
1:R:105:GLN:O	1:R:106:TRP:CD1	2.66	0.48
1:V:214:LYS:NZ	1:V:217:MET:HG3	2.29	0.48
1:H:324:SER:O	1:H:340:ARG:HA	2.14	0.48
1:I:213:THR:O	1:I:217:MET:HG2	2.14	0.48
1:K:226:ASP:OD1	1:K:228:THR:HG22	2.13	0.48
1:L:165:ARG:HD2	1:L:192:PRO:O	2.14	0.48
1:P:243:GLU:OE1	4:P:607:HOH:O	2.20	0.48
1:T:164:PRO:HB2	1:T:167:GLU:HG3	1.94	0.48
1:H:213:THR:O	1:H:217:MET:HG2	2.14	0.48
1:B:250:HIS:HE1	1:B:255:SER:CB	2.26	0.48
1:C:179:LYS:HB2	1:C:181:GLU:HG3	1.95	0.48
1:L:183:ILE:HG13	1:L:184:ASP:N	2.28	0.48
1:L:316:GLU:H	1:L:316:GLU:CD	2.17	0.48
1:M:348:ARG:HB2	1:M:365:LYS:HD3	1.96	0.48
1:N:318:ARG:HB3	1:N:352:TYR:CE2	2.48	0.48
1:B:177:ASN:HD21	1:B:181:GLU:HB2	1.79	0.48
1:C:43[B]:LYS:NZ	1:D:306:THR:O	2.25	0.48
1:J:179:LYS:NZ	4:J:628:HOH:O	2.47	0.48
1:B:206:MET:HE1	1:B:252:PHE:CD1	2.48	0.47
1:A:88:ARG:HD2	1:C:55:LEU:HD21	1.96	0.47
1:J:179:LYS:HG2	1:J:181:GLU:CG	2.43	0.47
1:T:166:MET:HE3	4:T:922:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:242:VAL:HG22	1:T:247:ALA:CB	2.44	0.47
1:D:242:VAL:HG12	1:D:247:ALA:HB3	1.96	0.47
1:I:25:ARG:NH1	4:I:612:HOH:O	2.27	0.47
1:M:340:ARG:NH2	4:M:621:HOH:O	2.43	0.47
1:A:312:VAL:CG2	1:B:286:VAL:HG22	2.45	0.47
1:B:149:HIS:NE2	1:B:212:HIS:CD2	2.82	0.47
1:D:312:VAL:CG1	1:D:334:SER:HA	2.44	0.47
1:N:399:ARG:NH2	4:N:607:HOH:O	2.23	0.47
1:O:140:MET:HE1	1:O:208:ILE:HD13	1.96	0.47
1:Q:145:MET:HG3	1:Q:250:HIS:CD2	2.49	0.47
1:T:43[B]:LYS:HA	4:T:695:HOH:O	2.14	0.47
1:T:104:THR:HB	1:T:106:TRP:CD1	2.49	0.47
1:T:108:GLU:HG2	1:T:112:ASP:OD2	2.14	0.47
1:G:304:GLU:O	1:G:306:THR:HG23	2.14	0.47
1:I:55:LEU:HD21	1:M:88:ARG:HD2	1.97	0.47
1:K:176:VAL:HG22	3:K:503:GOL:O1	2.14	0.47
1:O:394:GLU:CD	1:O:394:GLU:H	2.17	0.47
1:C:243:GLU:HG3	1:C:244:HIS:CD2	2.49	0.47
1:J:68:TYR:CE2	1:J:69:LYS:HG2	2.49	0.47
1:H:56:ASN:HB2	1:H:297:TRP:CZ2	2.50	0.47
1:B:149:HIS:NE2	1:B:212:HIS:HD2	2.13	0.47
1:C:416:ASP:OD1	4:C:609:HOH:O	2.20	0.47
1:K:177:ASN:OD1	1:K:179:LYS:N	2.39	0.47
4:K:607:HOH:O	1:V:120:ILE:HD12	2.14	0.47
1:L:160:CYS:N	1:L:206:MET:CE	2.77	0.47
1:L:175:HIS:CG	1:L:183:ILE:HD11	2.50	0.47
1:M:165:ARG:O	1:M:192:PRO:HG2	2.15	0.47
1:N:43[B]:LYS:NZ	4:N:602:HOH:O	2.06	0.47
1:V:304:GLU:O	1:V:306:THR:HG23	2.15	0.47
1:V:324:SER:O	1:V:340:ARG:HA	2.14	0.47
1:V:365:LYS:CD	4:V:629:HOH:O	2.63	0.47
1:D:400:PHE:O	4:D:607:HOH:O	2.20	0.47
1:F:159:GLY:HA3	1:F:206:MET:CE	2.40	0.47
1:T:318:ARG:HB3	1:T:352:TYR:CE2	2.50	0.47
1:D:105:GLN:O	1:D:106:TRP:HD1	1.98	0.47
1:F:221:ARG:NH1	4:F:610:HOH:O	2.32	0.47
1:L:116[B]:GLN:OE1	4:L:613:HOH:O	2.20	0.47
1:M:286:VAL:HG22	1:N:312:VAL:CG2	2.44	0.47
1:T:182:ILE:O	1:T:242:VAL:HG13	2.14	0.47
1:A:176:VAL:CG2	1:A:180:ASP:HA	2.45	0.47
1:D:356:GLU:O	1:D:373:SER:HA	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:214:LYS:HA	1:E:214:LYS:CE	2.45	0.47
1:J:18:LEU:HD13	1:J:110:THR:HB	1.97	0.47
1:K:411[A]:THR:HG22	4:K:785:HOH:O	2.15	0.47
1:L:213:THR:O	1:L:217:MET:HG2	2.15	0.47
1:P:354:ARG:CZ	4:P:605:HOH:O	2.59	0.47
1:R:350:ASN:HD21	1:R:365:LYS:HE3	1.79	0.47
1:H:239:PRO:HA	1:H:242:VAL:HG22	1.97	0.46
1:A:324:SER:O	1:A:340:ARG:HA	2.14	0.46
1:E:95:LEU:CD2	1:E:120:ILE:HD12	2.45	0.46
1:L:101:VAL:HG12	4:L:613:HOH:O	2.15	0.46
1:N:385:GLU:HG3	4:N:772:HOH:O	2.15	0.46
1:O:213:THR:O	1:O:217:MET:HG2	2.15	0.46
1:P:312:VAL:HG23	1:P:334:SER:HA	1.97	0.46
1:B:165:ARG:NH2	4:B:636:HOH:O	2.49	0.46
1:I:149:HIS:HD2	1:I:157:THR:OG1	1.98	0.46
1:L:140:MET:CE	1:L:142:TYR:CE1	2.98	0.46
1:V:394:GLU:CD	1:V:394:GLU:H	2.18	0.46
1:T:323:VAL:HG23	1:T:339:ASN:HD22	1.80	0.46
1:F:368:ARG:HH21	2:F:501:CIT:H41	1.79	0.46
1:J:235:LYS:HB2	1:J:235:LYS:HE2	1.79	0.46
1:J:397:ALA:O	3:J:503:GOL:H12	2.15	0.46
1:M:166:MET:H	1:M:166:MET:HG3	1.39	0.46
1:T:344:PHE:HB3	1:T:362:PRO:HG3	1.98	0.46
1:E:14:MET:HE3	1:E:129:MET:SD	2.55	0.46
1:G:12:ASP:HB3	1:G:127:GLU:HB2	1.98	0.46
1:I:99:GLN:HA	1:I:102:SER:O	2.16	0.46
1:A:111:ALA:HB2	1:A:237:ILE:CD1	2.45	0.46
1:C:32:ARG:NH2	4:C:610:HOH:O	2.21	0.46
1:G:273:ILE:HG13	4:G:675:HOH:O	2.14	0.46
1:I:312:VAL:O	1:I:313:HIS:HB2	2.16	0.46
1:P:313:HIS:HE2	2:P:501:CIT:C5	2.28	0.46
1:E:44[A]:ALA:HB1	1:E:48:ASP:HB2	1.97	0.46
1:L:312:VAL:HG23	1:L:334:SER:HA	1.98	0.46
1:V:184:ASP:N	4:V:603:HOH:O	2.10	0.46
1:A:111:ALA:CB	1:A:237:ILE:HD11	2.45	0.46
1:J:178:GLU:CD	1:J:178:GLU:H	2.19	0.46
1:N:411[A]:THR:HG21	4:N:624:HOH:O	2.16	0.46
1:E:39:TYR:HB3	1:E:344:PHE:CE2	2.51	0.46
1:G:39:TYR:HB3	1:G:344:PHE:CE2	2.50	0.46
1:N:267:TRP:CH2	1:N:269:ASP:HB3	2.50	0.46
1:Q:213:THR:O	1:Q:217:MET:HG2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:39:TYR:HB3	1:T:344:PHE:CE1	2.50	0.46
1:B:365:LYS:NZ	4:B:626:HOH:O	2.41	0.46
1:E:99:GLN:HA	1:E:102:SER:O	2.15	0.46
1:I:32:ARG:HB3	1:I:32:ARG:CZ	2.46	0.46
1:I:174:MET:HE2	1:I:182:ILE:HD13	1.97	0.46
1:O:238:ILE:O	1:O:242:VAL:HG13	2.16	0.46
1:O:267:TRP:CH2	1:O:269:ASP:HB3	2.50	0.46
1:B:250:HIS:CE1	1:B:255:SER:HB3	2.51	0.46
1:F:222:ARG:NH2	4:F:603:HOH:O	2.20	0.46
1:N:348:ARG:HG3	1:N:365:LYS:NZ	2.31	0.46
1:O:75:ARG:HD3	1:O:75:ARG:O	2.16	0.46
1:H:400:PHE:HA	1:H:414:MET:CE	2.46	0.45
1:H:411[A]:THR:HG22	4:H:792:HOH:O	2.16	0.45
1:N:312:VAL:O	1:N:313:HIS:HB2	2.16	0.45
1:O:306:THR:O	1:P:43[B]:LYS:NZ	2.35	0.45
1:R:159:GLY:HA3	1:R:206:MET:CE	2.46	0.45
1:C:43[B]:LYS:HG2	4:C:718:HOH:O	2.15	0.45
1:I:88:ARG:HD2	1:M:55:LEU:HD21	1.98	0.45
1:I:149:HIS:CE1	1:I:212:HIS:CD2	3.04	0.45
1:O:368:ARG:HH21	2:O:501:CIT:H41	1.81	0.45
1:T:182:ILE:HG22	1:T:242:VAL:HG11	1.98	0.45
1:B:39:TYR:HB3	1:B:344:PHE:CE2	2.51	0.45
1:G:238:ILE:O	1:G:242:VAL:HG13	2.16	0.45
1:I:79:ARG:NH2	4:I:638:HOH:O	2.49	0.45
1:I:372:LEU:HD11	1:I:383:ILE:HG21	1.99	0.45
1:L:310:LYS:NZ	4:L:629:HOH:O	2.46	0.45
1:N:39:TYR:HB3	1:N:344:PHE:CE2	2.51	0.45
1:O:326:VAL:HG22	1:P:310:LYS:HG3	1.98	0.45
1:V:286:VAL:HG22	1:T:312:VAL:CG2	2.46	0.45
1:D:79:ARG:HG2	1:D:79:ARG:HH11	1.82	0.45
1:G:200:GLY:N	4:G:603:HOH:O	2.05	0.45
1:J:175:HIS:O	1:J:183:ILE:HG12	2.16	0.45
1:K:43[B]:LYS:HE2	4:L:749:HOH:O	2.17	0.45
1:M:413:SER:O	1:M:417:LYS:HE2	2.16	0.45
1:N:221:ARG:NH1	4:N:604:HOH:O	2.16	0.45
1:H:316:GLU:HG3	1:H:317:ASP:N	2.30	0.45
1:F:235:LYS:HB2	1:F:235:LYS:HE2	1.67	0.45
1:L:69:LYS:HD3	4:L:790:HOH:O	2.16	0.45
1:L:145:MET:HG3	1:L:250:HIS:CD2	2.51	0.45
1:L:159:GLY:CA	1:L:206:MET:HE3	2.46	0.45
1:B:165:ARG:HD3	1:B:199:GLU:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:176:VAL:CG1	1:B:180:ASP:HA	2.46	0.45
1:B:214:LYS:HB2	4:B:676:HOH:O	2.17	0.45
1:E:145:MET:HG3	1:E:250:HIS:CD2	2.51	0.45
1:G:262:GLU:HG2	1:G:289:ASP:O	2.17	0.45
1:I:43[A]:LYS:NZ	4:I:642:HOH:O	2.49	0.45
1:K:55:LEU:HD21	1:T:88:ARG:HD2	1.99	0.45
1:K:310:LYS:HE2	4:L:636:HOH:O	2.16	0.45
1:M:221:ARG:HD2	4:M:608:HOH:O	2.15	0.45
1:M:344:PHE:HB3	1:M:362:PRO:HG3	1.99	0.45
1:D:164:PRO:HD2	1:D:167:GLU:HG3	1.97	0.45
1:D:324:SER:O	1:D:340:ARG:HA	2.16	0.45
1:E:318:ARG:HB3	1:E:352:TYR:CE2	2.51	0.45
1:F:324:SER:O	1:F:340:ARG:HA	2.16	0.45
1:J:154:ALA:HA	1:J:246:LYS:HD3	1.99	0.45
1:K:159:GLY:HA3	1:K:206:MET:HE3	1.98	0.45
4:M:783:HOH:O	1:N:43[B]:LYS:HE3	2.16	0.45
1:R:165:ARG:HD2	1:R:192:PRO:O	2.16	0.45
1:D:39:TYR:HB3	1:D:344:PHE:CE1	2.51	0.45
1:N:43[B]:LYS:HG2	1:N:301:THR:HG21	1.98	0.45
1:F:404:GLU:CD	1:F:404:GLU:H	2.21	0.45
1:Q:158:ILE:HD12	1:Q:174:MET:HE1	1.99	0.45
1:V:69:LYS:NZ	4:V:606:HOH:O	2.09	0.45
1:C:165:ARG:NH1	4:C:628:HOH:O	2.49	0.45
1:V:340:ARG:HH11	1:V:357:ASN:HD21	1.65	0.45
1:T:354:ARG:HD3	4:T:605:HOH:O	2.17	0.45
1:H:234:GLY:HA2	4:H:606:HOH:O	2.17	0.44
1:A:67:GLN:HE22	1:A:107:TYR:H	1.63	0.44
1:B:226:ASP:OD2	1:B:228:THR:N	2.43	0.44
1:C:310:LYS:HE2	4:D:713:HOH:O	2.18	0.44
1:G:145:MET:HG3	1:G:250:HIS:CD2	2.52	0.44
1:L:165:ARG:HG3	4:L:609:HOH:O	2.17	0.44
1:O:313:HIS:NE2	2:O:501:CIT:O3	2.29	0.44
1:R:18:LEU:HD13	1:R:110:THR:HB	1.99	0.44
1:T:104:THR:HB	1:T:106:TRP:HD1	1.83	0.44
1:C:148:GLN:OE1	1:C:250:HIS:HD2	2.01	0.44
1:D:67:GLN:OE1	1:D:107:TYR:N	2.45	0.44
1:F:267:TRP:CH2	1:F:269:ASP:HB3	2.52	0.44
1:I:138:TYR:CE2	1:I:140:MET:HE1	2.52	0.44
1:J:324:SER:O	1:J:340:ARG:HA	2.16	0.44
1:L:104:THR:HB	1:L:106:TRP:CD1	2.52	0.44
1:L:165:ARG:HG2	1:L:202:ALA:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:149:HIS:HE1	1:R:212:HIS:CD2	2.36	0.44
1:H:165:ARG:O	1:H:192:PRO:HG2	2.17	0.44
1:A:55:LEU:HD21	1:C:88:ARG:HD2	2.00	0.44
1:A:332:ILE:HB	1:A:348:ARG:HD3	1.99	0.44
1:J:274:ASP:OD1	1:J:405:SER:OG	2.34	0.44
1:M:339:ASN:O	1:M:356:GLU:HA	2.17	0.44
1:N:187:GLU:HG2	1:N:235:LYS:HZ1	1.81	0.44
1:H:75:ARG:NH2	4:H:628:HOH:O	2.50	0.44
1:E:95:LEU:HD22	1:E:120:ILE:HD12	1.99	0.44
1:G:206:MET:CE	1:G:252:PHE:HB2	2.47	0.44
1:G:214:LYS:H	1:G:214:LYS:HE2	1.82	0.44
1:M:304:GLU:O	1:M:306:THR:HG23	2.17	0.44
1:O:295:LYS:HB2	1:O:295:LYS:HE2	1.54	0.44
1:H:143:GLU:O	1:H:147:GLN:HG2	2.17	0.44
1:H:221:ARG:NH1	4:H:612:HOH:O	2.27	0.44
1:A:159:GLY:CA	1:A:206:MET:HE3	2.46	0.44
1:A:313:HIS:NE2	2:A:501:CIT:O3	2.35	0.44
1:L:104:THR:HB	1:L:106:TRP:HD1	1.83	0.44
1:N:214:LYS:NZ	4:N:615:HOH:O	2.30	0.44
1:O:31:ASP:O	1:O:69:LYS:NZ	2.37	0.44
1:K:267:TRP:CH2	1:K:269:ASP:HB3	2.53	0.44
1:D:238:ILE:HB	1:D:239:PRO:HD3	1.98	0.44
1:N:206:MET:HE2	1:N:206:MET:HB2	1.72	0.44
1:Q:217:MET:O	1:Q:221:ARG:HG3	2.17	0.44
1:G:213:THR:O	1:G:217:MET:HG2	2.18	0.44
1:E:43[B]:LYS:NZ	1:F:306:THR:O	2.29	0.44
1:E:69:LYS:HE3	1:E:69:LYS:HB3	1.72	0.44
1:E:418:LEU:HD23	1:E:418:LEU:HA	1.84	0.44
1:I:98:SER:OG	4:I:608:HOH:O	2.21	0.44
1:I:159:GLY:HA3	1:I:206:MET:HE3	2.00	0.44
1:I:368:ARG:HB2	1:I:385:GLU:OE2	2.18	0.44
1:K:160:CYS:N	1:K:206:MET:HE2	2.32	0.44
1:K:213:THR:O	1:K:217:MET:HG2	2.18	0.44
1:L:97:ALA:O	4:L:614:HOH:O	2.21	0.44
1:P:43[B]:LYS:HA	4:P:642:HOH:O	2.18	0.44
1:P:401:ARG:NH1	4:P:616:HOH:O	2.30	0.44
1:Q:400:PHE:CD1	1:Q:414:MET:CE	2.97	0.44
1:R:154:ALA:HA	1:R:246:LYS:HG2	2.00	0.44
1:R:368:ARG:HH21	2:R:501:CIT:H41	1.82	0.44
1:A:69:LYS:HE3	1:A:69:LYS:HB3	1.73	0.43
1:R:69:LYS:O	4:R:612:HOH:O	2.21	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:160:CYS:N	1:H:206:MET:HE3	2.33	0.43
1:C:159:GLY:HA3	1:C:206:MET:HE3	1.97	0.43
1:E:67:GLN:OE1	1:E:106:TRP:HA	2.18	0.43
1:E:238:ILE:O	1:E:242:VAL:HG23	2.18	0.43
1:E:319:ARG:HD2	1:F:323:VAL:HG11	2.00	0.43
1:P:156:VAL:HG12	1:P:241:ILE:HG21	2.00	0.43
1:R:39:TYR:HB3	1:R:344:PHE:CE1	2.53	0.43
1:H:160:CYS:N	1:H:206:MET:HE2	2.33	0.43
1:G:189:PRO:HG2	1:G:192:PRO:HG3	2.00	0.43
1:K:226:ASP:OD2	1:K:226:ASP:C	2.57	0.43
1:P:224:ALA:O	4:P:609:HOH:O	2.21	0.43
1:Q:165:ARG:O	1:Q:192:PRO:HG2	2.19	0.43
