



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 09:26 PM UTC

PDB ID : 3ANZ / pdb_00003anz
Title : Crystal Structure of alpha-hemolysin
Authors : Yamashita, K.; Kawauchi, H.; Tanaka, Y.; Yao, M.; Tanaka, I.
Deposited on : 2010-09-16
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

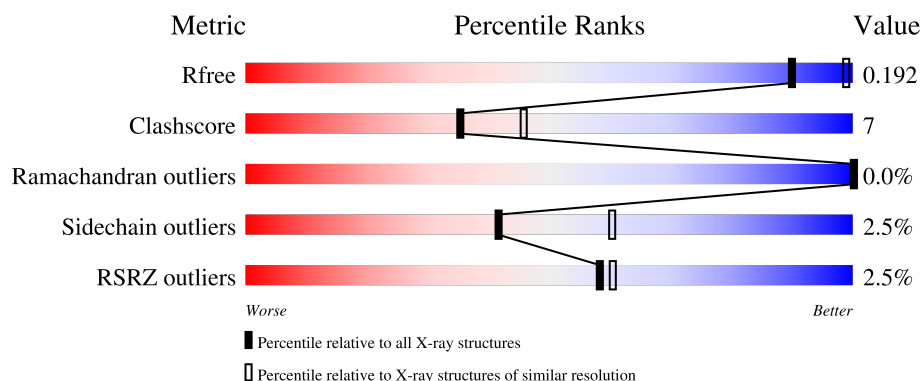
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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

Mol	Chain	Length	Quality of chain
1	A	302	0% 85% 12% ..
1	B	302	3% 83% 14% .
1	C	302	2% 81% 16% ..
1	D	302	2% 82% 14% ..
1	E	302	8% 79% 17% ..

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Mol	Chain	Length	Quality of chain
1	F	302	
1	G	302	
1	H	302	
1	I	302	
1	J	302	
1	K	302	
1	L	302	
1	M	302	
1	N	302	
1	O	302	
1	P	302	
1	Q	302	
1	R	302	
1	S	302	
1	T	302	
1	U	302	
1	V	302	
1	W	302	
1	X	302	
1	Y	302	
1	Z	302	
1	a	302	
1	b	302	

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
2	MPD	E	302	-	-	X	-
2	MPD	F	303	-	-	X	-
2	MPD	K	303	-	-	X	-
2	MPD	N	302	-	-	X	-
2	MPD	Q	302	-	-	X	-
2	MPD	R	303	-	-	X	-
2	MPD	V	303	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 68607 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 Alpha-hemolysin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	295	Total	C	N	O	S	0	0	0
			2362	1484	403	468	7			
1	B	295	Total	C	N	O	S	0	0	0
			2362	1484	403	468	7			
1	C	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	D	295	Total	C	N	O	S	0	0	0
			2362	1484	403	468	7			
1	E	295	Total	C	N	O	S	0	0	0
			2362	1484	403	468	7			
1	F	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	G	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	H	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	I	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	J	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	K	295	Total	C	N	O	S	0	0	0
			2362	1484	403	468	7			
1	L	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	M	295	Total	C	N	O	S	0	0	0
			2362	1484	403	468	7			
1	N	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	O	295	Total	C	N	O	S	0	0	0
			2362	1484	403	468	7			
1	P	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	R	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	S	295	Total	C	N	O	S	0	0	0
			2362	1484	403	468	7			
1	T	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	U	295	Total	C	N	O	S	0	0	0
			2362	1484	403	468	7			
1	V	295	Total	C	N	O	S	0	0	0
			2362	1484	403	468	7			
1	W	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	X	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	Y	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	Z	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	a	294	Total	C	N	O	S	0	0	0
			2353	1479	402	465	7			
1	b	295	Total	C	N	O	S	0	0	0
			2362	1484	403	468	7			

There are 252 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	expression tag	UNP Q99UU6
A	294	LEU	-	expression tag	UNP Q99UU6
A	295	GLU	-	expression tag	UNP Q99UU6
A	296	HIS	-	expression tag	UNP Q99UU6
A	297	HIS	-	expression tag	UNP Q99UU6
A	298	HIS	-	expression tag	UNP Q99UU6
A	299	HIS	-	expression tag	UNP Q99UU6
A	300	HIS	-	expression tag	UNP Q99UU6
A	301	HIS	-	expression tag	UNP Q99UU6
B	0	MET	-	expression tag	UNP Q99UU6
B	294	LEU	-	expression tag	UNP Q99UU6
B	295	GLU	-	expression tag	UNP Q99UU6
B	296	HIS	-	expression tag	UNP Q99UU6
B	297	HIS	-	expression tag	UNP Q99UU6
B	298	HIS	-	expression tag	UNP Q99UU6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	299	HIS	-	expression tag	UNP Q99UU6
B	300	HIS	-	expression tag	UNP Q99UU6
B	301	HIS	-	expression tag	UNP Q99UU6
C	0	MET	-	expression tag	UNP Q99UU6
C	294	LEU	-	expression tag	UNP Q99UU6
C	295	GLU	-	expression tag	UNP Q99UU6
C	296	HIS	-	expression tag	UNP Q99UU6
C	297	HIS	-	expression tag	UNP Q99UU6
C	298	HIS	-	expression tag	UNP Q99UU6
C	299	HIS	-	expression tag	UNP Q99UU6
C	300	HIS	-	expression tag	UNP Q99UU6
C	301	HIS	-	expression tag	UNP Q99UU6
D	0	MET	-	expression tag	UNP Q99UU6
D	294	LEU	-	expression tag	UNP Q99UU6
D	295	GLU	-	expression tag	UNP Q99UU6
D	296	HIS	-	expression tag	UNP Q99UU6
D	297	HIS	-	expression tag	UNP Q99UU6
D	298	HIS	-	expression tag	UNP Q99UU6
D	299	HIS	-	expression tag	UNP Q99UU6
D	300	HIS	-	expression tag	UNP Q99UU6
D	301	HIS	-	expression tag	UNP Q99UU6
E	0	MET	-	expression tag	UNP Q99UU6
E	294	LEU	-	expression tag	UNP Q99UU6
E	295	GLU	-	expression tag	UNP Q99UU6
E	296	HIS	-	expression tag	UNP Q99UU6
E	297	HIS	-	expression tag	UNP Q99UU6
E	298	HIS	-	expression tag	UNP Q99UU6
E	299	HIS	-	expression tag	UNP Q99UU6
E	300	HIS	-	expression tag	UNP Q99UU6
E	301	HIS	-	expression tag	UNP Q99UU6
F	0	MET	-	expression tag	UNP Q99UU6
F	294	LEU	-	expression tag	UNP Q99UU6
F	295	GLU	-	expression tag	UNP Q99UU6
F	296	HIS	-	expression tag	UNP Q99UU6
F	297	HIS	-	expression tag	UNP Q99UU6
F	298	HIS	-	expression tag	UNP Q99UU6
F	299	HIS	-	expression tag	UNP Q99UU6
F	300	HIS	-	expression tag	UNP Q99UU6
F	301	HIS	-	expression tag	UNP Q99UU6
G	0	MET	-	expression tag	UNP Q99UU6
G	294	LEU	-	expression tag	UNP Q99UU6
G	295	GLU	-	expression tag	UNP Q99UU6

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Chain	Residue	Modelled	Actual	Comment	Reference
G	296	HIS	-	expression tag	UNP Q99UU6
G	297	HIS	-	expression tag	UNP Q99UU6
G	298	HIS	-	expression tag	UNP Q99UU6
G	299	HIS	-	expression tag	UNP Q99UU6
G	300	HIS	-	expression tag	UNP Q99UU6
G	301	HIS	-	expression tag	UNP Q99UU6
H	0	MET	-	expression tag	UNP Q99UU6
H	294	LEU	-	expression tag	UNP Q99UU6
H	295	GLU	-	expression tag	UNP Q99UU6
H	296	HIS	-	expression tag	UNP Q99UU6
H	297	HIS	-	expression tag	UNP Q99UU6
H	298	HIS	-	expression tag	UNP Q99UU6
H	299	HIS	-	expression tag	UNP Q99UU6
H	300	HIS	-	expression tag	UNP Q99UU6
H	301	HIS	-	expression tag	UNP Q99UU6
I	0	MET	-	expression tag	UNP Q99UU6
I	294	LEU	-	expression tag	UNP Q99UU6
I	295	GLU	-	expression tag	UNP Q99UU6
I	296	HIS	-	expression tag	UNP Q99UU6
I	297	HIS	-	expression tag	UNP Q99UU6
I	298	HIS	-	expression tag	UNP Q99UU6
I	299	HIS	-	expression tag	UNP Q99UU6
I	300	HIS	-	expression tag	UNP Q99UU6
I	301	HIS	-	expression tag	UNP Q99UU6
J	0	MET	-	expression tag	UNP Q99UU6
J	294	LEU	-	expression tag	UNP Q99UU6
J	295	GLU	-	expression tag	UNP Q99UU6
J	296	HIS	-	expression tag	UNP Q99UU6
J	297	HIS	-	expression tag	UNP Q99UU6
J	298	HIS	-	expression tag	UNP Q99UU6
J	299	HIS	-	expression tag	UNP Q99UU6
J	300	HIS	-	expression tag	UNP Q99UU6
J	301	HIS	-	expression tag	UNP Q99UU6
K	0	MET	-	expression tag	UNP Q99UU6
K	294	LEU	-	expression tag	UNP Q99UU6
K	295	GLU	-	expression tag	UNP Q99UU6
K	296	HIS	-	expression tag	UNP Q99UU6
K	297	HIS	-	expression tag	UNP Q99UU6
K	298	HIS	-	expression tag	UNP Q99UU6
K	299	HIS	-	expression tag	UNP Q99UU6
K	300	HIS	-	expression tag	UNP Q99UU6
K	301	HIS	-	expression tag	UNP Q99UU6

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Chain	Residue	Modelled	Actual	Comment	Reference
L	0	MET	-	expression tag	UNP Q99UU6
L	294	LEU	-	expression tag	UNP Q99UU6
L	295	GLU	-	expression tag	UNP Q99UU6
L	296	HIS	-	expression tag	UNP Q99UU6
L	297	HIS	-	expression tag	UNP Q99UU6
L	298	HIS	-	expression tag	UNP Q99UU6
L	299	HIS	-	expression tag	UNP Q99UU6
L	300	HIS	-	expression tag	UNP Q99UU6
L	301	HIS	-	expression tag	UNP Q99UU6
M	0	MET	-	expression tag	UNP Q99UU6
M	294	LEU	-	expression tag	UNP Q99UU6
M	295	GLU	-	expression tag	UNP Q99UU6
M	296	HIS	-	expression tag	UNP Q99UU6
M	297	HIS	-	expression tag	UNP Q99UU6
M	298	HIS	-	expression tag	UNP Q99UU6
M	299	HIS	-	expression tag	UNP Q99UU6
M	300	HIS	-	expression tag	UNP Q99UU6
M	301	HIS	-	expression tag	UNP Q99UU6
N	0	MET	-	expression tag	UNP Q99UU6
N	294	LEU	-	expression tag	UNP Q99UU6
N	295	GLU	-	expression tag	UNP Q99UU6
N	296	HIS	-	expression tag	UNP Q99UU6
N	297	HIS	-	expression tag	UNP Q99UU6
N	298	HIS	-	expression tag	UNP Q99UU6
N	299	HIS	-	expression tag	UNP Q99UU6
N	300	HIS	-	expression tag	UNP Q99UU6
N	301	HIS	-	expression tag	UNP Q99UU6
O	0	MET	-	expression tag	UNP Q99UU6
O	294	LEU	-	expression tag	UNP Q99UU6
O	295	GLU	-	expression tag	UNP Q99UU6
O	296	HIS	-	expression tag	UNP Q99UU6
O	297	HIS	-	expression tag	UNP Q99UU6
O	298	HIS	-	expression tag	UNP Q99UU6
O	299	HIS	-	expression tag	UNP Q99UU6
O	300	HIS	-	expression tag	UNP Q99UU6
O	301	HIS	-	expression tag	UNP Q99UU6
P	0	MET	-	expression tag	UNP Q99UU6
P	294	LEU	-	expression tag	UNP Q99UU6
P	295	GLU	-	expression tag	UNP Q99UU6
P	296	HIS	-	expression tag	UNP Q99UU6
P	297	HIS	-	expression tag	UNP Q99UU6
P	298	HIS	-	expression tag	UNP Q99UU6

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Chain	Residue	Modelled	Actual	Comment	Reference
P	299	HIS	-	expression tag	UNP Q99UU6
P	300	HIS	-	expression tag	UNP Q99UU6
P	301	HIS	-	expression tag	UNP Q99UU6
Q	0	MET	-	expression tag	UNP Q99UU6
Q	294	LEU	-	expression tag	UNP Q99UU6
Q	295	GLU	-	expression tag	UNP Q99UU6
Q	296	HIS	-	expression tag	UNP Q99UU6
Q	297	HIS	-	expression tag	UNP Q99UU6
Q	298	HIS	-	expression tag	UNP Q99UU6
Q	299	HIS	-	expression tag	UNP Q99UU6
Q	300	HIS	-	expression tag	UNP Q99UU6
Q	301	HIS	-	expression tag	UNP Q99UU6
R	0	MET	-	expression tag	UNP Q99UU6
R	294	LEU	-	expression tag	UNP Q99UU6
R	295	GLU	-	expression tag	UNP Q99UU6
R	296	HIS	-	expression tag	UNP Q99UU6
R	297	HIS	-	expression tag	UNP Q99UU6
R	298	HIS	-	expression tag	UNP Q99UU6
R	299	HIS	-	expression tag	UNP Q99UU6
R	300	HIS	-	expression tag	UNP Q99UU6
R	301	HIS	-	expression tag	UNP Q99UU6
S	0	MET	-	expression tag	UNP Q99UU6
S	294	LEU	-	expression tag	UNP Q99UU6
S	295	GLU	-	expression tag	UNP Q99UU6
S	296	HIS	-	expression tag	UNP Q99UU6
S	297	HIS	-	expression tag	UNP Q99UU6
S	298	HIS	-	expression tag	UNP Q99UU6
S	299	HIS	-	expression tag	UNP Q99UU6
S	300	HIS	-	expression tag	UNP Q99UU6
S	301	HIS	-	expression tag	UNP Q99UU6
T	0	MET	-	expression tag	UNP Q99UU6
T	294	LEU	-	expression tag	UNP Q99UU6
T	295	GLU	-	expression tag	UNP Q99UU6
T	296	HIS	-	expression tag	UNP Q99UU6
T	297	HIS	-	expression tag	UNP Q99UU6
T	298	HIS	-	expression tag	UNP Q99UU6
T	299	HIS	-	expression tag	UNP Q99UU6
T	300	HIS	-	expression tag	UNP Q99UU6
T	301	HIS	-	expression tag	UNP Q99UU6
U	0	MET	-	expression tag	UNP Q99UU6
U	294	LEU	-	expression tag	UNP Q99UU6
U	295	GLU	-	expression tag	UNP Q99UU6

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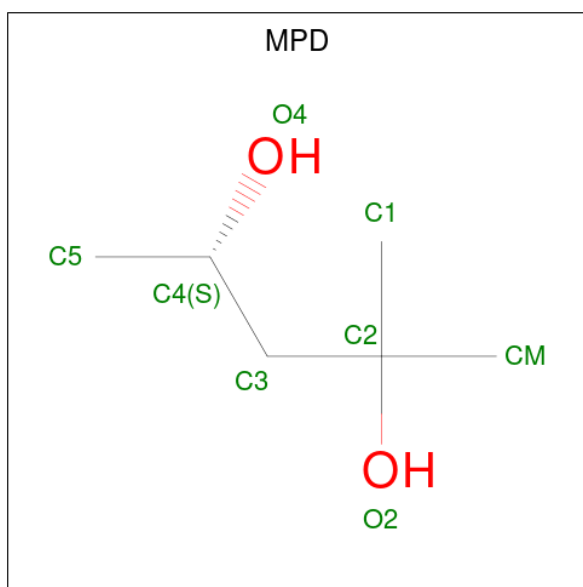
Chain	Residue	Modelled	Actual	Comment	Reference
U	296	HIS	-	expression tag	UNP Q99UU6
U	297	HIS	-	expression tag	UNP Q99UU6
U	298	HIS	-	expression tag	UNP Q99UU6
U	299	HIS	-	expression tag	UNP Q99UU6
U	300	HIS	-	expression tag	UNP Q99UU6
U	301	HIS	-	expression tag	UNP Q99UU6
V	0	MET	-	expression tag	UNP Q99UU6
V	294	LEU	-	expression tag	UNP Q99UU6
V	295	GLU	-	expression tag	UNP Q99UU6
V	296	HIS	-	expression tag	UNP Q99UU6
V	297	HIS	-	expression tag	UNP Q99UU6
V	298	HIS	-	expression tag	UNP Q99UU6
V	299	HIS	-	expression tag	UNP Q99UU6
V	300	HIS	-	expression tag	UNP Q99UU6
V	301	HIS	-	expression tag	UNP Q99UU6
W	0	MET	-	expression tag	UNP Q99UU6
W	294	LEU	-	expression tag	UNP Q99UU6
W	295	GLU	-	expression tag	UNP Q99UU6
W	296	HIS	-	expression tag	UNP Q99UU6
W	297	HIS	-	expression tag	UNP Q99UU6
W	298	HIS	-	expression tag	UNP Q99UU6
W	299	HIS	-	expression tag	UNP Q99UU6
W	300	HIS	-	expression tag	UNP Q99UU6
W	301	HIS	-	expression tag	UNP Q99UU6
X	0	MET	-	expression tag	UNP Q99UU6
X	294	LEU	-	expression tag	UNP Q99UU6
X	295	GLU	-	expression tag	UNP Q99UU6
X	296	HIS	-	expression tag	UNP Q99UU6
X	297	HIS	-	expression tag	UNP Q99UU6
X	298	HIS	-	expression tag	UNP Q99UU6
X	299	HIS	-	expression tag	UNP Q99UU6
X	300	HIS	-	expression tag	UNP Q99UU6
X	301	HIS	-	expression tag	UNP Q99UU6
Y	0	MET	-	expression tag	UNP Q99UU6
Y	294	LEU	-	expression tag	UNP Q99UU6
Y	295	GLU	-	expression tag	UNP Q99UU6
Y	296	HIS	-	expression tag	UNP Q99UU6
Y	297	HIS	-	expression tag	UNP Q99UU6
Y	298	HIS	-	expression tag	UNP Q99UU6
Y	299	HIS	-	expression tag	UNP Q99UU6
Y	300	HIS	-	expression tag	UNP Q99UU6
Y	301	HIS	-	expression tag	UNP Q99UU6

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Chain	Residue	Modelled	Actual	Comment	Reference
Z	0	MET	-	expression tag	UNP Q99UU6
Z	294	LEU	-	expression tag	UNP Q99UU6
Z	295	GLU	-	expression tag	UNP Q99UU6
Z	296	HIS	-	expression tag	UNP Q99UU6
Z	297	HIS	-	expression tag	UNP Q99UU6
Z	298	HIS	-	expression tag	UNP Q99UU6
Z	299	HIS	-	expression tag	UNP Q99UU6
Z	300	HIS	-	expression tag	UNP Q99UU6
Z	301	HIS	-	expression tag	UNP Q99UU6
a	0	MET	-	expression tag	UNP Q99UU6
a	294	LEU	-	expression tag	UNP Q99UU6
a	295	GLU	-	expression tag	UNP Q99UU6
a	296	HIS	-	expression tag	UNP Q99UU6
a	297	HIS	-	expression tag	UNP Q99UU6
a	298	HIS	-	expression tag	UNP Q99UU6
a	299	HIS	-	expression tag	UNP Q99UU6
a	300	HIS	-	expression tag	UNP Q99UU6
a	301	HIS	-	expression tag	UNP Q99UU6
b	0	MET	-	expression tag	UNP Q99UU6
b	294	LEU	-	expression tag	UNP Q99UU6
b	295	GLU	-	expression tag	UNP Q99UU6
b	296	HIS	-	expression tag	UNP Q99UU6
b	297	HIS	-	expression tag	UNP Q99UU6
b	298	HIS	-	expression tag	UNP Q99UU6
b	299	HIS	-	expression tag	UNP Q99UU6
b	300	HIS	-	expression tag	UNP Q99UU6
b	301	HIS	-	expression tag	UNP Q99UU6

- Molecule 2 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			8	6	2		
2	A	1	Total	C	O	0	0
			8	6	2		
2	B	1	Total	C	O	0	0
			8	6	2		
2	B	1	Total	C	O	0	0
			8	6	2		
2	C	1	Total	C	O	0	0
			8	6	2		
2	C	1	Total	C	O	0	0
			8	6	2		
2	D	1	Total	C	O	0	0
			8	6	2		
2	D	1	Total	C	O	0	0
			8	6	2		
2	E	1	Total	C	O	0	0
			8	6	2		
2	E	1	Total	C	O	0	0
			8	6	2		
2	F	1	Total	C	O	0	0
			8	6	2		
2	F	1	Total	C	O	0	0
			8	6	2		
2	G	1	Total	C	O	0	0
			8	6	2		
2	G	1	Total	C	O	0	0
			8	6	2		

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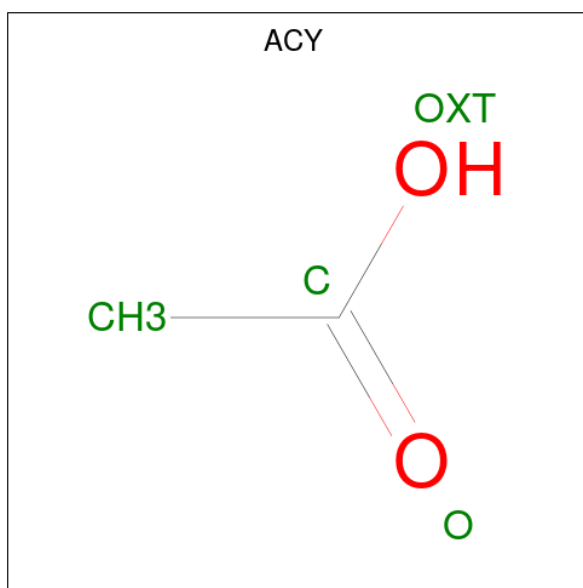
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
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			8	6	2		
2	H	1	Total	C	O	0	0
			8	6	2		
2	I	1	Total	C	O	0	0
			8	6	2		
2	I	1	Total	C	O	0	0
			8	6	2		
2	J	1	Total	C	O	0	0
			8	6	2		
2	J	1	Total	C	O	0	0
			8	6	2		
2	K	1	Total	C	O	0	0
			8	6	2		
2	K	1	Total	C	O	0	0
			8	6	2		
2	L	1	Total	C	O	0	0
			8	6	2		
2	L	1	Total	C	O	0	0
			8	6	2		
2	M	1	Total	C	O	0	0
			8	6	2		
2	M	1	Total	C	O	0	0
			8	6	2		
2	N	1	Total	C	O	0	0
			8	6	2		
2	N	1	Total	C	O	0	0
			8	6	2		
2	O	1	Total	C	O	0	0
			8	6	2		
2	O	1	Total	C	O	0	0
			8	6	2		
2	P	1	Total	C	O	0	0
			8	6	2		
2	P	1	Total	C	O	0	0
			8	6	2		
2	Q	1	Total	C	O	0	0
			8	6	2		
2	Q	1	Total	C	O	0	0
			8	6	2		
2	R	1	Total	C	O	0	0
			8	6	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	R	1	Total	C	O	0	0
			8	6	2		
2	S	1	Total	C	O	0	0
			8	6	2		
2	S	1	Total	C	O	0	0
			8	6	2		
2	T	1	Total	C	O	0	0
			8	6	2		
2	T	1	Total	C	O	0	0
			8	6	2		
2	U	1	Total	C	O	0	0
			8	6	2		
2	U	1	Total	C	O	0	0
			8	6	2		
2	V	1	Total	C	O	0	0
			8	6	2		
2	V	1	Total	C	O	0	0
			8	6	2		
2	W	1	Total	C	O	0	0
			8	6	2		
2	W	1	Total	C	O	0	0
			8	6	2		
2	X	1	Total	C	O	0	0
			8	6	2		
2	X	1	Total	C	O	0	0
			8	6	2		
2	X	1	Total	C	O	0	0
			8	6	2		
2	Y	1	Total	C	O	0	0
			8	6	2		
2	Y	1	Total	C	O	0	0
			8	6	2		
2	Z	1	Total	C	O	0	0
			8	6	2		
2	a	1	Total	C	O	0	0
			8	6	2		
2	a	1	Total	C	O	0	0
			8	6	2		
2	b	1	Total	C	O	0	0
			8	6	2		
2	b	1	Total	C	O	0	0
			8	6	2		

