



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 03:34 AM UTC

PDB ID : 6NAH / pdb_00006nah
Title : Crystal structure of Neisseria meningitidis ClpP protease in complex with Acyldepsipeptide-14 (ADEP-14)
Authors : Mabanglo, M.F.; Houry, W.A.
Deposited on : 2018-12-05
Resolution : 2.70 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

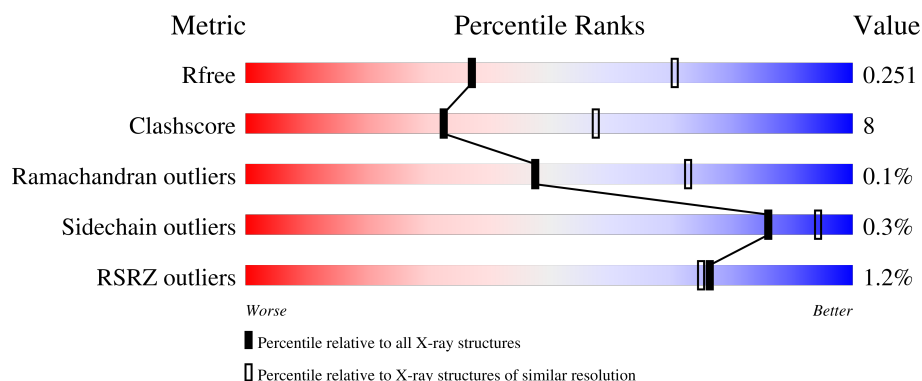
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	217	% 67% 14% 19%
1	B	217	% 66% 13% 20%
1	C	217	% 70% 9% 20%
1	D	217	67% 12% 20%
1	E	217	% 70% 10% 20%

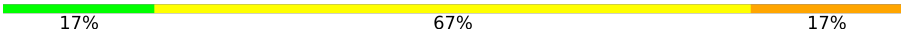
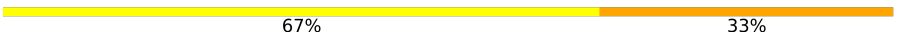
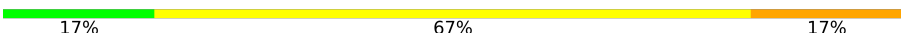
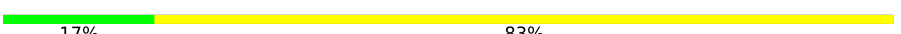
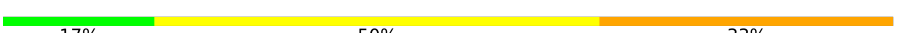
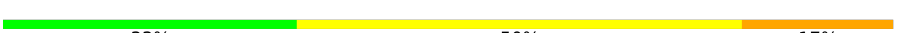
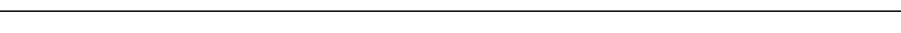
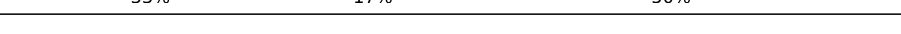
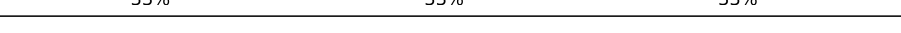
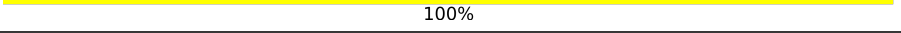
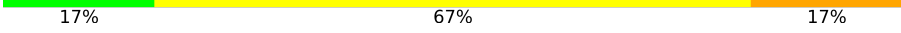
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Mol	Chain	Length	Quality of chain
1	F	217	% 64% 15% 21%
1	G	217	69% 11% 20%
1	H	217	% 62% 16% 20%
1	I	217	% 65% 13% 21%
1	J	217	65% 13% 21%
1	K	217	2% 66% 12% 20%
1	L	217	% 62% 18% 20%
1	M	217	% 65% 13% 21%
1	N	217	% 68% 12% 20%
1	O	217	60% 19% 21%
1	P	217	64% 15% 20%
1	Q	217	% 66% 13% 20%
1	R	217	65% 16% 19%
1	S	217	% 62% 17% 20%
1	T	217	% 66% 12% 21%
1	U	217	2% 60% 20% 20%
1	V	217	% 64% 14% 21%
1	W	217	% 67% 12% 20%
1	X	217	% 67% 12% 20%
1	Y	217	% 65% 14% 21%
1	Z	217	% 64% 15% 21%
1	a	217	% 67% 12% 20%
1	b	217	65% 14% 21%
2	0	6	33% 67%
2	1	6	33% 50% 17%

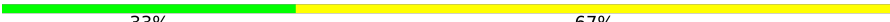
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Mol	Chain	Length	Quality of chain
2	2	6	
2	3	6	
2	4	6	
2	c	6	
2	e	6	
2	f	6	
2	g	6	
2	h	6	
2	i	6	
2	j	6	
2	k	6	
2	l	6	
2	m	6	
2	n	6	
2	o	6	
2	p	6	
2	q	6	
2	r	6	
2	s	6	
2	t	6	
2	u	6	
2	v	6	
2	w	6	
2	x	6	
2	y	6	

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Mol	Chain	Length	Quality of chain
2	z	6	 33% 67%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 40076 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 ATP-dependent Clp protease proteolytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	175	Total	C	N	O	S	0	0	0
			1359	858	230	263	8			
1	B	173	Total	C	N	O	S	0	0	0
			1347	850	228	261	8			
1	C	173	Total	C	N	O	S	0	0	0
			1346	851	228	259	8			
1	D	173	Total	C	N	O	S	0	0	0
			1351	852	229	262	8			
1	E	174	Total	C	N	O	S	0	0	0
			1359	858	230	263	8			
1	F	171	Total	C	N	O	S	0	0	0
			1339	846	226	259	8			
1	G	173	Total	C	N	O	S	0	0	0
			1351	854	228	261	8			
1	H	174	Total	C	N	O	S	0	0	0
			1359	858	230	263	8			
1	I	171	Total	C	N	O	S	0	0	0
			1339	846	226	259	8			
1	J	172	Total	C	N	O	S	0	0	0
			1347	850	228	261	8			
1	K	174	Total	C	N	O	S	0	0	0
			1359	858	230	263	8			
1	L	174	Total	C	N	O	S	0	0	0
			1355	856	229	262	8			
1	M	171	Total	C	N	O	S	0	0	0
			1339	846	226	259	8			
1	N	173	Total	C	N	O	S	0	0	0
			1347	850	228	261	8			
1	O	172	Total	C	N	O	S	0	0	0
			1347	850	228	261	8			
1	P	173	Total	C	N	O	S	0	0	0
			1347	850	228	261	8			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	173	Total	C	N	O	S	0	0	0
			1351	854	228	261	8			
1	R	175	Total	C	N	O	S	0	0	0
			1363	860	231	264	8			
1	S	173	Total	C	N	O	S	0	0	0
			1351	852	229	262	8			
1	T	171	Total	C	N	O	S	0	0	0
			1339	846	227	258	8			
1	U	174	Total	C	N	O	S	0	0	0
			1350	853	229	260	8			
1	V	171	Total	C	N	O	S	0	0	0
			1341	847	227	259	8			
1	W	173	Total	C	N	O	S	0	0	0
			1347	850	228	261	8			
1	X	174	Total	C	N	O	S	0	0	0
			1355	856	229	262	8			
1	Y	172	Total	C	N	O	S	0	0	0
			1347	850	228	261	8			
1	Z	172	Total	C	N	O	S	0	0	0
			1343	848	227	260	8			
1	a	173	Total	C	N	O	S	0	0	0
			1351	852	229	262	8			
1	b	172	Total	C	N	O	S	0	0	0
			1345	851	227	259	8			

There are 364 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-12	HIS	-	expression tag	UNP I4E574
A	-11	HIS	-	expression tag	UNP I4E574
A	-10	HIS	-	expression tag	UNP I4E574
A	-9	HIS	-	expression tag	UNP I4E574
A	-8	HIS	-	expression tag	UNP I4E574
A	-7	HIS	-	expression tag	UNP I4E574
A	-6	GLU	-	expression tag	UNP I4E574
A	-5	ASN	-	expression tag	UNP I4E574
A	-4	LEU	-	expression tag	UNP I4E574
A	-3	TYR	-	expression tag	UNP I4E574
A	-2	PHE	-	expression tag	UNP I4E574
A	-1	GLN	-	expression tag	UNP I4E574
A	0	GLY	-	expression tag	UNP I4E574
B	-12	HIS	-	expression tag	UNP I4E574
B	-11	HIS	-	expression tag	UNP I4E574

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	HIS	-	expression tag	UNP I4E574
B	-9	HIS	-	expression tag	UNP I4E574
B	-8	HIS	-	expression tag	UNP I4E574
B	-7	HIS	-	expression tag	UNP I4E574
B	-6	GLU	-	expression tag	UNP I4E574
B	-5	ASN	-	expression tag	UNP I4E574
B	-4	LEU	-	expression tag	UNP I4E574
B	-3	TYR	-	expression tag	UNP I4E574
B	-2	PHE	-	expression tag	UNP I4E574
B	-1	GLN	-	expression tag	UNP I4E574
B	0	GLY	-	expression tag	UNP I4E574
C	-12	HIS	-	expression tag	UNP I4E574
C	-11	HIS	-	expression tag	UNP I4E574
C	-10	HIS	-	expression tag	UNP I4E574
C	-9	HIS	-	expression tag	UNP I4E574
C	-8	HIS	-	expression tag	UNP I4E574
C	-7	HIS	-	expression tag	UNP I4E574
C	-6	GLU	-	expression tag	UNP I4E574
C	-5	ASN	-	expression tag	UNP I4E574
C	-4	LEU	-	expression tag	UNP I4E574
C	-3	TYR	-	expression tag	UNP I4E574
C	-2	PHE	-	expression tag	UNP I4E574
C	-1	GLN	-	expression tag	UNP I4E574
C	0	GLY	-	expression tag	UNP I4E574
D	-12	HIS	-	expression tag	UNP I4E574
D	-11	HIS	-	expression tag	UNP I4E574
D	-10	HIS	-	expression tag	UNP I4E574
D	-9	HIS	-	expression tag	UNP I4E574
D	-8	HIS	-	expression tag	UNP I4E574
D	-7	HIS	-	expression tag	UNP I4E574
D	-6	GLU	-	expression tag	UNP I4E574
D	-5	ASN	-	expression tag	UNP I4E574
D	-4	LEU	-	expression tag	UNP I4E574
D	-3	TYR	-	expression tag	UNP I4E574
D	-2	PHE	-	expression tag	UNP I4E574
D	-1	GLN	-	expression tag	UNP I4E574
D	0	GLY	-	expression tag	UNP I4E574
E	-12	HIS	-	expression tag	UNP I4E574
E	-11	HIS	-	expression tag	UNP I4E574
E	-10	HIS	-	expression tag	UNP I4E574
E	-9	HIS	-	expression tag	UNP I4E574
E	-8	HIS	-	expression tag	UNP I4E574

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-7	HIS	-	expression tag	UNP I4E574
E	-6	GLU	-	expression tag	UNP I4E574
E	-5	ASN	-	expression tag	UNP I4E574
E	-4	LEU	-	expression tag	UNP I4E574
E	-3	TYR	-	expression tag	UNP I4E574
E	-2	PHE	-	expression tag	UNP I4E574
E	-1	GLN	-	expression tag	UNP I4E574
E	0	GLY	-	expression tag	UNP I4E574
F	-12	HIS	-	expression tag	UNP I4E574
F	-11	HIS	-	expression tag	UNP I4E574
F	-10	HIS	-	expression tag	UNP I4E574
F	-9	HIS	-	expression tag	UNP I4E574
F	-8	HIS	-	expression tag	UNP I4E574
F	-7	HIS	-	expression tag	UNP I4E574
F	-6	GLU	-	expression tag	UNP I4E574
F	-5	ASN	-	expression tag	UNP I4E574
F	-4	LEU	-	expression tag	UNP I4E574
F	-3	TYR	-	expression tag	UNP I4E574
F	-2	PHE	-	expression tag	UNP I4E574
F	-1	GLN	-	expression tag	UNP I4E574
F	0	GLY	-	expression tag	UNP I4E574
G	-12	HIS	-	expression tag	UNP I4E574
G	-11	HIS	-	expression tag	UNP I4E574
G	-10	HIS	-	expression tag	UNP I4E574
G	-9	HIS	-	expression tag	UNP I4E574
G	-8	HIS	-	expression tag	UNP I4E574
G	-7	HIS	-	expression tag	UNP I4E574
G	-6	GLU	-	expression tag	UNP I4E574
G	-5	ASN	-	expression tag	UNP I4E574
G	-4	LEU	-	expression tag	UNP I4E574
G	-3	TYR	-	expression tag	UNP I4E574
G	-2	PHE	-	expression tag	UNP I4E574
G	-1	GLN	-	expression tag	UNP I4E574
G	0	GLY	-	expression tag	UNP I4E574
H	-12	HIS	-	expression tag	UNP I4E574
H	-11	HIS	-	expression tag	UNP I4E574
H	-10	HIS	-	expression tag	UNP I4E574
H	-9	HIS	-	expression tag	UNP I4E574
H	-8	HIS	-	expression tag	UNP I4E574
H	-7	HIS	-	expression tag	UNP I4E574
H	-6	GLU	-	expression tag	UNP I4E574
H	-5	ASN	-	expression tag	UNP I4E574

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Chain	Residue	Modelled	Actual	Comment	Reference
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H	-3	TYR	-	expression tag	UNP I4E574
H	-2	PHE	-	expression tag	UNP I4E574
H	-1	GLN	-	expression tag	UNP I4E574
H	0	GLY	-	expression tag	UNP I4E574
I	-12	HIS	-	expression tag	UNP I4E574
I	-11	HIS	-	expression tag	UNP I4E574
I	-10	HIS	-	expression tag	UNP I4E574
I	-9	HIS	-	expression tag	UNP I4E574
I	-8	HIS	-	expression tag	UNP I4E574
I	-7	HIS	-	expression tag	UNP I4E574
I	-6	GLU	-	expression tag	UNP I4E574
I	-5	ASN	-	expression tag	UNP I4E574
I	-4	LEU	-	expression tag	UNP I4E574
I	-3	TYR	-	expression tag	UNP I4E574
I	-2	PHE	-	expression tag	UNP I4E574
I	-1	GLN	-	expression tag	UNP I4E574
I	0	GLY	-	expression tag	UNP I4E574
J	-12	HIS	-	expression tag	UNP I4E574
J	-11	HIS	-	expression tag	UNP I4E574
J	-10	HIS	-	expression tag	UNP I4E574
J	-9	HIS	-	expression tag	UNP I4E574
J	-8	HIS	-	expression tag	UNP I4E574
J	-7	HIS	-	expression tag	UNP I4E574
J	-6	GLU	-	expression tag	UNP I4E574
J	-5	ASN	-	expression tag	UNP I4E574
J	-4	LEU	-	expression tag	UNP I4E574
J	-3	TYR	-	expression tag	UNP I4E574
J	-2	PHE	-	expression tag	UNP I4E574
J	-1	GLN	-	expression tag	UNP I4E574
J	0	GLY	-	expression tag	UNP I4E574
K	-12	HIS	-	expression tag	UNP I4E574
K	-11	HIS	-	expression tag	UNP I4E574
K	-10	HIS	-	expression tag	UNP I4E574
K	-9	HIS	-	expression tag	UNP I4E574
K	-8	HIS	-	expression tag	UNP I4E574
K	-7	HIS	-	expression tag	UNP I4E574
K	-6	GLU	-	expression tag	UNP I4E574
K	-5	ASN	-	expression tag	UNP I4E574
K	-4	LEU	-	expression tag	UNP I4E574
K	-3	TYR	-	expression tag	UNP I4E574
K	-2	PHE	-	expression tag	UNP I4E574

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Chain	Residue	Modelled	Actual	Comment	Reference
K	-1	GLN	-	expression tag	UNP I4E574
K	0	GLY	-	expression tag	UNP I4E574
L	-12	HIS	-	expression tag	UNP I4E574
L	-11	HIS	-	expression tag	UNP I4E574
L	-10	HIS	-	expression tag	UNP I4E574
L	-9	HIS	-	expression tag	UNP I4E574
L	-8	HIS	-	expression tag	UNP I4E574
L	-7	HIS	-	expression tag	UNP I4E574
L	-6	GLU	-	expression tag	UNP I4E574
L	-5	ASN	-	expression tag	UNP I4E574
L	-4	LEU	-	expression tag	UNP I4E574
L	-3	TYR	-	expression tag	UNP I4E574
L	-2	PHE	-	expression tag	UNP I4E574
L	-1	GLN	-	expression tag	UNP I4E574
L	0	GLY	-	expression tag	UNP I4E574
M	-12	HIS	-	expression tag	UNP I4E574
M	-11	HIS	-	expression tag	UNP I4E574
M	-10	HIS	-	expression tag	UNP I4E574
M	-9	HIS	-	expression tag	UNP I4E574
M	-8	HIS	-	expression tag	UNP I4E574
M	-7	HIS	-	expression tag	UNP I4E574
M	-6	GLU	-	expression tag	UNP I4E574
M	-5	ASN	-	expression tag	UNP I4E574
M	-4	LEU	-	expression tag	UNP I4E574
M	-3	TYR	-	expression tag	UNP I4E574
M	-2	PHE	-	expression tag	UNP I4E574
M	-1	GLN	-	expression tag	UNP I4E574
M	0	GLY	-	expression tag	UNP I4E574
N	-12	HIS	-	expression tag	UNP I4E574
N	-11	HIS	-	expression tag	UNP I4E574
N	-10	HIS	-	expression tag	UNP I4E574
N	-9	HIS	-	expression tag	UNP I4E574
N	-8	HIS	-	expression tag	UNP I4E574
N	-7	HIS	-	expression tag	UNP I4E574
N	-6	GLU	-	expression tag	UNP I4E574
N	-5	ASN	-	expression tag	UNP I4E574
N	-4	LEU	-	expression tag	UNP I4E574
N	-3	TYR	-	expression tag	UNP I4E574
N	-2	PHE	-	expression tag	UNP I4E574
N	-1	GLN	-	expression tag	UNP I4E574
N	0	GLY	-	expression tag	UNP I4E574
O	-12	HIS	-	expression tag	UNP I4E574