1:B:18:LEU:HD13	1:B:110:THR:HB	2.01	0.43
1:E:304:GLU:O	1:E:306:THR:HG23	2.17	0.43
1:J:32:ARG:HH12	1:J:411[A]:THR:HG21	1.84	0.43
1:L:126:PRO:HA	4:L:690:HOH:O	2.17	0.43
1:M:338:LEU:HD22	1:M:355:LEU:HD12	2.00	0.43
1:P:118:ILE:HG12	1:P:217:MET:SD	2.58	0.43
1:A:160:CYS:N	1:A:206:MET:HE2	2.33	0.43
1:E:186:ILE:HG21	1:E:189:PRO:HB3	1.99	0.43
1:G:221:ARG:HD2	4:G:622:HOH:O	2.18	0.43
1:N:348:ARG:HG3	1:N:365:LYS:HE2	2.00	0.43
1:V:356:GLU:O	1:V:373:SER:HA	2.18	0.43
1:T:345[B]:THR:HG23	4:T:611:HOH:O	2.18	0.43
1:H:18:LEU:HD13	1:H:110:THR:HB	2.00	0.43
1:H:242:VAL:HG12	1:H:247:ALA:CB	2.48	0.43
1:A:213:THR:O	1:A:217:MET:HG2	2.18	0.43
1:B:225:ALA:C	4:B:606:HOH:O	2.51	0.43
1:B:374:ASN:HB3	1:B:407:ILE:HD13	2.00	0.43
1:D:143:GLU:O	1:D:147:GLN:HG2	2.18	0.43
1:E:214:LYS:HE3	4:E:828:HOH:O	2.18	0.43
1:G:43[A]:LYS:NZ	4:G:627:HOH:O	2.44	0.43
1:G:274:ASP:OD2	1:G:403:THR:HG21	2.18	0.43
1:L:12:ASP:HB3	1:L:127:GLU:HB2	2.00	0.43
1:M:418:LEU:HD23	1:M:418:LEU:HA	1.84	0.43
1:C:160:CYS:N	1:C:206:MET:CE	2.82	0.43
1:C:368:ARG:HH21	2:C:501:CIT:C4	2.31	0.43
1:D:171:PHE:CG	1:D:204:ALA:HB1	2.54	0.43
1:G:221:ARG:NH1	4:G:622:HOH:O	2.41	0.43
1:I:372:LEU:HD23	1:I:389:VAL:HB	2.01	0.43
1:J:414[A]:MET:HE2	1:J:414[A]:MET:HB2	1.93	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:132:LEU:HD23	1:O:208:ILE:HG23	1.99	0.43
1:V:310:LYS:HE2	4:T:724:HOH:O	2.18	0.43
1:T:310:LYS:NZ	4:T:631:HOH:O	2.48	0.43
1:B:75:ARG:HD3	4:B:880:HOH:O	2.18	0.43
1:T:69:LYS:HD3	4:T:678:HOH:O	2.18	0.43
1:T:145:MET:HG3	1:T:250:HIS:CD2	2.53	0.43
1:J:411[A]:THR:HG22	4:J:706:HOH:O	2.17	0.43
1:O:39:TYR:HB3	1:O:344:PHE:CE2	2.54	0.43
1:V:286:VAL:HG22	1:T:312:VAL:HG23	2.00	0.43
1:A:186:ILE:HG21	1:A:189:PRO:HB3	2.01	0.43
1:F:155:ASP:HB3	1:F:215:PHE:CZ	2.54	0.43
1:J:88:ARG:HD2	1:N:55:LEU:HD21	2.01	0.43
1:J:400:PHE:HA	1:J:414[A]:MET:HE3	2.01	0.43
1:L:39:TYR:HB3	1:L:344:PHE:CE2	2.54	0.43
1:M:140:MET:CE	1:M:208:ILE:HD13	2.49	0.43
1:M:324:SER:O	1:M:340:ARG:HA	2.18	0.43
1:D:192:PRO:HA	1:D:193:PRO:HD3	1.96	0.42
1:E:199:GLU:CD	1:E:199:GLU:H	2.22	0.42
1:K:148:GLN:NE2	1:K:250:HIS:ND1	2.67	0.42
1:L:312:VAL:O	1:L:313:HIS:HB2	2.19	0.42
1:N:32:ARG:NH1	1:N:32:ARG:HB2	2.34	0.42
1:O:393:PRO:HD2	1:O:394:GLU:OE1	2.18	0.42
1:T:257:VAL:HG23	1:T:257:VAL:O	2.18	0.42
1:H:43[B]:LYS:HE3	4:G:849:HOH:O	2.19	0.42
1:H:67:GLN:HB3	1:H:68:TYR:H	1.64	0.42
1:A:165:ARG:O	1:A:192:PRO:HG2	2.19	0.42
1:A:214:LYS:HA	1:A:214:LYS:HD2	1.78	0.42
1:A:286:VAL:HG13	1:B:312:VAL:HG21	2.01	0.42
1:O:32:ARG:HH21	1:O:69:LYS:NZ	2.17	0.42
1:V:277:TRP:CD2	1:V:407:ILE:HD12	2.54	0.42
1:A:319:ARG:HD2	1:B:323:VAL:HG11	2.01	0.42
1:C:357:ASN:HD22	1:C:374:ASN:ND2	2.15	0.42
1:D:34:ALA:HB2	1:D:68:TYR:O	2.19	0.42
1:G:156:VAL:HG12	1:G:241:ILE:HG21	2.01	0.42
1:L:317:ASP:HB2	4:L:711:HOH:O	2.19	0.42
1:L:214:LYS:HG2	4:L:738:HOH:O	2.20	0.42
1:N:340:ARG:HH11	1:N:357:ASN:ND2	2.18	0.42
1:T:140:MET:HG2	1:T:142:TYR:CD1	2.55	0.42
1:H:43[B]:LYS:NZ	1:G:306:THR:O	2.42	0.42
1:C:324:SER:O	1:C:340:ARG:HA	2.19	0.42
1:I:312:VAL:HG23	1:I:334:SER:HA	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:399:ARG:HB3	1:L:418:LEU:HD21	2.01	0.42
1:N:353:SER:HA	1:N:370:ALA:O	2.18	0.42
1:A:145:MET:HG3	1:A:250:HIS:CD2	2.54	0.42
1:A:368:ARG:HH21	2:A:501:CIT:H41	1.85	0.42
1:G:159:GLY:CA	1:G:206:MET:HE3	2.50	0.42
1:G:418:LEU:HD23	1:G:418:LEU:HA	1.83	0.42
1:J:338:LEU:HD22	1:J:355:LEU:HD12	2.01	0.42
1:K:393:PRO:HB3	1:K:402:ARG:CZ	2.50	0.42
1:Q:409:LEU:O	1:Q:414:MET:HE1	2.19	0.42
1:R:369:HIS:HE1	4:R:852:HOH:O	2.01	0.42
1:H:400:PHE:HA	1:H:414:MET:HE2	2.01	0.42
1:A:165:ARG:HH11	1:A:165:ARG:HD2	1.71	0.42
1:F:344:PHE:HB3	1:F:362:PRO:HG3	2.00	0.42
1:F:369:HIS:HE1	4:F:858:HOH:O	2.02	0.42
1:I:162:GLU:OE1	4:I:609:HOH:O	2.22	0.42
1:P:6:VAL:N	4:P:635:HOH:O	2.52	0.42
1:V:213:THR:O	1:V:217:MET:HG2	2.20	0.42
1:V:250:HIS:CE1	1:V:255:SER:OG	2.71	0.42
1:A:350:ASN:HB3	2:A:501:CIT:O1	2.20	0.42
1:B:199:GLU:HG2	1:F:391:GLU:OE1	2.19	0.42
1:F:350:ASN:HB3	2:F:501:CIT:O1	2.20	0.42
1:J:318:ARG:HB3	1:J:352:TYR:CE2	2.55	0.42
1:O:418:LEU:HD23	1:O:418:LEU:HA	1.90	0.42
1:B:324:SER:O	1:B:340:ARG:HA	2.19	0.42
1:E:324:SER:O	1:E:340:ARG:HA	2.19	0.42
1:E:340:ARG:HD3	1:E:357:ASN:OD1	2.20	0.42
1:F:313:HIS:NE2	2:F:501:CIT:O3	2.25	0.42
1:M:75:ARG:HD2	1:M:75:ARG:O	2.19	0.42
1:N:74:ILE:O	1:N:78:GLN:HG3	2.19	0.42
1:P:165:ARG:HH11	1:P:165:ARG:HD2	1.70	0.42
1:H:338:LEU:HD22	1:H:355:LEU:HD12	2.02	0.42
1:A:154:ALA:HA	1:A:246:LYS:HD2	2.02	0.42
1:A:318:ARG:HB3	1:A:352:TYR:CE1	2.55	0.42
1:B:356:GLU:O	1:B:373:SER:HA	2.20	0.42
1:D:105:GLN:C	1:D:106:TRP:CD1	2.93	0.42
1:M:75:ARG:CZ	1:M:79:ARG:HD3	2.50	0.42
1:M:235:LYS:HG2	1:M:236:ASP:OD2	2.19	0.42
1:M:282:ASP:HB3	4:M:806:HOH:O	2.20	0.42
1:P:103:GLU:O	1:P:105:GLN:N	2.53	0.42
1:P:418:LEU:HD23	1:P:418:LEU:HA	1.78	0.42
1:Q:105:GLN:C	1:Q:106:TRP:HD1	2.23	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:165:ARG:NH1	1:R:191:ASP:OD2	2.35	0.42
1:H:267:TRP:CH2	1:H:269:ASP:HB3	2.55	0.41
1:F:318:ARG:HB3	1:F:352:TYR:CE1	2.55	0.41
1:M:312:VAL:CG2	1:N:286:VAL:HG22	2.50	0.41
1:N:374:ASN:HB3	1:N:407:ILE:HD13	2.01	0.41
1:P:266:TYR:CZ	1:P:268:ARG:HG3	2.55	0.41
1:R:165:ARG:O	1:R:192:PRO:HG2	2.20	0.41
1:T:221:ARG:NH1	4:T:626:HOH:O	2.44	0.41
1:C:414[A]:MET:HE2	1:C:414[A]:MET:HB2	1.82	0.41
1:E:313:HIS:NE2	2:E:501:CIT:O4	2.27	0.41
1:M:419:ASP:HA	4:M:703:HOH:O	2.20	0.41
1:N:340:ARG:HH11	1:N:357:ASN:HD21	1.68	0.41
1:Q:324:SER:O	1:Q:340:ARG:HA	2.20	0.41
1:Q:410:ILE:HA	1:Q:414:MET:HE1	2.00	0.41
1:H:37:ALA:HB2	1:H:73:LEU:HB2	2.02	0.41
1:C:339:ASN:O	1:C:356:GLU:HA	2.21	0.41
1:F:192:PRO:HA	1:F:193:PRO:HD3	1.94	0.41
1:I:226:ASP:OD2	1:I:228:THR:HB	2.20	0.41
1:K:222:ARG:HG2	1:K:222:ARG:H	1.66	0.41
1:N:166:MET:H	1:N:166:MET:HG3	1.46	0.41
1:O:25:ARG:NH1	4:O:645:HOH:O	2.54	0.41
1:O:192:PRO:HA	1:O:193:PRO:HD3	1.93	0.41
1:P:372:LEU:HD11	1:P:383:ILE:HG21	2.02	0.41
1:V:365:LYS:NZ	4:V:629:HOH:O	2.46	0.41
1:A:349:ALA:HA	1:A:366:ILE:HB	2.03	0.41
1:G:419:ASP:CG	1:G:420:LEU:N	2.72	0.41
1:J:165:ARG:HD2	1:J:192:PRO:O	2.20	0.41
1:J:213:THR:O	1:J:217:MET:HG2	2.20	0.41
1:Q:79:ARG:NE	4:Q:647:HOH:O	2.54	0.41
1:Q:313:HIS:HE1	2:Q:501:CIT:C5	2.33	0.41
1:V:12:ASP:HB3	1:V:127:GLU:HB2	2.02	0.41
1:V:110:THR:HG23	1:V:233:PHE:H	1.79	0.41
1:T:213:THR:O	1:T:217:MET:HG2	2.20	0.41
1:H:165:ARG:NH1	1:H:192:PRO:O	2.53	0.41
1:B:246:LYS:N	4:B:605:HOH:O	2.52	0.41
1:D:33:ARG:NH1	4:D:642:HOH:O	2.53	0.41
1:L:43[A]:LYS:HG3	4:L:827:HOH:O	2.19	0.41
1:L:55:LEU:HD21	1:V:88:ARG:HD2	2.03	0.41
1:O:239:PRO:O	1:O:242:VAL:HG22	2.20	0.41
1:R:404:GLU:HB3	4:R:613:HOH:O	2.20	0.41
1:V:165:ARG:CD	1:V:199:GLU:O	2.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:242:VAL:HG12	1:H:247:ALA:HB3	2.03	0.41
1:A:39:TYR:HB3	1:A:344:PHE:CE2	2.56	0.41
1:B:74:ILE:O	1:B:78:GLN:HG3	2.20	0.41
1:B:399:ARG:HG2	1:B:418:LEU:HD21	2.01	0.41
1:F:165:ARG:O	1:F:192:PRO:HG2	2.20	0.41
1:K:88:ARG:HD2	1:T:55:LEU:HD21	2.02	0.41
1:K:203:LEU:N	3:K:503:GOL:O3	2.46	0.41
1:K:356:GLU:O	1:K:373:SER:HA	2.20	0.41
1:L:127:GLU:OE2	4:L:615:HOH:O	2.22	0.41
1:M:143:GLU:OE2	1:M:147:GLN:NE2	2.54	0.41
1:O:331[A]:CYS:SG	1:O:347:VAL:HB	2.61	0.41
1:P:267:TRP:CH2	1:P:269:ASP:HB3	2.55	0.41
1:V:310:LYS:NZ	4:V:628:HOH:O	2.46	0.41
1:E:340:ARG:HD3	1:E:340:ARG:HH11	1.74	0.41
1:F:39:TYR:HB3	1:F:344:PHE:CE1	2.56	0.41
1:G:239:PRO:O	1:G:242:VAL:HG22	2.20	0.41
2:J:501:CIT:H42	4:J:604:HOH:O	2.21	0.41
1:K:238:ILE:O	1:K:242:VAL:HG13	2.21	0.41
1:M:156:VAL:HG12	1:M:241:ILE:HG21	2.02	0.41
1:Q:326:VAL:HG22	1:R:310:LYS:HG3	2.02	0.41
1:Q:400:PHE:HA	1:Q:414:MET:HE3	2.03	0.41
1:V:53:ASN:HB3	1:V:142:TYR:CE2	2.56	0.41
1:H:316:GLU:HG3	1:H:317:ASP:CG	2.40	0.41
1:L:165:ARG:HD2	1:L:165:ARG:HH11	1.72	0.41
1:O:232:ASP:HB3	1:O:235:LYS:HG2	2.02	0.41
1:Q:318:ARG:HB3	1:Q:352:TYR:CE2	2.56	0.41
1:T:106:TRP:HD1	1:T:106:TRP:O	2.03	0.41
1:T:165:ARG:HD2	1:T:192:PRO:O	2.20	0.41
1:T:312:VAL:O	1:T:313:HIS:HB2	2.21	0.41
1:A:44[A]:ALA:HB1	1:A:48:ASP:HB2	2.03	0.41
1:A:192:PRO:HA	1:A:193:PRO:HD3	1.94	0.41
1:D:333:ILE:HD12	1:D:338:LEU:HD11	2.03	0.41
1:F:330:ASP:O	1:F:346:GLY:HA2	2.20	0.41
1:G:368:ARG:HH21	2:G:501:CIT:C4	2.33	0.41
1:I:295:LYS:NZ	4:I:616:HOH:O	2.31	0.41
1:J:132:LEU:HD23	1:J:208:ILE:HG23	2.02	0.41
1:L:186:ILE:HA	4:L:823:HOH:O	2.21	0.41
1:M:145:MET:HG2	1:M:250:HIS:CG	2.56	0.41
1:O:69:LYS:HG2	4:O:772:HOH:O	2.21	0.41
1:P:277:TRP:CE2	1:P:281:ILE:HG13	2.56	0.41
1:R:399:ARG:HH21	1:R:420:LEU:C	2.24	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:37:ALA:HB2	1:T:73:LEU:HB2	2.01	0.41
1:T:79:ARG:NH2	4:T:633:HOH:O	2.50	0.41
1:A:217:MET:O	1:A:221:ARG:HG3	2.21	0.41
1:A:286:VAL:HG22	1:B:312:VAL:CG2	2.51	0.41
1:D:61:ARG:HG3	4:D:844:HOH:O	2.20	0.41
1:F:28:GLU:HB3	1:F:411[B]:THR:HG21	2.03	0.41
1:F:415:ILE:O	1:F:418:LEU:HG	2.21	0.41
1:G:178:GLU:HG2	1:G:179:LYS:HG3	2.03	0.41
1:M:140:MET:CE	1:M:142:TYR:HE1	2.34	0.41
1:O:25:ARG:NH1	4:O:646:HOH:O	2.54	0.41
1:O:417:LYS:NZ	4:O:614:HOH:O	2.33	0.41
1:Q:178:GLU:H	1:Q:178:GLU:CD	2.24	0.41
1:H:173:VAL:HG11	1:H:193:PRO:HD2	2.02	0.40
1:A:12:ASP:OD2	4:A:607:HOH:O	2.22	0.40
1:D:148:GLN:HE22	1:D:250:HIS:HB2	1.86	0.40
1:I:323:VAL:HG23	1:I:339:ASN:HD22	1.86	0.40
1:L:281:ILE:HD11	1:L:340:ARG:O	2.21	0.40
1:A:28:GLU:HB3	1:A:411[B]:THR:HG21	2.02	0.40
1:K:214:LYS:NZ	4:K:632:HOH:O	2.49	0.40
1:K:353:SER:HA	1:K:370:ALA:O	2.21	0.40
1:Q:304:GLU:O	1:Q:306:THR:HG23	2.21	0.40
1:H:217:MET:O	1:H:221:ARG:HG3	2.21	0.40
1:E:187:GLU:OE2	1:E:188:LYS:NZ	2.53	0.40
1:G:186:ILE:HD13	1:G:193:PRO:HG2	2.02	0.40
1:I:174:MET:HE3	1:I:182:ILE:CG2	2.51	0.40
1:L:192:PRO:HA	1:L:193:PRO:HD3	1.95	0.40
1:L:226:ASP:OD2	1:L:228:THR:HB	2.21	0.40
1:N:166:MET:HG3	4:N:879:HOH:O	2.21	0.40
1:P:44[A]:ALA:HB1	1:P:48:ASP:HB2	2.03	0.40
1:V:231:ARG:HG3	1:V:231:ARG:HH11	1.86	0.40
1:H:345[B]:THR:HG23	4:H:611:HOH:O	2.20	0.40
1:E:333:ILE:HD12	1:E:338:LEU:HD11	2.04	0.40
1:F:70:ALA:O	1:F:74:ILE:HG12	2.21	0.40
2:J:501:CIT:C6	4:J:604:HOH:O	2.65	0.40
1:K:239:PRO:O	1:K:242:VAL:HG22	2.22	0.40
1:M:75:ARG:HD2	1:M:79:ARG:HG2	2.02	0.40
1:R:45:ARG:HH11	1:R:45:ARG:HD3	1.77	0.40
1:C:118:ILE:HG12	1:C:217:MET:SD	2.62	0.40
1:D:206:MET:HE2	1:D:206:MET:HB2	1.85	0.40
1:E:313:HIS:HE2	2:E:501:CIT:C5	2.25	0.40
1:G:312:VAL:O	1:G:313:HIS:HB2	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:235:LYS:O	1:I:239:PRO:HG2	2.21	0.40
1:M:155:ASP:HB3	1:M:215:PHE:CZ	2.56	0.40
1:R:365:LYS:CD	4:R:609:HOH:O	2.67	0.40
1:V:222:ARG:NH2	4:V:605:HOH:O	2.08	0.40