- Molecule 3 is ACETIC ACID (CCD ID: ACY) (formula: $C_2H_4O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	E	1	Total	C	O	0	0
			4	2	2		
3	F	1	Total	C	O	0	0
			4	2	2		
3	H	1	Total	C	O	0	0
			4	2	2		
3	H	1	Total	C	O	0	0
			4	2	2		
3	J	1	Total	C	O	0	0
			4	2	2		
3	K	1	Total	C	O	0	0
			4	2	2		
3	K	1	Total	C	O	0	0
			4	2	2		
3	L	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	M	1	Total C O 4 2 2	0	0
3	O	1	Total C O 4 2 2	0	0
3	O	1	Total C O 4 2 2	0	0
3	P	1	Total C O 4 2 2	0	0
3	Q	1	Total C O 4 2 2	0	0
3	S	1	Total C O 4 2 2	0	0
3	S	1	Total C O 4 2 2	0	0
3	T	1	Total C O 4 2 2	0	0
3	V	1	Total C O 4 2 2	0	0
3	W	1	Total C O 4 2 2	0	0
3	X	1	Total C O 4 2 2	0	0
3	Y	1	Total C O 4 2 2	0	0
3	Z	1	Total C O 4 2 2	0	0
3	a	1	Total C O 4 2 2	0	0
3	b	1	Total C O 4 2 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	94	Total O 94 94	0	0
4	B	79	Total O 79 79	0	0
4	C	73	Total O 73 73	0	0
4	D	66	Total O 66 66	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	E	54	Total O 54 54	0	0
4	F	46	Total O 46 46	0	0
4	G	70	Total O 70 70	0	0
4	H	89	Total O 89 89	0	0
4	I	87	Total O 87 87	0	0
4	J	86	Total O 86 86	0	0
4	K	97	Total O 97 97	0	0
4	L	71	Total O 71 71	0	0
4	M	61	Total O 61 61	0	0
4	N	73	Total O 73 73	0	0
4	O	91	Total O 91 91	0	0
4	P	73	Total O 73 73	0	0
4	Q	61	Total O 61 61	0	0
4	R	73	Total O 73 73	0	0
4	S	69	Total O 69 69	0	0
4	T	82	Total O 82 82	0	0
4	U	85	Total O 85 85	0	0
4	V	82	Total O 82 82	0	0
4	W	50	Total O 50 50	0	0
4	X	46	Total O 46 46	0	0
4	Y	53	Total O 53 53	0	0

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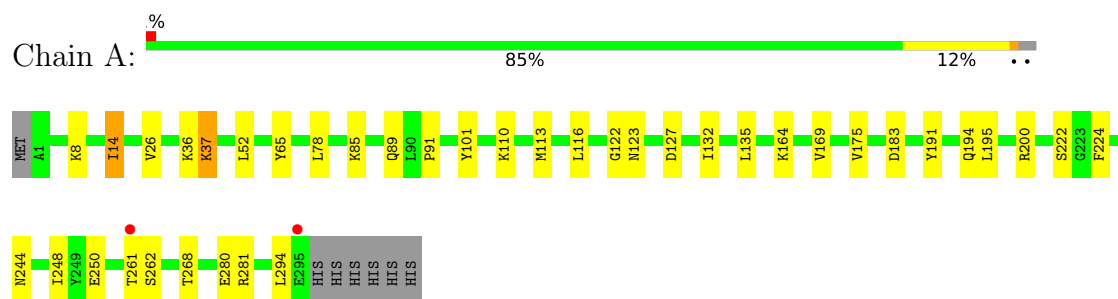
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	Z	82	Total 82	O 82	0	0
4	a	99	Total 99	O 99	0	0
4	b	72	Total 72	O 72	0	0

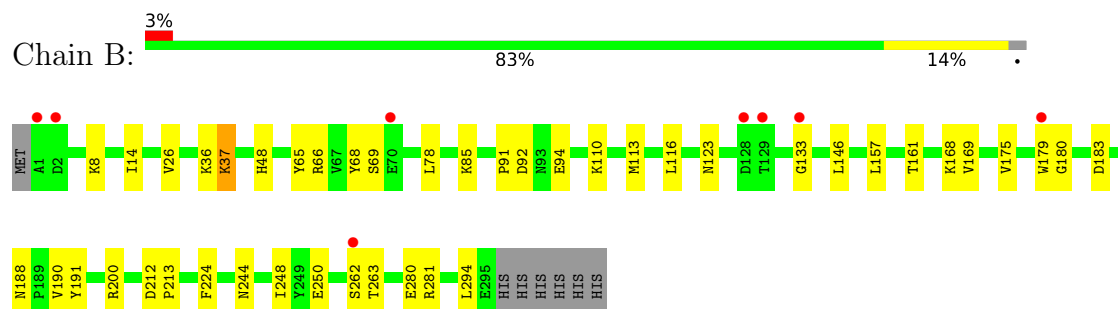
3 Residue-property plots [i](#)

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.

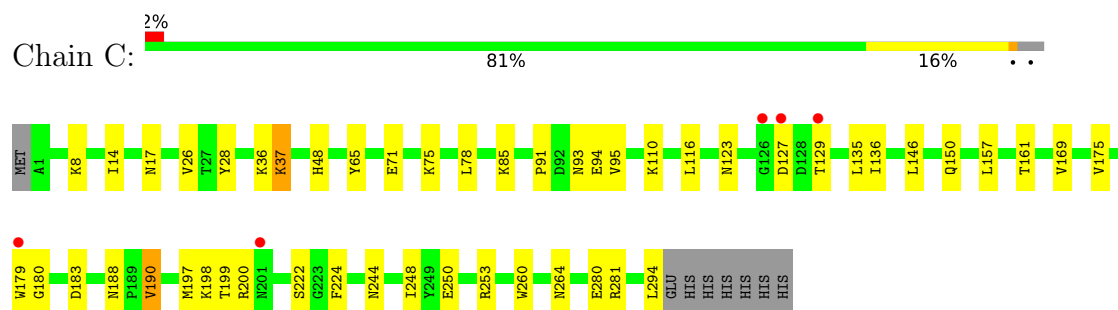
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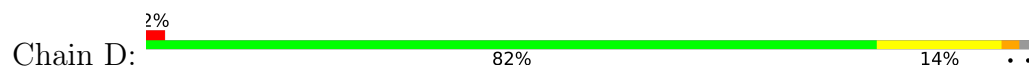
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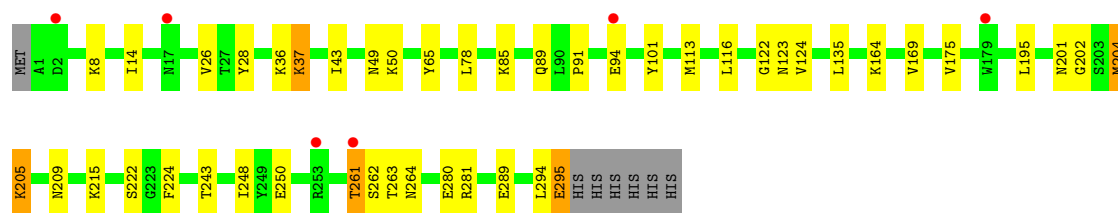


• Molecule 1: Alpha-hemolysin

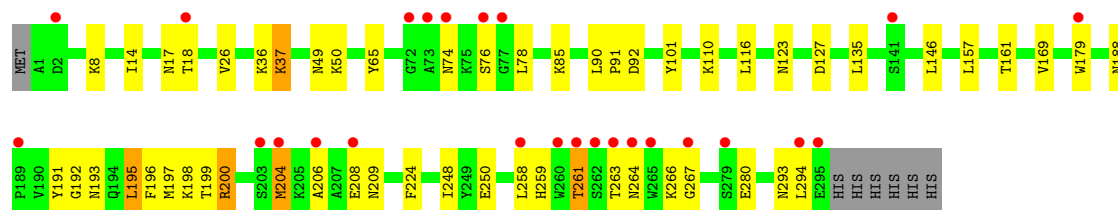
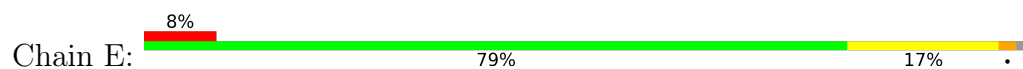


• Molecule 1: Alpha-hemolysin

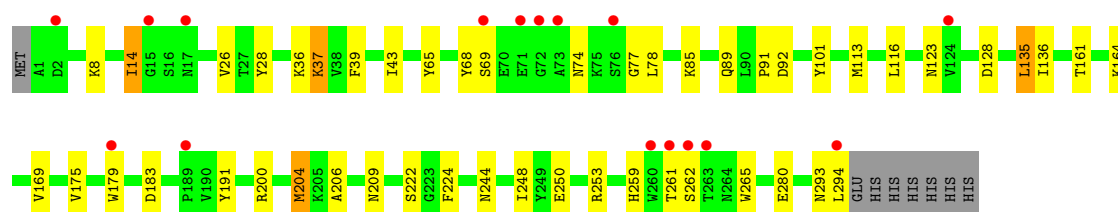
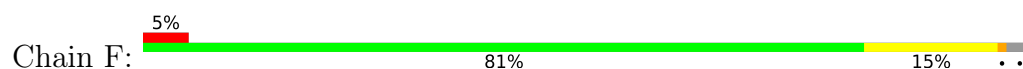




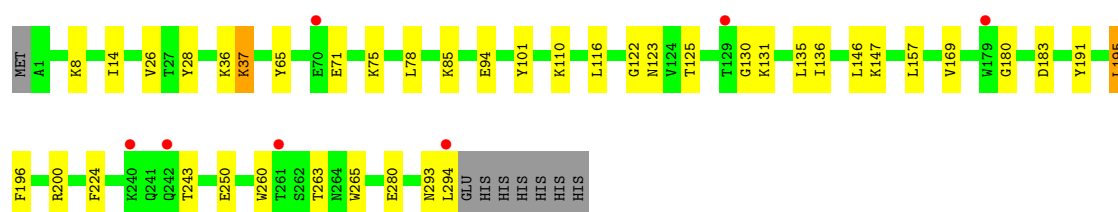
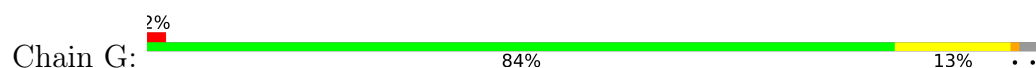
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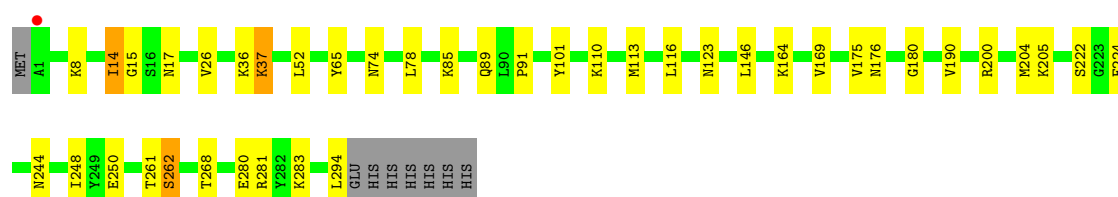
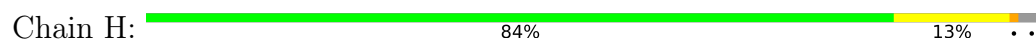
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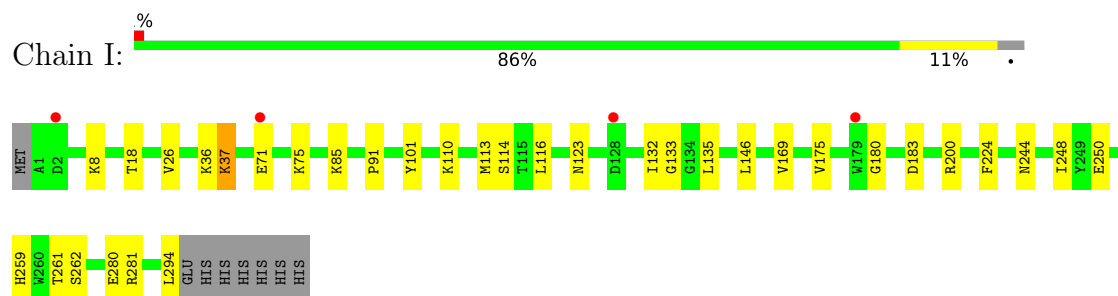
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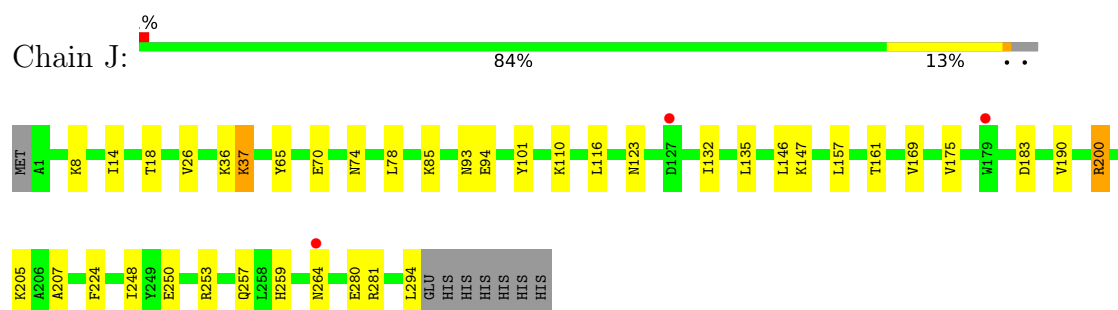
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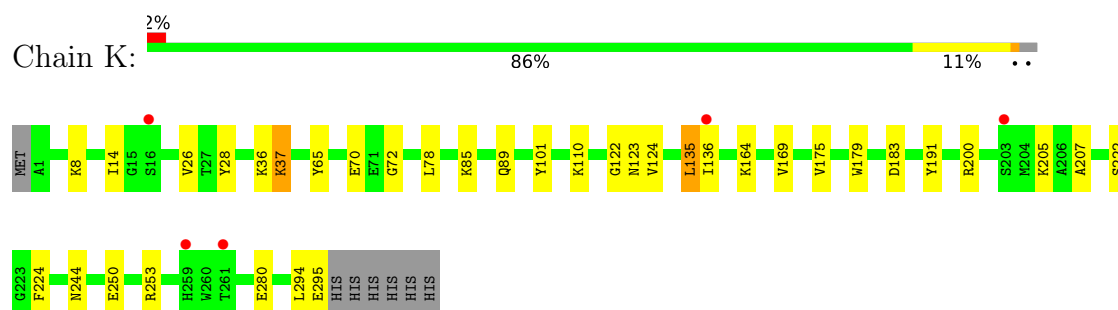
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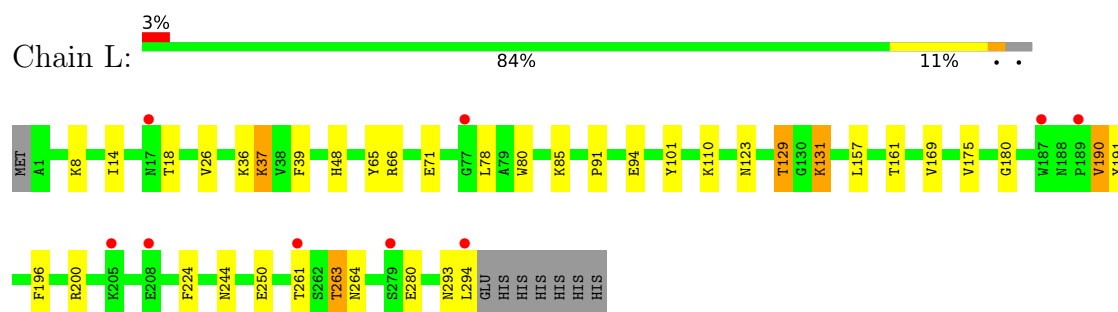
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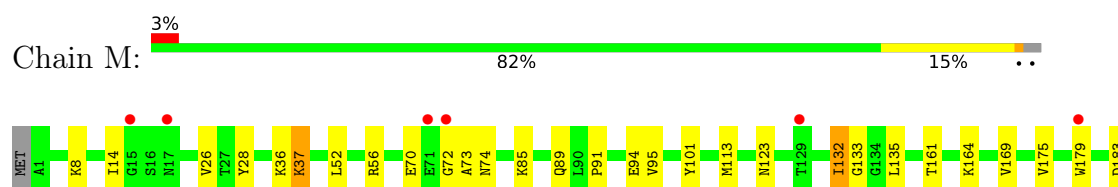
- Molecule 1: Alpha-hemolysin



- Molecule 1: Alpha-hemolysin

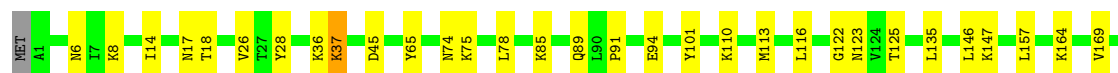
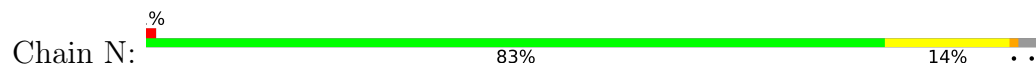


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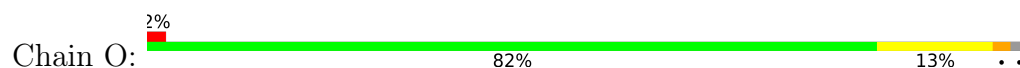




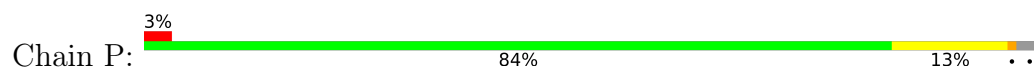
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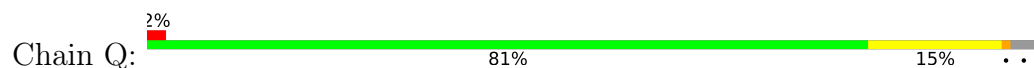
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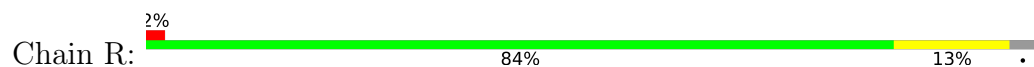
• Molecule 1: Alpha-hemolysin

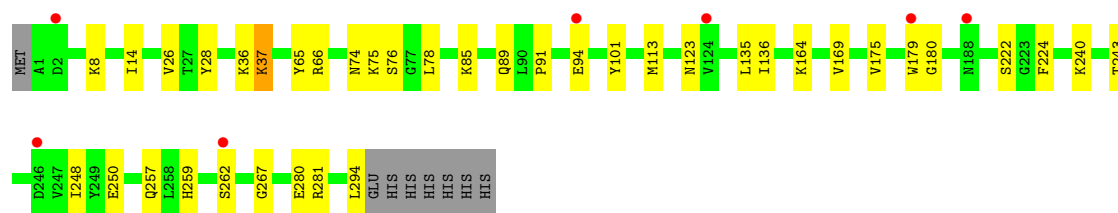


• Molecule 1: Alpha-hemolysin

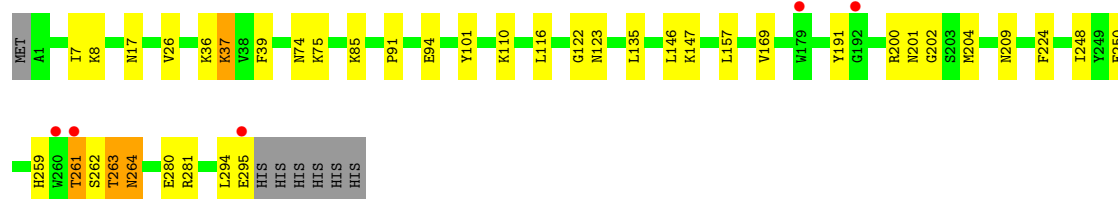
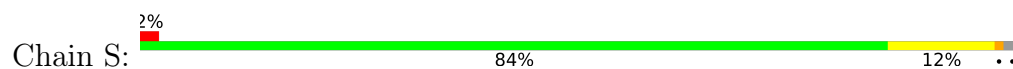


• Molecule 1: Alpha-hemolysin

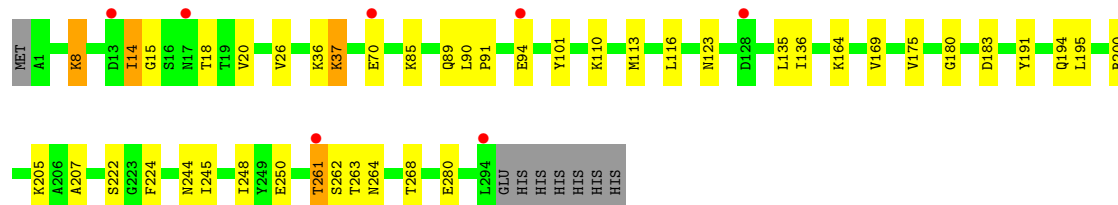
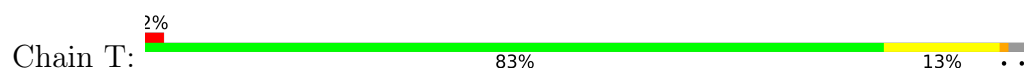




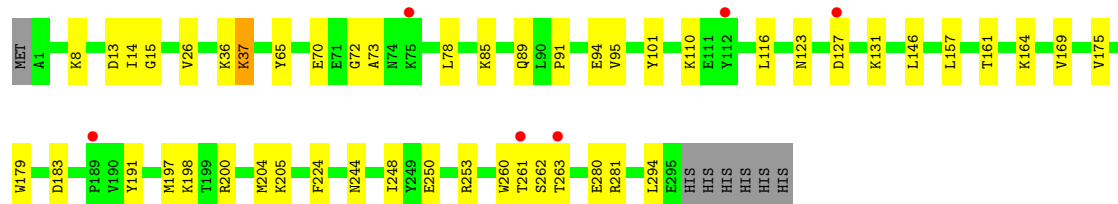
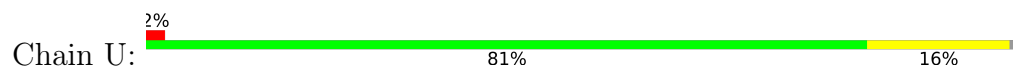
• Molecule 1: Alpha-hemolysin