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Chain	Residue	Modelled	Actual	Comment	Reference
O	-11	HIS	-	expression tag	UNP I4E574
O	-10	HIS	-	expression tag	UNP I4E574
O	-9	HIS	-	expression tag	UNP I4E574
O	-8	HIS	-	expression tag	UNP I4E574
O	-7	HIS	-	expression tag	UNP I4E574
O	-6	GLU	-	expression tag	UNP I4E574
O	-5	ASN	-	expression tag	UNP I4E574
O	-4	LEU	-	expression tag	UNP I4E574
O	-3	TYR	-	expression tag	UNP I4E574
O	-2	PHE	-	expression tag	UNP I4E574
O	-1	GLN	-	expression tag	UNP I4E574
O	0	GLY	-	expression tag	UNP I4E574
P	-12	HIS	-	expression tag	UNP I4E574
P	-11	HIS	-	expression tag	UNP I4E574
P	-10	HIS	-	expression tag	UNP I4E574
P	-9	HIS	-	expression tag	UNP I4E574
P	-8	HIS	-	expression tag	UNP I4E574
P	-7	HIS	-	expression tag	UNP I4E574
P	-6	GLU	-	expression tag	UNP I4E574
P	-5	ASN	-	expression tag	UNP I4E574
P	-4	LEU	-	expression tag	UNP I4E574
P	-3	TYR	-	expression tag	UNP I4E574
P	-2	PHE	-	expression tag	UNP I4E574
P	-1	GLN	-	expression tag	UNP I4E574
P	0	GLY	-	expression tag	UNP I4E574
Q	-12	HIS	-	expression tag	UNP I4E574
Q	-11	HIS	-	expression tag	UNP I4E574
Q	-10	HIS	-	expression tag	UNP I4E574
Q	-9	HIS	-	expression tag	UNP I4E574
Q	-8	HIS	-	expression tag	UNP I4E574
Q	-7	HIS	-	expression tag	UNP I4E574
Q	-6	GLU	-	expression tag	UNP I4E574
Q	-5	ASN	-	expression tag	UNP I4E574
Q	-4	LEU	-	expression tag	UNP I4E574
Q	-3	TYR	-	expression tag	UNP I4E574
Q	-2	PHE	-	expression tag	UNP I4E574
Q	-1	GLN	-	expression tag	UNP I4E574
Q	0	GLY	-	expression tag	UNP I4E574
R	-12	HIS	-	expression tag	UNP I4E574
R	-11	HIS	-	expression tag	UNP I4E574
R	-10	HIS	-	expression tag	UNP I4E574
R	-9	HIS	-	expression tag	UNP I4E574

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Chain	Residue	Modelled	Actual	Comment	Reference
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R	-7	HIS	-	expression tag	UNP I4E574
R	-6	GLU	-	expression tag	UNP I4E574
R	-5	ASN	-	expression tag	UNP I4E574
R	-4	LEU	-	expression tag	UNP I4E574
R	-3	TYR	-	expression tag	UNP I4E574
R	-2	PHE	-	expression tag	UNP I4E574
R	-1	GLN	-	expression tag	UNP I4E574
R	0	GLY	-	expression tag	UNP I4E574
S	-12	HIS	-	expression tag	UNP I4E574
S	-11	HIS	-	expression tag	UNP I4E574
S	-10	HIS	-	expression tag	UNP I4E574
S	-9	HIS	-	expression tag	UNP I4E574
S	-8	HIS	-	expression tag	UNP I4E574
S	-7	HIS	-	expression tag	UNP I4E574
S	-6	GLU	-	expression tag	UNP I4E574
S	-5	ASN	-	expression tag	UNP I4E574
S	-4	LEU	-	expression tag	UNP I4E574
S	-3	TYR	-	expression tag	UNP I4E574
S	-2	PHE	-	expression tag	UNP I4E574
S	-1	GLN	-	expression tag	UNP I4E574
S	0	GLY	-	expression tag	UNP I4E574
T	-12	HIS	-	expression tag	UNP I4E574
T	-11	HIS	-	expression tag	UNP I4E574
T	-10	HIS	-	expression tag	UNP I4E574
T	-9	HIS	-	expression tag	UNP I4E574
T	-8	HIS	-	expression tag	UNP I4E574
T	-7	HIS	-	expression tag	UNP I4E574
T	-6	GLU	-	expression tag	UNP I4E574
T	-5	ASN	-	expression tag	UNP I4E574
T	-4	LEU	-	expression tag	UNP I4E574
T	-3	TYR	-	expression tag	UNP I4E574
T	-2	PHE	-	expression tag	UNP I4E574
T	-1	GLN	-	expression tag	UNP I4E574
T	0	GLY	-	expression tag	UNP I4E574
U	-12	HIS	-	expression tag	UNP I4E574
U	-11	HIS	-	expression tag	UNP I4E574
U	-10	HIS	-	expression tag	UNP I4E574
U	-9	HIS	-	expression tag	UNP I4E574
U	-8	HIS	-	expression tag	UNP I4E574
U	-7	HIS	-	expression tag	UNP I4E574
U	-6	GLU	-	expression tag	UNP I4E574

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Chain	Residue	Modelled	Actual	Comment	Reference
U	-5	ASN	-	expression tag	UNP I4E574
U	-4	LEU	-	expression tag	UNP I4E574
U	-3	TYR	-	expression tag	UNP I4E574
U	-2	PHE	-	expression tag	UNP I4E574
U	-1	GLN	-	expression tag	UNP I4E574
U	0	GLY	-	expression tag	UNP I4E574
V	-12	HIS	-	expression tag	UNP I4E574
V	-11	HIS	-	expression tag	UNP I4E574
V	-10	HIS	-	expression tag	UNP I4E574
V	-9	HIS	-	expression tag	UNP I4E574
V	-8	HIS	-	expression tag	UNP I4E574
V	-7	HIS	-	expression tag	UNP I4E574
V	-6	GLU	-	expression tag	UNP I4E574
V	-5	ASN	-	expression tag	UNP I4E574
V	-4	LEU	-	expression tag	UNP I4E574
V	-3	TYR	-	expression tag	UNP I4E574
V	-2	PHE	-	expression tag	UNP I4E574
V	-1	GLN	-	expression tag	UNP I4E574
V	0	GLY	-	expression tag	UNP I4E574
W	-12	HIS	-	expression tag	UNP I4E574
W	-11	HIS	-	expression tag	UNP I4E574
W	-10	HIS	-	expression tag	UNP I4E574
W	-9	HIS	-	expression tag	UNP I4E574
W	-8	HIS	-	expression tag	UNP I4E574
W	-7	HIS	-	expression tag	UNP I4E574
W	-6	GLU	-	expression tag	UNP I4E574
W	-5	ASN	-	expression tag	UNP I4E574
W	-4	LEU	-	expression tag	UNP I4E574
W	-3	TYR	-	expression tag	UNP I4E574
W	-2	PHE	-	expression tag	UNP I4E574
W	-1	GLN	-	expression tag	UNP I4E574
W	0	GLY	-	expression tag	UNP I4E574
X	-12	HIS	-	expression tag	UNP I4E574
X	-11	HIS	-	expression tag	UNP I4E574
X	-10	HIS	-	expression tag	UNP I4E574
X	-9	HIS	-	expression tag	UNP I4E574
X	-8	HIS	-	expression tag	UNP I4E574
X	-7	HIS	-	expression tag	UNP I4E574
X	-6	GLU	-	expression tag	UNP I4E574
X	-5	ASN	-	expression tag	UNP I4E574
X	-4	LEU	-	expression tag	UNP I4E574
X	-3	TYR	-	expression tag	UNP I4E574

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Chain	Residue	Modelled	Actual	Comment	Reference
X	-2	PHE	-	expression tag	UNP I4E574
X	-1	GLN	-	expression tag	UNP I4E574
X	0	GLY	-	expression tag	UNP I4E574
Y	-12	HIS	-	expression tag	UNP I4E574
Y	-11	HIS	-	expression tag	UNP I4E574
Y	-10	HIS	-	expression tag	UNP I4E574
Y	-9	HIS	-	expression tag	UNP I4E574
Y	-8	HIS	-	expression tag	UNP I4E574
Y	-7	HIS	-	expression tag	UNP I4E574
Y	-6	GLU	-	expression tag	UNP I4E574
Y	-5	ASN	-	expression tag	UNP I4E574
Y	-4	LEU	-	expression tag	UNP I4E574
Y	-3	TYR	-	expression tag	UNP I4E574
Y	-2	PHE	-	expression tag	UNP I4E574
Y	-1	GLN	-	expression tag	UNP I4E574
Y	0	GLY	-	expression tag	UNP I4E574
Z	-12	HIS	-	expression tag	UNP I4E574
Z	-11	HIS	-	expression tag	UNP I4E574
Z	-10	HIS	-	expression tag	UNP I4E574
Z	-9	HIS	-	expression tag	UNP I4E574
Z	-8	HIS	-	expression tag	UNP I4E574
Z	-7	HIS	-	expression tag	UNP I4E574
Z	-6	GLU	-	expression tag	UNP I4E574
Z	-5	ASN	-	expression tag	UNP I4E574
Z	-4	LEU	-	expression tag	UNP I4E574
Z	-3	TYR	-	expression tag	UNP I4E574
Z	-2	PHE	-	expression tag	UNP I4E574
Z	-1	GLN	-	expression tag	UNP I4E574
Z	0	GLY	-	expression tag	UNP I4E574
a	-12	HIS	-	expression tag	UNP I4E574
a	-11	HIS	-	expression tag	UNP I4E574
a	-10	HIS	-	expression tag	UNP I4E574
a	-9	HIS	-	expression tag	UNP I4E574
a	-8	HIS	-	expression tag	UNP I4E574
a	-7	HIS	-	expression tag	UNP I4E574
a	-6	GLU	-	expression tag	UNP I4E574
a	-5	ASN	-	expression tag	UNP I4E574
a	-4	LEU	-	expression tag	UNP I4E574
a	-3	TYR	-	expression tag	UNP I4E574
a	-2	PHE	-	expression tag	UNP I4E574
a	-1	GLN	-	expression tag	UNP I4E574
a	0	GLY	-	expression tag	UNP I4E574

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Chain	Residue	Modelled	Actual	Comment	Reference
b	-12	HIS	-	expression tag	UNP I4E574
b	-11	HIS	-	expression tag	UNP I4E574
b	-10	HIS	-	expression tag	UNP I4E574
b	-9	HIS	-	expression tag	UNP I4E574
b	-8	HIS	-	expression tag	UNP I4E574
b	-7	HIS	-	expression tag	UNP I4E574
b	-6	GLU	-	expression tag	UNP I4E574
b	-5	ASN	-	expression tag	UNP I4E574
b	-4	LEU	-	expression tag	UNP I4E574
b	-3	TYR	-	expression tag	UNP I4E574
b	-2	PHE	-	expression tag	UNP I4E574
b	-1	GLN	-	expression tag	UNP I4E574
b	0	GLY	-	expression tag	UNP I4E574

- Molecule 2 is a protein called Acyldepsipeptide-14.

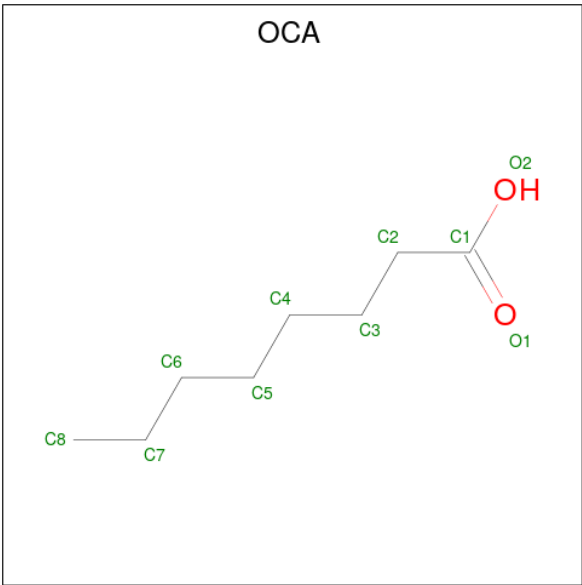
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	c	6	Total	C	F	N	O	0	0	0
			48	33	2	6	7			
2	e	6	Total	C	F	N	O	0	0	0
			48	33	2	6	7			
2	f	6	Total	C	F	N	O	0	0	0
			48	33	2	6	7			
2	g	6	Total	C	F	N	O	0	0	0
			48	33	2	6	7			
2	h	6	Total	C	F	N	O	0	0	0
			48	33	2	6	7			
2	i	6	Total	C	F	N	O	0	0	0
			48	33	2	6	7			
2	j	6	Total	C	F	N	O	0	0	0
			48	33	2	6	7			
2	k	6	Total	C	F	N	O	0	0	0
			48	33	2	6	7			
2	l	6	Total	C	F	N	O	0	0	0
			48	33	2	6	7			
2	m	6	Total	C	F	N	O	0	0	0
			48	33	2	6	7			
2	n	6	Total	C	F	N	O	0	0	0
			48	33	2	6	7			
2	o	6	Total	C	F	N	O	0	0	0
			48	33	2	6	7			
2	p	6	Total	C	F	N	O	0	0	0
			48	33	2	6	7			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	q	6	Total 48	C 33	F 2	N 6	O 7	0	0	0
2	r	6	Total 48	C 33	F 2	N 6	O 7	0	0	0
2	s	6	Total 48	C 33	F 2	N 6	O 7	0	0	0
2	t	6	Total 48	C 33	F 2	N 6	O 7	0	0	0
2	u	6	Total 48	C 33	F 2	N 6	O 7	0	0	0
2	v	6	Total 48	C 33	F 2	N 6	O 7	0	0	0
2	w	6	Total 48	C 33	F 2	N 6	O 7	0	0	0
2	x	6	Total 48	C 33	F 2	N 6	O 7	0	0	0
2	y	6	Total 48	C 33	F 2	N 6	O 7	0	0	0
2	z	6	Total 48	C 33	F 2	N 6	O 7	0	0	0
2	0	6	Total 48	C 33	F 2	N 6	O 7	0	0	0
2	1	6	Total 48	C 33	F 2	N 6	O 7	0	0	0
2	2	6	Total 48	C 33	F 2	N 6	O 7	0	0	0
2	3	6	Total 48	C 33	F 2	N 6	O 7	0	0	0
2	4	6	Total 48	C 33	F 2	N 6	O 7	0	0	0

- Molecule 3 is OCTANOIC ACID (CAPRYLIC ACID) (CCD ID: OCA) (formula: C₈H₁₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	c	1	Total	C	O	0	0
			9	8	1		
3	e	1	Total	C	O	0	0
			9	8	1		
3	f	1	Total	C	O	0	0
			9	8	1		
3	g	1	Total	C	O	0	0
			9	8	1		
3	h	1	Total	C	O	0	0
			9	8	1		
3	i	1	Total	C	O	0	0
			9	8	1		
3	j	1	Total	C	O	0	0
			9	8	1		
3	k	1	Total	C	O	0	0
			9	8	1		
3	l	1	Total	C	O	0	0
			9	8	1		
3	m	1	Total	C	O	0	0
			9	8	1		
3	n	1	Total	C	O	0	0
			9	8	1		
3	o	1	Total	C	O	0	0
			9	8	1		
3	p	1	Total	C	O	0	0
			9	8	1		
3	q	1	Total	C	O	0	0
			9	8	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	r	1	Total C O 9 8 1	0	0
3	s	1	Total C O 9 8 1	0	0
3	t	1	Total C O 9 8 1	0	0
3	u	1	Total C O 9 8 1	0	0
3	v	1	Total C O 9 8 1	0	0
3	w	1	Total C O 9 8 1	0	0
3	x	1	Total C O 9 8 1	0	0
3	y	1	Total C O 9 8 1	0	0
3	z	1	Total C O 9 8 1	0	0
3	0	1	Total C O 9 8 1	0	0
3	1	1	Total C O 9 8 1	0	0
3	2	1	Total C O 9 8 1	0	0
3	3	1	Total C O 9 8 1	0	0
3	4	1	Total C O 9 8 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	25	Total O 25 25	0	0
4	B	27	Total O 27 27	0	0
4	C	21	Total O 21 21	0	0
4	D	17	Total O 17 17	0	0
4	E	24	Total O 24 24	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	F	22	Total O 22 22	0	0
4	G	26	Total O 26 26	0	0
4	H	19	Total O 19 19	0	0
4	I	20	Total O 20 20	0	0
4	J	21	Total O 21 21	0	0
4	K	32	Total O 32 32	0	0
4	L	26	Total O 26 26	0	0
4	M	31	Total O 31 31	0	0
4	N	35	Total O 35 35	0	0
4	O	22	Total O 22 22	0	0
4	P	30	Total O 30 30	0	0
4	Q	22	Total O 22 22	0	0
4	R	27	Total O 27 27	0	0
4	S	17	Total O 17 17	0	0
4	T	24	Total O 24 24	0	0
4	U	25	Total O 25 25	0	0
4	V	29	Total O 29 29	0	0
4	W	24	Total O 24 24	0	0
4	X	24	Total O 24 24	0	0
4	Y	17	Total O 17 17	0	0
4	Z	32	Total O 32 32	0	0