All (19) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:864:HOH:O	4:P:949:HOH:O[1_455]	1.80	0.40
4:F:803:HOH:O	4:R:826:HOH:O[1_565]	1.84	0.36
4:H:794:HOH:O	4:P:722:HOH:O[1_565]	1.87	0.33
4:F:940:HOH:O	4:Q:1021:HOH:O[1_565]	1.89	0.31
4:E:826:HOH:O	4:Q:671:HOH:O[1_565]	2.01	0.19
4:G:922:HOH:O	4:O:939:HOH:O[1_565]	2.01	0.19
1:F:75:ARG:NH1	1:Q:93:ASP:OD2[1_565]	2.04	0.16
4:H:940:HOH:O	4:P:948:HOH:O[1_565]	2.07	0.13
4:F:889:HOH:O	4:Q:973:HOH:O[1_565]	2.08	0.12
4:F:920:HOH:O	4:R:977:HOH:O[1_565]	2.08	0.12
4:C:830:HOH:O	4:Q:928:HOH:O[1_455]	2.10	0.10
4:F:722:HOH:O	4:R:617:HOH:O[1_565]	2.11	0.09
4:I:965:HOH:O	4:T:861:HOH:O[1_556]	2.11	0.09
4:D:865:HOH:O	4:P:896:HOH:O[1_455]	2.12	0.08
4:D:881:HOH:O	4:P:705:HOH:O[1_455]	2.12	0.08
4:F:959:HOH:O	4:Q:1021:HOH:O[1_565]	2.12	0.08
4:G:885:HOH:O	4:O:806:HOH:O[1_565]	2.14	0.06
4:E:953:HOH:O	4:R:997:HOH:O[1_565]	2.17	0.03
4:D:614:HOH:O	4:P:891:HOH:O[1_455]	2.19	0.01