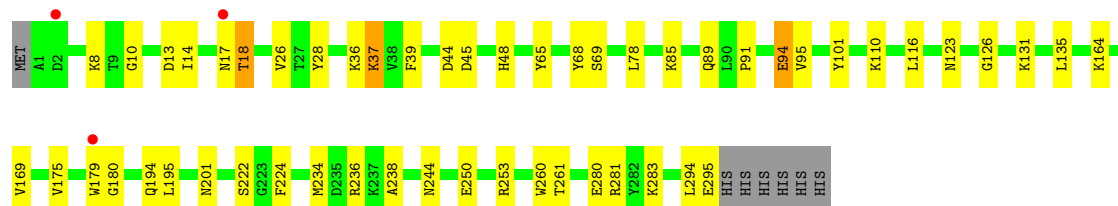
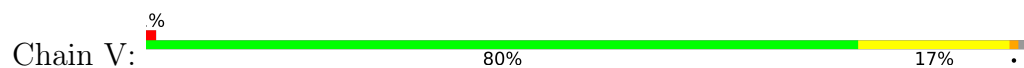
• Molecule 1: Alpha-hemolysin



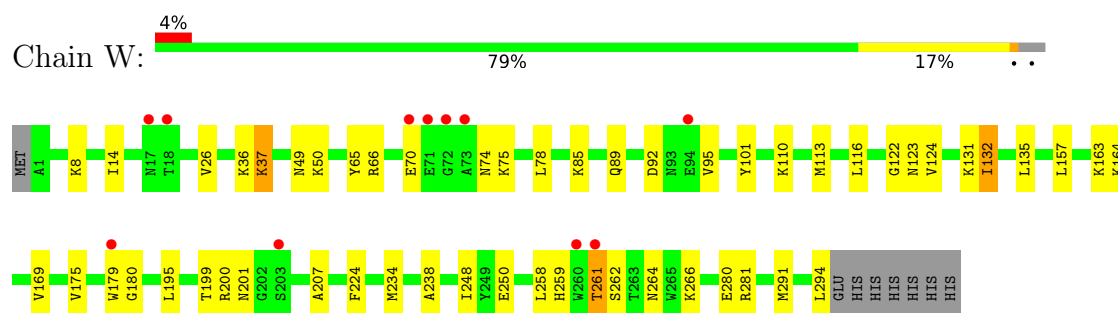
• Molecule 1: Alpha-hemolysin



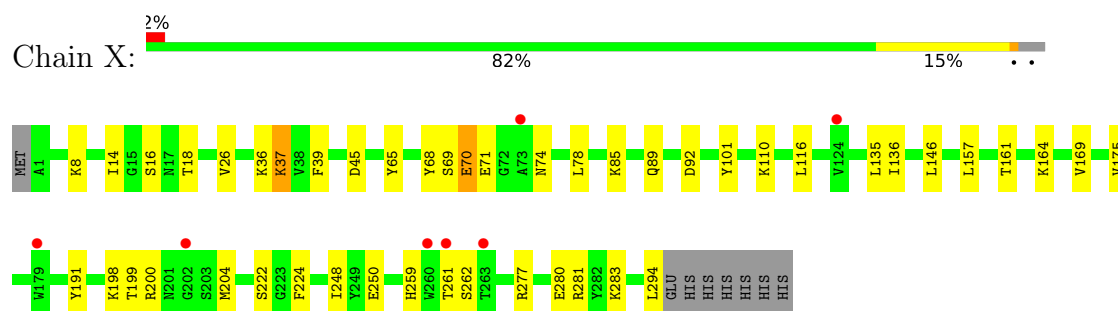
• Molecule 1: Alpha-hemolysin



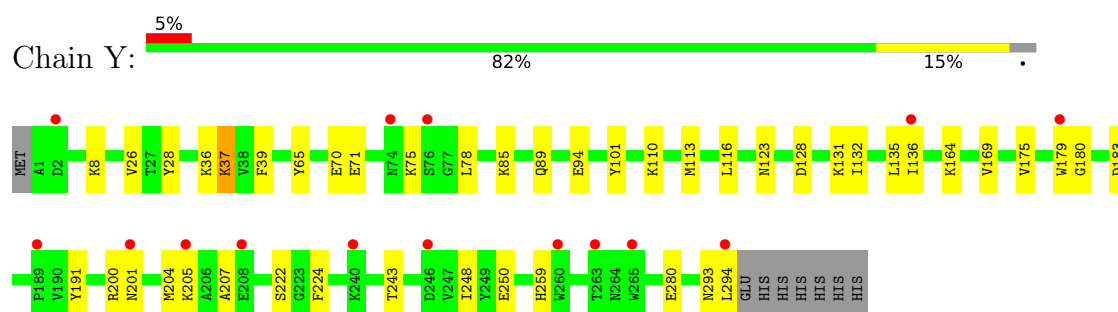
- Molecule 1: Alpha-hemolysin



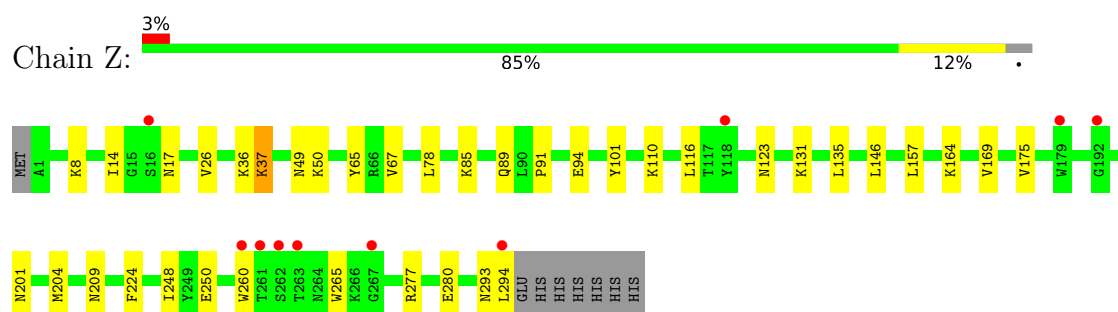
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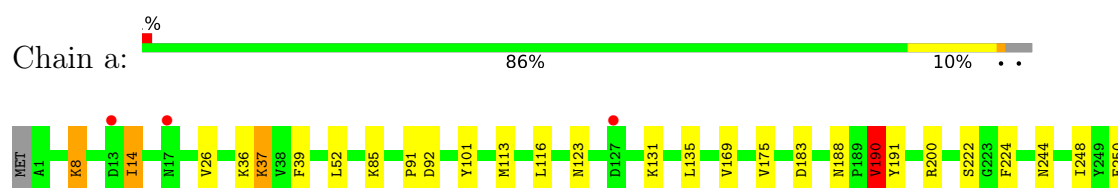
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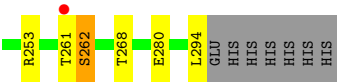


- Molecule 1: Alpha-hemolysin

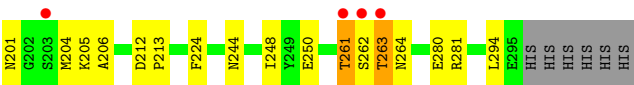
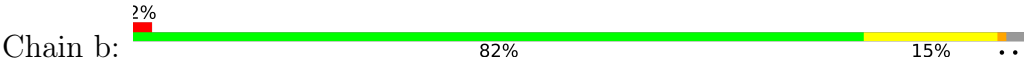


- Molecule 1: Alpha-hemolysin





● Molecule 1: Alpha-hemolysin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	102.82Å 293.88Å 170.54Å 90.00° 92.42° 90.00°	Depositor
Resolution (Å)	19.98 – 2.30 19.98 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.8 (19.98-2.30) 97.9 (19.98-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.77 (at 2.30Å)	Xtriage
Refinement program	PHENIX 1.6.4_486	Depositor
R, R_{free}	0.193 , 0.226 0.194 , 0.192	Depositor DCC
R_{free} test set	31076 reflections (7.15%)	wwPDB-VP
Wilson B-factor (Å ²)	36.5	Xtriage
Anisotropy	0.672	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 56.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.010 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	68607	wwPDB-VP
Average B, all atoms (Å ²)	52.0	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 61.01 % 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 1.3976e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.42	0/2414	0.66	1/3267 (0.0%)
1	B	0.41	0/2414	0.62	0/3267
1	C	0.40	0/2405	0.63	0/3255
1	D	0.41	0/2414	0.63	0/3267
1	E	0.39	0/2414	0.62	0/3267
1	F	0.38	0/2405	0.63	0/3255
1	G	0.40	0/2405	0.64	1/3255 (0.0%)
1	H	0.43	0/2405	0.67	2/3255 (0.1%)
1	I	0.41	0/2405	0.63	0/3255
1	J	0.42	0/2405	0.62	0/3255
1	K	0.42	0/2414	0.64	0/3267
1	L	0.42	0/2405	0.63	0/3255
1	M	0.39	0/2414	0.62	0/3267
1	N	0.40	0/2405	0.62	0/3255
1	O	0.41	0/2414	0.64	0/3267
1	P	0.41	0/2405	0.64	1/3255 (0.0%)
1	Q	0.39	0/2405	0.62	0/3255
1	R	0.41	0/2405	0.64	0/3255
1	S	0.38	0/2414	0.62	1/3267 (0.0%)
1	T	0.42	0/2405	0.64	1/3255 (0.0%)
1	U	0.40	0/2414	0.62	0/3267
1	V	0.40	0/2414	0.63	0/3267
1	W	0.39	0/2405	0.63	0/3255
1	X	0.37	0/2405	0.62	0/3255
1	Y	0.39	0/2405	0.65	0/3255
1	Z	0.39	0/2405	0.63	0/3255
1	a	0.42	0/2405	0.64	2/3255 (0.1%)
1	b	0.41	0/2414	0.63	0/3267
All	All	0.40	0/67439	0.63	9/91272 (0.0%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	261	THR	N-CA-C	-10.17	96.82	110.55
1	A	261	THR	N-CA-C	-10.05	96.98	110.55
1	P	261	THR	N-CA-C	-6.08	101.51	110.52
1	H	190	VAL	N-CA-C	6.00	117.03	111.45
1	T	261	THR	N-CA-C	-5.89	102.20	110.50
1	a	261	THR	N-CA-C	-5.49	102.75	110.50
1	G	130	GLY	N-CA-C	-5.21	108.08	115.43
1	S	264	ASN	N-CA-C	5.17	114.97	108.45
1	a	190	VAL	N-CA-C	5.04	116.15	111.81