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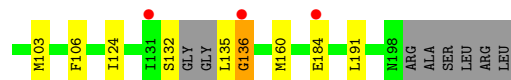
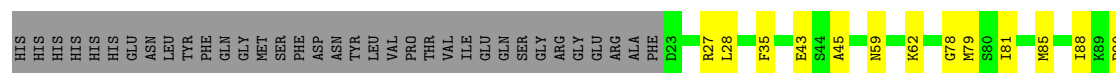
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	a	29	Total 29	O 29	0	0
4	b	37	Total 37	O 37	0	0
4	0	1	Total 1	O 1	0	0

- Molecule 1: ATP-dependent Clp protease proteolytic subunit

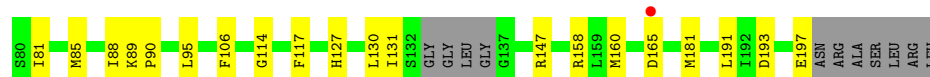
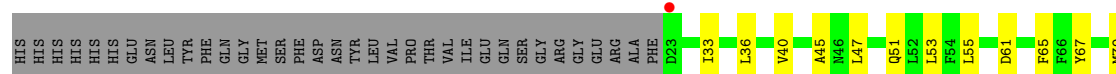




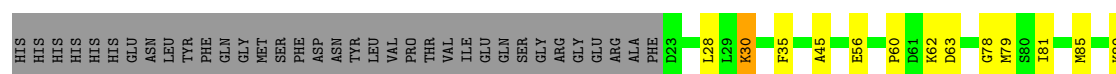
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



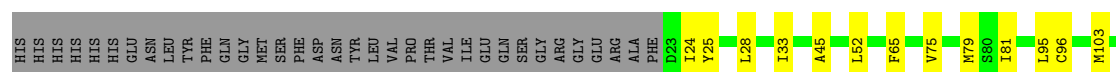
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit

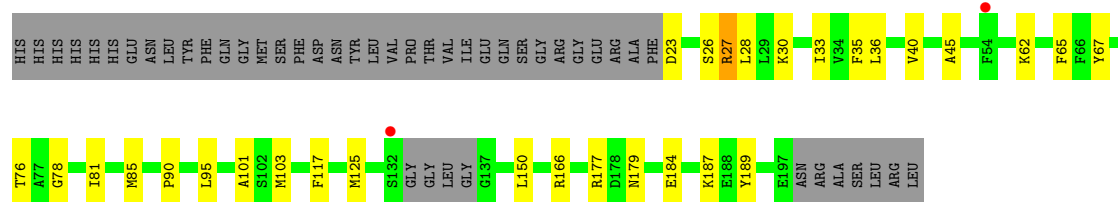


- Molecule 1: ATP-dependent Clp protease proteolytic subunit



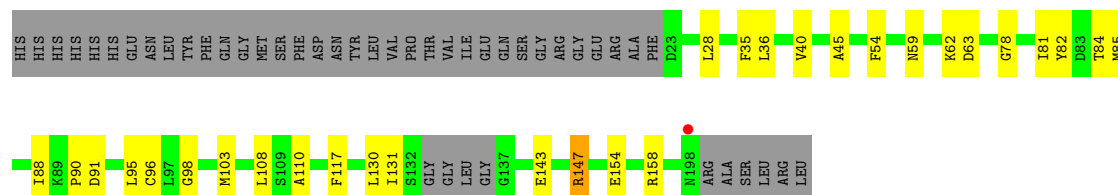
- Molecule 1: ATP-dependent Clp protease proteolytic subunit





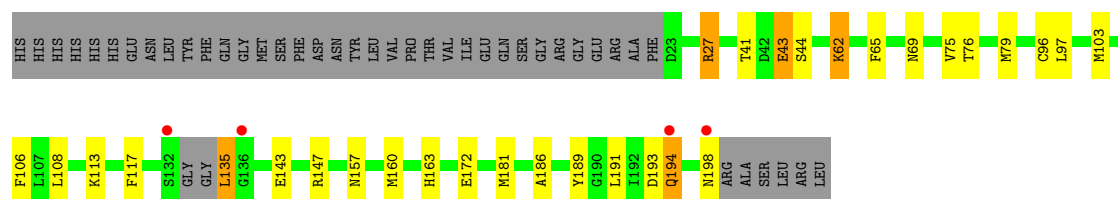
- Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain J: 65% 13% 21%



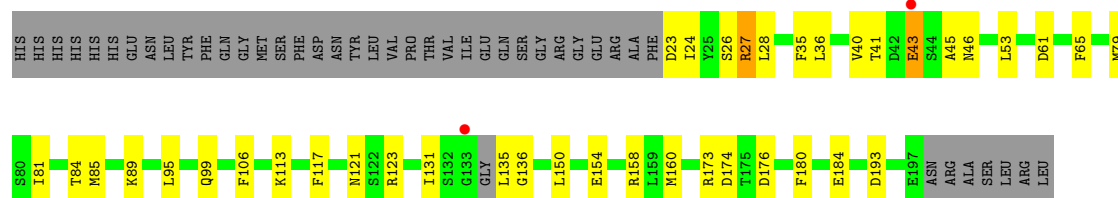
- Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain K: 66% 12% 20%



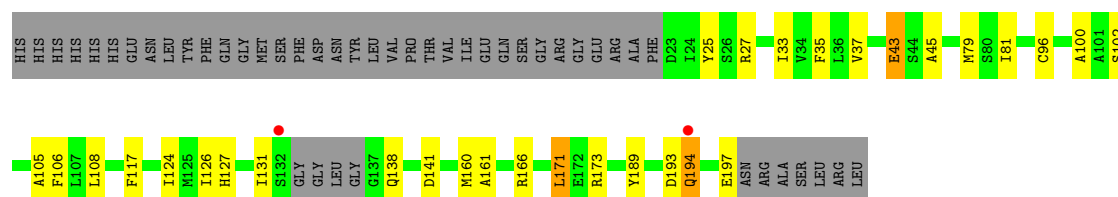
- Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain L: 62% 18% 20%

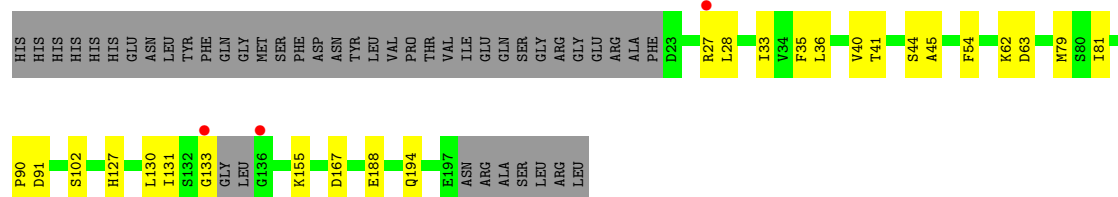


- Molecule 1: ATP-dependent Clp protease proteolytic subunit

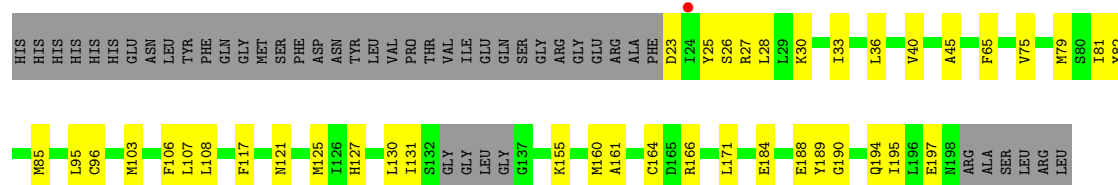
Chain M: 65% 13% 21%



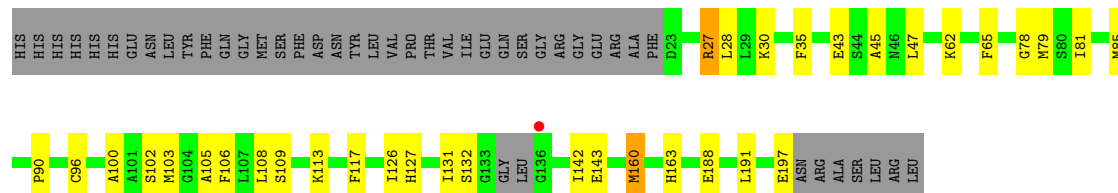
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



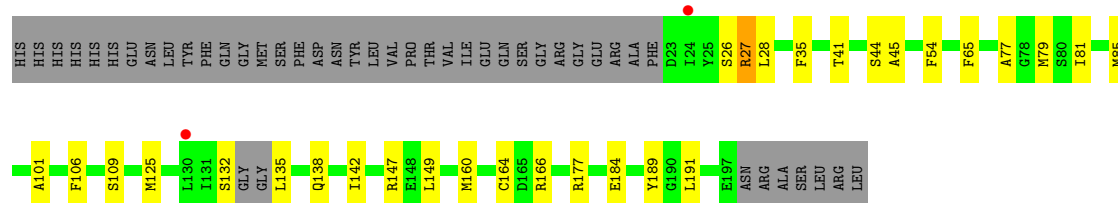
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



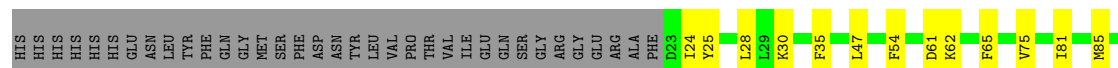
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit

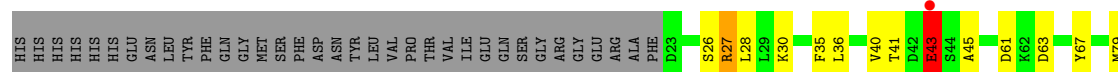


- Molecule 1: ATP-dependent Clp protease proteolytic subunit

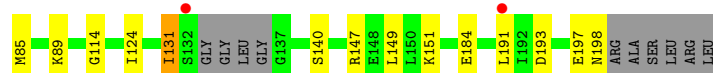
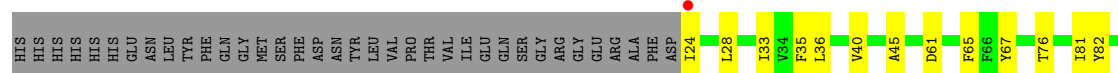




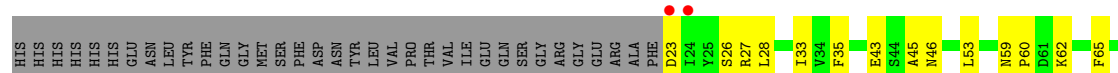
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



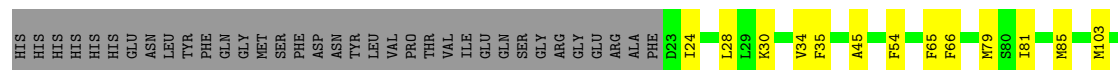
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit

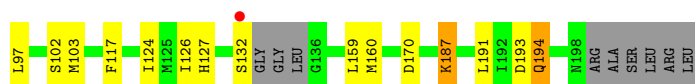


- Chain X:
-
- | Residue | Category |
|---------|----------|
| HIS | Grey |
| HIS | Grey |
| HIS | Grey |
| HIS | Grey |
| HIS | Grey |
| GLU | Grey |
| ASN | Grey |
| LEU | Grey |
| TVR | Grey |
| PHE | Grey |
| GLN | Grey |
| GLY | Grey |
| MET | Grey |
| SER | Grey |
| PHE | Grey |
| ASP | Grey |
| ASN | Grey |
| TYR | Grey |
| LEU | Grey |
| VAL | Grey |
| PRO | Grey |
| THR | Grey |
| VAL | Grey |
| ILE | Grey |
| GLU | Grey |
| GLN | Grey |
| SER | Grey |
| GLY | Grey |
| ARG | Grey |
| GLY | Grey |
| GLU | Grey |
| ARG | Grey |
| ALA | Grey |
| PHE | Grey |
| D23 | Green |
| R27 | Yellow |
| L28 | Yellow |
| F35 | Yellow |
| L36 | Yellow |
| V40 | Yellow |
| T41 | Yellow |
| S44 | Yellow |
| A45 | Yellow |
| M46 | Yellow |
| N59 | Yellow |
| K62 | Yellow |
| I81 | Yellow |
| S102 | Yellow |
| M103 | Yellow |
| T124 | Yellow |
-
- | Residue | Category |
|---------|----------|
| H127 | Green |
| G133 | Green |
| GLY | Green |
| L135 | Green |
| G136 | Green |
| G137 | Green |
| Q138 | Green |
| E154 | Green |
| K155 | Green |
| R158 | Green |
| L159 | Green |
| R173 | Green |
| D174 | Green |
| R177 | Green |
| E184 | Green |
| L191 | Green |
| E197 | Green |
| ASN | Green |
| ARG | Green |
| ALA | Green |
| ALA | Green |
| SER | Green |
| LEU | Green |
| ARG | Green |
| LEU | Green |

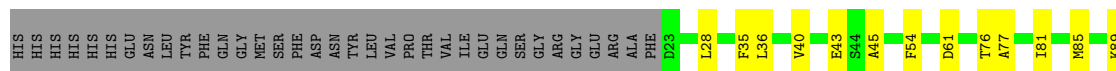
- Chain Y: 

- [illegible]

- Chain a: 



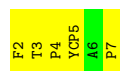
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



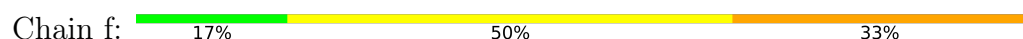
- Molecule 2: Acyldepsipeptide-14



- Molecule 2: Acyldepsipeptide-14



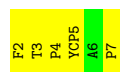
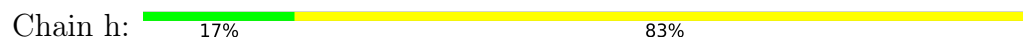
- Molecule 2: Acyldepsipeptide-14



- Molecule 2: Acyldepsipeptide-14



- Molecule 2: Acyldepsipeptide-14



- Molecule 2: Acyldepsipeptide-14





- Molecule 2: Acyldepsipeptide-14



- Molecule 2: Acyldepsipeptide-14



- Molecule 2: Acyldepsipeptide-14



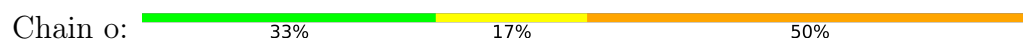
- Molecule 2: Acyldepsipeptide-14



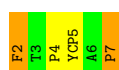
- Molecule 2: Acyldepsipeptide-14



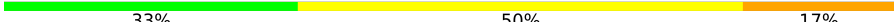
- Molecule 2: Acyldepsipeptide-14



- Molecule 2: Acyldepsipeptide-14

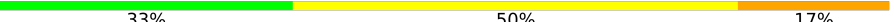


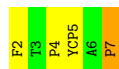
- Molecule 2: Acyldepsipeptide-14

Chain q:  33% 50% 17%



- Molecule 2: Acyldepsipeptide-14

Chain r:  33% 50% 17%

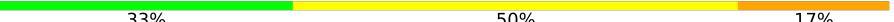


- Molecule 2: Acyldepsipeptide-14

Chain s:  50% 50%

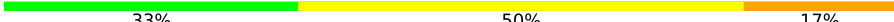


- Molecule 2: Acyldepsipeptide-14

Chain t:  33% 50% 17%

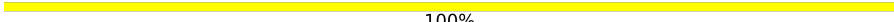


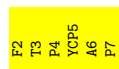
- Molecule 2: Acyldepsipeptide-14

Chain u:  33% 50% 17%

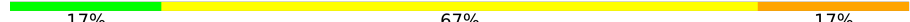


- Molecule 2: Acyldepsipeptide-14

Chain v:  100%

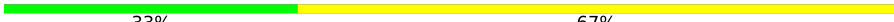


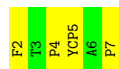
- Molecule 2: Acyldepsipeptide-14

Chain w:  17% 67% 17%

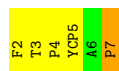


- Molecule 2: Acyldepsipeptide-14

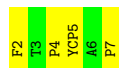
Chain x:  33% 67%



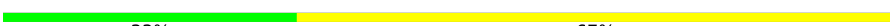
● Molecule 2: Acyldepsipeptide-14

Chain y:  17% 67% 17%

● Molecule 2: Acyldepsipeptide-14

Chain z:  33% 67%

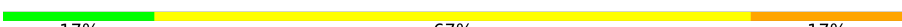
● Molecule 2: Acyldepsipeptide-14

Chain 0:  33% 67%

● Molecule 2: Acyldepsipeptide-14

Chain 1:  33% 50% 17%

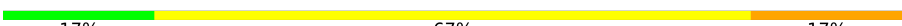
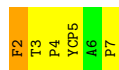
● Molecule 2: Acyldepsipeptide-14

Chain 2:  17% 67% 17%

● Molecule 2: Acyldepsipeptide-14

Chain 3:  67% 33%

● Molecule 2: Acyldepsipeptide-14

Chain 4:  17% 67% 17%

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	117.24Å 195.98Å 139.89Å 90.00° 97.42° 90.00°	Depositor
Resolution (Å)	49.63 – 2.70 49.63 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.6 (49.63-2.70) 98.5 (49.63-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.49 (at 2.69Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575)	Depositor
R, R_{free}	0.192 , 0.248 0.196 , 0.251	Depositor DCC
R_{free} test set	2000 reflections (1.17%)	wwPDB-VP
Wilson B-factor (Å ²)	43.2	Xtriage
Anisotropy	0.211	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 44.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	40076	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: OCA, WFP, MP8, ALO, YCP