## 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	419/440 (95%)	403 (96%)	13 (3%)	3 (1%)	22	10
1	B	419/440 (95%)	398 (95%)	18 (4%)	3 (1%)	22	10
1	C	414/440 (94%)	398 (96%)	14 (3%)	2 (0%)	29	15
1	D	419/440 (95%)	402 (96%)	15 (4%)	2 (0%)	29	15
1	E	419/440 (95%)	397 (95%)	19 (4%)	3 (1%)	22	10
1	F	419/440 (95%)	393 (94%)	25 (6%)	1 (0%)	47	33
1	G	420/440 (96%)	400 (95%)	16 (4%)	4 (1%)	15	5
1	H	419/440 (95%)	399 (95%)	16 (4%)	4 (1%)	15	5
1	I	420/440 (96%)	403 (96%)	14 (3%)	3 (1%)	22	10
1	J	411/440 (93%)	395 (96%)	15 (4%)	1 (0%)	47	33
1	K	419/440 (95%)	399 (95%)	16 (4%)	4 (1%)	15	5
1	L	420/440 (96%)	396 (94%)	20 (5%)	4 (1%)	15	5
1	M	419/440 (95%)	398 (95%)	18 (4%)	3 (1%)	22	10
1	N	419/440 (95%)	400 (96%)	16 (4%)	3 (1%)	22	10
1	O	421/440 (96%)	398 (94%)	21 (5%)	2 (0%)	29	15
1	P	421/440 (96%)	400 (95%)	17 (4%)	4 (1%)	15	5
1	Q	420/440 (96%)	401 (96%)	16 (4%)	3 (1%)	22	10
1	R	420/440 (96%)	403 (96%)	16 (4%)	1 (0%)	47	33
1	T	420/440 (96%)	398 (95%)	18 (4%)	4 (1%)	15	5
1	V	419/440 (95%)	399 (95%)	19 (4%)	1 (0%)	47	33
All	All	8377/8800 (95%)	7980 (95%)	342 (4%)	55 (1%)	22	10