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2362	0	2289	31	0
1	B	2362	0	2289	38	0
1	C	2353	0	2283	38	0
1	D	2362	0	2289	35	0
1	E	2362	0	2289	44	0
1	F	2353	0	2283	44	0
1	G	2353	0	2283	33	0
1	H	2353	0	2283	32	0
1	I	2353	0	2283	29	0
1	J	2353	0	2283	28	0
1	K	2362	0	2289	33	0
1	L	2353	0	2283	29	0
1	M	2362	0	2289	33	0
1	N	2353	0	2283	42	0
1	O	2362	0	2289	38	0
1	P	2353	0	2283	30	0
1	Q	2353	0	2283	39	0
1	R	2353	0	2283	32	0
1	S	2362	0	2289	32	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	T	2353	0	2283	38	0
1	U	2362	0	2289	38	0
1	V	2362	0	2289	45	0
1	W	2353	0	2283	42	0
1	X	2353	0	2283	41	0
1	Y	2353	0	2283	34	0
1	Z	2353	0	2283	29	0
1	a	2353	0	2283	30	0
1	b	2362	0	2289	44	0
2	A	16	0	28	4	0
2	B	16	0	28	9	0
2	C	16	0	28	6	0
2	D	16	0	28	2	0
2	E	16	0	28	9	0
2	F	16	0	28	10	0
2	G	16	0	28	7	0
2	H	16	0	28	5	0
2	I	16	0	28	7	0
2	J	16	0	28	5	0
2	K	16	0	28	11	0
2	L	16	0	28	5	0
2	M	16	0	28	4	0
2	N	16	0	28	11	0
2	O	16	0	28	10	0
2	P	16	0	28	4	0
2	Q	16	0	28	11	0
2	R	16	0	28	6	0
2	S	16	0	28	4	0
2	T	16	0	28	4	0
2	U	16	0	28	5	0
2	V	16	0	28	8	0
2	W	16	0	28	9	0
2	X	24	0	42	6	0
2	Y	16	0	28	6	0
2	Z	8	0	14	1	0
2	a	16	0	28	8	0
2	b	16	0	28	6	0
3	A	8	0	6	0	0
3	C	4	0	3	0	0
3	D	8	0	6	0	0
3	E	4	0	3	0	0
3	F	4	0	3	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	8	0	6	0	0
3	J	4	0	3	0	0
3	K	8	0	6	0	0
3	L	4	0	3	0	0
3	M	4	0	3	0	0
3	O	8	0	6	0	0
3	P	4	0	3	0	0
3	Q	4	0	3	0	0
3	S	8	0	6	0	0
3	T	4	0	3	0	0
3	V	4	0	3	0	0
3	W	4	0	3	0	0
3	X	4	0	3	0	0
3	Y	4	0	3	0	0
3	Z	4	0	3	0	0
3	a	4	0	3	0	0
3	b	4	0	3	0	0
4	A	94	0	0	1	0
4	B	79	0	0	2	0
4	C	73	0	0	6	0
4	D	66	0	0	1	0
4	E	54	0	0	1	0
4	F	46	0	0	2	0
4	G	70	0	0	1	0
4	H	89	0	0	1	0
4	I	87	0	0	2	0
4	J	86	0	0	3	0
4	K	97	0	0	4	0
4	L	71	0	0	0	0
4	M	61	0	0	2	0
4	N	73	0	0	3	0
4	O	91	0	0	1	0
4	P	73	0	0	1	0
4	Q	61	0	0	2	0
4	R	73	0	0	0	0
4	S	69	0	0	0	0
4	T	82	0	0	3	0
4	U	85	0	0	2	0
4	V	82	0	0	7	0
4	W	50	0	0	1	0
4	X	46	0	0	1	0
4	Y	53	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	Z	82	0	0	2	0
4	a	99	0	0	2	0
4	b	72	0	0	2	0
All	All	68607	0	64858	968	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (968) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:303:MPD:H11	2:B:303:MPD:H53	1.27	1.09
1:R:180:GLY:O	2:R:303:MPD:H11	1.52	1.06
1:E:199:THR:HA	2:E:302:MPD:H51	1.38	1.04
1:I:180:GLY:O	2:I:303:MPD:H12	1.60	1.01
2:S:303:MPD:H53	2:S:303:MPD:H12	1.44	0.98
2:Y:303:MPD:H12	2:Y:303:MPD:H53	1.49	0.93
2:H:303:MPD:H52	2:H:303:MPD:H13	1.51	0.91
2:A:302:MPD:HM3	2:A:302:MPD:H52	1.58	0.86
1:F:179:TRP:CE3	2:F:303:MPD:C5	2.61	0.83
1:P:180:GLY:O	2:P:303:MPD:H13	1.79	0.82
1:O:91:PRO:HG2	1:O:94:GLU:CG	2.08	0.82
1:F:179:TRP:HA	2:F:303:MPD:H53	1.61	0.81
1:U:91:PRO:HB2	1:U:94:GLU:HG2	1.61	0.81
1:M:179:TRP:CE3	2:M:303:MPD:H51	2.16	0.80
1:C:281:ARG:HG2	1:C:294:LEU:HD12	1.61	0.80
1:W:200:ARG:HG3	2:W:302:MPD:H52	1.62	0.80
1:D:91:PRO:HG2	1:D:94:GLU:HB2	1.63	0.79
2:A:303:MPD:H12	2:A:303:MPD:H53	1.65	0.79
2:G:303:MPD:H12	2:G:303:MPD:H53	1.63	0.79
1:V:131:LYS:HE2	1:b:127:ASP:HB2	1.67	0.76
1:F:179:TRP:CD2	2:F:303:MPD:C5	2.69	0.76
1:T:91:PRO:HB2	1:T:94:GLU:HG2	1.67	0.75
1:E:199:THR:CA	2:E:302:MPD:H51	2.14	0.75
1:G:180:GLY:O	2:G:303:MPD:H13	1.86	0.75
1:O:91:PRO:HG2	1:O:94:GLU:HG2	1.68	0.75
2:I:302:MPD:O4	2:I:302:MPD:H11	1.86	0.75
2:A:302:MPD:H52	2:A:302:MPD:CM	2.17	0.74
1:B:179:TRP:CE3	2:B:303:MPD:H51	2.23	0.74
1:F:179:TRP:CE3	2:F:303:MPD:H52	2.22	0.74
1:b:91:PRO:HB2	1:b:94:GLU:HG2	1.68	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:b:303:MPD:H12	2:b:303:MPD:H52	1.69	0.74
1:Q:199:THR:CA	2:Q:302:MPD:H51	2.18	0.74
1:Z:91:PRO:HG2	1:Z:94:GLU:CG	2.18	0.74
1:O:179:TRP:HA	2:O:303:MPD:H53	1.71	0.72
2:W:303:MPD:H12	2:W:303:MPD:H53	1.71	0.72
1:K:179:TRP:HA	2:K:303:MPD:H52	1.70	0.72
1:E:179:TRP:CG	2:E:302:MPD:HM3	2.24	0.72
1:F:179:TRP:CE3	2:F:303:MPD:H51	2.24	0.72
2:G:302:MPD:H11	2:G:302:MPD:O4	1.89	0.71
1:V:179:TRP:HA	2:V:303:MPD:H53	1.70	0.71
2:a:302:MPD:O4	2:a:302:MPD:H11	1.90	0.71
1:T:91:PRO:HG2	1:T:94:GLU:HG3	1.73	0.71
4:C:1969:HOH:O	1:D:215:LYS:HE2	1.89	0.71
2:L:303:MPD:H12	2:L:303:MPD:H53	1.73	0.70
1:O:91:PRO:HG2	1:O:94:GLU:HG3	1.73	0.70
1:U:91:PRO:HG2	1:U:94:GLU:HG3	1.74	0.70
1:H:262:SER:HB2	1:a:268:THR:HG21	1.74	0.69
1:F:179:TRP:CD2	2:F:303:MPD:H51	2.27	0.69
2:R:303:MPD:H52	2:R:303:MPD:H12	1.74	0.69
1:N:180:GLY:O	2:N:303:MPD:H13	1.92	0.69
1:K:200:ARG:CZ	2:K:302:MPD:C1	2.71	0.69
2:O:302:MPD:O4	2:O:302:MPD:H11	1.91	0.68
1:K:124:VAL:HG11	1:b:262:SER:HB3	1.74	0.68
2:K:303:MPD:H52	2:K:303:MPD:H11	1.75	0.68
2:B:303:MPD:H11	2:B:303:MPD:C5	2.13	0.68
2:N:303:MPD:H12	2:N:303:MPD:H53	1.76	0.68
1:C:179:TRP:CE3	2:C:303:MPD:H51	2.29	0.67
1:V:281:ARG:HG2	1:V:294:LEU:HD12	1.76	0.67
1:J:281:ARG:HG2	1:J:294:LEU:HD12	1.77	0.67
2:J:302:MPD:HM1	2:J:302:MPD:O4	1.95	0.67
2:K:303:MPD:H11	2:K:303:MPD:C5	2.23	0.67
2:M:303:MPD:H12	2:M:303:MPD:H53	1.76	0.67
1:M:72:GLY:C	1:M:205:LYS:HZ1	2.02	0.66
1:S:91:PRO:HB2	1:S:94:GLU:HG2	1.77	0.66
1:X:116:LEU:HB2	2:X:304:MPD:HM1	1.76	0.66
1:b:200:ARG:HE	2:b:302:MPD:C1	2.09	0.66
2:B:302:MPD:H11	2:B:302:MPD:O4	1.95	0.66
1:J:200:ARG:HD2	2:J:302:MPD:H32	1.77	0.66
1:E:65:TYR:CE1	1:E:78:LEU:HD21	2.31	0.65
1:Y:180:GLY:O	2:Y:303:MPD:H13	1.97	0.65
1:W:179:TRP:CE3	2:W:303:MPD:H51	2.32	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Y:116:LEU:HB2	2:a:303:MPD:H11	1.79	0.65
1:E:204:MET:N	4:E:1879:HOH:O	2.28	0.65
1:S:75:LYS:HE3	1:S:259:HIS:HB2	1.78	0.65
1:W:261:THR:HG22	1:W:264:ASN:O	1.96	0.64
1:B:91:PRO:HG2	1:B:94:GLU:HG3	1.78	0.64
1:S:281:ARG:HG2	1:S:294:LEU:HD12	1.80	0.64
1:B:281:ARG:HG2	1:B:294:LEU:HD12	1.79	0.64
1:S:191:TYR:CE1	1:S:200:ARG:HB3	2.32	0.64
1:Q:14:ILE:HD11	1:Q:52:LEU:HD22	1.78	0.64
1:E:74:ASN:O	1:E:259:HIS:HA	1.97	0.63
1:T:180:GLY:O	2:T:303:MPD:H12	1.98	0.63
1:b:261:THR:HG22	1:b:264:ASN:H	1.63	0.63
1:G:200:ARG:HE	2:G:302:MPD:C1	2.11	0.63
1:Z:91:PRO:HG2	1:Z:94:GLU:HG3	1.80	0.63
1:W:261:THR:HG23	1:W:262:SER:N	2.13	0.63
2:B:303:MPD:H53	2:B:303:MPD:C1	2.15	0.63
1:I:262:SER:HB3	1:W:124:VAL:HG11	1.79	0.63
1:D:94:GLU:HG2	1:D:243:THR:HG23	1.80	0.63
1:O:179:TRP:CE3	2:O:303:MPD:H51	2.34	0.63
1:b:180:GLY:O	2:b:303:MPD:H13	1.98	0.63
1:G:200:ARG:HE	2:G:302:MPD:H13	1.63	0.62
1:N:293:ASN:O	1:N:294:LEU:HB2	1.99	0.62
1:Q:198:LYS:C	2:Q:302:MPD:H53	2.25	0.62
1:Q:14:ILE:CD1	1:Q:52:LEU:HD22	2.30	0.62
1:S:200:ARG:HE	2:S:302:MPD:H13	1.64	0.62
2:E:303:MPD:H52	2:E:303:MPD:H12	1.82	0.62
1:Z:94:GLU:HB2	4:Z:738:HOH:O	1.99	0.62
1:F:65:TYR:CE2	1:F:78:LEU:HD21	2.34	0.62
1:Q:199:THR:C	2:Q:302:MPD:H51	2.24	0.62
2:P:302:MPD:H13	2:P:302:MPD:O4	1.98	0.62
1:Q:18:THR:HG21	1:Q:45:ASP:HA	1.82	0.62
1:Q:199:THR:HA	2:Q:302:MPD:H51	1.82	0.62
2:V:302:MPD:O4	2:V:302:MPD:HM3	1.97	0.62
2:T:302:MPD:HM1	2:T:302:MPD:O4	1.99	0.62
1:D:261:THR:HG22	1:D:263:THR:H	1.65	0.61
2:Q:303:MPD:H12	2:Q:303:MPD:H52	1.82	0.61
1:J:183:ASP:HB2	4:J:309:HOH:O	2.01	0.61
1:C:36:LYS:HZ2	1:C:280:GLU:HG2	1.66	0.61
1:A:268:THR:HG21	1:T:262:SER:HB2	1.83	0.61
1:C:150:GLN:HG2	4:C:508:HOH:O	1.99	0.61
1:H:14:ILE:HG22	1:N:6:ASN:O	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:200:ARG:NH2	2:K:302:MPD:H11	2.16	0.60
1:Y:94:GLU:CD	1:Y:243:THR:HG23	2.25	0.60
1:b:206:ALA:HB3	4:b:1684:HOH:O	2.00	0.60
1:D:36:LYS:HZ2	1:D:280:GLU:HG2	1.66	0.60
1:O:110:LYS:HD2	1:P:175:VAL:HG23	1.84	0.60
1:G:195:LEU:HD13	1:G:196:PHE:CE2	2.36	0.60
2:C:302:MPD:O4	2:C:302:MPD:H11	2.02	0.60
1:N:191:TYR:CZ	1:N:200:ARG:HD3	2.37	0.60
1:D:281:ARG:HG2	1:D:294:LEU:HD12	1.84	0.60
1:P:14:ILE:HD11	1:Q:39:PHE:HE1	1.66	0.60
1:X:281:ARG:HG2	1:X:294:LEU:HD12	1.83	0.60
1:a:294:LEU:HD12	1:a:294:LEU:C	2.27	0.60
1:H:268:THR:HG21	1:a:262:SER:HB2	1.84	0.59
1:Q:200:ARG:HE	2:Q:302:MPD:CM	2.14	0.59
1:Q:179:TRP:CG	2:Q:302:MPD:HO2	2.20	0.59
2:O:302:MPD:O4	2:O:302:MPD:HM3	2.02	0.59
1:X:65:TYR:CE2	1:X:78:LEU:HD21	2.37	0.59
1:b:91:PRO:HG2	1:b:94:GLU:HG3	1.85	0.59
1:L:261:THR:HG22	1:L:264:ASN:O	2.02	0.59
2:K:303:MPD:H53	2:K:303:MPD:HM2	1.85	0.59
1:C:179:TRP:HA	2:C:303:MPD:H53	1.84	0.59
1:I:200:ARG:CZ	2:I:302:MPD:C1	2.81	0.59
1:C:197:MET:HE3	1:C:200:ARG:HG2	1.85	0.59
1:T:205:LYS:HG2	4:T:1509:HOH:O	2.03	0.59
1:L:261:THR:HG23	1:L:263:THR:H	1.67	0.58
1:N:18:THR:HG22	4:N:1336:HOH:O	2.03	0.58
1:O:200:ARG:CZ	2:O:302:MPD:C1	2.81	0.58
1:V:85:LYS:HB2	1:V:250:GLU:HB3	1.85	0.58
1:A:14:ILE:CD1	1:A:52:LEU:HD22	2.32	0.58
1:B:85:LYS:HB2	1:B:250:GLU:HB3	1.85	0.58
1:E:192:GLY:N	1:E:266:LYS:HG3	2.17	0.58
1:M:85:LYS:HB2	1:M:250:GLU:HB3	1.86	0.58
1:F:200:ARG:HE	2:F:302:MPD:C1	2.15	0.58
1:S:91:PRO:HG2	1:S:94:GLU:HG3	1.85	0.58
1:W:281:ARG:HG2	1:W:294:LEU:HD12	1.84	0.58
1:E:204:MET:HE2	1:E:204:MET:H	1.69	0.58
1:J:205:LYS:HD2	1:J:205:LYS:N	2.17	0.58
1:D:65:TYR:CE1	1:D:78:LEU:HD21	2.39	0.58
1:G:71:GLU:HB2	1:G:75:LYS:HB2	1.86	0.58
1:Y:191:TYR:CE1	1:Y:200:ARG:HB3	2.38	0.58
1:O:200:ARG:CZ	2:O:302:MPD:H13	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:135:LEU:C	1:F:135:LEU:HD12	2.28	0.58
1:I:259:HIS:ND1	4:I:1717:HOH:O	2.32	0.58
1:Q:179:TRP:CD1	2:Q:302:MPD:HO2	2.22	0.58
1:A:262:SER:HB2	1:T:268:THR:HG21	1.84	0.58
1:X:70:GLU:O	1:X:70:GLU:HG3	2.04	0.58
1:I:91:PRO:HD3	1:I:244:ASN:O	2.04	0.57
1:V:131:LYS:HE3	1:b:129:THR:HG23	1.86	0.57
1:F:91:PRO:HD3	1:F:244:ASN:O	2.05	0.57
1:L:65:TYR:CE2	1:L:78:LEU:HD21	2.40	0.57
2:P:303:MPD:H52	2:P:303:MPD:O2	2.03	0.57
1:Q:200:ARG:HE	2:Q:302:MPD:HM1	1.68	0.57
1:I:183:ASP:HB2	4:I:1915:HOH:O	2.04	0.57
1:Y:94:GLU:OE1	1:Y:243:THR:HG23	2.04	0.57
2:C:303:MPD:H52	2:C:303:MPD:O2	2.05	0.57
2:D:302:MPD:HM1	2:D:302:MPD:C5	2.35	0.57
1:F:293:ASN:O	1:F:294:LEU:HB2	2.02	0.57
1:R:85:LYS:HB2	1:R:250:GLU:HB3	1.87	0.57
1:T:180:GLY:O	2:T:303:MPD:C1	2.53	0.57
1:W:258:LEU:HD12	1:W:258:LEU:O	2.05	0.57
1:C:14:ILE:HD11	1:C:48:HIS:CE1	2.40	0.57
1:T:183:ASP:HB2	4:T:1635:HOH:O	2.03	0.57
1:B:262:SER:HB3	1:P:124:VAL:HG11	1.85	0.57
1:N:91:PRO:HB2	1:N:94:GLU:HG2	1.86	0.57
1:R:91:PRO:HG2	1:R:94:GLU:HB2	1.87	0.57
1:E:36:LYS:HZ2	1:E:280:GLU:HG2	1.69	0.57
1:K:244:ASN:HB2	4:K:1399:HOH:O	2.05	0.57
1:b:261:THR:HG22	1:b:264:ASN:O	2.05	0.57
1:B:91:PRO:CG	1:B:94:GLU:HG3	2.35	0.56
2:O:303:MPD:C5	2:O:303:MPD:H12	2.35	0.56
1:P:71:GLU:O	1:P:71:GLU:HG2	2.05	0.56
1:U:85:LYS:HB2	1:U:250:GLU:HB3	1.87	0.56
1:Y:85:LYS:HB2	1:Y:250:GLU:HB3	1.87	0.56
1:I:85:LYS:HB2	1:I:250:GLU:HB3	1.87	0.56
1:b:85:LYS:HB2	1:b:250:GLU:HB3	1.87	0.56
1:A:14:ILE:HD13	1:A:52:LEU:HD22	1.87	0.56
1:F:85:LYS:HB2	1:F:250:GLU:HB3	1.88	0.56
1:b:281:ARG:HG2	1:b:294:LEU:HD12	1.86	0.56
1:G:191:TYR:CE2	1:G:200:ARG:HB3	2.41	0.56
1:V:179:TRP:CE3	2:V:303:MPD:H51	2.41	0.56
1:b:65:TYR:CE2	1:b:78:LEU:HD21	2.41	0.56
1:D:135:LEU:C	1:D:135:LEU:HD23	2.31	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:169:VAL:HG21	1:D:224:PHE:CZ	2.41	0.56
1:U:36:LYS:HZ2	1:U:280:GLU:HG2	1.71	0.56
1:b:169:VAL:HG21	1:b:224:PHE:CZ	2.41	0.56
1:O:169:VAL:HG21	1:O:224:PHE:CZ	2.41	0.56
1:U:65:TYR:CE1	1:U:78:LEU:HD21	2.41	0.56
1:V:169:VAL:HG21	1:V:224:PHE:CZ	2.41	0.56
1:Z:36:LYS:HZ2	1:Z:280:GLU:HG2	1.69	0.56
1:L:85:LYS:HB2	1:L:250:GLU:HB3	1.88	0.56
1:F:169:VAL:HG21	1:F:224:PHE:CZ	2.41	0.56
1:G:85:LYS:HB2	1:G:250:GLU:HB3	1.88	0.56
1:X:14:ILE:HD11	1:Y:39:PHE:HE1	1.70	0.56
1:X:85:LYS:HB2	1:X:250:GLU:HB3	1.88	0.56
1:L:180:GLY:O	2:L:303:MPD:H12	2.06	0.56
1:Q:191:TYR:CZ	1:Q:200:ARG:HD2	2.40	0.55
1:H:85:LYS:HB2	1:H:250:GLU:HB3	1.87	0.55
1:N:169:VAL:HG21	1:N:224:PHE:CZ	2.41	0.55
1:V:135:LEU:HD23	1:V:135:LEU:C	2.32	0.55
1:E:200:ARG:CZ	2:E:302:MPD:C1	2.84	0.55
1:F:74:ASN:O	1:F:259:HIS:HA	2.06	0.55
1:N:200:ARG:CG	2:N:302:MPD:HM3	2.37	0.55
1:N:200:ARG:HG3	2:N:302:MPD:HM3	1.88	0.55
1:Q:85:LYS:HB2	1:Q:250:GLU:HB3	1.89	0.55
1:D:294:LEU:O	1:D:295:GLU:HB2	2.07	0.55
1:O:180:GLY:O	2:O:303:MPD:H13	2.07	0.55
1:G:94:GLU:CD	1:G:243:THR:HG23	2.32	0.55
1:I:281:ARG:HG2	1:I:294:LEU:HD12	1.87	0.55
1:N:85:LYS:HB2	1:N:250:GLU:HB3	1.88	0.55
1:R:14:ILE:HD11	1:S:39:PHE:HE1	1.71	0.55
2:Y:303:MPD:H12	2:Y:303:MPD:C5	2.28	0.55
1:F:191:TYR:CZ	1:F:200:ARG:HD3	2.41	0.55
1:H:169:VAL:HG21	1:H:224:PHE:CZ	2.42	0.55
1:G:260:TRP:HB2	1:G:265:TRP:CZ3	2.42	0.55
1:N:91:PRO:HG2	1:N:94:GLU:HG3	1.89	0.55
1:S:85:LYS:HB2	1:S:250:GLU:HB3	1.88	0.55
1:W:65:TYR:CE1	1:W:78:LEU:HD21	2.42	0.55
1:O:85:LYS:HB2	1:O:250:GLU:HB3	1.89	0.54
1:I:180:GLY:O	2:I:303:MPD:C1	2.45	0.54
1:M:183:ASP:HB2	4:M:869:HOH:O	2.07	0.54
1:N:179:TRP:CE3	2:N:303:MPD:H51	2.42	0.54
1:W:14:ILE:HD11	1:X:39:PHE:HE1	1.72	0.54
1:W:261:THR:HG22	1:W:264:ASN:H	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:169:VAL:HG21	1:I:224:PHE:CZ	2.41	0.54
1:P:85:LYS:HB2	1:P:250:GLU:HB3	1.90	0.54
1:Y:65:TYR:CE1	1:Y:78:LEU:HD21	2.42	0.54
1:B:200:ARG:HE	2:B:302:MPD:H13	1.73	0.54
1:T:85:LYS:HB2	1:T:250:GLU:HB3	1.89	0.54
1:Z:110:LYS:HD2	1:a:175:VAL:HG23	1.90	0.54
1:P:169:VAL:HG21	1:P:224:PHE:CZ	2.42	0.54
1:R:169:VAL:HG21	1:R:224:PHE:CZ	2.43	0.54
1:U:91:PRO:CG	1:U:94:GLU:HG3	2.36	0.54
1:a:85:LYS:HB2	1:a:250:GLU:HB3	1.89	0.54
1:A:135:LEU:HD13	1:A:135:LEU:C	2.32	0.54
1:A:169:VAL:HG21	1:A:224:PHE:CZ	2.43	0.54
1:Q:169:VAL:HG21	1:Q:224:PHE:CZ	2.43	0.54
1:U:253:ARG:NH2	4:U:702:HOH:O	2.34	0.54
1:W:85:LYS:HB2	1:W:250:GLU:HB3	1.90	0.54
1:C:180:GLY:O	2:C:303:MPD:H13	2.08	0.54
1:J:190:VAL:HG12	1:J:264:ASN:ND2	2.23	0.54
1:L:190:VAL:HG13	1:L:191:TYR:CD2	2.42	0.54
1:V:253:ARG:CZ	4:V:731:HOH:O	2.55	0.54
1:Y:169:VAL:HG21	1:Y:224:PHE:CZ	2.43	0.54
1:E:261:THR:HG22	1:E:264:ASN:O	2.08	0.54
1:G:169:VAL:HG21	1:G:224:PHE:CZ	2.42	0.54
1:E:293:ASN:OD1	1:E:294:LEU:N	2.42	0.53
2:T:303:MPD:H52	2:T:303:MPD:O2	2.08	0.53
1:F:200:ARG:HE	2:F:302:MPD:H13	1.72	0.53
1:W:199:THR:HA	2:W:302:MPD:H53	1.90	0.53
1:E:85:LYS:HB2	1:E:250:GLU:HB3	1.91	0.53
1:M:74:ASN:O	1:M:259:HIS:HA	2.08	0.53
1:N:135:LEU:C	1:N:135:LEU:HD23	2.33	0.53
1:D:204:MET:HE3	1:D:209:ASN:HB2	1.90	0.53
1:F:36:LYS:HZ2	1:F:280:GLU:HG2	1.73	0.53
1:P:135:LEU:C	1:P:135:LEU:HD23	2.34	0.53
2:U:302:MPD:O4	2:U:302:MPD:HM3	2.08	0.53
1:H:262:SER:HB2	1:a:268:THR:CG2	2.39	0.53
1:X:169:VAL:HG21	1:X:224:PHE:CZ	2.44	0.53
2:Y:302:MPD:O4	2:Y:302:MPD:H11	2.08	0.53
1:Z:85:LYS:HB2	1:Z:250:GLU:HB3	1.90	0.53
1:G:293:ASN:O	1:G:294:LEU:HB2	2.09	0.53
2:Y:302:MPD:O4	2:Y:302:MPD:HM3	2.07	0.53
1:C:110:LYS:HD2	1:D:175:VAL:HG23	1.90	0.53
1:E:76:SER:HB2	1:E:206:ALA:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:169:VAL:HG21	1:K:224:PHE:CZ	2.44	0.53
1:M:179:TRP:HA	2:M:303:MPD:H53	1.90	0.53
2:U:303:MPD:H12	2:U:303:MPD:C5	2.39	0.53
1:Z:169:VAL:HG21	1:Z:224:PHE:CZ	2.43	0.53
1:b:36:LYS:HZ2	1:b:280:GLU:HG2	1.74	0.53
1:C:169:VAL:HG21	1:C:224:PHE:CZ	2.43	0.53
1:L:169:VAL:HG21	1:L:224:PHE:CZ	2.44	0.53
2:N:303:MPD:O2	2:N:303:MPD:H52	2.09	0.53
1:P:94:GLU:OE1	1:P:243:THR:HG23	2.09	0.53
2:F:302:MPD:O4	2:F:302:MPD:HM3	2.09	0.53
1:I:75:LYS:HD3	1:I:259:HIS:HB3	1.91	0.53
1:M:91:PRO:HD3	1:M:244:ASN:O	2.09	0.53
1:M:169:VAL:HG21	1:M:224:PHE:CZ	2.45	0.53
1:B:91:PRO:HB2	1:B:94:GLU:CG	2.39	0.52
1:B:169:VAL:HG21	1:B:224:PHE:CZ	2.43	0.52
1:N:293:ASN:OD1	1:N:294:LEU:HD23	2.09	0.52
1:W:169:VAL:HG21	1:W:224:PHE:CZ	2.44	0.52
1:F:135:LEU:HD12	1:F:136:ILE:N	2.24	0.52
1:T:70:GLU:OE2	1:T:205:LYS:HE3	2.09	0.52
1:B:180:GLY:O	2:B:303:MPD:H12	2.10	0.52
1:C:85:LYS:HB2	1:C:250:GLU:HB3	1.90	0.52
1:J:85:LYS:HB2	1:J:250:GLU:HB3	1.91	0.52