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	0.40	0/1379	0.77	3/1856 (0.2%)
1	B	0.49	1/1366 (0.1%)	0.79	7/1837 (0.4%)
1	C	0.42	0/1365	0.70	2/1836 (0.1%)
1	D	0.45	0/1370	0.76	5/1843 (0.3%)
1	E	0.46	0/1378	0.82	5/1854 (0.3%)
1	F	0.46	0/1358	0.68	1/1827 (0.1%)
1	G	0.45	0/1370	0.70	3/1843 (0.2%)
1	H	0.43	0/1378	0.78	5/1854 (0.3%)
1	I	0.43	0/1358	0.72	4/1827 (0.2%)
1	J	0.42	0/1366	0.70	4/1838 (0.2%)
1	K	0.50	0/1378	0.89	9/1854 (0.5%)
1	L	0.46	0/1374	0.71	3/1848 (0.2%)
1	M	0.46	0/1358	0.80	8/1827 (0.4%)
1	N	0.39	0/1366	0.74	4/1837 (0.2%)
1	O	0.46	0/1366	0.77	2/1838 (0.1%)
1	P	0.48	1/1366 (0.1%)	0.83	5/1837 (0.3%)
1	Q	0.41	0/1370	0.79	5/1843 (0.3%)
1	R	0.43	0/1382	0.73	2/1859 (0.1%)
1	S	0.47	0/1370	0.81	5/1843 (0.3%)
1	T	0.43	1/1358 (0.1%)	0.69	2/1827 (0.1%)
1	U	0.47	0/1370	0.92	8/1844 (0.4%)
1	V	0.43	0/1360	0.66	0/1830
1	W	0.41	0/1366	0.72	2/1837 (0.1%)
1	X	0.41	0/1374	0.71	3/1848 (0.2%)
1	Y	0.40	0/1366	0.77	5/1838 (0.3%)
1	Z	0.40	0/1362	0.65	2/1832 (0.1%)
1	a	0.41	0/1370	0.72	3/1843 (0.2%)
1	b	0.52	0/1364	0.74	3/1835 (0.2%)
2	0	4.24	2/11 (18.2%)	1.29	0/12
2	1	4.18	2/11 (18.2%)	1.43	0/12
2	2	4.32	2/11 (18.2%)	1.98	0/12
2	3	4.23	2/11 (18.2%)	1.47	0/12

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
2	4	4.28	2/11 (18.2%)	1.30	0/12
2	c	4.19	2/11 (18.2%)	1.46	0/12
2	e	4.37	2/11 (18.2%)	1.26	0/12
2	f	4.17	2/11 (18.2%)	1.56	0/12
2	g	4.37	2/11 (18.2%)	1.50	0/12
2	h	4.41	2/11 (18.2%)	1.59	0/12
2	i	4.33	2/11 (18.2%)	1.51	0/12
2	j	4.26	2/11 (18.2%)	1.72	0/12
2	k	4.24	2/11 (18.2%)	1.41	0/12
2	l	4.28	2/11 (18.2%)	1.45	0/12
2	m	4.26	2/11 (18.2%)	2.26	0/12
2	n	4.24	2/11 (18.2%)	1.70	0/12
2	o	4.21	2/11 (18.2%)	1.55	0/12
2	p	4.14	2/11 (18.2%)	1.50	0/12
2	q	4.25	2/11 (18.2%)	1.37	0/12
2	r	4.38	2/11 (18.2%)	1.47	0/12
2	s	4.31	2/11 (18.2%)	1.65	0/12
2	t	4.40	2/11 (18.2%)	1.59	0/12
2	u	4.16	2/11 (18.2%)	2.11	0/12
2	v	4.18	2/11 (18.2%)	1.89	0/12
2	w	4.30	2/11 (18.2%)	1.35	0/12
2	x	4.21	2/11 (18.2%)	1.81	0/12
2	y	4.13	2/11 (18.2%)	1.59	0/12
2	z	4.29	2/11 (18.2%)	1.62	0/12
All	All	0.58	59/38616 (0.2%)	0.76	110/51871 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	S	0	1
1	W	0	1
All	All	0	2

All (59) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	i	4	PRO	N-CD	-11.09	1.32	1.47
2	h	4	PRO	N-CD	-11.09	1.32	1.47
2	e	4	PRO	N-CD	-11.08	1.32	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	r	4	PRO	N-CD	-11.02	1.32	1.47
2	s	4	PRO	N-CD	-10.95	1.32	1.47
2	2	4	PRO	N-CD	-10.95	1.32	1.47
2	0	4	PRO	N-CD	-10.92	1.32	1.47
2	4	4	PRO	N-CD	-10.89	1.32	1.47
2	j	4	PRO	N-CD	-10.74	1.32	1.47
2	t	4	PRO	N-CD	-10.74	1.32	1.47
2	3	4	PRO	N-CD	-10.67	1.32	1.47
2	w	4	PRO	N-CD	-10.66	1.32	1.47
2	n	4	PRO	N-CD	-10.65	1.32	1.47
2	o	4	PRO	N-CD	-10.64	1.32	1.47
2	1	4	PRO	N-CD	-10.62	1.32	1.47
2	c	4	PRO	N-CD	-10.59	1.32	1.47
2	f	4	PRO	N-CD	-10.55	1.32	1.47
2	g	4	PRO	N-CD	-10.54	1.32	1.47
2	z	4	PRO	N-CD	-10.52	1.33	1.47
2	m	4	PRO	N-CD	-10.50	1.33	1.47
2	l	4	PRO	N-CD	-10.45	1.33	1.47
2	q	4	PRO	N-CD	-10.44	1.33	1.47
2	x	4	PRO	N-CD	-10.41	1.33	1.47
2	p	4	PRO	N-CD	-10.40	1.33	1.47
2	u	4	PRO	N-CD	-10.39	1.33	1.47
2	k	4	PRO	N-CD	-10.24	1.33	1.47
2	v	4	PRO	N-CD	-10.23	1.33	1.47
2	y	4	PRO	N-CD	-10.13	1.33	1.47
2	t	4	PRO	N-CA	7.18	1.57	1.47
2	q	4	PRO	N-CA	6.79	1.57	1.47
1	P	27	ARG	CZ-NH2	6.79	1.42	1.33
2	l	4	PRO	N-CA	6.79	1.57	1.47
2	w	4	PRO	N-CA	6.75	1.57	1.47
2	g	4	PRO	N-CA	6.68	1.57	1.47
2	h	4	PRO	N-CA	6.61	1.56	1.47
2	e	4	PRO	N-CA	6.50	1.56	1.47
2	r	4	PRO	N-CA	6.49	1.56	1.47
2	z	4	PRO	N-CA	6.42	1.56	1.47
2	n	4	PRO	N-CA	6.41	1.56	1.47
2	v	4	PRO	N-CA	6.35	1.56	1.47
2	o	4	PRO	N-CA	6.35	1.56	1.47
2	s	4	PRO	N-CA	6.34	1.56	1.47
2	i	4	PRO	N-CA	6.33	1.56	1.47
2	k	4	PRO	N-CA	6.32	1.56	1.47
2	2	4	PRO	N-CA	6.30	1.56	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	4	4	PRO	N-CA	6.26	1.56	1.47
2	m	4	PRO	N-CA	6.22	1.56	1.47
2	x	4	PRO	N-CA	6.16	1.56	1.47
2	p	4	PRO	N-CA	6.08	1.56	1.47
2	j	4	PRO	N-CA	6.08	1.56	1.47
2	c	4	PRO	N-CA	5.99	1.56	1.47
2	f	4	PRO	N-CA	5.99	1.56	1.47
2	3	4	PRO	N-CA	5.92	1.55	1.47
2	0	4	PRO	N-CA	5.85	1.55	1.47
2	u	4	PRO	N-CA	5.73	1.55	1.47
2	y	4	PRO	N-CA	5.70	1.55	1.47
2	1	4	PRO	N-CA	5.48	1.55	1.47
1	T	131	ILE	C-O	-5.48	1.18	1.23
1	B	30	LYS	CB-CG	5.16	1.68	1.52

All (110) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	27	ARG	CA-CB-CG	12.51	139.13	114.10
1	A	147	ARG	CA-CB-CG	-12.36	89.38	114.10
1	P	188	GLU	CA-CB-CG	-11.55	91.00	114.10
1	K	27	ARG	CG-CD-NE	-11.21	87.34	112.00
1	S	43	GLU	CA-CB-CG	11.07	136.25	114.10
1	G	30	LYS	CB-CG-CD	-10.97	86.06	111.30
1	I	27	ARG	CB-CG-CD	-10.71	86.66	111.30
1	Q	184	GLU	CA-CB-CG	-10.68	92.75	114.10
1	U	43	GLU	CB-CA-C	-10.55	94.23	110.90
1	P	27	ARG	NE-CZ-NH2	10.28	128.45	119.20
1	T	147	ARG	CG-CD-NE	-10.09	89.81	112.00
1	U	43	GLU	CA-CB-CG	9.36	132.82	114.10
1	b	43	GLU	CA-CB-CG	-9.07	95.95	114.10
1	P	27	ARG	CG-CD-NE	8.89	131.56	112.00
1	K	27	ARG	CB-CG-CD	-8.74	91.19	111.30
1	K	194	GLN	CA-CB-CG	8.73	131.56	114.10
1	U	43	GLU	N-CA-CB	8.51	122.41	110.07
1	P	27	ARG	NE-CZ-NH1	-8.48	113.02	121.50
1	B	30	LYS	CG-CD-CE	8.40	130.63	111.30
1	I	27	ARG	CD-NE-CZ	8.08	135.72	124.40
1	I	184	GLU	CA-CB-CG	8.05	130.20	114.10
1	N	27	ARG	CB-CG-CD	-7.85	93.24	111.30
1	b	43	GLU	N-CA-CB	-7.78	98.62	110.13
1	N	27	ARG	CB-CA-C	-7.69	97.78	110.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	147	ARG	CB-CG-CD	-7.68	93.62	111.30
1	R	194	GLN	CA-CB-CG	7.62	129.34	114.10
1	U	184	GLU	CA-CB-CG	7.56	129.21	114.10
1	D	147	ARG	NE-CZ-NH2	7.54	125.99	119.20
1	U	43	GLU	CB-CG-CD	7.52	125.39	112.60
1	Y	43	GLU	CA-CB-CG	-7.51	99.08	114.10
1	S	43	GLU	CB-CA-C	-7.44	99.51	110.96
1	G	194	GLN	CA-CB-CG	7.41	128.92	114.10
1	E	27	ARG	CG-CD-NE	7.38	128.23	112.00
1	E	27	ARG	CB-CG-CD	-7.33	94.43	111.30
1	Q	27	ARG	N-CA-CB	7.18	121.42	110.28
1	F	147	ARG	CG-CD-NE	7.12	127.66	112.00
1	K	43	GLU	CB-CG-CD	7.09	124.65	112.60
1	Y	43	GLU	N-CA-CB	-7.04	99.71	110.13
1	X	184	GLU	CB-CG-CD	7.02	124.54	112.60
1	Y	173	ARG	CB-CG-CD	-6.97	95.27	111.30
1	M	173	ARG	CA-CB-CG	6.88	127.86	114.10
1	B	162	LYS	CA-CB-CG	6.87	127.84	114.10
1	B	30	LYS	CB-CG-CD	-6.75	95.78	111.30
1	Q	26	SER	CA-C-N	-6.71	110.11	120.31
1	Q	26	SER	C-N-CA	-6.71	110.11	120.31
1	J	147	ARG	CG-CD-NE	-6.67	97.33	112.00
1	J	147	ARG	NE-CZ-NH1	-6.55	114.95	121.50
1	R	30	LYS	CG-CD-CE	-6.54	96.26	111.30
1	X	184	GLU	CB-CA-C	-6.52	99.76	110.85
1	E	184	GLU	CA-CB-CG	6.51	127.13	114.10
1	W	43	GLU	CA-CB-CG	6.44	126.98	114.10
1	X	27	ARG	CB-CG-CD	6.40	126.02	111.30
1	B	43	GLU	CB-CG-CD	6.38	123.45	112.60
1	L	43	GLU	CA-CB-CG	6.34	126.79	114.10
1	D	147	ARG	NE-CZ-NH1	-6.33	115.17	121.50
1	A	147	ARG	CB-CG-CD	6.31	125.80	111.30
1	b	97	LEU	CA-CB-CG	6.29	138.31	116.30
1	S	27	ARG	CB-CA-C	-6.25	98.67	110.67
1	L	27	ARG	CG-CD-NE	6.17	125.57	112.00
1	J	147	ARG	CA-CB-CG	6.09	126.28	114.10
1	Z	184	GLU	CA-CB-CG	-6.06	101.98	114.10
1	H	113	LYS	CG-CD-CE	-6.03	97.44	111.30
1	U	184	GLU	CB-CG-CD	5.96	122.73	112.60
1	N	194	GLN	CA-CB-CG	5.93	125.95	114.10
1	Q	27	ARG	CB-CA-C	-5.92	99.30	110.67
1	L	27	ARG	CB-CA-C	5.91	122.01	110.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	30	LYS	CA-CB-CG	5.90	125.91	114.10
1	M	194	GLN	CA-CB-CG	5.89	125.89	114.10
1	D	188	GLU	CB-CA-C	-5.88	99.75	110.63
1	H	147	ARG	CA-CB-CG	-5.87	102.36	114.10
1	a	187	LYS	CB-CG-CD	5.82	124.69	111.30
1	D	27	ARG	CB-CA-C	5.80	121.81	110.67
1	N	27	ARG	CG-CD-NE	5.79	124.74	112.00
1	K	27	ARG	CD-NE-CZ	5.78	132.49	124.40
1	H	147	ARG	CB-CG-CD	5.77	124.57	111.30
1	P	43	GLU	N-CA-CB	5.74	118.39	110.07
1	W	43	GLU	CB-CA-C	-5.71	101.87	110.90
1	M	43	GLU	CB-CG-CD	-5.70	102.92	112.60
1	M	27	ARG	N-CA-CB	5.66	119.05	110.28
1	B	42	ASP	CA-C-N	-5.66	110.91	120.58
1	B	42	ASP	C-N-CA	-5.66	110.91	120.58
1	H	184	GLU	CB-CA-C	-5.60	101.33	110.85
1	K	62	LYS	CD-CE-NZ	-5.53	94.22	111.90
1	E	27	ARG	CB-CA-C	5.51	121.82	110.31
1	a	194	GLN	CA-CB-CG	5.50	125.11	114.10
1	S	26	SER	CA-C-N	-5.50	111.95	120.31
1	S	26	SER	C-N-CA	-5.50	111.95	120.31
1	U	188	GLU	N-CA-CB	-5.44	101.85	110.28
1	A	30	LYS	CG-CD-CE	-5.43	98.80	111.30
1	O	27	ARG	CD-NE-CZ	5.42	131.99	124.40
1	M	27	ARG	CB-CA-C	-5.42	100.27	110.67
1	K	27	ARG	NE-CZ-NH2	-5.37	114.37	119.20
1	M	193	ASP	CA-C-N	5.35	131.39	122.73
1	M	193	ASP	C-N-CA	5.35	131.39	122.73
1	D	147	ARG	CA-CB-CG	5.32	124.75	114.10
1	K	193	ASP	CA-C-N	5.32	131.35	122.73
1	K	193	ASP	C-N-CA	5.32	131.35	122.73
1	M	173	ARG	NE-CZ-NH1	-5.26	116.24	121.50
1	C	162	LYS	CD-CE-NZ	-5.22	95.20	111.90
1	T	184	GLU	CB-CG-CD	-5.18	103.80	112.60
1	G	194	GLN	CB-CA-C	5.17	118.04	109.72
1	I	27	ARG	CB-CA-C	-5.17	102.07	110.85
1	H	184	GLU	CB-CG-CD	5.14	121.34	112.60
1	Z	43	GLU	CA-CB-CG	-5.14	103.83	114.10
1	O	188	GLU	CA-CB-CG	5.12	124.35	114.10
1	C	162	LYS	CA-CB-CG	-5.11	103.88	114.10
1	U	27	ARG	CG-CD-NE	-5.08	100.81	112.00
1	Y	42	ASP	CA-C-N	5.02	129.09	120.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Y	42	ASP	C-N-CA	5.02	129.09	120.71
1	a	30	LYS	CB-CG-CD	5.02	122.85	111.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	S	43	GLU	Sidechain
1	W	193	ASP	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1359	0	1365	16	0
1	B	1347	0	1350	23	0
1	C	1346	0	1355	16	0
1	D	1351	0	1353	20	0
1	E	1359	0	1364	14	0
1	F	1339	0	1344	25	0
1	G	1351	0	1358	23	0
1	H	1359	0	1364	30	0
1	I	1339	0	1344	24	0
1	J	1347	0	1350	21	0
1	K	1359	0	1364	40	0
1	L	1355	0	1361	36	0
1	M	1339	0	1344	20	0
1	N	1347	0	1350	16	0
1	O	1347	0	1350	30	0
1	P	1347	0	1350	26	0
1	Q	1351	0	1358	25	0
1	R	1363	0	1367	26	0
1	S	1351	0	1353	31	0
1	T	1339	0	1346	22	0
1	U	1350	0	1359	25	0
1	V	1341	0	1345	24	0
1	W	1347	0	1350	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	X	1355	0	1361	25	0
1	Y	1347	0	1350	22	0
1	Z	1343	0	1347	22	0
1	a	1351	0	1353	24	0
1	b	1345	0	1353	22	0
2	0	48	0	42	0	0
2	1	48	0	41	1	0
2	2	48	0	41	3	0
2	3	48	0	41	3	0
2	4	48	0	41	3	0
2	c	48	0	42	1	0
2	e	48	0	42	1	0
2	f	48	0	41	4	0
2	g	48	0	42	1	0
2	h	48	0	42	1	0
2	i	48	0	41	1	0
2	j	48	0	41	0	0
2	k	48	0	42	2	0
2	l	48	0	42	3	0
2	m	48	0	41	0	0
2	n	48	0	41	2	0
2	o	48	0	41	2	0
2	p	48	0	41	2	0
2	q	48	0	42	1	0
2	r	48	0	41	2	0
2	s	48	0	42	4	0
2	t	48	0	41	1	0
2	u	48	0	42	1	0
2	v	48	0	41	1	0
2	w	48	0	41	2	0
2	x	48	0	41	0	0
2	y	48	0	42	2	0
2	z	48	0	41	0	0
3	0	9	0	15	0	0
3	1	9	0	15	0	0
3	2	9	0	15	1	0
3	3	9	0	15	3	0
3	4	9	0	15	2	0
3	c	9	0	15	0	0
3	e	9	0	15	2	0
3	f	9	0	15	1	0
3	g	9	0	15	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	h	9	0	15	2	0
3	i	9	0	15	0	0
3	j	9	0	15	1	0
3	k	9	0	15	2	0
3	l	9	0	15	1	0
3	m	9	0	15	0	0
3	n	9	0	15	0	0
3	o	9	0	15	0	0
3	p	9	0	15	2	0
3	q	9	0	15	1	0
3	r	9	0	15	1	0
3	s	9	0	15	1	0
3	t	9	0	15	2	0
3	u	9	0	15	0	0
3	v	9	0	15	0	0
3	w	9	0	15	0	0
3	x	9	0	15	1	0
3	y	9	0	15	1	0
3	z	9	0	15	0	0
4	O	1	0	0	0	0
4	A	25	0	0	0	0
4	B	27	0	0	3	0
4	C	21	0	0	0	0
4	D	17	0	0	0	0
4	E	24	0	0	2	0
4	F	22	0	0	3	0
4	G	26	0	0	1	0
4	H	19	0	0	0	0
4	I	20	0	0	1	0
4	J	21	0	0	0	0
4	K	32	0	0	2	0
4	L	26	0	0	3	0
4	M	31	0	0	2	0
4	N	35	0	0	2	0
4	O	22	0	0	0	0
4	P	30	0	0	3	0
4	Q	22	0	0	1	0
4	R	27	0	0	3	0
4	S	17	0	0	0	0
4	T	24	0	0	3	0
4	U	25	0	0	1	0
4	V	29	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	W	24	0	0	0	0
4	X	24	0	0	1	0
4	Y	17	0	0	1	0
4	Z	32	0	0	3	0
4	a	29	0	0	3	0
4	b	37	0	0	3	0
All	All	40076	0	39487	623	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 8.