All (55) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	104	THR
1	A	104	THR
1	B	104	THR
1	C	104	THR
1	D	104	THR
1	E	104	THR
1	G	104	THR
1	I	104	THR
1	L	104	THR
1	M	104	THR
1	O	104	THR

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Mol	Chain	Res	Type
1	P	104	THR
1	Q	104	THR
1	T	104	THR
1	J	69	LYS
1	K	104	THR
1	N	198	ASN
1	H	67	GLN
1	H	99	GLN
1	C	69	LYS
1	E	69	LYS
1	F	69	LYS
1	G	198	ASN
1	L	198	ASN
1	M	69	LYS
1	N	67	GLN
1	Q	69	LYS
1	T	69	LYS
1	H	69	LYS
1	A	67	GLN
1	B	69	LYS
1	B	313	HIS
1	D	67	GLN
1	E	67	GLN
1	K	67	GLN
1	K	100	ARG
1	P	67	GLN
1	P	69	LYS
1	P	100	ARG
1	Q	67	GLN
1	R	67	GLN
1	V	105	GLN
1	T	67	GLN
1	G	67	GLN
1	G	69	LYS
1	L	67	GLN
1	M	67	GLN
1	N	313	HIS
1	O	99	GLN
1	T	313	HIS
1	A	198	ASN
1	I	67	GLN
1	I	313	HIS

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Mol	Chain	Res	Type
1	K	69	LYS
1	L	270	VAL

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	345/369 (94%)	337 (98%)	8 (2%)	50	37
1	B	345/369 (94%)	341 (99%)	4 (1%)	71	65
1	C	343/369 (93%)	336 (98%)	7 (2%)	55	44
1	D	344/369 (93%)	340 (99%)	4 (1%)	71	65
1	E	345/369 (94%)	337 (98%)	8 (2%)	50	37
1	F	345/369 (94%)	341 (99%)	4 (1%)	71	65
1	G	346/369 (94%)	342 (99%)	4 (1%)	71	65
1	H	346/369 (94%)	337 (97%)	9 (3%)	46	32
1	I	346/369 (94%)	339 (98%)	7 (2%)	55	44
1	J	341/369 (92%)	335 (98%)	6 (2%)	59	48
1	K	345/369 (94%)	337 (98%)	8 (2%)	50	37
1	L	346/369 (94%)	342 (99%)	4 (1%)	71	65
1	M	345/369 (94%)	340 (99%)	5 (1%)	67	59
1	N	345/369 (94%)	338 (98%)	7 (2%)	55	44
1	O	347/369 (94%)	338 (97%)	9 (3%)	46	32
1	P	347/369 (94%)	342 (99%)	5 (1%)	67	59
1	Q	346/369 (94%)	341 (99%)	5 (1%)	67	59
1	R	346/369 (94%)	341 (99%)	5 (1%)	67	59
1	T	346/369 (94%)	339 (98%)	7 (2%)	55	44
1	V	345/369 (94%)	338 (98%)	7 (2%)	55	44
All	All	6904/7380 (94%)	6781 (98%)	123 (2%)	65	48

All (123) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	60	ARG
1	H	106	TRP
1	H	166	MET
1	H	175	HIS
1	H	214	LYS
1	H	345[A]	THR
1	H	345[B]	THR
1	H	411[A]	THR
1	H	411[B]	THR
1	A	67	GLN
1	A	75	ARG
1	A	106	TRP
1	A	230	SER
1	A	259	SER
1	A	365	LYS
1	A	411[A]	THR
1	A	411[B]	THR
1	B	75	ARG
1	B	342	LEU
1	B	411[A]	THR
1	B	411[B]	THR
1	C	75	ARG
1	C	175	HIS
1	C	259	SER
1	C	411[A]	THR
1	C	411[B]	THR
1	C	414[A]	MET
1	C	414[B]	MET
1	D	69	LYS
1	D	175	HIS
1	D	411[A]	THR
1	D	411[B]	THR
1	E	14	MET
1	E	75	ARG
1	E	222	ARG
1	E	235	LYS
1	E	354	ARG
1	E	411[A]	THR
1	E	411[B]	THR
1	E	420	LEU
1	F	175	HIS
1	F	251	ARG

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Mol	Chain	Res	Type
1	F	354	ARG
1	F	404	GLU
1	G	12	ASP
1	G	175	HIS
1	G	214	LYS
1	G	259	SER
1	I	43[A]	LYS
1	I	43[B]	LYS
1	I	106	TRP
1	I	175	HIS
1	I	342	LEU
1	I	411[A]	THR
1	I	411[B]	THR
1	J	175	HIS
1	J	259	SER
1	J	411[A]	THR
1	J	411[B]	THR
1	J	414[A]	MET
1	J	414[B]	MET
1	K	32	ARG
1	K	75	ARG
1	K	166	MET
1	K	175	HIS
1	K	235	LYS
1	K	259	SER
1	K	411[A]	THR
1	K	411[B]	THR
1	L	106	TRP
1	L	166	MET
1	L	175	HIS
1	L	417	LYS
1	M	175	HIS
1	M	188	LYS
1	M	411[A]	THR
1	M	411[B]	THR
1	M	417	LYS
1	N	75	ARG
1	N	175	HIS
1	N	179	LYS
1	N	259	SER
1	N	404	GLU
1	N	411[A]	THR

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Mol	Chain	Res	Type
1	N	411[B]	THR
1	O	75	ARG
1	O	106	TRP
1	O	175	HIS
1	O	244	HIS
1	O	259	SER
1	O	340	ARG
1	O	342	LEU
1	O	411[A]	THR
1	O	411[B]	THR
1	P	75	ARG
1	P	214	LYS
1	P	259	SER
1	P	411[A]	THR
1	P	411[B]	THR
1	Q	75	ARG
1	Q	259	SER
1	Q	404	GLU
1	Q	411[A]	THR
1	Q	411[B]	THR
1	R	175	HIS
1	R	179	LYS
1	R	259	SER
1	R	411[A]	THR
1	R	411[B]	THR
1	V	12	ASP
1	V	43[A]	LYS
1	V	43[B]	LYS
1	V	69	LYS
1	V	342	LEU
1	V	411[A]	THR
1	V	411[B]	THR
1	T	43[A]	LYS
1	T	43[B]	LYS
1	T	106	TRP
1	T	175	HIS
1	T	259	SER
1	T	411[A]	THR
1	T	411[B]	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (79) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	56	ASN
1	H	89	ASN
1	H	313	HIS
1	H	371	GLN
1	A	175	HIS
1	A	357	ASN
1	B	78	GLN
1	B	148	GLN
1	B	212	HIS
1	B	250	HIS
1	B	357	ASN
1	B	412	GLN
1	C	76	HIS
1	C	78	GLN
1	C	250	HIS
1	C	280	ASN
1	C	313	HIS
1	C	374	ASN
1	D	53	ASN
1	D	148	GLN
1	D	313	HIS
1	D	357	ASN
1	F	369	HIS
1	F	412	GLN
1	G	374	ASN
1	I	148	GLN
1	I	149	HIS
1	I	212	HIS
1	I	280	ASN
1	I	339	ASN
1	I	357	ASN
1	J	56	ASN
1	J	78	GLN
1	J	148	GLN
1	K	53	ASN
1	K	78	GLN
1	K	148	GLN
1	K	280	ASN
1	K	339	ASN
1	L	53	ASN
1	L	78	GLN
1	L	89	ASN
1	L	212	HIS

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Mol	Chain	Res	Type
1	L	313	HIS
1	M	78	GLN
1	M	148	GLN
1	M	357	ASN
1	N	148	GLN
1	N	244	HIS
1	N	357	ASN
1	O	357	ASN
1	P	53	ASN
1	P	280	ASN
1	P	371	GLN
1	Q	53	ASN
1	Q	56	ASN
1	Q	67	GLN
1	Q	313	HIS
1	Q	350	ASN
1	Q	371	GLN
1	R	149	HIS
1	R	339	ASN
1	R	357	ASN
1	R	369	HIS
1	R	374	ASN
1	V	53	ASN
1	V	147	GLN
1	V	198	ASN
1	V	244	HIS
1	V	250	HIS
1	V	350	ASN
1	V	357	ASN
1	V	371	GLN
1	T	89	ASN
1	T	212	HIS
1	T	313	HIS
1	T	339	ASN
1	T	350	ASN
1	T	371	GLN

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

42 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	$\# Z  > 2$
2	CIT	K	501	-	12,12,12	0.94	0	17,17,17	1.88	2 (11%)
2	CIT	T	501	-	12,12,12	1.01	0	17,17,17	1.93	5 (29%)
2	CIT	D	501	-	12,12,12	0.99	0	17,17,17	1.95	4 (23%)
3	GOL	P	502	-	5,5,5	0.38	0	5,5,5	0.30	0
3	GOL	R	502	-	5,5,5	0.32	0	5,5,5	0.87	0
2	CIT	O	501	-	12,12,12	1.03	0	17,17,17	1.90	5 (29%)
2	CIT	Q	501	-	12,12,12	0.99	0	17,17,17	1.92	6 (35%)
2	CIT	V	501	-	12,12,12	1.16	0	17,17,17	1.82	4 (23%)
3	GOL	Q	502	-	5,5,5	0.36	0	5,5,5	0.71	0
3	GOL	C	502	-	5,5,5	0.50	0	5,5,5	0.56	0
2	CIT	J	501	-	12,12,12	1.14	1 (8%)	17,17,17	2.08	5 (29%)
2	CIT	P	501	-	12,12,12	1.13	0	17,17,17	2.09	6 (35%)
2	CIT	C	501	-	12,12,12	1.20	1 (8%)	17,17,17	1.94	5 (29%)
3	GOL	J	502	-	5,5,5	0.34	0	5,5,5	0.28	0
3	GOL	O	502	-	5,5,5	0.40	0	5,5,5	0.62	0
3	GOL	H	502	-	5,5,5	0.41	0	5,5,5	0.26	0
2	CIT	L	501	-	12,12,12	1.12	0	17,17,17	2.01	5 (29%)
3	GOL	B	502	-	5,5,5	0.26	0	5,5,5	0.63	0
2	CIT	N	501	-	12,12,12	0.98	0	17,17,17	1.75	3 (17%)
3	GOL	V	502	-	5,5,5	0.22	0	5,5,5	0.85	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	GOL	F	502	-	5,5,5	0.37	0	5,5,5	0.64	0
3	GOL	K	503	-	5,5,5	0.45	0	5,5,5	0.48	0
2	CIT	B	501	-	12,12,12	1.14	1 (8%)	17,17,17	1.81	5 (29%)
2	CIT	H	501	-	12,12,12	0.98	0	17,17,17	1.89	7 (41%)
2	CIT	M	501	-	12,12,12	1.17	1 (8%)	17,17,17	1.94	5 (29%)
2	CIT	G	501	-	12,12,12	1.19	1 (8%)	17,17,17	1.73	4 (23%)
3	GOL	A	502	-	5,5,5	0.27	0	5,5,5	1.28	1 (20%)
3	GOL	M	502	-	5,5,5	0.36	0	5,5,5	0.55	0
3	GOL	E	502	-	5,5,5	0.42	0	5,5,5	0.34	0
3	GOL	G	502	-	5,5,5	0.58	0	5,5,5	0.87	0
3	GOL	J	503	-	5,5,5	0.51	0	5,5,5	1.70	2 (40%)
3	GOL	T	502	-	5,5,5	0.40	0	5,5,5	0.77	0
3	GOL	D	502	-	5,5,5	0.44	0	5,5,5	0.78	0
2	CIT	E	501	-	12,12,12	1.02	0	17,17,17	1.77	4 (23%)
3	GOL	K	502	-	5,5,5	0.39	0	5,5,5	0.98	0
2	CIT	A	501	-	12,12,12	1.00	0	17,17,17	1.97	5 (29%)
3	GOL	I	502	-	5,5,5	0.42	0	5,5,5	0.55	0
2	CIT	F	501	-	12,12,12	1.10	0	17,17,17	1.81	5 (29%)
2	CIT	R	501	-	12,12,12	1.14	0	17,17,17	1.97	7 (41%)
2	CIT	I	501	-	12,12,12	1.08	0	17,17,17	1.80	3 (17%)
3	GOL	N	502	-	5,5,5	0.19	0	5,5,5	1.02	0
3	GOL	L	502	-	5,5,5	0.46	0	5,5,5	0.47	0