1:L:129:THR:HG22	1:L:131:LYS:H	1.74	0.52
1:S:200:ARG:NE	2:S:302:MPD:H13	2.24	0.52
1:b:70:GLU:OE2	1:b:205:LYS:HE3	2.10	0.52
1:E:14:ILE:HD11	1:F:39:PHE:HE1	1.74	0.52
1:G:65:TYR:CE1	1:G:78:LEU:HD21	2.44	0.52
1:U:169:VAL:HG21	1:U:224:PHE:CZ	2.45	0.52
2:J:303:MPD:C1	2:J:303:MPD:H52	2.39	0.52
1:O:191:TYR:CZ	1:O:200:ARG:HD3	2.44	0.52
1:R:281:ARG:HG2	1:R:294:LEU:HD12	1.90	0.52
1:U:179:TRP:HA	2:U:303:MPD:H53	1.91	0.52
1:S:169:VAL:HG21	1:S:224:PHE:CZ	2.44	0.52
2:J:302:MPD:O4	2:J:302:MPD:CM	2.58	0.52
1:B:14:ILE:HD11	1:B:48:HIS:CE1	2.44	0.52
1:B:68:TYR:CE2	1:B:69:SER:HB2	2.44	0.52
1:E:169:VAL:HG21	1:E:224:PHE:CZ	2.44	0.52
1:H:36:LYS:HZ2	1:H:280:GLU:HG2	1.74	0.52
1:A:127:ASP:HA	1:B:133:GLY:HA2	1.92	0.52
1:D:85:LYS:HB2	1:D:250:GLU:HB3	1.91	0.52
1:L:36:LYS:HZ2	1:L:280:GLU:HG2	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:293:ASN:O	1:Q:294:LEU:HB2	2.09	0.52
1:X:198:LYS:HG2	1:X:199:THR:HG23	1.92	0.52
1:a:200:ARG:NH2	2:a:302:MPD:H11	2.25	0.52
1:E:188:ASN:HB3	1:E:191:TYR:O	2.10	0.52
1:N:125:THR:HG23	4:N:1671:HOH:O	2.10	0.52
2:W:303:MPD:H12	2:W:303:MPD:C5	2.39	0.51
1:X:68:TYR:CG	1:X:69:SER:N	2.77	0.51
1:a:169:VAL:HG21	1:a:224:PHE:CZ	2.44	0.51
1:A:36:LYS:HZ2	1:A:280:GLU:HG2	1.74	0.51
1:B:26:VAL:HG22	1:B:37:LYS:HG2	1.92	0.51
1:H:200:ARG:NH2	2:H:302:MPD:C1	2.73	0.51
1:N:200:ARG:HE	2:N:302:MPD:CM	2.24	0.51
1:A:91:PRO:HD3	1:A:244:ASN:O	2.09	0.51
1:U:91:PRO:HB2	1:U:94:GLU:CG	2.38	0.51
1:V:294:LEU:O	1:V:295:GLU:CB	2.59	0.51
1:D:261:THR:HB	1:D:264:ASN:O	2.10	0.51
1:E:204:MET:HE3	1:E:209:ASN:HB2	1.93	0.51
1:I:36:LYS:HZ2	1:I:280:GLU:HG2	1.74	0.51
1:O:198:LYS:O	2:O:302:MPD:H52	2.10	0.51
1:T:169:VAL:HG21	1:T:224:PHE:CZ	2.44	0.51
1:T:135:LEU:HD23	1:T:136:ILE:N	2.26	0.51
1:Z:91:PRO:CG	1:Z:94:GLU:HG3	2.41	0.51
1:C:188:ASN:OD1	1:C:190:VAL:HG13	2.11	0.51
1:P:179:TRP:CE3	2:P:303:MPD:H51	2.46	0.51
1:b:261:THR:HG23	1:b:263:THR:H	1.75	0.51
1:M:261:THR:HG22	1:M:262:SER:N	2.26	0.51
1:Q:91:PRO:HD3	1:Q:244:ASN:O	2.11	0.51
1:R:257:GLN:O	1:R:267:GLY:HA2	2.11	0.51
2:U:303:MPD:H12	2:U:303:MPD:H53	1.93	0.51
1:A:85:LYS:HB2	1:A:250:GLU:HB3	1.92	0.50
1:R:94:GLU:OE2	1:R:240:LYS:HE3	2.10	0.50
1:X:200:ARG:HG3	2:X:302:MPD:H13	1.92	0.50
2:V:303:MPD:HM1	1:a:116:LEU:HB3	1.93	0.50
1:B:190:VAL:HG23	1:B:191:TYR:CD1	2.45	0.50
1:D:124:VAL:HG11	1:U:262:SER:HB3	1.93	0.50
1:J:36:LYS:HZ2	1:J:280:GLU:HG2	1.77	0.50
1:K:135:LEU:HD13	1:K:136:ILE:N	2.26	0.50
2:L:302:MPD:H52	2:L:302:MPD:O2	2.11	0.50
1:N:191:TYR:CE1	1:N:200:ARG:HB3	2.47	0.50
1:U:73:ALA:HB2	1:U:205:LYS:HE3	1.94	0.50
1:V:180:GLY:O	2:V:303:MPD:H12	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:263:THR:HG23	1:S:264:ASN:OD1	2.12	0.50
1:V:94:GLU:HG3	4:V:1657:HOH:O	2.11	0.50
1:F:113:MET:HA	1:G:146:LEU:O	2.12	0.50
1:S:74:ASN:C	1:S:75:LYS:HD2	2.37	0.50
1:V:28:TYR:CZ	1:b:161:THR:HG22	2.47	0.50
1:Q:198:LYS:HE3	1:Q:210:PHE:O	2.12	0.50
1:R:36:LYS:HZ2	1:R:280:GLU:HG2	1.76	0.50
1:W:95:VAL:HG11	1:W:238:ALA:HB1	1.94	0.50
1:V:44:ASP:HB3	4:V:2007:HOH:O	2.11	0.50
1:b:26:VAL:HG22	1:b:37:LYS:HG2	1.94	0.50
1:K:110:LYS:HD2	1:L:175:VAL:HG23	1.92	0.50
1:S:261:THR:O	1:S:262:SER:C	2.54	0.50
1:U:281:ARG:HG2	1:U:294:LEU:HD12	1.92	0.50
1:W:74:ASN:O	1:W:259:HIS:HA	2.12	0.50
1:Y:71:GLU:HB3	1:Y:75:LYS:HB2	1.94	0.50
1:R:113:MET:HA	1:S:146:LEU:O	2.12	0.50
1:S:26:VAL:HG22	1:S:37:LYS:HG2	1.93	0.50
1:Y:36:LYS:HZ2	1:Y:280:GLU:HG2	1.76	0.50
1:G:36:LYS:NZ	1:G:280:GLU:HG2	2.27	0.49
1:J:253:ARG:NH2	4:J:1212:HOH:O	2.44	0.49
1:P:66:ARG:C	1:P:78:LEU:HD12	2.37	0.49
1:P:73:ALA:HB1	1:P:205:LYS:HG3	1.94	0.49
1:V:253:ARG:HD3	4:V:731:HOH:O	2.11	0.49
1:b:135:LEU:HD23	1:b:136:ILE:N	2.27	0.49
1:E:293:ASN:O	1:E:294:LEU:HB2	2.12	0.49
1:H:281:ARG:HG2	1:H:294:LEU:HD12	1.94	0.49
1:J:132:ILE:HG21	1:R:262:SER:HB3	1.94	0.49
1:J:169:VAL:HG21	1:J:224:PHE:CZ	2.46	0.49
1:V:222:SER:HB3	1:b:157:LEU:O	2.12	0.49
1:I:114:SER:CB	2:K:303:MPD:HM2	2.42	0.49
1:O:8:LYS:HD2	1:P:13:ASP:HB2	1.93	0.49
1:F:68:TYR:CE2	1:F:69:SER:HB3	2.47	0.49
1:L:200:ARG:NH2	2:L:302:MPD:C3	2.75	0.49
1:K:85:LYS:HB2	1:K:250:GLU:HB3	1.94	0.49
1:N:91:PRO:HG2	1:N:94:GLU:CG	2.42	0.49
1:N:191:TYR:CE1	1:N:200:ARG:HD3	2.47	0.49
1:b:191:TYR:CZ	1:b:200:ARG:HD3	2.47	0.49
1:B:36:LYS:HZ2	1:B:280:GLU:HG2	1.76	0.49
1:E:26:VAL:HG22	1:E:37:LYS:HG2	1.94	0.49
1:I:114:SER:HB3	2:K:303:MPD:HM2	1.93	0.49
1:J:110:LYS:HD2	1:K:175:VAL:HG23	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:113:MET:HA	1:N:146:LEU:O	2.12	0.49
1:Q:110:LYS:HD2	1:R:175:VAL:HG23	1.94	0.49
1:U:191:TYR:CZ	1:U:200:ARG:HD3	2.47	0.49
1:Y:293:ASN:O	1:Y:294:LEU:HB2	2.12	0.49
1:I:36:LYS:NZ	1:I:280:GLU:HG2	2.28	0.49
1:P:65:TYR:CE2	1:P:78:LEU:HD21	2.47	0.49
1:T:36:LYS:NZ	1:T:280:GLU:HG2	2.28	0.49
1:X:261:THR:O	1:X:262:SER:C	2.55	0.49
1:a:8:LYS:HD2	1:b:13:ASP:HB2	1.95	0.49
1:A:268:THR:CG2	1:T:262:SER:HB2	2.43	0.49
1:H:110:LYS:HD2	1:I:175:VAL:HG23	1.95	0.49
1:J:65:TYR:CE1	1:J:78:LEU:HD21	2.48	0.49
1:N:125:THR:CG2	4:N:1671:HOH:O	2.60	0.49
1:R:135:LEU:C	1:R:135:LEU:HD23	2.38	0.49
1:S:204:MET:HE3	1:S:209:ASN:OD1	2.12	0.49
1:V:175:VAL:HG23	1:b:110:LYS:HD2	1.94	0.49
1:X:68:TYR:CE1	1:X:69:SER:HB2	2.48	0.49
1:b:191:TYR:CE2	1:b:200:ARG:HB3	2.48	0.49
1:C:260:TRP:CZ3	1:C:264:ASN:HA	2.48	0.49
1:R:94:GLU:OE2	1:R:243:THR:HA	2.13	0.49
1:V:26:VAL:HG22	1:V:37:LYS:HG2	1.94	0.49
1:B:65:TYR:CE1	1:B:78:LEU:HD21	2.48	0.49
1:E:110:LYS:HD2	1:F:175:VAL:HG23	1.94	0.49
1:K:253:ARG:NH2	4:K:1974:HOH:O	2.45	0.49
1:R:135:LEU:HD23	1:R:136:ILE:N	2.28	0.49
1:T:205:LYS:HE2	1:T:207:ALA:HB3	1.94	0.49
1:C:71:GLU:OE1	1:C:75:LYS:HE3	2.13	0.48
1:L:200:ARG:NH2	2:L:302:MPD:H32	2.28	0.48
1:L:261:THR:HG22	1:L:264:ASN:H	1.78	0.48
1:W:74:ASN:C	1:W:75:LYS:HD2	2.37	0.48
1:X:110:LYS:HD2	1:Y:175:VAL:HG23	1.95	0.48
1:A:175:VAL:HG23	1:G:110:LYS:HD2	1.95	0.48
1:B:183:ASP:HB2	4:B:1231:HOH:O	2.13	0.48
1:L:91:PRO:HD3	1:L:244:ASN:O	2.13	0.48
1:O:26:VAL:HG22	1:O:37:LYS:HG2	1.94	0.48
1:V:110:LYS:HD2	1:W:175:VAL:HG23	1.95	0.48
1:A:222:SER:HB3	1:G:157:LEU:O	2.14	0.48
1:E:188:ASN:O	1:E:191:TYR:O	2.31	0.48
1:F:36:LYS:NZ	1:F:280:GLU:HG2	2.29	0.48
1:P:281:ARG:HG2	1:P:294:LEU:HD12	1.95	0.48
1:Q:69:SER:OG	1:Q:71:GLU:HG2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:135:LEU:HD23	1:X:135:LEU:C	2.38	0.48
1:G:135:LEU:HD23	1:G:135:LEU:C	2.38	0.48
1:K:183:ASP:HB2	4:K:309:HOH:O	2.13	0.48
2:N:302:MPD:H52	2:N:302:MPD:H13	1.96	0.48
1:U:36:LYS:NZ	1:U:280:GLU:HG2	2.29	0.48
1:X:70:GLU:O	1:X:71:GLU:C	2.55	0.48
1:O:175:VAL:HG23	1:U:110:LYS:HD2	1.95	0.48
2:O:303:MPD:H53	2:O:303:MPD:H12	1.95	0.48
1:R:26:VAL:HG22	1:R:37:LYS:HG2	1.95	0.48
1:V:37:LYS:HD2	1:V:37:LYS:C	2.38	0.48
1:W:75:LYS:HE3	1:W:259:HIS:HB2	1.96	0.48
1:X:283:LYS:HB2	1:X:294:LEU:HD21	1.96	0.48
1:L:110:LYS:HD2	1:M:175:VAL:HG23	1.94	0.48
1:S:202:GLY:HA3	1:S:204:MET:HE2	1.95	0.48
1:T:26:VAL:HG22	1:T:37:LYS:HG2	1.94	0.48
1:T:91:PRO:HB2	1:T:94:GLU:CG	2.41	0.48
1:V:36:LYS:HZ2	1:V:280:GLU:HG2	1.79	0.48
1:V:236:ARG:O	4:V:374:HOH:O	2.20	0.48
1:Y:26:VAL:HG22	1:Y:37:LYS:HG2	1.94	0.48
1:C:183:ASP:HB2	4:C:1668:HOH:O	2.13	0.48
1:P:113:MET:HA	1:Q:146:LEU:O	2.13	0.48
1:S:191:TYR:HE1	1:S:200:ARG:HB3	1.77	0.48
1:W:26:VAL:HG22	1:W:37:LYS:HG2	1.95	0.48
1:X:16:SER:O	1:X:18:THR:HG23	2.13	0.48
1:Y:116:LEU:CB	2:a:303:MPD:H11	2.43	0.48
1:a:113:MET:HA	1:b:146:LEU:O	2.12	0.48
1:Z:135:LEU:C	1:Z:135:LEU:HD23	2.38	0.48
1:a:36:LYS:NZ	1:a:280:GLU:HG2	2.29	0.48
1:b:36:LYS:NZ	1:b:280:GLU:HG2	2.29	0.48
1:H:222:SER:HB3	1:N:157:LEU:O	2.14	0.48
1:a:91:PRO:HD3	1:a:244:ASN:O	2.13	0.48
1:H:268:THR:CG2	1:a:262:SER:HB2	2.44	0.48
1:W:36:LYS:NZ	1:W:280:GLU:HG2	2.29	0.48
1:A:26:VAL:HG22	1:A:37:LYS:HG2	1.96	0.47
1:K:36:LYS:HZ2	1:K:280:GLU:HG2	1.78	0.47
1:K:124:VAL:HG11	1:b:262:SER:CB	2.43	0.47
1:V:68:TYR:O	1:V:69:SER:HB3	2.14	0.47
1:H:175:VAL:HG23	1:N:110:LYS:HD2	1.95	0.47
1:K:26:VAL:HG22	1:K:37:LYS:HG2	1.95	0.47
1:M:26:VAL:HG22	1:M:37:LYS:HG2	1.95	0.47
1:N:26:VAL:HG22	1:N:37:LYS:HG2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:36:LYS:NZ	1:N:280:GLU:HG2	2.29	0.47
1:P:26:VAL:HG22	1:P:37:LYS:HG2	1.95	0.47
1:Y:131:LYS:C	1:Y:132:ILE:HD12	2.40	0.47
1:C:95:VAL:HG12	4:C:712:HOH:O	2.14	0.47
1:D:26:VAL:HG22	1:D:37:LYS:HG2	1.95	0.47
1:J:93:ASN:OD1	1:J:94:GLU:HG3	2.14	0.47
1:O:294:LEU:O	1:O:295:GLU:CB	2.61	0.47
1:Q:36:LYS:NZ	1:Q:280:GLU:HG2	2.30	0.47
1:U:260:TRP:NE1	1:U:262:SER:HA	2.29	0.47
1:Z:91:PRO:CG	1:Z:94:GLU:CG	2.91	0.47
1:Z:260:TRP:HB2	1:Z:265:TRP:CZ3	2.50	0.47
1:E:90:LEU:HB3	1:E:91:PRO:HD2	1.96	0.47
1:E:192:GLY:HA3	1:E:267:GLY:O	2.15	0.47
1:L:14:ILE:HD11	1:L:48:HIS:CE1	2.50	0.47
1:M:200:ARG:HE	2:M:302:MPD:HM2	1.79	0.47
1:O:65:TYR:CE2	1:O:78:LEU:HD21	2.49	0.47
1:Z:26:VAL:HG22	1:Z:37:LYS:HG2	1.95	0.47
1:a:253:ARG:NH2	4:a:565:HOH:O	2.31	0.47
1:B:91:PRO:HB2	1:B:94:GLU:HG3	1.97	0.47
1:E:261:THR:HG22	1:E:264:ASN:H	1.79	0.47
1:F:26:VAL:HG22	1:F:37:LYS:HG2	1.96	0.47
1:G:26:VAL:HG22	1:G:37:LYS:HG2	1.96	0.47
1:G:94:GLU:OE2	1:G:243:THR:HG23	2.14	0.47
1:M:56:ARG:NH2	4:M:315:HOH:O	2.47	0.47
1:O:135:LEU:C	1:O:135:LEU:HD23	2.39	0.47
1:W:113:MET:HA	1:X:146:LEU:O	2.15	0.47
1:Z:36:LYS:NZ	1:Z:280:GLU:HG2	2.30	0.47
1:B:36:LYS:NZ	1:B:280:GLU:HG2	2.30	0.47
1:C:129:THR:HG22	1:C:129:THR:O	2.14	0.47
1:D:116:LEU:HB2	2:F:303:MPD:HM3	1.96	0.47
2:D:302:MPD:HM1	2:D:302:MPD:H53	1.96	0.47
1:H:17:ASN:N	4:H:1998:HOH:O	2.45	0.47
1:L:26:VAL:HG22	1:L:37:LYS:HG2	1.97	0.47
1:L:37:LYS:HD2	1:L:37:LYS:C	2.39	0.47
1:N:75:LYS:HD2	1:N:259:HIS:CB	2.45	0.47
1:O:39:PHE:HE1	1:U:14:ILE:HD11	1.80	0.47
1:V:18:THR:HG21	1:V:45:ASP:HA	1.96	0.47
1:W:14:ILE:HD11	1:X:39:PHE:CE1	2.50	0.47
1:a:200:ARG:CZ	2:a:302:MPD:C1	2.93	0.47
1:Q:36:LYS:HZ2	1:Q:280:GLU:HG2	1.79	0.47
1:Q:135:LEU:HD23	1:Q:136:ILE:N	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:110:LYS:HD2	1:T:175:VAL:HG23	1.97	0.47
1:T:90:LEU:HD12	1:T:245:ILE:HG13	1.95	0.47
1:X:161:THR:HG22	1:Y:28:TYR:CZ	2.50	0.47
1:I:132:ILE:HG22	1:I:133:GLY:N	2.30	0.47
1:R:36:LYS:NZ	1:R:280:GLU:HG2	2.29	0.47
1:X:36:LYS:NZ	1:X:280:GLU:HG2	2.30	0.47
1:X:191:TYR:CZ	1:X:200:ARG:HD3	2.50	0.47
1:A:36:LYS:NZ	1:A:280:GLU:HG2	2.29	0.47
1:E:197:MET:O	2:E:302:MPD:H52	2.15	0.47
1:L:66:ARG:C	1:L:78:LEU:HD12	2.40	0.47
1:O:14:ILE:HD11	1:P:39:PHE:HE1	1.80	0.47
1:U:191:TYR:CE1	1:U:200:ARG:HD3	2.50	0.47
2:V:303:MPD:HM1	1:a:116:LEU:CB	2.45	0.47
1:P:227:ASP:OD1	4:P:2010:HOH:O	2.21	0.46
1:V:179:TRP:HA	2:V:303:MPD:C5	2.41	0.46
1:W:258:LEU:HA	1:W:266:LYS:O	2.15	0.46
1:A:110:LYS:HD2	1:B:175:VAL:HG23	1.98	0.46
1:B:188:ASN:OD1	1:B:190:VAL:HG22	2.14	0.46
1:D:37:LYS:C	1:D:37:LYS:HD2	2.40	0.46
1:D:294:LEU:O	1:D:295:GLU:CB	2.63	0.46
1:U:26:VAL:HG22	1:U:37:LYS:HG2	1.97	0.46
1:V:126:GLY:HA2	1:V:131:LYS:O	2.16	0.46
1:X:200:ARG:CZ	2:X:302:MPD:H11	2.45	0.46
1:B:37:LYS:HD2	1:B:37:LYS:C	2.41	0.46
1:C:26:VAL:HG22	1:C:37:LYS:HG2	1.95	0.46
1:D:36:LYS:NZ	1:D:280:GLU:HG2	2.29	0.46
1:C:36:LYS:NZ	1:C:280:GLU:HG2	2.28	0.46
1:C:157:LEU:O	1:D:222:SER:HB3	2.15	0.46
1:I:26:VAL:HG22	1:I:37:LYS:HG2	1.97	0.46
1:J:36:LYS:HD2	4:J:527:HOH:O	2.16	0.46
1:K:37:LYS:C	1:K:37:LYS:HD2	2.41	0.46
1:M:36:LYS:NZ	1:M:280:GLU:HG2	2.31	0.46
1:N:263:THR:HG23	1:N:264:ASN:OD1	2.16	0.46
1:P:36:LYS:NZ	1:P:280:GLU:HG2	2.31	0.46
1:P:37:LYS:C	1:P:37:LYS:HD2	2.40	0.46
1:Q:26:VAL:HG22	1:Q:37:LYS:HG2	1.96	0.46
1:Q:183:ASP:HB2	4:Q:1535:HOH:O	2.13	0.46
1:A:123:ASN:OD1	1:A:123:ASN:C	2.59	0.46
1:E:36:LYS:NZ	1:E:280:GLU:HG2	2.30	0.46
1:L:36:LYS:NZ	1:L:280:GLU:HG2	2.31	0.46
1:V:39:PHE:HE1	1:b:14:ILE:HD11	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:192:GLY:H	1:E:266:LYS:HG3	1.80	0.46
2:Q:302:MPD:HM1	2:Q:302:MPD:H4	1.69	0.46
1:R:179:TRP:CE3	2:R:303:MPD:H51	2.51	0.46
2:R:303:MPD:H12	2:R:303:MPD:C5	2.41	0.46
1:V:253:ARG:NE	4:V:731:HOH:O	2.48	0.46
1:W:66:ARG:C	1:W:78:LEU:HD12	2.41	0.46
1:A:183:ASP:HB2	4:A:313:HOH:O	2.15	0.46
1:F:37:LYS:HD2	1:F:37:LYS:C	2.41	0.46
1:G:36:LYS:HZ2	1:G:280:GLU:HG2	1.80	0.46
1:H:36:LYS:NZ	1:H:280:GLU:HG2	2.30	0.46
1:K:36:LYS:NZ	1:K:280:GLU:HG2	2.31	0.46
1:X:37:LYS:HD2	1:X:37:LYS:C	2.41	0.46
1:X:277:ARG:NE	4:X:728:HOH:O	2.45	0.46
1:a:26:VAL:HG22	1:a:37:LYS:HG2	1.97	0.46
1:a:188:ASN:OD1	1:a:190:VAL:HG13	2.15	0.46
1:b:91:PRO:HB2	1:b:94:GLU:CG	2.42	0.46
1:A:122:GLY:HA2	1:A:135:LEU:O	2.16	0.46
2:A:302:MPD:HM3	2:A:302:MPD:C5	2.37	0.46
1:Q:161:THR:HG22	1:R:28:TYR:CZ	2.51	0.46
1:T:135:LEU:HD23	1:T:135:LEU:C	2.40	0.46
1:W:37:LYS:HD2	1:W:37:LYS:C	2.40	0.46
1:X:74:ASN:O	1:X:259:HIS:HA	2.15	0.46
1:a:123:ASN:C	1:a:123:ASN:OD1	2.59	0.46
1:D:123:ASN:C	1:D:123:ASN:OD1	2.59	0.46
1:H:14:ILE:HD13	1:H:52:LEU:HD22	1.97	0.46
1:H:26:VAL:HG22	1:H:37:LYS:HG2	1.98	0.46
1:J:26:VAL:HG22	1:J:37:LYS:HG2	1.96	0.46
1:V:253:ARG:CD	4:V:731:HOH:O	2.64	0.46
1:W:49:ASN:OD1	1:W:50:LYS:HG2	2.16	0.46
1:X:36:LYS:HZ2	1:X:280:GLU:HG2	1.80	0.46
1:D:49:ASN:OD1	1:D:50:LYS:HG2	2.15	0.46
1:P:161:THR:HG22	1:Q:28:TYR:CZ	2.51	0.46
2:S:303:MPD:H12	2:S:303:MPD:C5	2.27	0.46
1:Z:37:LYS:C	1:Z:37:LYS:HD2	2.41	0.46
1:Z:65:TYR:CE2	1:Z:78:LEU:HD21	2.51	0.46
1:H:14:ILE:HG12	1:H:15:GLY:N	2.31	0.45
1:Q:157:LEU:O	1:R:222:SER:HB3	2.16	0.45
2:W:303:MPD:H52	2:W:303:MPD:O2	2.16	0.45
1:H:65:TYR:CE1	1:H:78:LEU:HD21	2.52	0.45
2:N:302:MPD:H52	2:N:302:MPD:C1	2.46	0.45
1:S:37:LYS:C	1:S:37:LYS:HD2	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:W:122:GLY:HA2	1:W:135:LEU:O	2.15	0.45
1:Y:179:TRP:CZ3	1:Y:200:ARG:NH1	2.84	0.45
1:B:191:TYR:CZ	1:B:200:ARG:HD3	2.51	0.45
1:E:248:ILE:N	1:E:248:ILE:HD12	2.32	0.45
1:Q:37:LYS:HD2	1:Q:37:LYS:C	2.42	0.45
1:Y:36:LYS:NZ	1:Y:280:GLU:HG2	2.31	0.45
2:a:302:MPD:O4	2:a:302:MPD:HM3	2.17	0.45
1:B:113:MET:HA	1:C:146:LEU:O	2.16	0.45
1:C:65:TYR:CE1	1:C:78:LEU:HD21	2.51	0.45
1:D:113:MET:HA	1:E:146:LEU:O	2.16	0.45
1:H:176:ASN:ND2	2:H:302:MPD:HM1	2.32	0.45
1:P:157:LEU:O	1:Q:222:SER:HB3	2.16	0.45
1:S:157:LEU:O	1:T:222:SER:HB3	2.16	0.45
1:T:123:ASN:OD1	1:T:123:ASN:C	2.60	0.45
1:V:10:GLY:HA2	1:V:13:ASP:OD2	2.16	0.45
1:a:135:LEU:C	1:a:135:LEU:HD23	2.41	0.45
1:H:91:PRO:HD3	1:H:244:ASN:O	2.16	0.45
1:I:135:LEU:C	1:I:135:LEU:HD23	2.42	0.45
1:K:205:LYS:HG2	1:K:207:ALA:H	1.81	0.45
1:N:200:ARG:HE	2:N:302:MPD:HM3	1.81	0.45
1:X:26:VAL:HG22	1:X:37:LYS:HG2	1.97	0.45
1:b:91:PRO:HD3	1:b:244:ASN:O	2.16	0.45
1:B:123:ASN:OD1	1:B:123:ASN:C	2.60	0.45
1:J:70:GLU:HG3	1:J:207:ALA:HB2	1.99	0.45
1:M:73:ALA:HA	1:M:205:LYS:CE	2.46	0.45
1:W:74:ASN:O	1:W:75:LYS:HD2	2.16	0.45
1:Y:110:LYS:HD2	1:Z:175:VAL:HG23	1.99	0.45
1:Z:67:VAL:HG23	1:Z:67:VAL:O	2.16	0.45
1:A:65:TYR:CE2	1:A:78:LEU:HD21	2.52	0.45
1:I:123:ASN:OD1	1:I:123:ASN:C	2.60	0.45
2:J:303:MPD:H52	2:J:303:MPD:H12	1.98	0.45
1:O:261:THR:O	1:O:262:SER:C	2.60	0.45
1:Z:14:ILE:HD11	1:a:39:PHE:HE1	1.82	0.45
2:Z:302:MPD:H52	2:Z:302:MPD:H13	1.99	0.45
1:J:161:THR:HG22	1:K:28:TYR:CZ	2.51	0.45
1:L:293:ASN:O	1:L:294:LEU:HB2	2.16	0.45
1:S:7:ILE:HA	1:T:14:ILE:O	2.17	0.45
1:T:261:THR:HG22	1:T:263:THR:H	1.81	0.45
1:D:14:ILE:CD1	1:D:43:ILE:HG21	2.47	0.45
1:M:70:GLU:OE2	1:M:205:LYS:HE2	2.17	0.45
1:N:18:THR:HG21	1:N:45:ASP:HA	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:294:LEU:O	1:O:295:GLU:HB2	2.17	0.45
1:V:36:LYS:NZ	1:V:280:GLU:HG2	2.32	0.45
1:V:283:LYS:HB2	1:V:294:LEU:HD21	1.99	0.45
1:X:14:ILE:HD11	1:Y:39:PHE:CE1	2.49	0.45