All (623) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:W:54:PHE:CE1	1:X:27:ARG:NH1	2.08	1.20
1:L:41:THR:OG1	1:L:43:GLU:CD	1.95	1.09
1:I:150:LEU:HD11	1:S:143:GLU:HG3	1.41	1.03
1:L:41:THR:OG1	1:L:43:GLU:OE1	1.78	1.00
1:S:113:LYS:NZ	1:S:163:HIS:O	1.99	0.94
1:L:41:THR:HG1	1:L:43:GLU:CD	1.76	0.89
1:W:54:PHE:HE1	1:X:27:ARG:HH11	1.03	0.88
1:X:137:GLY:O	4:X:301:HOH:O	1.93	0.85
1:W:54:PHE:HE1	1:X:27:ARG:NH1	1.60	0.84
1:K:117:PHE:HE1	1:K:194:GLN:HG3	1.42	0.84
1:O:23:ASP:HB3	1:O:26:SER:HB2	1.56	0.84
1:L:41:THR:CB	1:L:43:GLU:OE1	2.26	0.83
1:X:155:LYS:HE2	1:X:159:LEU:HG	1.61	0.82
1:D:27:ARG:NH1	1:D:30:LYS:HG2	1.94	0.82
1:O:81:ILE:HG22	1:O:85:MET:HE2	1.63	0.81
1:a:27:ARG:HA	1:a:30:LYS:HG3	1.64	0.80
1:b:181:MET:SD	4:b:330:HOH:O	2.39	0.79
1:F:181:MET:HE1	1:F:191:LEU:HD12	1.63	0.79
1:B:160:MET:HE3	1:B:191:LEU:HD11	1.65	0.79
1:Q:138:GLN:OE1	1:R:177:ARG:NH2	2.16	0.79
1:G:101:ALA:HB1	1:G:125:MET:HE3	1.65	0.79
1:B:194:GLN:NE2	4:B:301:HOH:O	2.16	0.78
1:B:28:LEU:HD13	1:B:35:PHE:HE2	1.48	0.78
1:F:33:ILE:HD11	2:i:7:MP8:HDA	1.66	0.77
1:b:181:MET:HE1	1:b:191:LEU:HD12	1.67	0.77
1:J:143:GLU:HG2	1:J:147:ARG:HH22	1.51	0.76
1:R:117:PHE:HD2	1:R:194:GLN:HB3	1.49	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:146:ALA:O	1:V:150:LEU:HD22	1.85	0.76
1:J:82:TYR:HA	1:J:85:MET:HE2	1.68	0.76
1:U:131:ILE:O	4:U:301:HOH:O	2.04	0.76
1:E:78:GLY:HA3	1:E:103:MET:HE2	1.67	0.75
1:U:117:PHE:HB3	1:U:196:LEU:HD11	1.69	0.74
1:Y:102:SER:OG	1:Y:127:HIS:NE2	2.20	0.74
1:X:155:LYS:HZ3	1:X:159:LEU:HD21	1.53	0.74
1:G:30:LYS:NZ	1:K:62:LYS:NZ	2.35	0.74
1:F:130:LEU:O	1:F:131:ILE:HD13	1.87	0.74
1:K:117:PHE:CE1	1:K:194:GLN:HG3	2.20	0.74
1:b:102:SER:OG	1:b:127:HIS:NE2	2.21	0.74
1:D:27:ARG:HD2	1:D:27:ARG:O	1.86	0.74
1:N:45:ALA:HA	1:N:81:ILE:HD11	1.70	0.74
1:J:143:GLU:HG2	1:J:147:ARG:NH2	2.03	0.73
1:S:27:ARG:HG2	1:S:30:LYS:HE2	1.70	0.73
1:F:106:PHE:HD1	1:F:160:MET:HE2	1.52	0.73
1:S:106:PHE:HD1	1:S:160:MET:HE2	1.52	0.73
1:J:59:ASN:ND2	1:J:62:LYS:HG3	2.04	0.72
1:T:131:ILE:HD12	1:T:149:LEU:HD22	1.69	0.72
1:O:82:TYR:HA	1:O:85:MET:HE3	1.72	0.72
1:T:82:TYR:HA	1:T:85:MET:HE2	1.70	0.72
1:Z:173:ARG:NH2	4:Z:301:HOH:O	2.22	0.72
1:N:102:SER:OG	1:N:127:HIS:NE2	2.23	0.72
1:Y:61:ASP:OD1	1:Y:89:LYS:NZ	2.16	0.72
1:S:28:LEU:HD13	1:S:35:PHE:HE2	1.54	0.72
1:T:33:ILE:HD11	2:w:7:MP8:HDA	1.71	0.71
1:K:181:MET:HE1	1:K:191:LEU:HD13	1.72	0.71
1:R:28:LEU:HD13	1:R:35:PHE:HE2	1.54	0.71
1:J:59:ASN:HD22	1:J:62:LYS:HG3	1.56	0.71
1:H:28:LEU:HD22	3:k:101:OCA:H82	1.74	0.70
1:P:197:GLU:OE1	4:P:301:HOH:O	2.07	0.70
1:V:187:LYS:NZ	4:V:302:HOH:O	2.23	0.70
1:b:184:GLU:OE2	4:b:301:HOH:O	2.09	0.70
1:E:132:SER:OG	4:E:302:HOH:O	2.10	0.70
1:L:28:LEU:HD13	1:L:35:PHE:HE2	1.57	0.70
1:S:102:SER:OG	1:S:127:HIS:NE2	2.24	0.70
1:V:146:ALA:O	1:V:150:LEU:CD2	2.39	0.70
1:b:28:LEU:HD13	1:b:35:PHE:HE2	1.57	0.70
1:H:184:GLU:HG2	1:H:185:GLU:H	1.56	0.70
1:F:158:ARG:HH12	1:b:167:ASP:CG	2.00	0.70
1:N:28:LEU:HD13	1:N:35:PHE:HE2	1.55	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:33:ILE:HD13	1:T:65:PHE:HB2	1.74	0.69
1:X:155:LYS:HZ3	1:X:159:LEU:HD11	1.55	0.69
1:a:187:LYS:HD2	1:a:193:ASP:O	1.92	0.69
1:L:45:ALA:HA	1:L:81:ILE:HD11	1.74	0.69
1:W:138:GLN:OE1	1:X:177:ARG:NH2	2.25	0.69
1:X:28:LEU:HD13	1:X:35:PHE:HE2	1.55	0.69
1:M:45:ALA:HA	1:M:81:ILE:HD11	1.74	0.69
1:B:79:MET:HE2	1:B:155:LYS:HD3	1.75	0.69
1:S:61:ASP:OD1	1:S:89:LYS:NZ	2.21	0.69
1:U:45:ALA:HA	1:U:81:ILE:HD11	1.74	0.69
1:M:25:TYR:OH	4:M:301:HOH:O	2.11	0.68
1:G:30:LYS:HZ3	1:K:62:LYS:HZ1	1.42	0.68
1:U:117:PHE:CD1	1:U:194:GLN:HB2	2.29	0.68
1:b:181:MET:HE1	1:b:191:LEU:CD1	2.24	0.68
1:I:187:LYS:NZ	4:I:301:HOH:O	2.25	0.68
1:R:62:LYS:NZ	4:R:303:HOH:O	2.27	0.67
1:Z:143:GLU:O	1:Z:147:ARG:HG2	1.93	0.67
1:I:33:ILE:HD11	2:I:7:MP8:HDA	1.76	0.67
1:K:198:ASN:N	4:K:301:HOH:O	2.27	0.67
1:L:41:THR:HB	1:L:43:GLU:OE1	1.94	0.67
1:Q:132:SER:O	4:Q:301:HOH:O	2.12	0.67
1:T:24:ILE:N	4:T:301:HOH:O	2.27	0.67
1:X:155:LYS:NZ	1:X:159:LEU:HD11	2.09	0.67
1:I:28:LEU:HD13	1:I:35:PHE:HE2	1.60	0.66
1:O:25:TYR:OH	1:U:46:ASN:OD1	2.09	0.66
1:R:154:GLU:OE1	1:R:158:ARG:NH1	2.29	0.66
1:B:79:MET:HE3	1:C:121:ASN:HB3	1.78	0.66
1:a:27:ARG:HG3	3:3:101:OCA:H83	1.78	0.65
1:N:188:GLU:OE2	4:N:301:HOH:O	2.13	0.65
1:G:30:LYS:HZ3	1:K:62:LYS:NZ	1.94	0.65
1:Z:136:GLY:N	4:Z:303:HOH:O	2.29	0.65
1:F:197:GLU:OE2	4:F:301:HOH:O	2.14	0.65
1:Z:28:LEU:HD13	1:Z:35:PHE:HE2	1.60	0.65
1:S:197:GLU:CD	1:S:198:ASN:H	2.05	0.64
1:E:135:LEU:HD12	1:E:136:GLY:H	1.61	0.64
1:Y:55:LEU:HD13	1:Y:64:ILE:HD12	1.79	0.64
1:E:45:ALA:HA	1:E:81:ILE:HD11	1.78	0.64
1:a:28:LEU:HD13	1:a:35:PHE:HE2	1.61	0.64
1:a:45:ALA:HA	1:a:81:ILE:HD11	1.79	0.64
1:a:102:SER:OG	1:a:127:HIS:NE2	2.30	0.64
1:M:102:SER:OG	1:M:127:HIS:NE2	2.31	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:79:MET:HE2	1:O:155:LYS:HE3	1.80	0.64
1:D:27:ARG:NH1	1:D:31:GLU:OE2	2.32	0.63
1:N:131:ILE:HG22	1:N:133:GLY:HA3	1.79	0.63
1:G:79:MET:HE1	1:G:106:PHE:CE2	2.34	0.63
1:V:138:GLN:OE1	1:W:177:ARG:NH2	2.32	0.63
1:H:95:LEU:HD11	1:H:119:LEU:HD13	1.80	0.63
1:D:101:ALA:HB1	1:D:125:MET:HE3	1.81	0.63
1:V:125:MET:HE1	1:V:127:HIS:CG	2.34	0.63
1:K:117:PHE:CD1	1:K:194:GLN:HB2	2.34	0.63
1:R:135:LEU:N	4:R:304:HOH:O	2.31	0.63
1:b:173:ARG:HD2	1:b:174:ASP:OD2	1.99	0.62
1:V:117:PHE:HD1	1:V:194:GLN:HB3	1.64	0.62
1:H:184:GLU:HG2	1:H:185:GLU:N	2.14	0.62
1:Y:28:LEU:HD13	1:Y:35:PHE:HE2	1.63	0.62
1:D:27:ARG:HD2	1:D:27:ARG:C	2.25	0.62
1:W:27:ARG:HG2	1:W:30:LYS:HE2	1.82	0.62
1:T:24:ILE:N	4:T:302:HOH:O	2.34	0.61
1:Z:102:SER:HB2	1:Z:127:HIS:CE1	2.35	0.61
1:D:27:ARG:HH11	1:D:30:LYS:HG2	1.64	0.61
1:J:36:LEU:HD21	1:J:40:VAL:HG22	1.82	0.61
1:R:117:PHE:CD2	1:R:194:GLN:HB3	2.34	0.61
1:H:79:MET:HE3	1:H:155:LYS:HE3	1.82	0.61
1:T:124:ILE:HG21	1:T:191:LEU:HD22	1.83	0.61
1:O:75:VAL:HG22	1:O:103:MET:HE3	1.82	0.61
1:P:126:ILE:HG21	1:P:160:MET:HE1	1.83	0.61
1:b:96:CYS:HB2	1:b:108:LEU:HD22	1.82	0.61
1:Y:184:GLU:OE1	1:Y:184:GLU:HA	2.01	0.60
1:a:78:GLY:HA3	1:a:103:MET:HE2	1.83	0.60
1:D:160:MET:HE2	1:D:171:LEU:HD11	1.83	0.60
1:A:126:ILE:HG21	1:A:160:MET:HE1	1.84	0.60
1:M:117:PHE:CD1	1:M:194:GLN:HB2	2.36	0.60
1:A:184:GLU:HA	1:A:195:ILE:HD11	1.84	0.60
1:B:27:ARG:O	1:B:30:LYS:HB3	2.00	0.60
1:K:189:TYR:HB3	1:K:191:LEU:HD12	1.82	0.60
1:P:28:LEU:HD13	1:P:35:PHE:HE2	1.65	0.60
1:I:101:ALA:HB1	1:I:125:MET:HE3	1.84	0.60
1:I:150:LEU:HD11	1:S:143:GLU:CG	2.26	0.60
1:O:45:ALA:HA	1:O:81:ILE:HD11	1.84	0.59
1:T:28:LEU:HD13	1:T:35:PHE:HE2	1.68	0.59
1:Z:194:GLN:NE2	4:Z:305:HOH:O	2.32	0.59
1:Y:130:LEU:HD23	1:Y:131:ILE:H	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:25:TYR:OH	4:R:301:HOH:O	2.16	0.59
1:U:53:LEU:HD22	3:r:101:OCA:H32	1.85	0.59
1:X:36:LEU:HD21	1:X:40:VAL:HG22	1.84	0.59
1:G:28:LEU:HD13	1:G:35:PHE:HE2	1.66	0.59
1:S:28:LEU:HD13	1:S:35:PHE:CE2	2.37	0.59
1:W:79:MET:HE1	1:W:106:PHE:CE2	2.38	0.59
1:Q:45:ALA:HA	1:Q:81:ILE:HD11	1.84	0.59
1:M:43:GLU:N	1:M:43:GLU:OE1	2.35	0.59
1:P:78:GLY:HA3	1:P:103:MET:HE2	1.84	0.59
1:Z:189:TYR:HB3	1:Z:191:LEU:HD12	1.86	0.58
1:K:135:LEU:O	1:Q:135:LEU:N	2.36	0.58
1:V:28:LEU:HD13	1:V:35:PHE:HE2	1.68	0.58
1:W:63:ASP:OD2	1:W:115:LYS:NZ	2.28	0.58
1:B:63:ASP:OD2	1:B:115:LYS:NZ	2.28	0.58
1:C:86:ASN:HB3	1:D:198:ASN:ND2	2.19	0.58
1:M:79:MET:HE1	1:M:106:PHE:CE2	2.39	0.58
1:O:130:LEU:O	1:O:131:ILE:HD12	2.03	0.58
1:U:117:PHE:HD1	1:U:194:GLN:HB2	1.66	0.58
1:I:76:THR:HG23	1:J:98:GLY:C	2.29	0.58
1:U:102:SER:OG	1:U:127:HIS:NE2	2.36	0.57
1:H:45:ALA:HA	1:H:81:ILE:HD11	1.84	0.57
1:P:102:SER:OG	1:P:127:HIS:NE2	2.37	0.57
1:R:189:TYR:HB3	1:R:191:LEU:HD12	1.86	0.57
1:L:150:LEU:HD11	1:P:143:GLU:HB2	1.87	0.57
1:L:79:MET:HE1	1:L:106:PHE:CE2	2.39	0.57
1:S:41:THR:OG1	1:S:43:GLU:OE1	2.20	0.57
1:X:46:ASN:HB3	4:Y:301:HOH:O	2.04	0.57
1:K:75:VAL:HG22	1:K:103:MET:HE3	1.86	0.57
1:P:45:ALA:HA	1:P:81:ILE:HD11	1.86	0.57
1:a:170:ASP:OD2	4:a:301:HOH:O	2.18	0.57
1:S:45:ALA:HA	1:S:81:ILE:HD11	1.87	0.56
1:Z:81:ILE:O	1:Z:85:MET:HG3	2.05	0.56
1:K:147:ARG:HH12	1:Q:147:ARG:NH1	2.02	0.56
1:O:30:LYS:O	1:O:30:LYS:HG2	2.05	0.56
1:V:130:LEU:O	1:V:131:ILE:HD12	2.05	0.56
1:W:28:LEU:HD13	1:W:35:PHE:HE2	1.71	0.56
1:T:61:ASP:OD1	1:T:89:LYS:NZ	2.24	0.56
1:V:125:MET:HE1	1:V:127:HIS:ND1	2.20	0.56
1:A:61:ASP:OD1	1:A:89:LYS:NZ	2.39	0.56
1:L:154:GLU:O	1:L:158:ARG:HG3	2.05	0.56
1:T:81:ILE:O	1:T:85:MET:HG3	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Y:124:ILE:HG21	1:Y:191:LEU:HD22	1.87	0.56
1:A:36:LEU:HD21	1:A:40:VAL:HG22	1.87	0.55
1:L:53:LEU:HD22	3:p:101:OCA:H32	1.89	0.55
1:P:62:LYS:NZ	4:P:303:HOH:O	2.38	0.55
1:Z:79:MET:HE1	1:Z:106:PHE:CE2	2.42	0.55
1:G:188:GLU:O	1:G:188:GLU:HG2	2.06	0.55
1:a:27:ARG:O	1:a:30:LYS:HB2	2.06	0.55
1:K:143:GLU:O	1:K:147:ARG:HG2	2.06	0.55
1:U:28:LEU:HD13	1:U:35:PHE:HE2	1.72	0.55
1:U:60:PRO:HB2	1:U:89:LYS:HD3	1.87	0.55
1:H:197:GLU:OE2	1:H:198:ASN:N	2.40	0.55
1:Q:106:PHE:HA	1:Q:160:MET:HE3	1.90	0.55
1:O:33:ILE:HD11	2:r:7:MP8:HDA	1.87	0.54
1:C:81:ILE:O	1:C:85:MET:HG3	2.08	0.54
1:H:182:SER:OG	1:H:184:GLU:HG2	2.07	0.54
1:D:109:SER:HB3	1:D:191:LEU:HD12	1.90	0.54
1:b:36:LEU:HD21	1:b:40:VAL:HG22	1.89	0.54
1:H:65:PHE:CE2	2:k:7:MP8:HB	2.42	0.54
1:L:23:ASP:O	1:L:26:SER:OG	2.21	0.54
1:Q:135:LEU:CD1	1:Q:142:ILE:HG13	2.37	0.54
1:F:158:ARG:NH1	1:b:167:ASP:OD1	2.41	0.54
1:W:81:ILE:O	1:W:85:MET:HG3	2.07	0.54
1:J:131:ILE:HD12	1:R:136:GLY:HA2	1.89	0.54
1:L:184:GLU:OE2	4:L:301:HOH:O	2.19	0.54
1:P:30:LYS:HE2	1:P:30:LYS:HA	1.90	0.54
1:W:45:ALA:HA	1:W:81:ILE:HD11	1.89	0.54
1:Z:33:ILE:HD11	2:2:7:MP8:HDA	1.88	0.54
1:L:81:ILE:O	1:L:85:MET:HG3	2.08	0.54
1:O:85:MET:CE	1:O:107:LEU:HD22	2.38	0.54
1:B:173:ARG:HD2	1:B:174:ASP:OD1	2.08	0.54
1:J:78:GLY:HA3	1:J:103:MET:HE2	1.90	0.54
1:a:67:TYR:OH	2:3:6:ALA:O	2.20	0.54
1:S:113:LYS:HZ1	1:S:163:HIS:HA	1.71	0.53
1:F:79:MET:HE1	1:F:106:PHE:CE2	2.43	0.53
1:G:106:PHE:HA	1:G:160:MET:HE2	1.90	0.53
1:O:117:PHE:CD1	1:O:194:GLN:HB3	2.42	0.53
1:X:155:LYS:HZ3	1:X:159:LEU:CD2	2.21	0.53
1:D:32:ARG:NH1	1:D:55:LEU:O	2.40	0.53
1:Q:79:MET:HE3	1:R:121:ASN:HB3	1.90	0.53
1:W:27:ARG:O	1:W:30:LYS:HG2	2.08	0.53
1:D:61:ASP:OD2	1:D:89:LYS:NZ	2.27	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Q:27:ARG:HD3	3:t:101:OCA:H83	1.91	0.53
1:U:28:LEU:HG	3:x:101:OCA:H82	1.89	0.53
1:A:78:GLY:HA3	1:A:103:MET:HE2	1.90	0.53
1:G:30:LYS:NZ	1:K:62:LYS:HZ1	1.99	0.53
1:L:173:ARG:HD2	1:L:174:ASP:OD1	2.08	0.53
1:a:45:ALA:HB2	1:a:77:ALA:HB1	1.90	0.53
1:b:81:ILE:O	1:b:85:MET:HG3	2.08	0.53
1:F:33:ILE:HD13	1:F:65:PHE:HB2	1.89	0.53
1:U:96:CYS:HB2	1:U:108:LEU:HD22	1.90	0.53
1:X:155:LYS:HZ3	1:X:159:LEU:CD1	2.22	0.53
1:H:135:LEU:HD22	1:H:142:ILE:HD13	1.91	0.53
1:Y:181:MET:CE	1:Y:186:ALA:HA	2.39	0.52
1:T:28:LEU:HD13	1:T:35:PHE:CE2	2.44	0.52
2:3:7:MP8:HD	3:3:101:OCA:H22	1.91	0.52
1:Z:46:ASN:HB3	4:a:303:HOH:O	2.09	0.52
1:C:33:ILE:HD11	2:f:7:MP8:HDA	1.92	0.52
1:O:117:PHE:HD1	1:O:194:GLN:HB3	1.74	0.52
1:U:170:ASP:HA	1:U:173:ARG:NH1	2.24	0.52
1:K:43:GLU:HG3	1:K:44:SER:N	2.23	0.52
1:K:157:ASN:HD22	1:K:172:GLU:HG3	1.75	0.52
1:O:106:PHE:HA	1:O:160:MET:HE2	1.92	0.52
1:F:53:LEU:HB3	3:j:101:OCA:H32	1.91	0.52
1:K:106:PHE:HA	1:K:160:MET:HE3	1.92	0.52
1:Y:184:GLU:HA	1:Y:195:ILE:HD11	1.92	0.52
1:B:197:GLU:O	4:B:302:HOH:O	2.19	0.52
1:H:135:LEU:HD23	1:H:136:GLY:N	2.25	0.52
1:V:81:ILE:O	1:V:85:MET:HG3	2.10	0.52
1:X:155:LYS:NZ	1:X:159:LEU:HD21	2.23	0.51
1:a:81:ILE:O	1:a:85:MET:HG3	2.11	0.51
1:B:159:LEU:O	1:B:162:LYS:HB2	2.11	0.51
1:Y:55:LEU:HB3	1:Y:64:ILE:HD11	1.92	0.51
1:Y:130:LEU:HD23	1:Y:131:ILE:N	2.25	0.51
1:K:79:MET:HE3	1:L:121:ASN:HB3	1.92	0.51
1:P:100:ALA:O	1:P:105:ALA:HB2	2.10	0.51
1:K:189:TYR:HB3	1:K:191:LEU:CD1	2.41	0.51
1:N:36:LEU:HD21	1:N:40:VAL:HG22	1.92	0.51
1:Q:101:ALA:HB1	1:Q:125:MET:HE3	1.92	0.51
1:U:59:ASN:HD21	1:U:62:LYS:HE2	1.75	0.51
1:V:166:ARG:HD2	1:V:189:TYR:O	2.10	0.51
1:I:23:ASP:OD1	1:I:23:ASP:N	2.43	0.51
1:J:131:ILE:HD11	1:R:142:ILE:HD11	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:117:PHE:HD1	1:M:194:GLN:HB2	1.75	0.51
1:O:121:ASN:HB3	1:U:79:MET:HE3	1.93	0.51
1:S:79:MET:HE1	1:S:106:PHE:CE2	2.46	0.51
1:A:81:ILE:O	1:A:85:MET:HG3	2.11	0.50
1:G:81:ILE:O	1:G:85:MET:HG3	2.12	0.50
1:Q:166:ARG:HD2	1:Q:189:TYR:O	2.10	0.50
1:X:124:ILE:HG21	1:X:191:LEU:HD13	1.92	0.50
1:a:117:PHE:HD1	1:a:194:GLN:HB2	1.76	0.50
1:E:79:MET:HE1	1:E:106:PHE:CE2	2.47	0.50
1:T:65:PHE:CE2	2:w:7:MP8:HB	2.46	0.50
1:C:154:GLU:O	1:C:158:ARG:HG3	2.12	0.50
1:B:109:SER:HB3	1:B:191:LEU:HD12	1.93	0.50
1:Z:45:ALA:HA	1:Z:81:ILE:HD11	1.94	0.50
1:Z:102:SER:HB2	1:Z:127:HIS:HE1	1.75	0.50
1:I:45:ALA:HA	1:I:81:ILE:HD11	1.94	0.49
1:C:45:ALA:HA	1:C:81:ILE:HD11	1.95	0.49
1:I:166:ARG:HD2	1:I:189:TYR:O	2.12	0.49
1:F:36:LEU:HD21	1:F:40:VAL:HG22	1.94	0.49
1:I:78:GLY:HA3	1:I:103:MET:HE2	1.94	0.49
1:S:81:ILE:O	1:S:85:MET:HG3	2.13	0.49
1:S:184:GLU:N	1:S:184:GLU:OE1	2.45	0.49
1:T:131:ILE:HD12	1:T:149:LEU:CD2	2.39	0.49
1:b:154:GLU:O	1:b:158:ARG:HG3	2.12	0.49
1:E:160:MET:HE1	1:E:191:LEU:HD21	1.94	0.49
1:X:59:ASN:ND2	1:X:62:LYS:HE3	2.28	0.49
1:I:65:PHE:CE2	2:l:7:MP8:HB	2.48	0.49
1:b:184:GLU:O	1:b:188:GLU:HB2	2.13	0.49
1:J:45:ALA:HA	1:J:81:ILE:HD11	1.94	0.49
1:W:113:LYS:HE2	1:W:163:HIS:O	2.13	0.49
1:X:41:THR:H	1:X:44:SER:HG	1.61	0.49
1:Y:65:PHE:CE2	2:1:7:MP8:HB	2.48	0.49
1:H:195:ILE:HG22	1:H:195:ILE:O	2.12	0.49
1:T:76:THR:HG21	1:U:99:GLN:HG2	1.95	0.49
1:I:67:TYR:CE2	2:l:2:WFP:HD2	2.48	0.48
1:L:176:ASP:CG	1:Q:177:ARG:HH12	2.21	0.48
1:K:181:MET:HE1	1:K:191:LEU:CD1	2.42	0.48
1:P:81:ILE:O	1:P:85:MET:HG3	2.12	0.48
1:L:65:PHE:CE2	2:o:7:MP8:HB	2.49	0.48
1:b:28:LEU:HD13	1:b:35:PHE:CE2	2.43	0.48
1:B:164:CYS:SG	1:B:191:LEU:HD13	2.53	0.48
1:N:79:MET:HE2	1:N:155:LYS:HE2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:131:ILE:HG22	1:P:132:SER:H	1.78	0.48
1:S:113:LYS:NZ	1:S:163:HIS:HA	2.28	0.48
1:X:154:GLU:O	1:X:158:ARG:HG3	2.12	0.48
1:A:23:ASP:OD1	1:A:24:ILE:N	2.47	0.48
1:S:157:ASN:HB3	1:S:168:LEU:HD12	1.96	0.48
1:Z:120:PRO:HG3	1:Z:197:GLU:HG2	1.95	0.48
1:B:138:GLN:OE1	1:C:177:ARG:NH2	2.45	0.48
1:G:45:ALA:HA	1:G:81:ILE:HD11	1.95	0.48
1:N:130:LEU:HA	1:U:137:GLY:O	2.13	0.48
1:R:65:PHE:CE2	2:u:7:MP8:HB	2.49	0.48
1:G:56:GLU:HB3	4:G:302:HOH:O	2.13	0.48
1:G:78:GLY:HA3	1:G:103:MET:HE2	1.95	0.48
1:H:143:GLU:OE2	1:H:147:ARG:NE	2.45	0.48
1:P:117:PHE:CZ	2:s:5:YCP:HD	2.49	0.48
1:R:94:THR:HB	1:R:108:LEU:HD12	1.95	0.48
1:T:36:LEU:HD21	1:T:40:VAL:HG22	1.95	0.48
1:A:45:ALA:HA	1:A:81:ILE:HD11	1.96	0.48
1:G:28:LEU:HD13	1:G:35:PHE:CE2	2.48	0.48
1:H:113:LYS:HD3	1:H:116:ARG:CZ	2.43	0.48
1:O:166:ARG:HD2	1:O:189:TYR:O	2.14	0.48
1:L:113:LYS:HD2	1:L:113:LYS:N	2.29	0.48
1:L:123:ARG:HD3	1:L:180:PHE:HD2	1.78	0.48
2:4:3:ALO:N	3:4:101:OCA:O1	2.46	0.48
1:D:65:PHE:CE2	2:g:7:MP8:HB	2.49	0.47
1:J:63:ASP:OD1	1:J:91:ASP:HB2	2.14	0.47
1:T:33:ILE:HD12	1:T:67:TYR:CZ	2.50	0.47
1:H:75:VAL:HG22	1:H:103:MET:HE3	1.96	0.47
1:K:41:THR:OG1	1:K:43:GLU:HG2	2.14	0.47
1:P:79:MET:HE1	1:P:106:PHE:CE1	2.49	0.47
1:R:54:PHE:HE1	1:S:27:ARG:NH1	2.12	0.47
1:E:43:GLU:O	1:E:43:GLU:HG2	2.02	0.47
1:V:106:PHE:HA	1:V:160:MET:HE2	1.95	0.47