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
2	CIT	K	501	-	-	5/16/16/16	-
2	CIT	T	501	-	-	5/16/16/16	-
2	CIT	D	501	-	-	5/16/16/16	-
3	GOL	P	502	-	-	0/4/4/4	-
3	GOL	R	502	-	-	2/4/4/4	-
2	CIT	O	501	-	-	5/16/16/16	-
2	CIT	Q	501	-	-	5/16/16/16	-
2	CIT	V	501	-	-	5/16/16/16	-
3	GOL	Q	502	-	-	1/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	C	502	-	-	0/4/4/4	-
2	CIT	J	501	-	-	5/16/16/16	-
2	CIT	P	501	-	-	5/16/16/16	-
2	CIT	C	501	-	-	8/16/16/16	-
3	GOL	J	502	-	-	0/4/4/4	-
3	GOL	O	502	-	-	0/4/4/4	-
3	GOL	H	502	-	-	0/4/4/4	-
2	CIT	L	501	-	-	5/16/16/16	-
3	GOL	B	502	-	-	0/4/4/4	-
2	CIT	N	501	-	-	5/16/16/16	-
3	GOL	V	502	-	-	0/4/4/4	-
3	GOL	F	502	-	-	0/4/4/4	-
3	GOL	K	503	-	-	2/4/4/4	-
2	CIT	B	501	-	-	6/16/16/16	-
2	CIT	H	501	-	-	5/16/16/16	-
2	CIT	M	501	-	-	6/16/16/16	-
2	CIT	G	501	-	-	5/16/16/16	-
3	GOL	A	502	-	-	4/4/4/4	-
3	GOL	M	502	-	-	0/4/4/4	-
3	GOL	E	502	-	-	3/4/4/4	-
3	GOL	G	502	-	-	3/4/4/4	-
3	GOL	J	503	-	-	3/4/4/4	-
3	GOL	T	502	-	-	2/4/4/4	-
3	GOL	D	502	-	-	1/4/4/4	-
2	CIT	E	501	-	-	5/16/16/16	-
3	GOL	K	502	-	-	2/4/4/4	-
2	CIT	A	501	-	-	5/16/16/16	-
3	GOL	I	502	-	-	0/4/4/4	-
2	CIT	F	501	-	-	5/16/16/16	-
2	CIT	R	501	-	-	5/16/16/16	-
2	CIT	I	501	-	-	5/16/16/16	-
3	GOL	N	502	-	-	2/4/4/4	-
3	GOL	L	502	-	-	0/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	M	501	CIT	C3-C6	-2.40	1.50	1.53
2	C	501	CIT	C3-C6	-2.38	1.50	1.53
2	J	501	CIT	C3-C6	-2.19	1.51	1.53
2	G	501	CIT	C3-C6	-2.12	1.51	1.53
2	B	501	CIT	C3-C6	-2.06	1.51	1.53

All (98) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	P	501	CIT	C3-C2-C1	-5.11	101.44	113.81
2	T	501	CIT	C3-C2-C1	-4.89	101.97	113.81
2	K	501	CIT	O6-C6-C3	4.57	120.99	113.05
2	N	501	CIT	O6-C6-C3	4.44	120.77	113.05
2	J	501	CIT	O6-C6-C3	4.39	120.67	113.05
2	R	501	CIT	O6-C6-C3	4.32	120.56	113.05
2	A	501	CIT	O6-C6-C3	4.31	120.53	113.05
2	C	501	CIT	C3-C2-C1	-4.28	103.45	113.81
2	E	501	CIT	O6-C6-C3	4.24	120.41	113.05
2	M	501	CIT	C3-C2-C1	-4.22	103.61	113.81
2	A	501	CIT	C3-C2-C1	-4.18	103.70	113.81
2	L	501	CIT	C3-C2-C1	-4.07	103.95	113.81
2	H	501	CIT	O6-C6-C3	4.07	120.11	113.05
2	K	501	CIT	C3-C2-C1	-3.98	104.19	113.81
2	J	501	CIT	C3-C2-C1	-3.95	104.25	113.81
2	O	501	CIT	O6-C6-C3	3.92	119.85	113.05
2	Q	501	CIT	O6-C6-C3	3.90	119.83	113.05
2	I	501	CIT	O6-C6-C3	3.86	119.75	113.05
2	D	501	CIT	C3-C2-C1	-3.80	104.62	113.81
2	F	501	CIT	O6-C6-C3	3.80	119.64	113.05
2	P	501	CIT	O6-C6-C3	3.79	119.64	113.05
2	B	501	CIT	O6-C6-C3	3.78	119.61	113.05
2	G	501	CIT	O7-C3-C6	-3.77	103.57	108.86
2	L	501	CIT	O6-C6-C3	3.70	119.48	113.05
2	F	501	CIT	C3-C2-C1	-3.64	105.01	113.81
2	V	501	CIT	O6-C6-C3	3.62	119.33	113.05
2	I	501	CIT	C3-C2-C1	-3.61	105.08	113.81
2	V	501	CIT	O7-C3-C6	-3.60	103.81	108.86
2	H	501	CIT	C3-C2-C1	-3.59	105.12	113.81
2	D	501	CIT	O6-C6-C3	3.57	119.25	113.05
2	B	501	CIT	O7-C3-C6	-3.52	103.92	108.86
2	M	501	CIT	O6-C6-C3	3.48	119.10	113.05
2	Q	501	CIT	C3-C2-C1	-3.39	105.61	113.81
2	V	501	CIT	C3-C2-C1	-3.36	105.67	113.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	O	501	CIT	C3-C2-C1	-3.35	105.70	113.81
2	R	501	CIT	O7-C3-C6	-3.34	104.17	108.86
2	T	501	CIT	O6-C6-C3	3.25	118.69	113.05
2	J	501	CIT	O5-C6-C3	-3.07	117.90	122.25
2	L	501	CIT	O2-C1-C2	3.06	124.17	114.35
2	N	501	CIT	C3-C2-C1	-3.02	106.50	113.81
2	C	501	CIT	O7-C3-C6	-2.98	104.69	108.86
2	G	501	CIT	O6-C6-C3	2.96	118.19	113.05
2	E	501	CIT	C3-C2-C1	-2.96	106.65	113.81
2	C	501	CIT	O6-C6-C3	2.92	118.11	113.05
2	C	501	CIT	O4-C5-C4	2.90	123.67	114.35
2	R	501	CIT	C3-C2-C1	-2.83	106.97	113.81
2	D	501	CIT	O4-C5-C4	2.77	123.25	114.35
2	B	501	CIT	C3-C2-C1	-2.73	107.19	113.81
2	J	501	CIT	O4-C5-O3	-2.70	116.58	123.30
2	J	501	CIT	O4-C5-C4	2.67	122.94	114.35
2	O	501	CIT	O2-C1-C2	2.63	122.78	114.35
2	E	501	CIT	O4-C5-O3	-2.61	116.80	123.30
2	Q	501	CIT	O4-C5-C4	2.60	122.70	114.35
2	D	501	CIT	O2-C1-C2	2.58	122.63	114.35
2	G	501	CIT	C3-C2-C1	-2.55	107.65	113.81
3	A	502	GOL	C3-C2-C1	-2.51	101.93	111.70
2	T	501	CIT	O2-C1-C2	2.50	122.39	114.35
2	Q	501	CIT	O2-C1-C2	2.47	122.28	114.35
3	J	503	GOL	O2-C2-C1	2.46	119.96	109.12
2	P	501	CIT	O7-C3-C6	-2.45	105.42	108.86
2	R	501	CIT	O4-C5-O3	-2.44	117.22	123.30
2	M	501	CIT	O2-C1-C2	2.44	122.18	114.35
2	F	501	CIT	O4-C5-C4	2.38	122.00	114.35
2	O	501	CIT	O4-C5-C4	2.38	122.00	114.35
3	J	503	GOL	O1-C1-C2	2.37	121.55	110.20
2	L	501	CIT	O2-C1-O1	-2.35	117.45	123.30
2	E	501	CIT	O4-C5-C4	2.35	121.89	114.35
2	R	501	CIT	O5-C6-C3	-2.34	118.94	122.25
2	H	501	CIT	O2-C1-C2	2.31	121.76	114.35
2	A	501	CIT	O4-C5-C4	2.31	121.75	114.35
2	Q	501	CIT	O4-C5-O3	-2.30	117.57	123.30
2	F	501	CIT	O4-C5-O3	-2.28	117.61	123.30
2	M	501	CIT	O4-C5-C4	2.28	121.67	114.35
2	H	501	CIT	O4-C5-C4	2.27	121.64	114.35
2	M	501	CIT	O5-C6-C3	-2.27	119.04	122.25
2	V	501	CIT	O4-C5-O3	-2.26	117.66	123.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	501	CIT	O4-C5-O3	-2.26	117.68	123.30
2	N	501	CIT	O4-C5-O3	-2.22	117.77	123.30
2	F	501	CIT	O2-C1-C2	2.21	121.44	114.35
2	I	501	CIT	O5-C6-C3	-2.20	119.14	122.25
2	H	501	CIT	O2-C1-O1	-2.19	117.83	123.30
2	B	501	CIT	O4-C5-O3	-2.19	117.84	123.30
2	G	501	CIT	O2-C1-O1	-2.17	117.89	123.30
2	B	501	CIT	O5-C6-C3	-2.16	119.20	122.25
2	P	501	CIT	C4-C3-C6	2.14	114.70	110.11
2	T	501	CIT	C3-C4-C5	-2.10	108.72	113.81
2	C	501	CIT	O4-C5-O3	-2.10	118.06	123.30
2	O	501	CIT	O4-C5-O3	-2.09	118.09	123.30
2	L	501	CIT	O7-C3-C6	-2.09	105.93	108.86
2	P	501	CIT	O4-C5-C4	2.08	121.03	114.35
2	H	501	CIT	C3-C4-C5	-2.08	108.78	113.81
2	A	501	CIT	O4-C5-O3	-2.08	118.12	123.30
2	A	501	CIT	O2-C1-C2	2.03	120.88	114.35
2	Q	501	CIT	O5-C6-C3	-2.03	119.37	122.25
2	R	501	CIT	O4-C5-C4	2.03	120.86	114.35
2	H	501	CIT	O4-C5-O3	-2.02	118.27	123.30
2	P	501	CIT	O4-C5-O3	-2.01	118.28	123.30
2	R	501	CIT	O2-C1-O1	-2.01	118.30	123.30

There are no chirality outliers.