1:A:89:GLN:HG3	1:A:164:LYS:HG2	1.99	0.45
1:K:200:ARG:NH2	2:K:302:MPD:C1	2.80	0.45
2:N:303:MPD:H12	2:N:303:MPD:C5	2.44	0.45
1:S:36:LYS:NZ	1:S:280:GLU:HG2	2.32	0.45
1:Y:113:MET:HA	1:Z:146:LEU:O	2.17	0.45
1:J:36:LYS:NZ	1:J:280:GLU:HG2	2.31	0.44
1:L:71:GLU:HG3	1:L:71:GLU:O	2.16	0.44
1:S:135:LEU:HD23	1:S:135:LEU:C	2.41	0.44
1:V:131:LYS:CE	1:b:127:ASP:HB2	2.44	0.44
1:Z:277:ARG:NE	4:Z:1846:HOH:O	2.50	0.44
1:E:261:THR:HG23	1:E:263:THR:H	1.83	0.44
1:M:261:THR:HG22	1:M:262:SER:H	1.80	0.44
1:O:18:THR:HG21	1:O:45:ASP:HA	1.98	0.44
1:C:95:VAL:HG22	4:C:1924:HOH:O	2.17	0.44
1:I:37:LYS:C	1:I:37:LYS:HD2	2.43	0.44
1:K:294:LEU:O	1:K:295:GLU:CB	2.65	0.44
1:S:116:LEU:C	1:S:116:LEU:HD13	2.42	0.44
1:T:37:LYS:HD2	1:T:37:LYS:C	2.42	0.44
1:O:28:TYR:CZ	1:U:161:THR:HG22	2.52	0.44
1:V:18:THR:HG21	1:V:44:ASP:O	2.18	0.44
1:A:281:ARG:HG2	1:A:294:LEU:HD12	2.00	0.44
1:B:157:LEU:O	1:C:222:SER:HB3	2.18	0.44
1:J:135:LEU:C	1:J:135:LEU:HD23	2.43	0.44
1:W:89:GLN:HG3	1:W:164:LYS:HG2	1.99	0.44
1:b:191:TYR:CE2	1:b:200:ARG:HD3	2.53	0.44
2:G:302:MPD:O4	2:G:302:MPD:HM3	2.17	0.44
1:O:37:LYS:HD2	1:O:37:LYS:C	2.42	0.44
1:R:74:ASN:O	1:R:259:HIS:HA	2.17	0.44
1:T:191:TYR:CZ	1:T:200:ARG:HD3	2.52	0.44
1:Y:205:LYS:HG3	1:Y:207:ALA:H	1.83	0.44
1:G:191:TYR:CZ	1:G:200:ARG:HD3	2.53	0.44
1:Q:200:ARG:NE	2:Q:302:MPD:HM1	2.33	0.44
1:R:91:PRO:CG	1:R:94:GLU:HG3	2.48	0.44
1:U:261:THR:O	1:U:263:THR:N	2.51	0.44
1:E:206:ALA:C	1:E:208:GLU:H	2.26	0.44
1:F:293:ASN:O	1:F:294:LEU:CB	2.65	0.44
1:J:74:ASN:O	1:J:259:HIS:HA	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:37:LYS:HD2	1:E:37:LYS:C	2.42	0.44
1:S:123:ASN:C	1:S:123:ASN:OD1	2.61	0.44
1:S:261:THR:HG22	1:S:264:ASN:H	1.83	0.44
1:W:70:GLU:HB3	1:W:207:ALA:HB2	2.00	0.44
1:b:135:LEU:HD23	1:b:135:LEU:C	2.43	0.44
1:A:135:LEU:HD23	1:G:125:THR:CG2	2.47	0.43
1:C:37:LYS:HD2	1:C:37:LYS:C	2.41	0.43
1:J:123:ASN:OD1	1:J:123:ASN:C	2.61	0.43
1:K:14:ILE:HD11	1:L:39:PHE:HE1	1.83	0.43
1:N:65:TYR:CE1	1:N:78:LEU:HD21	2.53	0.43
1:O:222:SER:HB3	1:U:157:LEU:O	2.18	0.43
2:a:303:MPD:H52	2:a:303:MPD:H13	2.00	0.43
1:A:37:LYS:HD2	1:A:37:LYS:C	2.42	0.43
1:A:262:SER:HB2	1:T:268:THR:CG2	2.47	0.43
1:B:68:TYR:CD2	1:B:69:SER:HB2	2.53	0.43
1:C:93:ASN:ND2	1:C:94:GLU:HG3	2.33	0.43
1:R:75:LYS:HG2	1:R:76:SER:N	2.33	0.43
1:V:123:ASN:OD1	1:V:123:ASN:C	2.60	0.43
1:X:18:THR:HG21	1:X:45:ASP:HA	2.00	0.43
1:B:91:PRO:HB2	1:B:94:GLU:HG2	1.99	0.43
1:D:289:GLU:O	4:D:311:HOH:O	2.21	0.43
1:F:14:ILE:CD1	1:F:43:ILE:HG21	2.48	0.43
1:T:20:VAL:O	4:T:611:HOH:O	2.21	0.43
1:Y:200:ARG:HE	2:Y:302:MPD:H11	1.82	0.43
1:Z:116:LEU:HD13	1:Z:116:LEU:C	2.43	0.43
1:a:37:LYS:C	1:a:37:LYS:HD2	2.43	0.43
1:a:131:LYS:HE2	1:a:131:LYS:HB3	1.88	0.43
1:C:123:ASN:OD1	1:C:123:ASN:C	2.61	0.43
1:C:135:LEU:HD23	1:C:136:ILE:N	2.33	0.43
1:F:253:ARG:NH2	4:F:741:HOH:O	2.44	0.43
2:I:303:MPD:HM1	2:I:303:MPD:H4	1.90	0.43
1:K:191:TYR:CZ	1:K:200:ARG:HD3	2.53	0.43
1:N:248:ILE:HD12	1:N:248:ILE:N	2.33	0.43
1:R:91:PRO:HG2	1:R:94:GLU:HG3	2.00	0.43
1:U:123:ASN:C	1:U:123:ASN:OD1	2.61	0.43
1:V:95:VAL:HG11	1:V:238:ALA:HB1	2.00	0.43
2:V:303:MPD:H52	2:V:303:MPD:O2	2.19	0.43
1:C:116:LEU:HD13	1:C:116:LEU:C	2.44	0.43
1:X:116:LEU:HB2	2:X:304:MPD:CM	2.45	0.43
1:Y:89:GLN:HG3	1:Y:164:LYS:HG2	1.99	0.43
1:b:37:LYS:HD2	1:b:37:LYS:C	2.42	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:49:ASN:OD1	1:E:50:LYS:HG2	2.18	0.43
1:G:122:GLY:HA2	1:G:135:LEU:O	2.17	0.43
1:H:283:LYS:HB2	1:H:294:LEU:HD21	2.01	0.43
1:J:116:LEU:C	1:J:116:LEU:HD13	2.44	0.43
1:K:135:LEU:HD13	1:K:135:LEU:C	2.43	0.43
1:R:113:MET:SD	1:S:147:LYS:HE2	2.58	0.43
1:S:122:GLY:HA2	1:S:135:LEU:O	2.19	0.43
1:U:248:ILE:HD12	1:U:248:ILE:N	2.34	0.43
1:V:17:ASN:CG	1:V:17:ASN:O	2.62	0.43
1:W:123:ASN:OD1	1:W:123:ASN:C	2.62	0.43
1:Z:123:ASN:C	1:Z:123:ASN:OD1	2.61	0.43
1:b:200:ARG:HE	2:b:302:MPD:H11	1.81	0.43
1:E:200:ARG:NH1	2:E:302:MPD:H12	2.34	0.43
1:F:204:MET:HE2	1:F:204:MET:H	1.83	0.43
1:T:261:THR:HB	1:T:264:ASN:O	2.18	0.43
1:D:261:THR:CG2	1:D:262:SER:N	2.82	0.43
1:G:123:ASN:C	1:G:123:ASN:OD1	2.61	0.43
1:I:116:LEU:HD13	1:I:116:LEU:C	2.43	0.43
1:N:75:LYS:HD2	1:N:259:HIS:HB3	2.00	0.43
1:O:116:LEU:C	1:O:116:LEU:HD13	2.44	0.43
1:Q:248:ILE:HD12	1:Q:248:ILE:N	2.34	0.43
1:U:116:LEU:HD13	1:U:116:LEU:C	2.44	0.43
1:U:197:MET:CE	1:U:200:ARG:HG2	2.49	0.43
1:Y:135:LEU:HD23	1:Y:136:ILE:N	2.33	0.43
1:A:248:ILE:N	1:A:248:ILE:HD12	2.33	0.43
1:C:198:LYS:HG2	1:C:199:THR:HG23	2.01	0.43
1:E:161:THR:HG22	1:F:28:TYR:CZ	2.54	0.43
1:F:248:ILE:N	1:F:248:ILE:HD12	2.34	0.43
1:H:146:LEU:O	1:N:113:MET:HA	2.19	0.43
1:H:248:ILE:N	1:H:248:ILE:HD12	2.33	0.43
1:K:123:ASN:OD1	1:K:123:ASN:C	2.62	0.43
1:R:65:TYR:CE1	1:R:78:LEU:HD21	2.53	0.43
1:R:180:GLY:O	2:R:303:MPD:C1	2.44	0.43
1:T:110:LYS:HD2	1:U:175:VAL:HG23	2.01	0.43
1:W:131:LYS:O	1:W:132:ILE:HD12	2.19	0.43
1:X:199:THR:HA	2:X:302:MPD:C4	2.48	0.43
1:a:14:ILE:HD11	1:a:52:LEU:HD22	2.00	0.43
1:C:253:ARG:NH2	4:C:1548:HOH:O	2.51	0.43
2:I:302:MPD:O4	2:I:302:MPD:HM3	2.19	0.43
1:K:70:GLU:HG2	1:K:72:GLY:O	2.19	0.43
1:K:244:ASN:ND2	4:K:1399:HOH:O	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:116:LEU:HD13	1:P:116:LEU:C	2.44	0.43
1:U:91:PRO:HD3	1:U:244:ASN:O	2.18	0.43
1:U:183:ASP:HB2	4:U:1063:HOH:O	2.18	0.43
1:W:180:GLY:O	2:W:303:MPD:H13	2.19	0.43
1:b:123:ASN:OD1	1:b:123:ASN:C	2.62	0.43
1:B:168:LYS:NZ	4:B:612:HOH:O	2.52	0.42
1:D:89:GLN:HG3	1:D:164:LYS:HG2	2.01	0.42
1:H:37:LYS:HD2	1:H:37:LYS:C	2.43	0.42
1:M:37:LYS:HD2	1:M:37:LYS:C	2.44	0.42
1:N:123:ASN:C	1:N:123:ASN:OD1	2.61	0.42
1:U:89:GLN:HG3	1:U:164:LYS:HG2	2.01	0.42
1:W:110:LYS:HD2	1:X:175:VAL:HG23	2.01	0.42
1:Y:123:ASN:OD1	1:Y:123:ASN:C	2.62	0.42
1:Z:204:MET:HE3	1:Z:209:ASN:OD1	2.19	0.42
1:G:183:ASP:HB2	4:G:1948:HOH:O	2.19	0.42
1:U:14:ILE:HG22	1:U:15:GLY:N	2.34	0.42
1:V:116:LEU:C	1:V:116:LEU:HD13	2.44	0.42
1:b:212:ASP:OD1	1:b:213:PRO:HD2	2.19	0.42
1:H:116:LEU:HD13	1:H:116:LEU:C	2.44	0.42
1:H:204:MET:HG2	1:H:205:LYS:O	2.19	0.42
1:L:91:PRO:HB2	1:L:94:GLU:HG3	2.00	0.42
1:Y:37:LYS:HD2	1:Y:37:LYS:C	2.44	0.42
1:Z:89:GLN:HG3	1:Z:164:LYS:HG2	2.02	0.42
1:b:116:LEU:C	1:b:116:LEU:HD13	2.44	0.42
1:D:204:MET:HE3	1:D:209:ASN:CB	2.48	0.42
1:E:179:TRP:CB	2:E:302:MPD:HM3	2.49	0.42
1:I:200:ARG:NH2	2:I:302:MPD:H11	2.33	0.42
1:K:89:GLN:HG3	1:K:164:LYS:HG2	2.02	0.42
1:K:179:TRP:CD1	2:K:302:MPD:HM3	2.54	0.42
1:M:14:ILE:HD12	1:M:52:LEU:HD22	2.01	0.42
1:M:255:ASP:HB3	1:M:270:THR:HB	2.00	0.42
1:R:248:ILE:N	1:R:248:ILE:HD12	2.35	0.42
2:B:302:MPD:O4	2:B:302:MPD:C1	2.62	0.42
1:I:261:THR:O	1:I:262:SER:C	2.62	0.42
1:M:89:GLN:HG3	1:M:164:LYS:HG2	2.01	0.42
1:R:123:ASN:OD1	1:R:123:ASN:C	2.62	0.42
1:T:91:PRO:HD3	1:T:244:ASN:O	2.20	0.42
1:T:248:ILE:N	1:T:248:ILE:HD12	2.35	0.42
1:U:37:LYS:C	1:U:37:LYS:HD2	2.44	0.42
1:W:95:VAL:HG13	1:W:234:MET:SD	2.59	0.42
1:W:200:ARG:NH2	2:W:302:MPD:C1	2.82	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Z:293:ASN:O	1:Z:294:LEU:HB2	2.19	0.42
1:B:66:ARG:C	1:B:78:LEU:HD12	2.45	0.42
1:C:161:THR:HG22	1:D:28:TYR:CZ	2.54	0.42
1:D:122:GLY:HA2	1:D:135:LEU:O	2.20	0.42
1:F:113:MET:SD	1:G:147:LYS:HE2	2.60	0.42
1:N:116:LEU:HD13	1:N:116:LEU:C	2.44	0.42
1:O:123:ASN:OD1	1:O:123:ASN:C	2.61	0.42
1:O:135:LEU:HD23	1:O:136:ILE:N	2.35	0.42
1:P:14:ILE:HD11	1:Q:39:PHE:CE1	2.50	0.42
1:X:248:ILE:HD12	1:X:248:ILE:N	2.35	0.42
1:Y:248:ILE:N	1:Y:248:ILE:HD12	2.35	0.42
2:C:303:MPD:H53	2:C:303:MPD:H12	2.00	0.42
1:F:204:MET:HE3	1:F:209:ASN:HB2	2.01	0.42
1:P:89:GLN:HG3	1:P:164:LYS:HG2	2.01	0.42
1:W:248:ILE:N	1:W:248:ILE:HD12	2.34	0.42
1:A:191:TYR:CZ	1:A:200:ARG:HD3	2.54	0.42
1:B:161:THR:HG22	1:C:28:TYR:CZ	2.55	0.42
1:E:123:ASN:OD1	1:E:123:ASN:C	2.63	0.42
2:H:302:MPD:H13	2:H:302:MPD:O4	2.19	0.42
1:M:95:VAL:HG23	1:M:234:MET:SD	2.60	0.42
1:N:74:ASN:O	1:N:259:HIS:HA	2.19	0.42
1:P:71:GLU:O	1:P:71:GLU:CG	2.67	0.42
1:Y:70:GLU:HG3	1:Y:70:GLU:O	2.19	0.42
1:Y:75:LYS:HG2	1:Y:259:HIS:HB2	2.00	0.42
1:a:248:ILE:HD12	1:a:248:ILE:N	2.35	0.42
1:B:91:PRO:HD3	1:B:244:ASN:O	2.20	0.42
1:M:132:ILE:HD12	1:M:133:GLY:N	2.35	0.42
1:W:157:LEU:O	1:X:222:SER:HB3	2.19	0.42
1:X:157:LEU:O	1:Y:222:SER:HB3	2.20	0.42
1:Z:248:ILE:HD12	1:Z:248:ILE:N	2.35	0.42
1:b:70:GLU:HG2	1:b:72:GLY:O	2.19	0.42
1:E:135:LEU:HD23	1:E:135:LEU:C	2.45	0.42
1:H:14:ILE:CD1	1:H:52:LEU:HD22	2.49	0.42
1:J:37:LYS:HD2	1:J:37:LYS:C	2.44	0.42
1:W:116:LEU:HD13	1:W:116:LEU:C	2.45	0.42
1:W:291:MET:O	4:W:1028:HOH:O	2.21	0.42
1:a:191:TYR:CZ	1:a:200:ARG:HD3	2.55	0.42
1:a:200:ARG:CZ	2:a:302:MPD:H11	2.50	0.42
1:A:116:LEU:HD13	1:A:116:LEU:C	2.45	0.41
1:B:248:ILE:HD12	1:B:248:ILE:N	2.34	0.41
1:F:116:LEU:HD13	1:F:116:LEU:C	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:116:LEU:C	1:G:116:LEU:HD13	2.45	0.41
1:O:36:LYS:NZ	1:O:280:GLU:HG2	2.35	0.41
1:P:126:GLY:HA2	1:P:131:LYS:O	2.20	0.41
1:R:37:LYS:HD2	1:R:37:LYS:C	2.45	0.41
1:b:56:ARG:NH2	4:b:384:HOH:O	2.52	0.41
1:A:135:LEU:HD23	1:G:125:THR:HG22	2.02	0.41
1:C:135:LEU:HD23	1:C:135:LEU:C	2.45	0.41
1:P:123:ASN:C	1:P:123:ASN:OD1	2.63	0.41
1:Q:205:LYS:HB2	1:Q:205:LYS:HE3	1.87	0.41
1:E:193:ASN:N	1:E:267:GLY:O	2.53	0.41
1:F:123:ASN:OD1	1:F:123:ASN:C	2.62	0.41
1:F:161:THR:HG22	1:G:28:TYR:CZ	2.56	0.41
1:K:122:GLY:HA2	1:K:135:LEU:O	2.19	0.41
1:N:89:GLN:HG3	1:N:164:LYS:HG2	2.02	0.41
1:O:91:PRO:HD3	1:O:244:ASN:O	2.20	0.41
1:P:248:ILE:HD12	1:P:248:ILE:N	2.35	0.41
1:V:260:TRP:CD1	1:V:261:THR:O	2.73	0.41
1:Z:49:ASN:OD1	1:Z:50:LYS:HG2	2.21	0.41
1:D:116:LEU:HD13	1:D:116:LEU:C	2.46	0.41
1:I:110:LYS:HD2	1:J:175:VAL:HG23	2.01	0.41
1:Q:65:TYR:CE2	1:Q:78:LEU:HD21	2.54	0.41
1:T:113:MET:HA	1:U:146:LEU:O	2.19	0.41
1:D:205:LYS:H	1:D:205:LYS:HG3	1.64	0.41
1:M:161:THR:HG22	1:N:28:TYR:CZ	2.56	0.41
1:Q:94:GLU:OE2	1:Q:240:LYS:HE3	2.21	0.41
1:b:200:ARG:HE	2:b:302:MPD:H13	1.82	0.41
1:b:248:ILE:HD12	1:b:248:ILE:N	2.36	0.41
1:A:113:MET:HA	1:B:146:LEU:O	2.21	0.41
1:C:248:ILE:N	1:C:248:ILE:HD12	2.35	0.41
1:H:123:ASN:OD1	1:H:123:ASN:C	2.64	0.41
1:I:248:ILE:HD12	1:I:248:ILE:N	2.34	0.41
1:K:135:LEU:C	1:K:135:LEU:CD1	2.94	0.41
1:L:261:THR:CG2	1:L:264:ASN:H	2.34	0.41
1:M:248:ILE:N	1:M:248:ILE:HD12	2.36	0.41
1:U:200:ARG:CZ	2:U:302:MPD:H13	2.51	0.41
1:A:194:GLN:O	1:A:195:LEU:C	2.63	0.41
1:B:212:ASP:OD1	1:B:213:PRO:HD2	2.20	0.41
1:C:91:PRO:HD3	1:C:244:ASN:O	2.20	0.41
1:F:261:THR:O	1:F:262:SER:C	2.63	0.41
1:H:180:GLY:O	2:H:303:MPD:O2	2.38	0.41
1:I:113:MET:HA	1:J:146:LEU:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:293:ASN:O	1:N:294:LEU:CB	2.68	0.41
1:O:39:PHE:CE1	1:U:14:ILE:HD11	2.55	0.41
1:Q:184:ARG:HG3	4:Q:1535:HOH:O	2.20	0.41
1:V:65:TYR:CE2	1:V:78:LEU:HD21	2.56	0.41
1:X:135:LEU:HD23	1:X:136:ILE:N	2.35	0.41
1:M:113:MET:SD	1:N:147:LYS:HE2	2.61	0.41
1:N:122:GLY:HA2	1:N:135:LEU:O	2.20	0.41
1:O:194:GLN:O	1:O:195:LEU:C	2.63	0.41
1:Q:89:GLN:HG3	1:Q:164:LYS:HG2	2.03	0.41
1:T:14:ILE:HA	1:T:15:GLY:HA2	1.85	0.41
1:T:89:GLN:HG3	1:T:164:LYS:HG2	2.01	0.41
1:V:89:GLN:HG3	1:V:164:LYS:HG2	2.02	0.41
1:V:194:GLN:O	1:V:195:LEU:C	2.63	0.41
1:D:248:ILE:HD12	1:D:248:ILE:N	2.34	0.41
1:E:116:LEU:C	1:E:116:LEU:HD13	2.46	0.41
1:F:206:ALA:HA	1:F:265:TRP:CH2	2.56	0.41
1:G:37:LYS:HD2	1:G:37:LYS:C	2.45	0.41
2:G:303:MPD:H12	2:G:303:MPD:C5	2.43	0.41
1:H:113:MET:HA	1:I:146:LEU:O	2.21	0.41
1:J:248:ILE:N	1:J:248:ILE:HD12	2.36	0.41
1:M:123:ASN:OD1	1:M:123:ASN:C	2.63	0.41
1:M:135:LEU:HD23	1:M:135:LEU:C	2.46	0.41
1:M:190:VAL:HG13	1:M:191:TYR:CD1	2.56	0.41
1:O:248:ILE:N	1:O:248:ILE:HD12	2.36	0.41
1:O:261:THR:HG22	1:O:263:THR:H	1.86	0.41
1:T:194:GLN:O	1:T:195:LEU:C	2.64	0.41
1:V:14:ILE:HD11	1:V:48:HIS:CE1	2.54	0.41
1:W:92:ASP:O	1:W:163:LYS:NZ	2.50	0.41
1:X:116:LEU:C	1:X:116:LEU:HD13	2.46	0.41
1:Y:116:LEU:HD13	1:Y:116:LEU:C	2.46	0.41
1:Z:91:PRO:HG2	1:Z:94:GLU:HG2	1.97	0.41
1:C:127:ASP:OD2	1:C:129:THR:HB	2.20	0.41
1:F:89:GLN:HG3	1:F:164:LYS:HG2	2.03	0.41
1:H:89:GLN:HG3	1:H:164:LYS:HG2	2.03	0.41
1:I:113:MET:SD	1:J:147:LYS:HE2	2.61	0.41
1:L:157:LEU:O	1:M:222:SER:HB3	2.21	0.41
1:R:180:GLY:H	2:R:303:MPD:H53	1.86	0.41
1:S:191:TYR:CZ	1:S:200:ARG:HD3	2.56	0.41
1:T:8:LYS:HD2	1:U:13:ASP:CB	2.51	0.41
1:X:89:GLN:HG3	1:X:164:LYS:HG2	2.03	0.41
1:X:199:THR:HA	2:X:302:MPD:H4	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Y:183:ASP:HB2	4:Y:858:HOH:O	2.21	0.41
1:B:110:LYS:HD2	1:C:175:VAL:HG23	2.02	0.40
1:B:116:LEU:HD13	1:B:116:LEU:C	2.45	0.40
2:B:302:MPD:O4	2:B:302:MPD:HM3	2.20	0.40
1:D:202:GLY:HA3	1:D:204:MET:HE1	2.03	0.40
1:E:157:LEU:O	1:F:222:SER:HB3	2.21	0.40
1:E:195:LEU:HD13	1:E:196:PHE:CE1	2.56	0.40
1:F:68:TYR:CE1	1:F:77:GLY:HA3	2.56	0.40
1:F:183:ASP:HB2	4:F:625:HOH:O	2.22	0.40
1:O:89:GLN:HG3	1:O:164:LYS:HG2	2.02	0.40
1:K:65:TYR:CE1	1:K:78:LEU:HD21	2.57	0.40
2:K:302:MPD:H11	2:K:302:MPD:H4	1.83	0.40
1:O:273:LYS:NZ	4:O:716:HOH:O	2.53	0.40
1:S:248:ILE:HD12	1:S:248:ILE:N	2.36	0.40
1:V:95:VAL:HG13	1:V:234:MET:SD	2.61	0.40
1:Z:157:LEU:O	1:a:222:SER:HB3	2.20	0.40
1:b:263:THR:HG23	1:b:264:ASN:OD1	2.21	0.40
1:F:68:TYR:CE1	1:F:77:GLY:C	3.00	0.40
1:G:294:LEU:HD23	1:G:294:LEU:HA	1.96	0.40
1:L:80:TRP:CE2	1:L:196:PHE:HE2	2.40	0.40
1:L:161:THR:HG22	1:M:28:TYR:CZ	2.56	0.40
1:M:294:LEU:O	1:M:295:GLU:HG2	2.21	0.40
1:R:89:GLN:HG3	1:R:164:LYS:HG2	2.02	0.40
1:S:36:LYS:HZ2	1:S:280:GLU:HG2	1.85	0.40
1:T:116:LEU:C	1:T:116:LEU:HD13	2.45	0.40
1:a:183:ASP:HB2	4:a:1308:HOH:O	2.21	0.40
1:E:198:LYS:C	2:E:302:MPD:C5	2.94	0.40
1:G:135:LEU:HD23	1:G:136:ILE:N	2.36	0.40
1:N:37:LYS:C	1:N:37:LYS:HD2	2.46	0.40
1:V:91:PRO:HD3	1:V:244:ASN:O	2.21	0.40
1:W:200:ARG:H	2:W:302:MPD:H52	1.86	0.40
1:F:128:ASP:OD1	1:G:131:LYS:HA	2.21	0.40
1:J:157:LEU:O	1:K:222:SER:HB3	2.21	0.40
1:L:123:ASN:OD1	1:L:123:ASN:C	2.65	0.40
1:U:70:GLU:HG2	1:U:72:GLY:O	2.22	0.40
2:b:303:MPD:H12	2:b:303:MPD:C5	2.45	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	293/302 (97%)	280 (96%)	13 (4%)	0	100	100
1	B	293/302 (97%)	278 (95%)	15 (5%)	0	100	100
1	C	292/302 (97%)	282 (97%)	10 (3%)	0	100	100
1	D	293/302 (97%)	282 (96%)	11 (4%)	0	100	100
1	E	293/302 (97%)	275 (94%)	18 (6%)	0	100	100
1	F	292/302 (97%)	280 (96%)	12 (4%)	0	100	100
1	G	292/302 (97%)	279 (96%)	13 (4%)	0	100	100
1	H	292/302 (97%)	278 (95%)	13 (4%)	1 (0%)	36	46
1	I	292/302 (97%)	280 (96%)	12 (4%)	0	100	100
1	J	292/302 (97%)	278 (95%)	14 (5%)	0	100	100
1	K	293/302 (97%)	284 (97%)	9 (3%)	0	100	100
1	L	292/302 (97%)	277 (95%)	15 (5%)	0	100	100
1	M	293/302 (97%)	282 (96%)	11 (4%)	0	100	100
1	N	292/302 (97%)	281 (96%)	11 (4%)	0	100	100
1	O	293/302 (97%)	278 (95%)	14 (5%)	1 (0%)	36	46
1	P	292/302 (97%)	281 (96%)	11 (4%)	0	100	100
1	Q	292/302 (97%)	282 (97%)	10 (3%)	0	100	100
1	R	292/302 (97%)	279 (96%)	13 (4%)	0	100	100
1	S	293/302 (97%)	280 (96%)	13 (4%)	0	100	100
1	T	292/302 (97%)	276 (94%)	16 (6%)	0	100	100
1	U	293/302 (97%)	283 (97%)	10 (3%)	0	100	100
1	V	293/302 (97%)	282 (96%)	11 (4%)	0	100	100
1	W	292/302 (97%)	279 (96%)	13 (4%)	0	100	100
1	X	292/302 (97%)	281 (96%)	11 (4%)	0	100	100
1	Y	292/302 (97%)	277 (95%)	15 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	Z	292/302 (97%)	280 (96%)	12 (4%)	0	100	100
1	a	292/302 (97%)	279 (96%)	12 (4%)	1 (0%)	36	46
1	b	293/302 (97%)	282 (96%)	11 (4%)	0	100	100
All	All	8187/8456 (97%)	7835 (96%)	349 (4%)	3 (0%)	100	100