1:X:173:ARG:HD2	1:X:174:ASP:OD1	2.14	0.47
1:B:81:ILE:O	1:B:85:MET:HG3	2.15	0.47
1:F:88:ILE:HD12	1:F:90:PRO:HG2	1.95	0.47
1:F:165:ASP:OD2	4:F:302:HOH:O	2.21	0.47
1:H:184:GLU:HA	1:H:195:ILE:HD11	1.95	0.47
1:J:28:LEU:HD13	1:J:35:PHE:CE2	2.49	0.47
1:K:117:PHE:HD1	1:K:194:GLN:HB2	1.76	0.47
1:Q:164:CYS:SG	1:Q:191:LEU:HD13	2.54	0.47
1:Z:65:PHE:CE2	2:2:7:MP8:HB	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:65:PHE:CE2	2:f:7:MP8:HB	2.49	0.47
1:C:106:PHE:HA	1:C:160:MET:HE2	1.97	0.47
1:F:95:LEU:HA	1:F:117:PHE:O	2.15	0.47
1:G:60:PRO:HB2	1:G:89:LYS:HD3	1.97	0.47
1:K:76:THR:HG21	1:L:99:GLN:HG2	1.95	0.47
1:K:198:ASN:ND2	4:K:304:HOH:O	2.43	0.47
1:R:28:LEU:HD13	1:R:35:PHE:CE2	2.42	0.47
1:R:47:LEU:HD12	1:R:47:LEU:HA	1.63	0.47
1:a:27:ARG:HG3	3:3:101:OCA:C8	2.42	0.47
2:y:3:ALO:H	3:y:101:OCA:C1	2.27	0.47
1:K:113:LYS:HE3	1:K:163:HIS:HA	1.96	0.47
1:K:157:ASN:ND2	1:K:172:GLU:HG3	2.30	0.47
1:M:33:ILE:HD11	2:p:7:MP8:HDA	1.96	0.47
1:D:81:ILE:O	1:D:85:MET:HG3	2.13	0.47
1:K:181:MET:HE3	1:K:186:ALA:HA	1.97	0.47
1:L:61:ASP:OD1	1:L:89:LYS:NZ	2.43	0.47
1:K:65:PHE:CE2	2:n:7:MP8:HB	2.49	0.46
1:T:45:ALA:HA	1:T:81:ILE:HD11	1.96	0.46
1:T:131:ILE:CD1	1:T:149:LEU:HD22	2.41	0.46
1:E:59:ASN:ND2	1:E:62:LYS:HE2	2.30	0.46
1:P:65:PHE:CE1	2:s:6:ALA:HA	2.50	0.46
1:P:197:GLU:H	1:P:197:GLU:CD	2.23	0.46
1:B:167:ASP:HB3	4:B:317:HOH:O	2.15	0.46
1:H:28:LEU:HD13	3:k:101:OCA:H61	1.97	0.46
1:S:162:LYS:HA	1:S:162:LYS:HD3	1.71	0.46
1:T:197:GLU:HG3	1:T:198:ASN:H	1.79	0.46
1:U:81:ILE:O	1:U:85:MET:HG3	2.16	0.46
1:Y:79:MET:CE	1:Y:155:LYS:HE3	2.45	0.46
1:B:36:LEU:HD21	1:B:40:VAL:HG22	1.97	0.46
1:D:147:ARG:NH2	1:D:151:LYS:NZ	2.63	0.46
1:F:197:GLU:HB3	4:F:301:HOH:O	2.16	0.46
1:O:96:CYS:HB2	1:O:108:LEU:HD22	1.97	0.46
1:U:75:VAL:O	1:U:79:MET:HG2	2.16	0.46
1:X:102:SER:OG	1:X:103:MET:N	2.43	0.46
1:Z:62:LYS:O	1:Z:90:PRO:HB3	2.15	0.46
2:4:3:ALO:HG23	3:4:101:OCA:O1	2.15	0.46
1:J:95:LEU:HD23	1:J:117:PHE:HB2	1.96	0.46
1:O:25:TYR:HA	1:O:28:LEU:HD12	1.97	0.46
1:Q:81:ILE:O	1:Q:85:MET:HG3	2.15	0.46
1:Q:135:LEU:HD11	1:Q:142:ILE:HG13	1.98	0.46
1:I:36:LEU:HD21	1:I:40:VAL:HG22	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:193:ASP:O	4:L:303:HOH:O	2.21	0.46
1:W:124:ILE:HG21	1:W:191:LEU:HD22	1.98	0.46
1:A:197:GLU:HA	1:A:197:GLU:OE1	2.16	0.46
1:M:100:ALA:O	1:M:105:ALA:HB2	2.16	0.46
1:Q:28:LEU:HD22	1:Q:35:PHE:HE2	1.82	0.46
1:a:45:ALA:CB	1:a:77:ALA:HB1	2.46	0.46
1:V:146:ALA:O	1:V:150:LEU:HD23	2.15	0.45
1:D:126:ILE:HG21	1:D:160:MET:HE1	1.97	0.45
1:G:89:LYS:HB3	1:G:90:PRO:HD3	1.98	0.45
1:P:109:SER:HB3	1:P:191:LEU:HG	1.97	0.45
1:A:65:PHE:CE2	2:c:7:MP8:HB	2.51	0.45
1:V:79:MET:HE2	1:V:103:MET:HE1	1.96	0.45
1:Y:114:GLY:N	1:Y:193:ASP:OD1	2.49	0.45
1:F:114:GLY:H	1:F:193:ASP:CG	2.25	0.45
1:I:33:ILE:HD12	1:I:67:TYR:CZ	2.51	0.45
1:L:135:LEU:HB3	1:L:136:GLY:H	1.54	0.45
1:H:95:LEU:HD11	1:H:119:LEU:CD1	2.46	0.45
1:L:131:ILE:HD11	1:P:142:ILE:HD13	1.98	0.45
1:L:95:LEU:HA	1:L:117:PHE:O	2.17	0.45
1:O:65:PHE:CE2	2:r:7:MP8:HB	2.52	0.45
1:P:28:LEU:HD13	1:P:35:PHE:CE2	2.49	0.45
1:F:61:ASP:OD1	1:F:89:LYS:NZ	2.48	0.45
1:G:30:LYS:NZ	1:K:62:LYS:CE	2.80	0.45
1:G:30:LYS:HB3	1:G:30:LYS:HE3	1.27	0.45
1:W:183:ALA:HB1	1:W:195:ILE:HG12	1.98	0.45
1:H:121:ASN:HB3	1:N:79:MET:HE3	1.99	0.45
1:Q:135:LEU:C	1:Q:135:LEU:HD12	2.42	0.45
1:Z:63:ASP:OD2	1:Z:115:LYS:NZ	2.44	0.45
1:B:54:PHE:CG	1:C:24:ILE:HD11	2.51	0.45
1:M:96:CYS:HB2	1:M:108:LEU:HD22	1.98	0.45
1:T:151:LYS:NZ	4:T:303:HOH:O	2.50	0.45
1:X:102:SER:HB2	1:X:127:HIS:NE2	2.32	0.45
2:h:3:ALO:H	3:h:101:OCA:C1	2.28	0.45
1:B:159:LEU:HD23	1:B:159:LEU:HA	1.79	0.45
1:S:113:LYS:HZ3	1:S:163:HIS:C	2.11	0.45
1:F:47:LEU:O	1:F:51:GLN:HG3	2.17	0.44
1:N:33:ILE:HD11	2:q:7:MP8:HDA	1.98	0.44
1:O:95:LEU:HA	1:O:117:PHE:O	2.16	0.44
1:V:181:MET:HE1	1:V:191:LEU:HD13	1.99	0.44
1:K:79:MET:HE3	1:L:121:ASN:CB	2.47	0.44
1:O:33:ILE:HD13	1:O:65:PHE:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:THR:HA	2:f:2:WFP:F1	2.08	0.44
1:J:28:LEU:HD13	1:J:35:PHE:HE2	1.82	0.44
1:J:54:PHE:HE1	1:K:27:ARG:HH11	1.66	0.44
1:J:88:ILE:HB	1:J:90:PRO:HD2	1.98	0.44
1:T:114:GLY:N	1:T:193:ASP:OD2	2.47	0.44
2:2:3:ALO:N	3:2:101:OCA:O1	2.51	0.44
1:L:160:MET:HE3	1:L:160:MET:HB2	1.84	0.44
1:I:28:LEU:HD13	1:I:35:PHE:CE2	2.47	0.44
1:P:131:ILE:HG22	1:P:132:SER:N	2.32	0.44
1:Q:41:THR:H	1:Q:44:SER:HG	1.65	0.44
1:C:95:LEU:HA	1:C:117:PHE:O	2.18	0.44
1:E:81:ILE:O	1:E:85:MET:HG3	2.17	0.44
1:M:166:ARG:HD2	1:M:189:TYR:O	2.18	0.44
1:U:161:ALA:HB1	1:U:166:ARG:O	2.17	0.44
1:H:141:ASP:OD1	1:I:177:ARG:NH1	2.50	0.44
1:H:161:ALA:HA	1:H:171:LEU:HD13	1.98	0.44
1:W:125:MET:HA	1:W:179:ASN:O	2.17	0.44
1:b:61:ASP:OD1	1:b:89:LYS:NZ	2.49	0.44
1:K:135:LEU:HD12	1:K:135:LEU:N	2.32	0.44
1:P:106:PHE:CD2	1:P:160:MET:HG3	2.52	0.44
1:R:81:ILE:O	1:R:85:MET:HG3	2.17	0.44
1:U:114:GLY:HA2	1:U:193:ASP:OD2	2.17	0.44
1:K:189:TYR:CD2	1:K:191:LEU:HD11	2.53	0.44
1:I:27:ARG:HB3	3:l:101:OCA:H83	2.00	0.43
1:I:33:ILE:HD13	1:I:65:PHE:HB2	2.00	0.43
1:O:130:LEU:C	1:O:131:ILE:HD12	2.43	0.43
1:M:126:ILE:HG21	1:M:160:MET:HE1	2.00	0.43
1:I:81:ILE:O	1:I:85:MET:HG3	2.17	0.43
1:L:46:ASN:ND2	1:M:37:VAL:HG21	2.33	0.43
1:M:138:GLN:HB2	1:M:141:ASP:OD2	2.18	0.43
1:U:126:ILE:HG21	1:U:160:MET:HE1	2.01	0.43
1:D:89:LYS:N	1:D:90:PRO:HD2	2.32	0.43
1:M:197:GLU:O	4:M:302:HOH:O	2.21	0.43
1:R:61:ASP:OD1	1:R:89:LYS:NZ	2.28	0.43
1:V:65:PHE:CE2	2:y:7:MP8:HB	2.53	0.43
1:A:66:PHE:HB3	1:A:94:THR:HG22	2.00	0.43
1:G:63:ASP:OD2	1:G:115:LYS:NZ	2.35	0.43
1:H:114:GLY:N	1:H:193:ASP:OD2	2.45	0.43
1:L:24:ILE:HA	1:L:27:ARG:HH11	1.83	0.43
1:W:126:ILE:HG21	1:W:160:MET:HE1	2.00	0.43
1:b:45:ALA:HA	1:b:81:ILE:HD11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:36:LEU:HD21	1:L:40:VAL:HG22	2.00	0.43
1:O:164:CYS:HA	1:O:190:GLY:O	2.19	0.43
1:R:154:GLU:O	1:R:158:ARG:HG3	2.18	0.43
1:Y:132:SER:O	1:Y:132:SER:OG	2.28	0.43
1:J:85:MET:HE1	1:J:110:ALA:HB3	2.00	0.43
1:S:63:ASP:CG	1:S:115:LYS:HZ3	2.24	0.43
1:H:131:ILE:HD12	1:H:149:LEU:HD22	1.99	0.43
1:H:195:ILE:O	1:H:195:ILE:CG2	2.67	0.43
1:O:160:MET:HB3	1:O:171:LEU:HD22	2.01	0.43
1:P:27:ARG:HH11	1:P:27:ARG:HD3	1.46	0.43
1:Q:27:ARG:HB3	3:t:101:OCA:H83	2.01	0.43
1:E:28:LEU:CD2	3:h:101:OCA:H82	2.49	0.43
1:G:30:LYS:HZ3	1:K:62:LYS:CE	2.31	0.43
1:I:125:MET:HA	1:I:179:ASN:O	2.19	0.43
1:S:67:TYR:OH	2:v:6:ALA:O	2.36	0.43
1:a:124:ILE:HG21	1:a:191:LEU:HD13	1.99	0.43
2:e:3:ALO:H	3:e:101:OCA:C1	2.31	0.43
1:a:88:ILE:HD12	1:a:90:PRO:HG2	1.99	0.43
1:a:102:SER:HB3	1:a:103:MET:H	1.58	0.43
1:A:117:PHE:HD1	1:A:194:GLN:HB2	1.84	0.42
1:H:128:GLN:OE1	1:T:140:SER:N	2.41	0.42
1:L:46:ASN:ND2	1:M:35:PHE:CD1	2.85	0.42
1:V:34:VAL:HG23	1:V:66:PHE:CD1	2.54	0.42
1:B:79:MET:CE	1:B:155:LYS:HD3	2.45	0.42
1:C:133:GLY:O	1:C:135:LEU:N	2.53	0.42
1:Q:135:LEU:HD12	1:Q:135:LEU:O	2.18	0.42
1:Q:149:LEU:HD12	1:Q:149:LEU:HA	1.85	0.42
1:V:123:ARG:HD2	1:b:148:GLU:OE1	2.19	0.42
1:Z:80:SER:HB2	1:a:97:LEU:HD12	2.00	0.42
1:F:33:ILE:HD12	1:F:67:TYR:CZ	2.54	0.42
1:N:41:THR:H	1:N:44:SER:HG	1.66	0.42
1:O:85:MET:HE1	1:O:107:LEU:HD22	2.01	0.42
1:S:131:ILE:HD13	1:S:131:ILE:HA	1.89	0.42
1:a:132:SER:N	4:a:308:HOH:O	2.52	0.42
1:M:108:LEU:CD2	1:M:124:ILE:HD11	2.49	0.42
1:M:124:ILE:HG23	1:M:124:ILE:HD12	1.77	0.42
1:R:164:CYS:SG	1:R:191:LEU:HG	2.59	0.42
1:Y:174:ASP:OD2	1:Y:189:TYR:OH	2.13	0.42
1:A:79:MET:HE3	1:B:121:ASN:HB3	2.01	0.42
1:D:154:GLU:O	1:D:158:ARG:HG3	2.19	0.42
1:E:88:ILE:HB	1:E:90:PRO:HD2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:160:MET:HB3	1:E:160:MET:HE3	1.83	0.42
1:G:154:GLU:O	1:G:158:ARG:HG3	2.20	0.42
1:P:96:CYS:HB2	1:P:108:LEU:HD22	2.00	0.42
1:D:46:ASN:HB3	4:E:305:HOH:O	2.19	0.42
1:Q:65:PHE:CE2	2:t:7:MP8:HB	2.54	0.42
1:J:84:THR:HA	2:n:2:WFP:F1	2.10	0.42
1:O:36:LEU:HD21	1:O:40:VAL:HG22	2.02	0.42
1:S:95:LEU:HA	1:S:117:PHE:O	2.20	0.42
1:E:124:ILE:HG21	1:E:191:LEU:HD13	2.01	0.42
1:H:24:ILE:HG23	1:H:25:TYR:N	2.35	0.42
1:R:148:GLU:OE2	1:S:123:ARG:HD2	2.19	0.42
1:U:23:ASP:OD1	1:U:26:SER:HB2	2.20	0.42
1:V:54:PHE:CE1	1:W:27:ARG:NH1	2.88	0.42
1:Y:79:MET:HE2	1:Y:155:LYS:HE3	2.01	0.42
1:F:47:LEU:HD12	1:F:47:LEU:HA	1.90	0.42
1:N:28:LEU:HG	3:q:101:OCA:H82	2.02	0.42
1:O:161:ALA:HB2	1:O:171:LEU:HD13	2.01	0.42
1:V:125:MET:CE	1:V:127:HIS:CG	3.01	0.42
1:B:79:MET:HE3	1:C:121:ASN:CB	2.48	0.42
1:D:30:LYS:O	1:D:30:LYS:HG3	2.20	0.42
1:X:138:GLN:OE1	1:Y:177:ARG:NH2	2.53	0.42
1:Y:41:THR:HB	1:Y:43:GLU:OE1	2.20	0.42
1:C:30:LYS:HD2	1:C:30:LYS:HA	1.77	0.41
1:M:131:ILE:HD13	1:M:131:ILE:HG21	1.80	0.41
1:M:161:ALA:HA	1:M:171:LEU:HD13	2.02	0.41
1:P:62:LYS:O	1:P:90:PRO:HB3	2.20	0.41
1:Q:54:PHE:CG	1:R:24:ILE:HD11	2.55	0.41
1:Z:55:LEU:HD23	1:Z:55:LEU:HA	1.88	0.41
1:B:31:GLU:HG3	3:e:101:OCA:H62	2.02	0.41
1:F:45:ALA:HA	1:F:81:ILE:HD11	2.01	0.41
1:H:96:CYS:HB2	1:H:108:LEU:HD22	2.02	0.41
1:N:167:ASP:HB3	4:N:328:HOH:O	2.20	0.41
1:P:113:LYS:HE2	1:P:163:HIS:O	2.20	0.41
1:S:36:LEU:HD21	1:S:40:VAL:HG22	2.01	0.41
1:I:95:LEU:HA	1:I:117:PHE:O	2.20	0.41
1:L:23:ASP:OD1	4:L:302:HOH:O	2.21	0.41
1:O:184:GLU:HG2	1:O:195:ILE:CD1	2.50	0.41
1:Q:109:SER:HB3	1:Q:191:LEU:HD12	2.01	0.41
1:V:164:CYS:HA	1:V:190:GLY:O	2.20	0.41
2:o:4:PRO:HA	2:o:5:YCP:HA	1.86	0.41
1:A:127:HIS:ND1	1:A:178:ASP:OD1	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:125:MET:HA	1:S:179:ASN:O	2.21	0.41
1:b:45:ALA:HB2	1:b:77:ALA:HB1	2.03	0.41
1:F:106:PHE:CD1	1:F:160:MET:HE2	2.41	0.41
1:O:184:GLU:HA	1:O:195:ILE:HD11	2.02	0.41
1:S:113:LYS:NZ	1:S:163:HIS:C	2.73	0.41
1:V:24:ILE:HD11	1:b:54:PHE:CG	2.54	0.41
1:Y:113:LYS:NZ	1:Y:165:ASP:OD2	2.49	0.41
1:a:117:PHE:CD1	1:a:194:GLN:HB2	2.54	0.41
1:H:24:ILE:HD11	1:N:54:PHE:HB2	2.02	0.41
1:K:76:THR:HG21	1:L:99:GLN:CG	2.50	0.41
1:N:62:LYS:O	1:N:90:PRO:HB3	2.21	0.41
1:S:117:PHE:CD1	1:S:194:GLN:HB2	2.56	0.41
1:V:45:ALA:HA	1:V:81:ILE:HD11	2.03	0.41
1:W:119:LEU:HD23	1:W:119:LEU:HA	1.84	0.41
1:Z:119:LEU:N	1:Z:119:LEU:CD1	2.83	0.41
1:b:159:LEU:HD23	1:b:159:LEU:HA	1.84	0.41
1:J:154:GLU:O	1:J:158:ARG:HG3	2.21	0.41
1:K:69:ASN:HB2	1:K:97:LEU:O	2.20	0.41
1:E:28:LEU:HD12	1:E:35:PHE:HE2	1.86	0.41
1:I:62:LYS:O	1:I:90:PRO:HB3	2.20	0.41
1:K:117:PHE:CD1	1:K:194:GLN:CB	3.03	0.41
1:N:63:ASP:OD1	1:N:91:ASP:HB2	2.20	0.41
1:V:30:LYS:NZ	4:V:303:HOH:O	2.33	0.41
1:X:155:LYS:HZ3	1:X:159:LEU:CG	2.33	0.41
1:Z:75:VAL:HG22	1:Z:103:MET:SD	2.60	0.41
1:F:81:ILE:O	1:F:85:MET:HG3	2.20	0.41
1:J:96:CYS:HB2	1:J:108:LEU:HD22	2.02	0.41
1:Q:45:ALA:HB2	1:Q:77:ALA:HB1	2.03	0.41
1:R:75:VAL:HG22	1:R:103:MET:HE3	2.03	0.41
1:R:145:HIS:NE2	1:S:178:ASP:OD2	2.50	0.41
1:X:45:ALA:HA	1:X:81:ILE:HD11	2.02	0.41
1:Z:184:GLU:HG3	1:Z:195:ILE:CD1	2.50	0.41
2:f:3:ALO:N	3:f:101:OCA:O1	2.53	0.41
1:A:125:MET:HG3	1:A:180:PHE:CE1	2.56	0.41
1:G:62:LYS:O	1:G:90:PRO:HB3	2.21	0.41
2:s:7:MP8:HD	3:s:101:OCA:H22	2.02	0.41
1:F:181:MET:HE1	1:F:191:LEU:CD1	2.42	0.40
1:G:30:LYS:NZ	1:K:62:LYS:HZ2	2.16	0.40
1:K:117:PHE:HD1	1:K:194:GLN:CB	2.35	0.40
1:a:84:THR:HA	2:4:2:WFP:F1	2.11	0.40
1:C:106:PHE:HD1	1:C:160:MET:HE3	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:117:PHE:CD1	1:C:194:GLN:HG3	2.57	0.40
1:F:55:LEU:HD23	1:F:55:LEU:HA	1.91	0.40
4:P:301:HOH:O	2:s:4:PRO:O	2.22	0.40
1:b:91:ASP:OD1	4:b:302:HOH:O	2.22	0.40
1:A:63:ASP:OD2	1:A:115:LYS:NZ	2.47	0.40
1:D:159:LEU:O	1:D:162:LYS:HB3	2.21	0.40
1:H:52:LEU:HD23	1:H:52:LEU:HA	1.86	0.40
1:K:96:CYS:HB2	1:K:108:LEU:HD22	2.04	0.40
1:O:125:MET:HE1	1:O:127:HIS:ND1	2.36	0.40
1:R:100:ALA:O	1:R:105:ALA:HB2	2.22	0.40
1:a:126:ILE:HG21	1:a:160:MET:HE1	2.04	0.40
1:H:33:ILE:HD11	2:k:7:MP8:HDA	2.03	0.40
1:L:53:LEU:HB3	3:p:101:OCA:H41	2.04	0.40
1:U:33:ILE:HA	1:U:65:PHE:O	2.22	0.40
1:Y:55:LEU:HA	1:Y:55:LEU:HD23	1.82	0.40
1:a:159:LEU:HD23	1:a:159:LEU:HA	1.84	0.40
2:3:3:ALO:HA	2:3:4:PRO:C	2.45	0.40
1:I:26:SER:O	1:I:30:LYS:HG3	2.21	0.40
1:L:84:THR:HA	2:p:2:WFP:F1	2.12	0.40
1:P:47:LEU:HA	1:P:47:LEU:HD23	1.85	0.40
1:U:113:LYS:HA	1:U:114:GLY:HA2	1.71	0.40
1:Y:102:SER:HB3	1:Y:103:MET:H	1.75	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	173/217 (80%)	168 (97%)	5 (3%)	0	100	100
1	B	169/217 (78%)	166 (98%)	3 (2%)	0	100	100
1	C	169/217 (78%)	166 (98%)	3 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	169/217 (78%)	165 (98%)	4 (2%)	0	100	100
1	E	170/217 (78%)	165 (97%)	4 (2%)	1 (1%)	21	44
1	F	167/217 (77%)	165 (99%)	2 (1%)	0	100	100
1	G	169/217 (78%)	165 (98%)	4 (2%)	0	100	100
1	H	170/217 (78%)	165 (97%)	3 (2%)	2 (1%)	10	27
1	I	167/217 (77%)	165 (99%)	2 (1%)	0	100	100
1	J	168/217 (77%)	166 (99%)	2 (1%)	0	100	100
1	K	170/217 (78%)	165 (97%)	5 (3%)	0	100	100
1	L	170/217 (78%)	166 (98%)	4 (2%)	0	100	100
1	M	167/217 (77%)	165 (99%)	2 (1%)	0	100	100
1	N	169/217 (78%)	167 (99%)	2 (1%)	0	100	100
1	O	168/217 (77%)	164 (98%)	3 (2%)	1 (1%)	21	44
1	P	169/217 (78%)	167 (99%)	2 (1%)	0	100	100
1	Q	169/217 (78%)	166 (98%)	3 (2%)	0	100	100
1	R	171/217 (79%)	168 (98%)	3 (2%)	0	100	100
1	S	169/217 (78%)	165 (98%)	4 (2%)	0	100	100
1	T	167/217 (77%)	165 (99%)	2 (1%)	0	100	100
1	U	172/217 (79%)	165 (96%)	5 (3%)	2 (1%)	10	27
1	V	167/217 (77%)	164 (98%)	2 (1%)	1 (1%)	21	44
1	W	169/217 (78%)	167 (99%)	2 (1%)	0	100	100
1	X	170/217 (78%)	165 (97%)	5 (3%)	0	100	100
1	Y	168/217 (77%)	165 (98%)	3 (2%)	0	100	100
1	Z	168/217 (77%)	165 (98%)	3 (2%)	0	100	100
1	a	169/217 (78%)	167 (99%)	2 (1%)	0	100	100
1	b	168/217 (77%)	166 (99%)	2 (1%)	0	100	100
2	0	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	1	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	2	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	3	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	4	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	c	2/6 (33%)	1 (50%)	1 (50%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	e	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	f	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	g	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	h	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	i	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	j	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	k	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	l	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	m	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	n	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	o	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	p	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	q	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	r	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	s	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	t	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	u	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	v	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	w	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	x	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	y	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
2	z	2/6 (33%)	1 (50%)	1 (50%)	0	100	100
All	All	4787/6244 (77%)	4666 (98%)	114 (2%)	7 (0%)	48	73