All (130) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	H	501	CIT	C2-C3-C6-O5
2	H	501	CIT	C2-C3-C6-O6
2	H	501	CIT	O7-C3-C6-O5
2	H	501	CIT	O7-C3-C6-O6
2	A	501	CIT	C2-C3-C6-O5
2	A	501	CIT	C2-C3-C6-O6
2	A	501	CIT	O7-C3-C6-O5
2	A	501	CIT	O7-C3-C6-O6
2	B	501	CIT	C2-C3-C6-O5
2	B	501	CIT	C2-C3-C6-O6
2	B	501	CIT	O7-C3-C6-O5
2	B	501	CIT	O7-C3-C6-O6
2	C	501	CIT	C2-C3-C6-O5
2	C	501	CIT	C2-C3-C6-O6
2	C	501	CIT	O7-C3-C6-O5

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Mol	Chain	Res	Type	Atoms
2	C	501	CIT	O7-C3-C6-O6
2	D	501	CIT	C2-C3-C6-O5
2	D	501	CIT	C2-C3-C6-O6
2	D	501	CIT	O7-C3-C6-O5
2	D	501	CIT	O7-C3-C6-O6
2	E	501	CIT	C2-C3-C6-O5
2	E	501	CIT	C2-C3-C6-O6
2	E	501	CIT	O7-C3-C6-O5
2	E	501	CIT	O7-C3-C6-O6
2	F	501	CIT	C2-C3-C6-O5
2	F	501	CIT	C2-C3-C6-O6
2	F	501	CIT	O7-C3-C6-O5
2	F	501	CIT	O7-C3-C6-O6
2	G	501	CIT	C2-C3-C6-O5
2	G	501	CIT	C2-C3-C6-O6
2	G	501	CIT	O7-C3-C6-O5
2	G	501	CIT	O7-C3-C6-O6
2	I	501	CIT	C2-C3-C6-O5
2	I	501	CIT	C2-C3-C6-O6
2	I	501	CIT	O7-C3-C6-O5
2	I	501	CIT	O7-C3-C6-O6
2	J	501	CIT	C2-C3-C6-O5
2	J	501	CIT	C2-C3-C6-O6
2	J	501	CIT	O7-C3-C6-O5
2	J	501	CIT	O7-C3-C6-O6
2	K	501	CIT	C2-C3-C6-O5
2	K	501	CIT	C2-C3-C6-O6
2	K	501	CIT	O7-C3-C6-O5
2	K	501	CIT	O7-C3-C6-O6
2	L	501	CIT	C2-C3-C6-O5
2	L	501	CIT	C2-C3-C6-O6
2	L	501	CIT	O7-C3-C6-O5
2	L	501	CIT	O7-C3-C6-O6
2	M	501	CIT	C2-C3-C6-O5
2	M	501	CIT	C2-C3-C6-O6
2	M	501	CIT	O7-C3-C6-O5
2	M	501	CIT	O7-C3-C6-O6
2	N	501	CIT	C2-C3-C6-O5
2	N	501	CIT	C2-C3-C6-O6
2	N	501	CIT	O7-C3-C6-O5
2	N	501	CIT	O7-C3-C6-O6
2	O	501	CIT	C2-C3-C6-O5

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Mol	Chain	Res	Type	Atoms
2	O	501	CIT	C2-C3-C6-O6
2	O	501	CIT	O7-C3-C6-O5
2	O	501	CIT	O7-C3-C6-O6
2	P	501	CIT	C2-C3-C6-O5
2	P	501	CIT	C2-C3-C6-O6
2	P	501	CIT	O7-C3-C6-O5
2	P	501	CIT	O7-C3-C6-O6
2	Q	501	CIT	C2-C3-C6-O5
2	Q	501	CIT	C2-C3-C6-O6
2	Q	501	CIT	O7-C3-C6-O5
2	Q	501	CIT	O7-C3-C6-O6
2	R	501	CIT	C2-C3-C6-O5
2	R	501	CIT	C2-C3-C6-O6
2	R	501	CIT	O7-C3-C6-O5
2	R	501	CIT	O7-C3-C6-O6
2	V	501	CIT	C2-C3-C6-O5
2	V	501	CIT	C2-C3-C6-O6
2	V	501	CIT	O7-C3-C6-O5
2	V	501	CIT	O7-C3-C6-O6
2	T	501	CIT	C2-C3-C6-O5
2	T	501	CIT	C2-C3-C6-O6
2	T	501	CIT	O7-C3-C6-O5
2	T	501	CIT	O7-C3-C6-O6
3	G	502	GOL	O1-C1-C2-C3
3	G	502	GOL	C1-C2-C3-O3
3	J	503	GOL	O1-C1-C2-C3
3	K	503	GOL	C1-C2-C3-O3
3	N	502	GOL	C1-C2-C3-O3
3	T	502	GOL	O1-C1-C2-C3
3	N	502	GOL	O2-C2-C3-O3
3	A	502	GOL	O1-C1-C2-C3
3	A	502	GOL	C1-C2-C3-O3
3	E	502	GOL	C1-C2-C3-O3
3	A	502	GOL	O2-C2-C3-O3
3	E	502	GOL	O2-C2-C3-O3
3	G	502	GOL	O1-C1-C2-O2
3	J	503	GOL	O1-C1-C2-O2
3	K	503	GOL	O2-C2-C3-O3
3	R	502	GOL	O1-C1-C2-C3
3	J	503	GOL	O2-C2-C3-O3
3	K	502	GOL	O1-C1-C2-O2
3	T	502	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	D	501	CIT	O7-C3-C4-C5
2	I	501	CIT	O7-C3-C4-C5
2	O	501	CIT	O7-C3-C4-C5
2	R	501	CIT	O7-C3-C4-C5
2	V	501	CIT	O7-C3-C4-C5
2	H	501	CIT	O7-C3-C4-C5
2	A	501	CIT	O7-C3-C4-C5
2	G	501	CIT	O7-C3-C4-C5
2	J	501	CIT	O7-C3-C4-C5
2	K	501	CIT	O7-C3-C4-C5
2	N	501	CIT	O7-C3-C4-C5
2	T	501	CIT	O7-C3-C4-C5
2	M	501	CIT	O7-C3-C4-C5
2	P	501	CIT	O7-C3-C4-C5
3	A	502	GOL	O1-C1-C2-O2
3	R	502	GOL	O1-C1-C2-O2
2	B	501	CIT	O7-C3-C4-C5
2	C	501	CIT	O7-C3-C4-C5
2	E	501	CIT	O7-C3-C4-C5
2	F	501	CIT	O7-C3-C4-C5
2	L	501	CIT	O7-C3-C4-C5
2	Q	501	CIT	O7-C3-C4-C5
2	C	501	CIT	C4-C3-C6-O6
2	M	501	CIT	C2-C3-C4-C5
3	K	502	GOL	O1-C1-C2-C3
2	C	501	CIT	C3-C4-C5-O4
2	C	501	CIT	C3-C4-C5-O3
3	D	502	GOL	O1-C1-C2-O2
3	E	502	GOL	O1-C1-C2-C3
2	B	501	CIT	C2-C3-C4-C5
3	Q	502	GOL	O1-C1-C2-O2

There are no ring outliers.

23 monomers are involved in 52 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	K	501	CIT	1	0
2	T	501	CIT	2	0
2	D	501	CIT	2	0
2	O	501	CIT	2	0
2	Q	501	CIT	2	0
2	V	501	CIT	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	J	501	CIT	4	0
2	P	501	CIT	1	0
2	C	501	CIT	2	0
2	L	501	CIT	4	0
2	N	501	CIT	1	0
3	K	503	GOL	5	0
2	B	501	CIT	2	0
2	H	501	CIT	1	0
2	M	501	CIT	2	0
2	G	501	CIT	2	0
3	G	502	GOL	1	0
3	J	503	GOL	4	0
2	E	501	CIT	2	0
2	A	501	CIT	3	0
2	F	501	CIT	3	0
2	R	501	CIT	3	0
2	I	501	CIT	1	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	415/440 (94%)	-0.29	13 (3%)	49	43	12, 28, 57, 181	0
1	B	415/440 (94%)	-0.17	12 (2%)	51	46	14, 29, 59, 177	0
1	C	411/440 (93%)	-0.41	10 (2%)	59	54	12, 24, 51, 142	0
1	D	415/440 (94%)	-0.19	17 (4%)	37	31	14, 30, 60, 165	0
1	E	415/440 (94%)	-0.30	11 (2%)	54	49	13, 27, 55, 168	0
1	F	415/440 (94%)	-0.33	14 (3%)	45	39	12, 27, 55, 136	0
1	G	415/440 (94%)	-0.20	17 (4%)	37	31	13, 28, 57, 166	0
1	H	415/440 (94%)	-0.28	12 (2%)	51	46	12, 27, 55, 172	0
1	I	415/440 (94%)	-0.24	12 (2%)	51	46	13, 28, 56, 159	0
1	J	408/440 (92%)	-0.49	4 (0%)	82	80	12, 24, 46, 150	0
1	K	415/440 (94%)	-0.28	14 (3%)	45	39	12, 27, 55, 171	0
1	L	415/440 (94%)	-0.24	17 (4%)	37	31	14, 29, 56, 159	0
1	M	415/440 (94%)	-0.21	13 (3%)	49	43	15, 30, 59, 184	0
1	N	415/440 (94%)	-0.27	13 (3%)	49	43	13, 27, 59, 152	0
1	O	415/440 (94%)	-0.25	16 (3%)	39	33	12, 28, 58, 167	0
1	P	415/440 (94%)	-0.31	13 (3%)	49	43	13, 26, 55, 163	0
1	Q	415/440 (94%)	-0.38	11 (2%)	54	49	12, 23, 48, 149	0
1	R	415/440 (94%)	-0.36	12 (2%)	51	46	11, 23, 48, 140	0
1	T	415/440 (94%)	-0.32	13 (3%)	49	43	12, 26, 55, 172	0
1	V	415/440 (94%)	-0.24	14 (3%)	45	39	11, 29, 61, 154	0
All	All	8289/8800 (94%)	-0.29	258 (3%)	49	43	11, 27, 56, 184	0

All (258) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	420	LEU	29.2
1	N	420	LEU	20.9
1	P	101	VAL	19.0
1	M	420	LEU	18.7
1	G	420	LEU	17.8
1	V	420	LEU	17.6
1	D	420	LEU	16.9
1	H	99	GLN	16.7
1	E	101	VAL	16.6
1	O	420	LEU	15.8
1	F	420	LEU	15.5
1	H	101	VAL	14.9
1	O	101	VAL	14.9
1	J	419	ASP	14.4
1	T	420	LEU	14.3
1	R	420	LEU	14.2
1	P	420	LEU	13.8
1	Q	420	LEU	13.7
1	Q	101	VAL	12.9
1	A	420	LEU	12.5
1	I	420	LEU	12.4
1	K	99	GLN	12.2
1	H	420	LEU	12.1
1	V	98	SER	12.0
1	D	101	VAL	12.0
1	J	420	LEU	11.5
1	L	98	SER	11.5
1	R	101	VAL	11.4
1	M	102	SER	11.3
1	E	420	LEU	11.2
1	T	101	VAL	11.1
1	L	102	SER	11.1
1	I	100	ARG	11.0
1	M	101	VAL	10.9
1	B	102	SER	10.8
1	I	419	ASP	10.7
1	A	101	VAL	10.7
1	H	102	SER	10.6
1	G	101	VAL	10.0
1	K	100	ARG	9.9
1	N	101	VAL	9.6
1	K	420	LEU	9.5
1	L	99	GLN	9.3