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	O	262	SER
1	a	262	SER
1	H	262	SER

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	261/268 (97%)	256 (98%)	5 (2%)	50	69
1	B	261/268 (97%)	257 (98%)	4 (2%)	57	75
1	C	260/268 (97%)	256 (98%)	4 (2%)	57	75
1	D	261/268 (97%)	252 (97%)	9 (3%)	32	49
1	E	261/268 (97%)	249 (95%)	12 (5%)	24	36
1	F	260/268 (97%)	253 (97%)	7 (3%)	39	58
1	G	260/268 (97%)	254 (98%)	6 (2%)	44	63
1	H	260/268 (97%)	255 (98%)	5 (2%)	50	69
1	I	260/268 (97%)	255 (98%)	5 (2%)	50	69
1	J	260/268 (97%)	253 (97%)	7 (3%)	39	58
1	K	261/268 (97%)	257 (98%)	4 (2%)	57	75
1	L	260/268 (97%)	252 (97%)	8 (3%)	35	52
1	M	261/268 (97%)	256 (98%)	5 (2%)	50	69
1	N	260/268 (97%)	253 (97%)	7 (3%)	39	58

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	O	261/268 (97%)	252 (97%)	9 (3%)	32	49
1	P	260/268 (97%)	255 (98%)	5 (2%)	50	69
1	Q	260/268 (97%)	253 (97%)	7 (3%)	39	58
1	R	260/268 (97%)	256 (98%)	4 (2%)	57	75
1	S	261/268 (97%)	253 (97%)	8 (3%)	35	52
1	T	260/268 (97%)	255 (98%)	5 (2%)	50	69
1	U	261/268 (97%)	253 (97%)	8 (3%)	35	52
1	V	261/268 (97%)	255 (98%)	6 (2%)	44	63
1	W	260/268 (97%)	253 (97%)	7 (3%)	39	58
1	X	260/268 (97%)	254 (98%)	6 (2%)	44	63
1	Y	260/268 (97%)	254 (98%)	6 (2%)	44	63
1	Z	260/268 (97%)	254 (98%)	6 (2%)	44	63
1	a	260/268 (97%)	254 (98%)	6 (2%)	44	63
1	b	261/268 (97%)	253 (97%)	8 (3%)	35	52
All	All	7291/7504 (97%)	7112 (98%)	179 (2%)	42	60

All (179) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LYS
1	A	14	ILE
1	A	37	LYS
1	A	101	TYR
1	A	132	ILE
1	B	8	LYS
1	B	37	LYS
1	B	92	ASP
1	B	263	THR
1	C	8	LYS
1	C	17	ASN
1	C	37	LYS
1	C	190	VAL
1	D	8	LYS
1	D	37	LYS
1	D	101	TYR
1	D	195	LEU
1	D	201	ASN

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Mol	Chain	Res	Type
1	D	204	MET
1	D	205	LYS
1	D	261	THR
1	D	295	GLU
1	E	8	LYS
1	E	17	ASN
1	E	18	THR
1	E	37	LYS
1	E	92	ASP
1	E	101	TYR
1	E	127	ASP
1	E	195	LEU
1	E	200	ARG
1	E	204	MET
1	E	258	LEU
1	E	261	THR
1	F	8	LYS
1	F	14	ILE
1	F	37	LYS
1	F	92	ASP
1	F	101	TYR
1	F	135	LEU
1	F	204	MET
1	G	8	LYS
1	G	14	ILE
1	G	37	LYS
1	G	101	TYR
1	G	195	LEU
1	G	263	THR
1	H	8	LYS
1	H	14	ILE
1	H	37	LYS
1	H	74	ASN
1	H	101	TYR
1	I	8	LYS
1	I	18	THR
1	I	37	LYS
1	I	71	GLU
1	I	101	TYR
1	J	8	LYS
1	J	14	ILE
1	J	18	THR

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Mol	Chain	Res	Type
1	J	37	LYS
1	J	101	TYR
1	J	200	ARG
1	J	257	GLN
1	K	8	LYS
1	K	37	LYS
1	K	101	TYR
1	K	135	LEU
1	L	8	LYS
1	L	18	THR
1	L	37	LYS
1	L	101	TYR
1	L	129	THR
1	L	131	LYS
1	L	190	VAL
1	L	263	THR
1	M	8	LYS
1	M	37	LYS
1	M	94	GLU
1	M	101	TYR
1	M	132	ILE
1	N	8	LYS
1	N	14	ILE
1	N	17	ASN
1	N	37	LYS
1	N	101	TYR
1	N	263	THR
1	N	294	LEU
1	O	8	LYS
1	O	18	THR
1	O	37	LYS
1	O	67	VAL
1	O	101	TYR
1	O	204	MET
1	O	258	LEU
1	O	261	THR
1	O	295	GLU
1	P	8	LYS
1	P	37	LYS
1	P	101	TYR
1	P	129	THR
1	P	205	LYS

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Mol	Chain	Res	Type
1	Q	8	LYS
1	Q	14	ILE
1	Q	17	ASN
1	Q	37	LYS
1	Q	101	TYR
1	Q	190	VAL
1	Q	201	ASN
1	R	8	LYS
1	R	37	LYS
1	R	66	ARG
1	R	101	TYR
1	S	8	LYS
1	S	17	ASN
1	S	37	LYS
1	S	101	TYR
1	S	201	ASN
1	S	261	THR
1	S	263	THR
1	S	295	GLU
1	T	8	LYS
1	T	14	ILE
1	T	18	THR
1	T	37	LYS
1	T	101	TYR
1	U	8	LYS
1	U	37	LYS
1	U	95	VAL
1	U	101	TYR
1	U	127	ASP
1	U	131	LYS
1	U	198	LYS
1	U	204	MET
1	V	8	LYS
1	V	18	THR
1	V	37	LYS
1	V	94	GLU
1	V	101	TYR
1	V	201	ASN
1	W	8	LYS
1	W	37	LYS
1	W	101	TYR
1	W	132	ILE

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Mol	Chain	Res	Type
1	W	195	LEU
1	W	201	ASN
1	W	261	THR
1	X	8	LYS
1	X	37	LYS
1	X	70	GLU
1	X	92	ASP
1	X	101	TYR
1	X	204	MET
1	Y	8	LYS
1	Y	37	LYS
1	Y	101	TYR
1	Y	128	ASP
1	Y	201	ASN
1	Y	204	MET
1	Z	8	LYS
1	Z	17	ASN
1	Z	37	LYS
1	Z	101	TYR
1	Z	131	LYS
1	Z	201	ASN
1	a	8	LYS
1	a	14	ILE
1	a	37	LYS
1	a	92	ASP
1	a	101	TYR
1	a	190	VAL
1	b	8	LYS
1	b	37	LYS
1	b	101	TYR
1	b	125	THR
1	b	201	ASN
1	b	204	MET
1	b	261	THR
1	b	263	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	121	ASN
1	A	177	GLN
1	B	74	ASN

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Mol	Chain	Res	Type
1	B	194	GLN
1	B	201	ASN
1	C	93	ASN
1	C	121	ASN
1	C	264	ASN
1	E	121	ASN
1	F	121	ASN
1	F	194	GLN
1	G	121	ASN
1	G	194	GLN
1	G	259	HIS
1	H	74	ASN
1	H	121	ASN
1	I	121	ASN
1	J	121	ASN
1	K	121	ASN
1	K	144	HIS
1	K	264	ASN
1	L	6	ASN
1	L	74	ASN
1	L	121	ASN
1	M	121	ASN
1	N	121	ASN
1	O	121	ASN
1	O	264	ASN
1	P	6	ASN
1	P	121	ASN
1	Q	121	ASN
1	Q	144	HIS
1	R	121	ASN
1	S	121	ASN
1	T	121	ASN
1	U	121	ASN
1	V	121	ASN
1	V	264	ASN
1	W	6	ASN
1	W	121	ASN
1	X	121	ASN
1	X	144	HIS
1	Y	121	ASN
1	Y	201	ASN
1	Z	121	ASN

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Mol	Chain	Res	Type
1	a	121	ASN
1	b	121	ASN
1	b	194	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

84 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	MPD	A	303	-	7,7,7	0.36	0	9,10,10	0.66	0
2	MPD	B	303	-	7,7,7	0.33	0	9,10,10	0.48	0
2	MPD	J	302	-	7,7,7	0.39	0	9,10,10	0.72	0
3	ACY	K	304	-	3,3,3	0.82	0	3,3,3	0.80	0
3	ACY	H	304	-	3,3,3	0.82	0	3,3,3	0.84	0
2	MPD	Q	302	-	7,7,7	0.26	0	9,10,10	0.59	0
3	ACY	C	304	-	3,3,3	0.84	0	3,3,3	0.73	0
2	MPD	O	303	-	7,7,7	0.35	0	9,10,10	0.59	0
2	MPD	b	303	-	7,7,7	0.34	0	9,10,10	0.49	0
2	MPD	F	303	-	7,7,7	0.25	0	9,10,10	0.47	0
2	MPD	D	303	-	7,7,7	0.31	0	9,10,10	0.24	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MPD	U	303	-	7,7,7	0.38	0	9,10,10	0.50	0
2	MPD	Y	303	-	7,7,7	0.37	0	9,10,10	0.64	0
3	ACY	D	305	-	3,3,3	0.82	0	3,3,3	0.79	0
3	ACY	O	304	-	3,3,3	0.81	0	3,3,3	0.76	0
3	ACY	X	305	-	3,3,3	0.83	0	3,3,3	0.86	0
3	ACY	A	305	-	3,3,3	0.79	0	3,3,3	0.89	0
2	MPD	L	303	-	7,7,7	0.29	0	9,10,10	0.59	0
3	ACY	D	304	-	3,3,3	0.81	0	3,3,3	0.75	0
3	ACY	S	305	-	3,3,3	0.82	0	3,3,3	0.77	0
2	MPD	I	302	-	7,7,7	0.31	0	9,10,10	0.26	0
2	MPD	G	303	-	7,7,7	0.36	0	9,10,10	0.55	0
3	ACY	a	304	-	3,3,3	0.83	0	3,3,3	0.82	0
3	ACY	O	305	-	3,3,3	0.86	0	3,3,3	0.69	0
2	MPD	M	303	-	7,7,7	0.36	0	9,10,10	0.59	0
2	MPD	C	302	-	7,7,7	0.30	0	9,10,10	0.33	0
3	ACY	H	305	-	3,3,3	0.80	0	3,3,3	0.83	0
2	MPD	P	302	-	7,7,7	0.27	0	9,10,10	0.25	0
3	ACY	Y	304	-	3,3,3	0.76	0	3,3,3	0.96	0
2	MPD	N	303	-	7,7,7	0.36	0	9,10,10	0.62	0
2	MPD	Z	302	-	7,7,7	0.30	0	9,10,10	0.42	0
2	MPD	S	303	-	7,7,7	0.35	0	9,10,10	0.48	0
2	MPD	B	302	-	7,7,7	0.28	0	9,10,10	0.36	0
2	MPD	F	302	-	7,7,7	0.30	0	9,10,10	0.34	0
3	ACY	V	304	-	3,3,3	0.82	0	3,3,3	0.83	0
2	MPD	X	302	-	7,7,7	0.28	0	9,10,10	0.57	0
2	MPD	V	302	-	7,7,7	0.28	0	9,10,10	0.34	0
3	ACY	P	304	-	3,3,3	0.81	0	3,3,3	0.88	0
2	MPD	Q	303	-	7,7,7	0.33	0	9,10,10	0.43	0
2	MPD	W	302	-	7,7,7	0.32	0	9,10,10	0.32	0
2	MPD	G	302	-	7,7,7	0.32	0	9,10,10	0.24	0
3	ACY	E	304	-	3,3,3	0.80	0	3,3,3	0.84	0
3	ACY	Q	304	-	3,3,3	0.77	0	3,3,3	0.96	0
2	MPD	L	302	-	7,7,7	0.32	0	9,10,10	0.54	0
2	MPD	R	302	-	7,7,7	0.29	0	9,10,10	0.62	0
3	ACY	M	304	-	3,3,3	0.85	0	3,3,3	0.73	0
2	MPD	N	302	-	7,7,7	0.28	0	9,10,10	0.43	0
2	MPD	b	302	-	7,7,7	0.30	0	9,10,10	0.50	0
2	MPD	W	303	-	7,7,7	0.33	0	9,10,10	0.52	0
2	MPD	X	304	-	7,7,7	0.32	0	9,10,10	0.52	0
2	MPD	H	302	-	7,7,7	0.30	0	9,10,10	0.48	0
2	MPD	K	302	-	7,7,7	0.29	0	9,10,10	0.36	0
2	MPD	I	303	-	7,7,7	0.30	0	9,10,10	0.48	0
3	ACY	W	304	-	3,3,3	0.87	0	3,3,3	0.69	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MPD	C	303	-	7,7,7	0.34	0	9,10,10	0.55	0
2	MPD	J	303	-	7,7,7	0.30	0	9,10,10	0.70	0
2	MPD	O	302	-	7,7,7	0.26	0	9,10,10	0.25	0
2	MPD	X	303	-	7,7,7	0.31	0	9,10,10	0.26	0
3	ACY	A	304	-	3,3,3	0.81	0	3,3,3	0.75	0
2	MPD	P	303	-	7,7,7	0.30	0	9,10,10	0.53	0
2	MPD	S	302	-	7,7,7	0.33	0	9,10,10	0.55	0
2	MPD	Y	302	-	7,7,7	0.30	0	9,10,10	0.41	0
3	ACY	J	304	-	3,3,3	0.79	0	3,3,3	0.84	0
3	ACY	S	304	-	3,3,3	0.82	0	3,3,3	0.68	0
2	MPD	U	302	-	7,7,7	0.33	0	9,10,10	0.29	0
3	ACY	K	305	-	3,3,3	0.82	0	3,3,3	0.85	0
2	MPD	E	303	-	7,7,7	0.35	0	9,10,10	0.58	0
2	MPD	a	302	-	7,7,7	0.31	0	9,10,10	0.33	0
2	MPD	a	303	-	7,7,7	0.27	0	9,10,10	0.64	0
2	MPD	V	303	-	7,7,7	0.30	0	9,10,10	0.71	0
3	ACY	F	304	-	3,3,3	0.82	0	3,3,3	0.78	0
3	ACY	L	304	-	3,3,3	0.81	0	3,3,3	0.80	0
2	MPD	T	302	-	7,7,7	0.34	0	9,10,10	0.27	0
2	MPD	M	302	-	7,7,7	0.29	0	9,10,10	0.24	0
2	MPD	A	302	-	7,7,7	0.28	0	9,10,10	0.91	0
2	MPD	T	303	-	7,7,7	0.29	0	9,10,10	0.89	0
3	ACY	Z	303	-	3,3,3	0.85	0	3,3,3	0.73	0
3	ACY	b	304	-	3,3,3	0.80	0	3,3,3	0.89	0
3	ACY	T	304	-	3,3,3	0.80	0	3,3,3	0.90	0
2	MPD	R	303	-	7,7,7	0.29	0	9,10,10	0.54	0
2	MPD	D	302	-	7,7,7	0.37	0	9,10,10	0.71	0
2	MPD	E	302	-	7,7,7	0.33	0	9,10,10	0.24	0
2	MPD	H	303	-	7,7,7	0.27	0	9,10,10	0.43	0
2	MPD	K	303	-	7,7,7	0.33	0	9,10,10	0.64	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	MPD	A	303	-	-	1/5/5/5	-
2	MPD	a	303	-	-	5/5/5/5	-
2	MPD	L	302	-	-	3/5/5/5	-
2	MPD	B	303	-	-	2/5/5/5	-
2	MPD	P	302	-	-	1/5/5/5	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MPD	J	302	-	-	2/5/5/5	-
2	MPD	V	303	-	-	4/5/5/5	-
2	MPD	R	302	-	-	3/5/5/5	-
2	MPD	N	303	-	-	2/5/5/5	-
2	MPD	N	302	-	-	4/5/5/5	-
2	MPD	Z	302	-	-	0/5/5/5	-
2	MPD	b	302	-	-	3/5/5/5	-
2	MPD	W	303	-	-	2/5/5/5	-
2	MPD	Q	302	-	-	1/5/5/5	-
2	MPD	X	304	-	-	2/5/5/5	-
2	MPD	S	303	-	-	1/5/5/5	-
2	MPD	O	303	-	-	2/5/5/5	-
2	MPD	H	302	-	-	0/5/5/5	-
2	MPD	K	302	-	-	1/5/5/5	-
2	MPD	b	303	-	-	1/5/5/5	-
2	MPD	I	303	-	-	2/5/5/5	-
2	MPD	B	302	-	-	3/5/5/5	-
2	MPD	F	303	-	-	3/5/5/5	-
2	MPD	F	302	-	-	1/5/5/5	-
2	MPD	D	303	-	-	1/5/5/5	-
2	MPD	T	302	-	-	2/5/5/5	-
2	MPD	M	302	-	-	1/5/5/5	-
2	MPD	U	303	-	-	4/5/5/5	-
2	MPD	C	303	-	-	4/5/5/5	-
2	MPD	X	302	-	-	3/5/5/5	-
2	MPD	Y	303	-	-	2/5/5/5	-
2	MPD	V	302	-	-	3/5/5/5	-
2	MPD	A	302	-	-	5/5/5/5	-
2	MPD	L	303	-	-	4/5/5/5	-
2	MPD	T	303	-	-	5/5/5/5	-
2	MPD	a	302	-	-	2/5/5/5	-
2	MPD	J	303	-	-	2/5/5/5	-
2	MPD	Q	303	-	-	3/5/5/5	-
2	MPD	O	302	-	-	2/5/5/5	-
2	MPD	X	303	-	-	2/5/5/5	-
2	MPD	P	303	-	-	4/5/5/5	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MPD	W	302	-	-	1/5/5/5	-
2	MPD	I	302	-	-	3/5/5/5	-
2	MPD	G	303	-	-	2/5/5/5	-
2	MPD	G	302	-	-	2/5/5/5	-
2	MPD	R	303	-	-	0/5/5/5	-
2	MPD	S	302	-	-	2/5/5/5	-
2	MPD	Y	302	-	-	3/5/5/5	-
2	MPD	U	302	-	-	1/5/5/5	-
2	MPD	M	303	-	-	1/5/5/5	-
2	MPD	C	302	-	-	3/5/5/5	-
2	MPD	D	302	-	-	2/5/5/5	-
2	MPD	E	303	-	-	2/5/5/5	-
2	MPD	E	302	-	-	2/5/5/5	-
2	MPD	H	303	-	-	3/5/5/5	-
2	MPD	K	303	-	-	2/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (127) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	302	MPD	O2-C2-C3-C4
2	A	302	MPD	CM-C2-C3-C4
2	A	303	MPD	C2-C3-C4-C5
2	B	302	MPD	C2-C3-C4-C5
2	B	303	MPD	C2-C3-C4-C5
2	C	303	MPD	C2-C3-C4-O4
2	C	303	MPD	C2-C3-C4-C5
2	D	302	MPD	C2-C3-C4-O4
2	D	302	MPD	C2-C3-C4-C5
2	E	303	MPD	C2-C3-C4-O4
2	E	303	MPD	C2-C3-C4-C5
2	F	303	MPD	CM-C2-C3-C4
2	G	302	MPD	C2-C3-C4-O4
2	G	303	MPD	C2-C3-C4-C5
2	I	302	MPD	C2-C3-C4-O4
2	J	303	MPD	C2-C3-C4-O4