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	136	GLY
1	H	136	GLY
1	H	196	LEU
1	U	132	SER
1	U	133	GLY
1	V	130	LEU
1	O	197	GLU

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	148/185 (80%)	148 (100%)	0	100	100
1	B	147/185 (80%)	147 (100%)	0	100	100
1	C	147/185 (80%)	145 (99%)	2 (1%)	59	82
1	D	148/185 (80%)	148 (100%)	0	100	100
1	E	149/185 (80%)	149 (100%)	0	100	100
1	F	147/185 (80%)	146 (99%)	1 (1%)	76	90
1	G	148/185 (80%)	148 (100%)	0	100	100
1	H	149/185 (80%)	149 (100%)	0	100	100
1	I	147/185 (80%)	147 (100%)	0	100	100
1	J	148/185 (80%)	147 (99%)	1 (1%)	76	90
1	K	149/185 (80%)	148 (99%)	1 (1%)	76	90
1	L	148/185 (80%)	148 (100%)	0	100	100
1	M	147/185 (80%)	146 (99%)	1 (1%)	76	90
1	N	147/185 (80%)	147 (100%)	0	100	100
1	O	148/185 (80%)	148 (100%)	0	100	100
1	P	147/185 (80%)	146 (99%)	1 (1%)	76	90
1	Q	148/185 (80%)	148 (100%)	0	100	100
1	R	149/185 (80%)	147 (99%)	2 (1%)	61	83
1	S	148/185 (80%)	148 (100%)	0	100	100
1	T	147/185 (80%)	147 (100%)	0	100	100
1	U	147/185 (80%)	146 (99%)	1 (1%)	76	90
1	V	147/185 (80%)	147 (100%)	0	100	100
1	W	147/185 (80%)	147 (100%)	0	100	100
1	X	148/185 (80%)	147 (99%)	1 (1%)	76	90
1	Y	148/185 (80%)	148 (100%)	0	100	100
1	Z	147/185 (80%)	147 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	a	148/185 (80%)	148 (100%)	0	100	100
1	b	147/185 (80%)	146 (99%)	1 (1%)	76	90
2	0	1/1 (100%)	1 (100%)	0	100	100
2	1	1/1 (100%)	1 (100%)	0	100	100
2	2	1/1 (100%)	1 (100%)	0	100	100
2	3	1/1 (100%)	1 (100%)	0	100	100
2	4	1/1 (100%)	1 (100%)	0	100	100
2	c	1/1 (100%)	1 (100%)	0	100	100
2	e	1/1 (100%)	1 (100%)	0	100	100
2	f	1/1 (100%)	1 (100%)	0	100	100
2	g	1/1 (100%)	1 (100%)	0	100	100
2	h	1/1 (100%)	1 (100%)	0	100	100
2	i	1/1 (100%)	1 (100%)	0	100	100
2	j	1/1 (100%)	1 (100%)	0	100	100
2	k	1/1 (100%)	1 (100%)	0	100	100
2	l	1/1 (100%)	1 (100%)	0	100	100
2	m	1/1 (100%)	1 (100%)	0	100	100
2	n	1/1 (100%)	1 (100%)	0	100	100
2	o	1/1 (100%)	1 (100%)	0	100	100
2	p	1/1 (100%)	1 (100%)	0	100	100
2	q	1/1 (100%)	1 (100%)	0	100	100
2	r	1/1 (100%)	1 (100%)	0	100	100
2	s	1/1 (100%)	1 (100%)	0	100	100
2	t	1/1 (100%)	1 (100%)	0	100	100
2	u	1/1 (100%)	1 (100%)	0	100	100
2	v	1/1 (100%)	1 (100%)	0	100	100
2	w	1/1 (100%)	1 (100%)	0	100	100
2	x	1/1 (100%)	1 (100%)	0	100	100
2	y	1/1 (100%)	1 (100%)	0	100	100
2	z	1/1 (100%)	1 (100%)	0	100	100
All	All	4163/5208 (80%)	4151 (100%)	12 (0%)	86	94