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Mol	Chain	Res	Type	RSRZ
1	G	98	SER	9.2
1	L	101	VAL	9.2
1	V	101	VAL	9.2
1	T	102	SER	9.1
1	C	420	LEU	9.1
1	D	419	ASP	9.0
1	I	101	VAL	9.0
1	F	101	VAL	8.9
1	N	419	ASP	8.8
1	K	102	SER	8.7
1	L	420	LEU	8.7
1	P	419	ASP	8.7
1	E	419	ASP	8.6
1	T	419	ASP	8.6
1	B	101	VAL	8.5
1	R	419	ASP	8.5
1	N	100	ARG	8.3
1	F	102	SER	8.2
1	O	98	SER	8.2
1	G	419	ASP	8.1
1	D	106	TRP	8.1
1	R	98	SER	8.0
1	A	99	GLN	8.0
1	D	102	SER	7.9
1	P	99	GLN	7.9
1	I	102	SER	7.8
1	O	99	GLN	7.8
1	H	419	ASP	7.8
1	C	419	ASP	7.8
1	P	100	ARG	7.7
1	Q	99	GLN	7.6
1	A	106	TRP	7.6
1	V	419	ASP	7.5
1	I	104	THR	7.3
1	N	99	GLN	7.3
1	T	106	TRP	7.3
1	D	99	GLN	7.2
1	D	104	THR	7.1
1	K	419	ASP	6.9
1	D	100	ARG	6.9
1	O	102	SER	6.8
1	R	99	GLN	6.7

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Mol	Chain	Res	Type	RSRZ
1	K	98	SER	6.6
1	R	104	THR	6.6
1	Q	419	ASP	6.6
1	L	100	ARG	6.6
1	K	101	VAL	6.6
1	G	102	SER	6.5
1	R	102	SER	6.5
1	B	419	ASP	6.5
1	T	99	GLN	6.4
1	A	102	SER	6.3
1	A	103	GLU	6.2
1	K	106	TRP	6.2
1	R	106	TRP	6.1
1	P	106	TRP	6.1
1	Q	102	SER	6.1
1	V	106	TRP	5.9
1	G	100	ARG	5.9
1	A	100	ARG	5.9
1	D	103	GLU	5.9
1	Q	100	ARG	5.9
1	Q	98	SER	5.8
1	M	99	GLN	5.8
1	B	98	SER	5.8
1	D	98	SER	5.8
1	V	100	ARG	5.7
1	M	419	ASP	5.7
1	N	102	SER	5.6
1	O	100	ARG	5.6
1	B	104	THR	5.6
1	G	99	GLN	5.5
1	B	6	VAL	5.4
1	K	104	THR	5.4
1	F	419	ASP	5.3
1	H	105	GLN	5.3
1	F	99	GLN	5.3
1	Q	106	TRP	5.3
1	H	104	THR	5.3
1	V	102	SER	5.3
1	R	100	ARG	5.2
1	E	98	SER	5.2
1	F	100	ARG	5.2
1	O	106	TRP	5.1

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Mol	Chain	Res	Type	RSRZ
1	M	100	ARG	5.1
1	M	104	THR	5.1
1	A	419	ASP	5.1
1	H	106	TRP	5.1
1	D	105	GLN	5.1
1	M	98	SER	5.1
1	T	104	THR	5.0
1	I	99	GLN	5.0
1	I	98	SER	5.0
1	R	105	GLN	4.9
1	L	6	VAL	4.8
1	E	104	THR	4.8
1	G	104	THR	4.8
1	O	419	ASP	4.8
1	B	99	GLN	4.7
1	T	68	TYR	4.7
1	B	100	ARG	4.7
1	H	100	ARG	4.7
1	A	98	SER	4.7
1	T	103	GLU	4.6
1	O	103	GLU	4.5
1	G	6	VAL	4.5
1	Q	103	GLU	4.4
1	E	100	ARG	4.4
1	P	104	THR	4.3
1	D	228	THR	4.3
1	H	98	SER	4.2
1	L	243	GLU	4.2
1	M	6	VAL	4.2
1	P	102	SER	4.2
1	V	99	GLN	4.2
1	C	104	THR	4.1
1	O	6	VAL	4.1
1	L	419	ASP	4.1
1	L	103	GLU	4.0
1	I	106	TRP	4.0
1	L	105	GLN	4.0
1	B	106	TRP	4.0
1	P	68	TYR	4.0
1	I	228	THR	4.0
1	V	104	THR	4.0
1	F	105	GLN	4.0

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Mol	Chain	Res	Type	RSRZ
1	L	68	TYR	3.9
1	F	104	THR	3.9
1	A	68	TYR	3.9
1	V	6	VAL	3.8
1	M	103	GLU	3.8
1	C	106	TRP	3.8
1	F	68	TYR	3.8
1	N	106	TRP	3.8
1	E	106	TRP	3.7
1	C	98	SER	3.7
1	G	68	TYR	3.6
1	L	104	THR	3.6
1	Q	104	THR	3.6
1	P	103	GLU	3.6
1	T	100	ARG	3.5
1	G	106	TRP	3.5
1	O	104	THR	3.5
1	A	105	GLN	3.5
1	M	106	TRP	3.5
1	A	104	THR	3.5
1	G	105	GLN	3.5
1	C	68	TYR	3.4
1	T	6	VAL	3.4
1	T	98	SER	3.4
1	E	105	GLN	3.3
1	J	106	TRP	3.3
1	L	106	TRP	3.3
1	I	105	GLN	3.3
1	D	68	TYR	3.3
1	G	103	GLU	3.2
1	C	103	GLU	3.2
1	N	103	GLU	3.2
1	E	102	SER	3.2
1	N	98	SER	3.2
1	H	103	GLU	3.1
1	N	68	TYR	3.1
1	L	197	GLY	3.1
1	K	68	TYR	3.1
1	O	105	GLN	3.0
1	O	68	TYR	3.0
1	E	68	TYR	3.0
1	C	228	THR	3.0

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Mol	Chain	Res	Type	RSRZ
1	O	242	VAL	3.0
1	K	103	GLU	2.9
1	D	6	VAL	2.9
1	H	68	TYR	2.8
1	K	105	GLN	2.8
1	J	68	TYR	2.7
1	G	228	THR	2.7
1	O	183	ILE	2.7
1	N	6	VAL	2.7
1	V	105	GLN	2.6
1	V	183	ILE	2.6
1	F	98	SER	2.6
1	M	68	TYR	2.6
1	R	103	GLU	2.5
1	A	6	VAL	2.5
1	N	104	THR	2.5
1	F	106	TRP	2.5
1	M	105	GLN	2.4
1	L	97	ALA	2.4
1	F	418	LEU	2.4
1	F	222	ARG	2.4
1	K	6	VAL	2.4
1	D	32	ARG	2.3
1	T	228	THR	2.3
1	N	105	GLN	2.3
1	D	69	LYS	2.3
1	B	103	GLU	2.2
1	R	6	VAL	2.2
1	O	240	TYR	2.2
1	C	105	GLN	2.2
1	G	240	TYR	2.2
1	L	240	TYR	2.2
1	C	6	VAL	2.2
1	P	6	VAL	2.2
1	K	228	THR	2.1
1	G	178	GLU	2.1
1	E	103	GLU	2.1
1	Q	6	VAL	2.1
1	V	228	THR	2.1
1	D	225	ALA	2.1
1	V	32	ARG	2.1
1	I	6	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	G	243	GLU	2.1
1	P	98	SER	2.1
1	B	105	GLN	2.0
1	F	6	VAL	2.0
1	P	105	GLN	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 monosaccharides 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. 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	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	GOL	J	503	6/6	0.79	0.19	37,41,49,51	0
2	CIT	B	501	13/13	0.81	0.19	31,60,71,72	0
2	CIT	F	501	13/13	0.83	0.15	44,72,79,80	0
2	CIT	V	501	13/13	0.84	0.18	29,49,66,67	0
2	CIT	N	501	13/13	0.84	0.16	41,69,76,76	0
2	CIT	J	501	13/13	0.85	0.16	23,42,51,53	0
2	CIT	L	501	13/13	0.87	0.20	26,57,70,71	0
2	CIT	O	501	13/13	0.88	0.14	27,59,68,69	0
2	CIT	Q	501	13/13	0.88	0.18	29,57,67,67	0
2	CIT	R	501	13/13	0.88	0.16	23,50,63,63	0
2	CIT	G	501	13/13	0.88	0.15	27,59,70,70	0
2	CIT	D	501	13/13	0.88	0.14	27,58,70,71	0
2	CIT	A	501	13/13	0.89	0.17	35,66,75,76	0
2	CIT	P	501	13/13	0.90	0.14	26,56,65,65	0
2	CIT	K	501	13/13	0.90	0.13	27,57,70,71	0
2	CIT	E	501	13/13	0.90	0.13	34,62,73,73	0
2	CIT	I	501	13/13	0.90	0.14	25,55,69,70	0
2	CIT	H	501	13/13	0.90	0.18	27,55,71,71	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	GOL	K	503	6/6	0.90	0.48	81,85,86,87	0
2	CIT	C	501	13/13	0.91	0.14	22,42,56,56	0
2	CIT	T	501	13/13	0.91	0.16	25,55,63,64	0
3	GOL	G	502	6/6	0.92	0.12	23,32,36,43	0
3	GOL	L	502	6/6	0.92	0.13	30,36,37,41	0
3	GOL	E	502	6/6	0.93	0.09	26,29,31,34	0
2	CIT	M	501	13/13	0.93	0.15	27,57,70,71	0
3	GOL	I	502	6/6	0.93	0.10	17,23,25,26	0
3	GOL	B	502	6/6	0.94	0.09	27,34,35,35	0
3	GOL	K	502	6/6	0.94	0.12	21,27,39,47	0
3	GOL	M	502	6/6	0.95	0.07	27,32,35,40	0
3	GOL	N	502	6/6	0.95	0.08	21,27,30,33	0
3	GOL	C	502	6/6	0.96	0.10	24,29,31,33	0
3	GOL	H	502	6/6	0.96	0.08	22,26,28,29	0
3	GOL	R	502	6/6	0.96	0.07	22,27,30,31	0
3	GOL	T	502	6/6	0.96	0.07	25,32,33,34	0
3	GOL	J	502	6/6	0.97	0.07	25,29,33,35	0
3	GOL	O	502	6/6	0.97	0.08	20,28,32,33	0
3	GOL	P	502	6/6	0.97	0.11	22,28,34,34	0
3	GOL	A	502	6/6	0.97	0.07	18,24,29,31	0
3	GOL	V	502	6/6	0.97	0.07	24,29,29,29	0
3	GOL	F	502	6/6	0.97	0.07	25,26,28,29	0
3	GOL	Q	502	6/6	0.98	0.07	18,26,29,29	0
3	GOL	D	502	6/6	0.98	0.06	17,23,27,29	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.