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Mol	Chain	Res	Type	Atoms
2	J	303	MPD	C2-C3-C4-C5
2	K	303	MPD	C2-C3-C4-C5
2	L	302	MPD	C2-C3-C4-O4
2	L	302	MPD	C2-C3-C4-C5
2	L	303	MPD	C1-C2-C3-C4
2	L	303	MPD	O2-C2-C3-C4
2	M	303	MPD	C2-C3-C4-C5
2	N	302	MPD	C2-C3-C4-O4
2	N	302	MPD	C2-C3-C4-C5
2	N	303	MPD	C2-C3-C4-C5
2	O	303	MPD	C2-C3-C4-O4
2	O	303	MPD	C2-C3-C4-C5
2	P	303	MPD	C2-C3-C4-O4
2	P	303	MPD	C2-C3-C4-C5
2	Q	303	MPD	C2-C3-C4-O4
2	Q	303	MPD	C2-C3-C4-C5
2	R	302	MPD	C1-C2-C3-C4
2	R	302	MPD	O2-C2-C3-C4
2	S	302	MPD	C1-C2-C3-C4
2	S	303	MPD	C2-C3-C4-C5
2	T	303	MPD	C2-C3-C4-O4
2	T	303	MPD	C2-C3-C4-C5
2	U	303	MPD	C2-C3-C4-O4
2	U	303	MPD	C2-C3-C4-C5
2	V	302	MPD	C2-C3-C4-C5
2	W	303	MPD	C2-C3-C4-C5
2	X	304	MPD	C2-C3-C4-O4
2	X	304	MPD	C2-C3-C4-C5
2	Y	302	MPD	C2-C3-C4-O4
2	Y	303	MPD	C2-C3-C4-C5
2	a	302	MPD	C2-C3-C4-O4
2	a	303	MPD	C1-C2-C3-C4
2	a	303	MPD	C2-C3-C4-O4
2	a	303	MPD	C2-C3-C4-C5
2	D	303	MPD	O2-C2-C3-C4
2	F	302	MPD	O2-C2-C3-C4
2	F	303	MPD	O2-C2-C3-C4
2	K	302	MPD	O2-C2-C3-C4
2	K	303	MPD	O2-C2-C3-C4
2	M	302	MPD	O2-C2-C3-C4
2	U	302	MPD	O2-C2-C3-C4
2	W	302	MPD	O2-C2-C3-C4

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Mol	Chain	Res	Type	Atoms
2	Y	302	MPD	O2-C2-C3-C4
2	a	303	MPD	O2-C2-C3-C4
2	A	302	MPD	C2-C3-C4-O4
2	N	303	MPD	C2-C3-C4-O4
2	O	302	MPD	C2-C3-C4-O4
2	W	303	MPD	C2-C3-C4-O4
2	Y	303	MPD	C2-C3-C4-O4
2	A	302	MPD	C1-C2-C3-C4
2	C	303	MPD	CM-C2-C3-C4
2	F	303	MPD	C1-C2-C3-C4
2	H	303	MPD	C1-C2-C3-C4
2	L	303	MPD	CM-C2-C3-C4
2	P	303	MPD	C1-C2-C3-C4
2	R	302	MPD	CM-C2-C3-C4
2	S	302	MPD	CM-C2-C3-C4
2	T	303	MPD	C1-C2-C3-C4
2	T	303	MPD	CM-C2-C3-C4
2	V	303	MPD	C1-C2-C3-C4
2	V	303	MPD	CM-C2-C3-C4
2	X	302	MPD	C1-C2-C3-C4
2	a	303	MPD	CM-C2-C3-C4
2	A	302	MPD	C2-C3-C4-C5
2	E	302	MPD	C2-C3-C4-C5
2	H	303	MPD	C2-C3-C4-C5
2	I	302	MPD	C2-C3-C4-C5
2	I	303	MPD	C2-C3-C4-C5
2	J	302	MPD	C2-C3-C4-C5
2	L	303	MPD	C2-C3-C4-C5
2	T	302	MPD	C2-C3-C4-C5
2	V	303	MPD	C2-C3-C4-C5
2	X	302	MPD	C2-C3-C4-C5
2	Y	302	MPD	C2-C3-C4-C5
2	b	303	MPD	C2-C3-C4-C5
2	B	302	MPD	O2-C2-C3-C4
2	C	302	MPD	O2-C2-C3-C4
2	G	302	MPD	O2-C2-C3-C4
2	I	302	MPD	O2-C2-C3-C4
2	O	302	MPD	O2-C2-C3-C4
2	T	303	MPD	O2-C2-C3-C4
2	U	303	MPD	O2-C2-C3-C4
2	V	303	MPD	O2-C2-C3-C4
2	X	303	MPD	O2-C2-C3-C4

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Mol	Chain	Res	Type	Atoms
2	a	302	MPD	O2-C2-C3-C4
2	b	302	MPD	O2-C2-C3-C4
2	B	302	MPD	C2-C3-C4-O4
2	C	302	MPD	C2-C3-C4-O4
2	E	302	MPD	C2-C3-C4-O4
2	G	303	MPD	C2-C3-C4-O4
2	H	303	MPD	C2-C3-C4-O4
2	I	303	MPD	C2-C3-C4-O4
2	J	302	MPD	C2-C3-C4-O4
2	P	302	MPD	C2-C3-C4-O4
2	Q	302	MPD	C2-C3-C4-O4
2	T	302	MPD	C2-C3-C4-O4
2	V	302	MPD	C2-C3-C4-O4
2	B	303	MPD	CM-C2-C3-C4
2	C	302	MPD	CM-C2-C3-C4
2	C	303	MPD	C1-C2-C3-C4
2	L	302	MPD	CM-C2-C3-C4
2	N	302	MPD	C1-C2-C3-C4
2	N	302	MPD	CM-C2-C3-C4
2	P	303	MPD	CM-C2-C3-C4
2	Q	303	MPD	C1-C2-C3-C4
2	U	303	MPD	C1-C2-C3-C4
2	V	302	MPD	C1-C2-C3-C4
2	X	302	MPD	CM-C2-C3-C4
2	X	303	MPD	CM-C2-C3-C4
2	b	302	MPD	C1-C2-C3-C4
2	b	302	MPD	CM-C2-C3-C4

There are no ring outliers.

53 monomers are involved in 183 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	303	MPD	1	0
2	B	303	MPD	5	0
2	J	302	MPD	3	0
2	Q	302	MPD	10	0
2	O	303	MPD	5	0
2	b	303	MPD	3	0
2	F	303	MPD	7	0
2	U	303	MPD	3	0
2	Y	303	MPD	3	0
2	L	303	MPD	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	I	302	MPD	4	0
2	G	303	MPD	3	0
2	M	303	MPD	3	0
2	C	302	MPD	1	0
2	P	302	MPD	1	0
2	N	303	MPD	5	0
2	Z	302	MPD	1	0
2	S	303	MPD	2	0
2	B	302	MPD	4	0
2	F	302	MPD	3	0
2	X	302	MPD	4	0
2	V	302	MPD	1	0
2	Q	303	MPD	1	0
2	W	302	MPD	4	0
2	G	302	MPD	4	0
2	L	302	MPD	3	0
2	N	302	MPD	6	0
2	b	302	MPD	3	0
2	W	303	MPD	5	0
2	X	304	MPD	2	0
2	H	302	MPD	3	0
2	K	302	MPD	5	0
2	I	303	MPD	3	0
2	C	303	MPD	5	0
2	J	303	MPD	2	0
2	O	302	MPD	5	0
2	P	303	MPD	3	0
2	S	302	MPD	2	0
2	Y	302	MPD	3	0
2	U	302	MPD	2	0
2	E	303	MPD	1	0
2	a	302	MPD	5	0
2	a	303	MPD	3	0
2	V	303	MPD	7	0
2	T	302	MPD	1	0
2	M	302	MPD	1	0
2	A	302	MPD	3	0
2	T	303	MPD	3	0
2	R	303	MPD	6	0
2	D	302	MPD	2	0
2	E	302	MPD	8	0
2	H	303	MPD	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	K	303	MPD	6	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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	295/302 (97%)	-0.03	2 (0%) 84 85	27, 45, 75, 108	0
1	B	295/302 (97%)	0.10	8 (2%) 56 58	27, 48, 82, 119	0
1	C	294/302 (97%)	0.18	5 (1%) 69 70	31, 50, 82, 108	0
1	D	295/302 (97%)	0.17	6 (2%) 65 66	33, 52, 83, 118	0
1	E	295/302 (97%)	0.45	25 (8%) 16 18	33, 54, 109, 151	0
1	F	294/302 (97%)	0.31	16 (5%) 31 33	30, 52, 95, 147	0
1	G	294/302 (97%)	0.16	7 (2%) 59 62	27, 50, 89, 127	0
1	H	294/302 (97%)	-0.02	1 (0%) 90 90	29, 46, 71, 100	0
1	I	294/302 (97%)	0.06	4 (1%) 73 75	27, 47, 75, 106	0
1	J	294/302 (97%)	0.05	3 (1%) 79 80	24, 46, 80, 122	0
1	K	295/302 (97%)	0.05	5 (1%) 69 70	27, 45, 76, 111	0
1	L	294/302 (97%)	0.11	9 (3%) 51 53	30, 47, 75, 131	0
1	M	295/302 (97%)	0.21	9 (3%) 51 53	29, 52, 92, 149	0
1	N	294/302 (97%)	0.16	3 (1%) 79 80	28, 51, 87, 117	0
1	O	295/302 (97%)	0.09	5 (1%) 69 70	30, 48, 78, 117	0
1	P	294/302 (97%)	0.13	8 (2%) 56 58	29, 48, 75, 131	0
1	Q	294/302 (97%)	0.13	6 (2%) 65 66	30, 51, 95, 140	0
1	R	294/302 (97%)	0.16	7 (2%) 59 62	32, 49, 77, 99	0
1	S	295/302 (97%)	0.19	5 (1%) 69 70	30, 52, 92, 118	0
1	T	294/302 (97%)	0.03	7 (2%) 59 62	30, 47, 76, 106	0
1	U	295/302 (97%)	0.08	6 (2%) 65 66	28, 48, 85, 128	0
1	V	295/302 (97%)	0.06	3 (1%) 79 80	29, 50, 81, 123	0
1	W	294/302 (97%)	0.25	11 (3%) 45 48	33, 53, 94, 155	0
1	X	294/302 (97%)	0.18	7 (2%) 59 62	33, 53, 96, 127	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	Y	294/302 (97%)	0.31	15 (5%) 33 35	29, 52, 102, 136	0
1	Z	294/302 (97%)	0.12	10 (3%) 48 50	28, 51, 87, 124	0
1	a	294/302 (97%)	-0.01	4 (1%) 73 75	28, 45, 74, 123	0
1	b	295/302 (97%)	0.05	6 (2%) 65 66	28, 48, 84, 151	0
All	All	8243/8456 (97%)	0.13	203 (2%) 58 60	24, 49, 86, 155	0

All (203) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	240	LYS	4.3
1	E	74	ASN	4.2
1	G	294	LEU	3.9
1	E	260	TRP	3.9
1	B	2	ASP	3.8
1	F	71	GLU	3.6
1	M	72	GLY	3.6
1	Y	294	LEU	3.6
1	Z	294	LEU	3.6
1	Z	262	SER	3.6
1	V	17	ASN	3.6
1	S	179	TRP	3.5
1	E	294	LEU	3.5
1	Z	179	TRP	3.5
1	M	261	THR	3.4
1	E	203	SER	3.4
1	E	73	ALA	3.4
1	F	17	ASN	3.4
1	P	93	ASN	3.4
1	E	206	ALA	3.4
1	a	261	THR	3.3
1	T	70	GLU	3.2
1	E	263	THR	3.2
1	Z	261	THR	3.2
1	E	265	TRP	3.2
1	A	295	GLU	3.2
1	E	76	SER	3.2
1	b	262	SER	3.2
1	O	17	ASN	3.2
1	R	179	TRP	3.2
1	Y	260	TRP	3.2
1	N	294	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
1	T	294	LEU	3.2
1	E	262	SER	3.1
1	B	179	TRP	3.1
1	J	127	ASP	3.1
1	C	179	TRP	3.1
1	E	72	GLY	3.1
1	W	17	ASN	3.1
1	J	179	TRP	3.0
1	W	72	GLY	3.0
1	W	73	ALA	3.0
1	F	2	ASP	3.0
1	Y	74	ASN	3.0
1	M	71	GLU	3.0
1	N	179	TRP	3.0
1	U	261	THR	2.9
1	R	2	ASP	2.9
1	S	192	GLY	2.9
1	S	261	THR	2.9
1	C	127	ASP	2.9
1	b	263	THR	2.9
1	E	77	GLY	2.8
1	Y	240	LYS	2.8
1	P	94	GLU	2.8
1	I	2	ASP	2.8
1	E	279	SER	2.8
1	M	15	GLY	2.8
1	R	124	VAL	2.8
1	b	261	THR	2.8
1	J	264	ASN	2.8
1	Y	2	ASP	2.8
1	L	205	LYS	2.8
1	D	261	THR	2.8
1	P	261	THR	2.8
1	F	260	TRP	2.8
1	M	17	ASN	2.8
1	U	263	THR	2.7
1	T	13	ASP	2.7
1	I	179	TRP	2.7
1	P	72	GLY	2.7
1	S	260	TRP	2.7
1	B	70	GLU	2.7
1	X	261	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	Y	246	ASP	2.7
1	F	72	GLY	2.7
1	E	179	TRP	2.7
1	S	295	GLU	2.7
1	F	76	SER	2.7
1	R	94	GLU	2.7
1	E	264	ASN	2.7
1	E	204	MET	2.7
1	T	261	THR	2.6
1	Q	294	LEU	2.6
1	E	208	GLU	2.6
1	P	71	GLU	2.6
1	B	262	SER	2.6
1	F	73	ALA	2.6
1	E	189	PRO	2.6
1	F	294	LEU	2.6
1	F	261	THR	2.6
1	W	261	THR	2.6
1	L	208	GLU	2.5
1	F	69	SER	2.5
1	W	203	SER	2.5
1	a	127	ASP	2.5
1	Z	192	GLY	2.5
1	W	18	THR	2.5
1	K	136	ILE	2.5
1	Y	136	ILE	2.5
1	D	94	GLU	2.5
1	L	17	ASN	2.5
1	L	77	GLY	2.5
1	G	261	THR	2.5
1	Q	148	TYR	2.5
1	C	201	ASN	2.5
1	H	1	ALA	2.5
1	X	73	ALA	2.5
1	F	15	GLY	2.5
1	P	77	GLY	2.5
1	F	262	SER	2.4
1	Q	262	SER	2.4
1	F	179	TRP	2.4
1	X	179	TRP	2.4
1	a	13	ASP	2.4
1	G	70	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	G	242	GLN	2.4
1	I	128	ASP	2.4
1	L	261	THR	2.4
1	W	179	TRP	2.4
1	B	128	ASP	2.4
1	M	295	GLU	2.4
1	O	71	GLU	2.4
1	F	189	PRO	2.4
1	L	279	SER	2.4
1	T	17	ASN	2.3
1	Y	189	PRO	2.3
1	Q	267	GLY	2.3
1	B	129	THR	2.3
1	K	261	THR	2.3
1	K	16	SER	2.3
1	E	295	GLU	2.3
1	U	127	ASP	2.3
1	Y	201	ASN	2.3
1	B	1	ALA	2.3
1	N	262	SER	2.3
1	C	126	GLY	2.3
1	X	260	TRP	2.3
1	X	263	THR	2.3
1	P	262	SER	2.3
1	W	71	GLU	2.3
1	W	94	GLU	2.3
1	E	2	ASP	2.3
1	Q	2	ASP	2.3
1	C	129	THR	2.3
1	M	129	THR	2.3
1	O	261	THR	2.3
1	Y	263	THR	2.3
1	U	189	PRO	2.2
1	Z	267	GLY	2.2
1	Y	179	TRP	2.2
1	Z	260	TRP	2.2
1	Y	205	LYS	2.2
1	D	179	TRP	2.2
1	V	179	TRP	2.2
1	T	128	ASP	2.2
1	Q	202	GLY	2.2
1	X	202	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	Y	76	SER	2.2
1	D	17	ASN	2.2
1	O	131	LYS	2.2
1	B	133	GLY	2.2
1	L	294	LEU	2.1
1	W	260	TRP	2.1
1	F	124	VAL	2.1
1	X	124	VAL	2.1
1	a	17	ASN	2.1
1	E	141	SER	2.1
1	F	263	THR	2.1
1	K	203	SER	2.1
1	Z	16	SER	2.1
1	b	1	ALA	2.1
1	U	112	TYR	2.1
1	R	246	ASP	2.1
1	W	70	GLU	2.1
1	Y	208	GLU	2.1
1	K	259	HIS	2.1
1	L	187	TRP	2.1
1	M	179	TRP	2.1
1	P	179	TRP	2.1
1	G	129	THR	2.1
1	b	203	SER	2.1
1	E	258	LEU	2.1
1	T	94	GLU	2.1
1	O	127	ASP	2.1
1	E	18	THR	2.1
1	E	261	THR	2.1
1	M	263	THR	2.1
1	R	262	SER	2.1
1	R	188	ASN	2.1
1	L	189	PRO	2.1
1	G	179	TRP	2.1
1	Y	265	TRP	2.1
1	A	261	THR	2.0
1	Z	263	THR	2.0
1	b	69	SER	2.0
1	I	71	GLU	2.0
1	D	2	ASP	2.0
1	Z	118	TYR	2.0
1	D	253	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	E	267	GLY	2.0
1	V	2	ASP	2.0
1	U	75	LYS	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 oligosaccharides 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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MPD	J	303	8/8	0.58	0.17	59,79,92,96	0
2	MPD	B	303	8/8	0.59	0.19	78,85,100,101	0
3	ACY	a	304	4/4	0.60	0.16	67,68,82,87	0
2	MPD	Q	303	8/8	0.61	0.17	60,78,87,88	0
2	MPD	F	303	8/8	0.61	0.19	68,77,84,91	0
3	ACY	J	304	4/4	0.62	0.20	61,65,78,90	0
3	ACY	X	305	4/4	0.64	0.18	58,74,85,90	0
2	MPD	Z	302	8/8	0.65	0.18	56,82,97,102	0
2	MPD	L	303	8/8	0.65	0.17	67,72,98,100	0
3	ACY	M	304	4/4	0.66	0.12	53,58,79,84	0
3	ACY	O	305	4/4	0.66	0.18	65,78,84,87	0
3	ACY	C	304	4/4	0.67	0.18	63,66,74,78	0
3	ACY	V	304	4/4	0.68	0.17	60,72,75,91	0
3	ACY	F	304	4/4	0.69	0.14	69,76,84,85	0
3	ACY	A	305	4/4	0.69	0.20	52,72,90,99	0
3	ACY	S	305	4/4	0.69	0.18	70,73,79,80	0
3	ACY	E	304	4/4	0.70	0.14	56,70,75,77	0
3	ACY	S	304	4/4	0.70	0.21	51,79,87,101	0
2	MPD	X	304	8/8	0.70	0.19	68,86,88,93	0
3	ACY	L	304	4/4	0.71	0.12	62,70,82,90	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MPD	X	302	8/8	0.71	0.15	70,85,88,89	0
2	MPD	F	302	8/8	0.72	0.14	71,79,91,103	0
3	ACY	W	304	4/4	0.72	0.16	49,65,70,83	0
2	MPD	U	303	8/8	0.72	0.15	52,76,87,93	0
3	ACY	Z	303	4/4	0.72	0.16	56,71,71,75	0
2	MPD	E	303	8/8	0.72	0.15	61,80,89,100	0
2	MPD	G	303	8/8	0.73	0.13	59,75,86,97	0
3	ACY	D	304	4/4	0.73	0.20	58,71,73,80	0
2	MPD	J	302	8/8	0.73	0.15	63,72,87,93	0
2	MPD	D	303	8/8	0.74	0.14	51,71,79,80	0
2	MPD	S	303	8/8	0.74	0.14	64,74,81,94	0
3	ACY	H	305	4/4	0.75	0.17	65,73,81,83	0
2	MPD	H	303	8/8	0.75	0.15	62,70,86,88	0
2	MPD	W	303	8/8	0.75	0.13	60,80,89,91	0
2	MPD	S	302	8/8	0.75	0.17	68,80,81,85	0
2	MPD	C	303	8/8	0.75	0.12	65,76,87,88	0
2	MPD	U	302	8/8	0.75	0.20	66,82,90,90	0
3	ACY	P	304	4/4	0.76	0.13	63,63,82,97	0
2	MPD	N	303	8/8	0.76	0.13	66,79,90,95	0
3	ACY	D	305	4/4	0.76	0.15	70,71,78,78	0
2	MPD	M	303	8/8	0.76	0.15	69,75,81,97	0
3	ACY	A	304	4/4	0.77	0.13	54,61,73,74	0
2	MPD	X	303	8/8	0.77	0.14	60,74,85,90	0
2	MPD	I	303	8/8	0.77	0.14	61,80,92,99	0
2	MPD	T	302	8/8	0.77	0.17	64,68,76,78	0
2	MPD	Q	302	8/8	0.78	0.15	69,76,88,103	0
2	MPD	P	303	8/8	0.78	0.12	59,79,86,88	0
3	ACY	b	304	4/4	0.78	0.13	66,69,91,92	0
2	MPD	R	302	8/8	0.79	0.17	60,73,83,86	0
2	MPD	T	303	8/8	0.80	0.11	56,72,85,89	0
2	MPD	Y	303	8/8	0.80	0.14	64,82,87,93	0
2	MPD	V	303	8/8	0.80	0.13	65,82,90,91	0
2	MPD	b	303	8/8	0.80	0.13	64,82,86,87	0
2	MPD	W	302	8/8	0.80	0.12	68,75,83,84	0
3	ACY	Q	304	4/4	0.81	0.12	60,64,80,87	0
2	MPD	E	302	8/8	0.81	0.12	72,81,85,87	0
2	MPD	R	303	8/8	0.81	0.14	64,77,82,87	0
3	ACY	T	304	4/4	0.81	0.12	57,65,76,81	0
3	ACY	H	304	4/4	0.81	0.13	53,60,76,79	0
2	MPD	a	302	8/8	0.82	0.17	60,74,83,89	0
2	MPD	K	303	8/8	0.82	0.11	63,78,80,88	0
2	MPD	C	302	8/8	0.83	0.13	67,73,79,82	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MPD	O	303	8/8	0.83	0.11	58,75,89,93	0
3	ACY	Y	304	4/4	0.83	0.13	64,75,79,82	0
2	MPD	O	302	8/8	0.84	0.13	68,74,80,82	0
2	MPD	N	302	8/8	0.84	0.14	64,77,82,92	0
2	MPD	K	302	8/8	0.84	0.15	66,73,83,90	0
2	MPD	a	303	8/8	0.85	0.10	55,80,89,93	0
2	MPD	A	303	8/8	0.86	0.11	59,72,85,90	0
3	ACY	O	304	4/4	0.86	0.12	59,68,77,86	0
2	MPD	A	302	8/8	0.86	0.16	58,69,77,81	0
3	ACY	K	304	4/4	0.86	0.12	52,67,72,73	0
2	MPD	P	302	8/8	0.86	0.11	54,76,79,80	0
2	MPD	D	302	8/8	0.87	0.13	65,73,83,85	0
2	MPD	V	302	8/8	0.87	0.10	49,61,75,80	0
2	MPD	M	302	8/8	0.87	0.12	64,73,79,82	0
2	MPD	H	302	8/8	0.87	0.16	58,64,75,78	0
2	MPD	Y	302	8/8	0.88	0.10	58,76,88,90	0
3	ACY	K	305	4/4	0.88	0.12	45,68,72,82	0
2	MPD	b	302	8/8	0.89	0.15	71,74,83,88	0
2	MPD	L	302	8/8	0.89	0.10	59,65,80,89	0
2	MPD	B	302	8/8	0.90	0.14	65,71,79,82	0
2	MPD	G	302	8/8	0.90	0.12	66,72,80,81	0
2	MPD	I	302	8/8	0.92	0.11	64,68,73,82	0

6.5 Other polymers [i](#)

There are no such residues in this entry.