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	43	GLU
1	C	135	LEU
1	F	127	HIS
1	J	130	LEU
1	K	135	LEU
1	M	171	LEU
1	P	160	MET
1	R	123	ARG
1	R	130	LEU
1	U	135	LEU
1	X	135	LEU
1	b	76	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	ASN
1	C	46	ASN
1	D	46	ASN
1	G	194	GLN
1	I	121	ASN
1	M	69	ASN
1	O	121	ASN
1	P	59	ASN
1	P	121	ASN
1	R	69	ASN
1	S	46	ASN
1	U	69	ASN
1	W	157	ASN
1	Y	46	ASN
1	Z	127	HIS
1	b	46	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

112 non-standard protein/DNA/RNA residues 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	MP8	i	7	2	6,8,9	4.12	4 (66%)	3,10,12	0.65	0
2	YCP	o	5	2	6,8,9	1.37	1 (16%)	7,9,11	1.78	3 (42%)
2	WFP	3	2	3,2	12,13,14	1.03	0	12,17,19	1.46	3 (25%)
2	WFP	s	2	3,2	12,13,14	1.31	0	12,17,19	1.87	4 (33%)
2	ALO	t	3	2	5,6,7	0.57	0	5,7,9	0.69	0
2	WFP	i	2	3,2	12,13,14	1.09	0	12,17,19	1.99	3 (25%)
2	ALO	o	3	2	5,6,7	0.58	0	5,7,9	0.72	0
2	ALO	w	3	2	5,6,7	0.72	0	5,7,9	1.17	1 (20%)
2	MP8	h	7	2	6,8,9	3.97	4 (66%)	3,10,12	1.43	1 (33%)
2	ALO	q	3	2	5,6,7	0.70	0	5,7,9	0.61	0
2	WFP	l	2	3,2	12,13,14	1.11	1 (8%)	12,17,19	1.49	3 (25%)
2	MP8	c	7	2	6,8,9	3.90	4 (66%)	3,10,12	1.25	0
2	YCP	2	5	2	6,8,9	1.42	1 (16%)	7,9,11	1.17	0
2	ALO	i	3	2	5,6,7	0.71	0	5,7,9	0.63	0
2	YCP	t	5	2	6,8,9	1.50	2 (33%)	7,9,11	1.55	2 (28%)
2	WFP	4	2	3,2	12,13,14	1.03	0	12,17,19	1.61	4 (33%)
2	MP8	3	7	2	6,8,9	3.92	4 (66%)	3,10,12	1.46	0
2	MP8	o	7	2	6,8,9	4.03	3 (50%)	3,10,12	1.37	0
2	YCP	p	5	2	6,8,9	1.52	2 (33%)	7,9,11	1.59	2 (28%)
2	ALO	s	3	2	5,6,7	0.69	0	5,7,9	1.00	1 (20%)
2	WFP	g	2	3,2	12,13,14	1.05	0	12,17,19	1.57	4 (33%)
2	WFP	n	2	3,2	12,13,14	1.11	0	12,17,19	1.92	5 (41%)
2	ALO	f	3	2	5,6,7	0.67	0	5,7,9	0.49	0
2	ALO	l	3	2	5,6,7	0.75	0	5,7,9	1.05	0
2	ALO	v	3	2	5,6,7	0.59	0	5,7,9	1.28	1 (20%)
2	WFP	j	2	3,2	12,13,14	1.15	1 (8%)	12,17,19	1.59	4 (33%)
2	YCP	y	5	2	6,8,9	1.47	2 (33%)	7,9,11	1.07	0
2	MP8	l	7	2	6,8,9	3.95	3 (50%)	3,10,12	1.01	0
2	MP8	x	7	2	6,8,9	4.06	3 (50%)	3,10,12	1.00	0
2	MP8	t	7	2	6,8,9	4.05	4 (66%)	3,10,12	1.42	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	WFP	0	2	3,2	12,13,14	1.08	0	12,17,19	1.66	3 (25%)
2	ALO	l	3	2	5,6,7	0.69	0	5,7,9	0.78	0
2	MP8	r	7	2	6,8,9	4.09	4 (66%)	3,10,12	1.35	0
2	YCP	s	5	2	6,8,9	1.54	2 (33%)	7,9,11	1.24	1 (14%)
2	YCP	z	5	2	6,8,9	1.46	1 (16%)	7,9,11	1.27	1 (14%)
2	MP8	f	7	2	6,8,9	4.09	4 (66%)	3,10,12	1.34	0
2	ALO	y	3	2	5,6,7	0.81	0	5,7,9	0.94	0
2	MP8	p	7	2	6,8,9	3.85	4 (66%)	3,10,12	0.74	0
2	YCP	w	5	2	6,8,9	1.41	2 (33%)	7,9,11	1.58	3 (42%)
2	MP8	m	7	2	6,8,9	3.95	4 (66%)	3,10,12	1.49	0
2	YCP	k	5	2	6,8,9	1.47	2 (33%)	7,9,11	1.30	1 (14%)
2	WFP	z	2	3,2	12,13,14	1.21	1 (8%)	12,17,19	1.89	4 (33%)
2	YCP	4	5	2	6,8,9	1.51	2 (33%)	7,9,11	1.24	0
2	WFP	y	2	3,2	12,13,14	0.99	0	12,17,19	1.71	4 (33%)
2	ALO	j	3	2	5,6,7	0.62	0	5,7,9	0.88	0
2	WFP	e	2	3,2	12,13,14	1.09	0	12,17,19	1.64	4 (33%)
2	YCP	i	5	2	6,8,9	1.34	1 (16%)	7,9,11	1.31	1 (14%)
2	YCP	0	5	2	6,8,9	1.61	2 (33%)	7,9,11	1.20	1 (14%)
2	MP8	s	7	2	6,8,9	4.06	4 (66%)	3,10,12	1.06	0
2	MP8	z	7	2	6,8,9	4.20	3 (50%)	3,10,12	1.46	1 (33%)
2	YCP	r	5	2	6,8,9	1.40	1 (16%)	7,9,11	1.32	1 (14%)
2	ALO	c	3	2	5,6,7	0.86	0	5,7,9	0.74	0
2	WFP	c	2	3,2	12,13,14	1.22	0	12,17,19	1.49	1 (8%)
2	ALO	0	3	2	5,6,7	0.70	0	5,7,9	0.71	0
2	YCP	l	5	2	6,8,9	1.42	2 (33%)	7,9,11	1.19	1 (14%)
2	MP8	4	7	2	6,8,9	3.88	3 (50%)	3,10,12	0.95	0
2	WFP	1	2	3,2	12,13,14	1.04	0	12,17,19	1.68	3 (25%)
2	YCP	n	5	2	6,8,9	1.43	2 (33%)	7,9,11	1.57	1 (14%)
2	WFP	m	2	3,2	12,13,14	1.17	0	12,17,19	1.77	3 (25%)
2	YCP	j	5	2	6,8,9	1.47	2 (33%)	7,9,11	1.44	1 (14%)
2	YCP	m	5	2	6,8,9	1.48	2 (33%)	7,9,11	1.47	1 (14%)
2	YCP	e	5	2	6,8,9	1.39	2 (33%)	7,9,11	0.98	0
2	YCP	1	5	2	6,8,9	1.30	1 (16%)	7,9,11	2.32	2 (28%)
2	WFP	h	2	3,2	12,13,14	1.10	0	12,17,19	1.49	3 (25%)
2	MP8	j	7	2	6,8,9	4.02	4 (66%)	3,10,12	1.61	1 (33%)
2	WFP	2	2	3,2	12,13,14	1.14	0	12,17,19	1.62	3 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	WFP	f	2	3,2	12,13,14	1.01	0	12,17,19	1.47	3 (25%)
2	MP8	0	7	2	6,8,9	4.08	4 (66%)	3,10,12	1.26	0
2	ALO	n	3	2	5,6,7	0.61	0	5,7,9	0.61	0
2	YCP	u	5	2	6,8,9	1.54	2 (33%)	7,9,11	0.80	0
2	MP8	1	7	2	6,8,9	3.96	4 (66%)	3,10,12	1.11	0
2	WFP	v	2	3,2	12,13,14	1.26	0	12,17,19	1.62	3 (25%)
2	ALO	k	3	2	5,6,7	0.67	0	5,7,9	1.01	0
2	YCP	v	5	2	6,8,9	1.58	2 (33%)	7,9,11	1.23	1 (14%)
2	MP8	g	7	2	6,8,9	4.01	4 (66%)	3,10,12	1.01	0
2	ALO	3	3	2	5,6,7	0.70	0	5,7,9	0.45	0
2	MP8	n	7	2	6,8,9	3.92	3 (50%)	3,10,12	1.45	0
2	YCP	x	5	2	6,8,9	1.48	2 (33%)	7,9,11	2.01	3 (42%)
2	WFP	w	2	3,2	12,13,14	1.17	1 (8%)	12,17,19	1.56	3 (25%)
2	MP8	k	7	2	6,8,9	3.86	4 (66%)	3,10,12	1.39	0
2	YCP	h	5	2	6,8,9	1.53	2 (33%)	7,9,11	1.41	1 (14%)
2	MP8	e	7	2	6,8,9	3.97	4 (66%)	3,10,12	1.53	0
2	ALO	m	3	2	5,6,7	0.68	0	5,7,9	0.56	0
2	ALO	x	3	2	5,6,7	0.55	0	5,7,9	0.85	0
2	WFP	u	2	3,2	12,13,14	1.09	1 (8%)	12,17,19	1.91	4 (33%)
2	ALO	4	3	2	5,6,7	0.71	0	5,7,9	0.94	0
2	ALO	z	3	2	5,6,7	0.82	0	5,7,9	0.80	0
2	MP8	2	7	2	6,8,9	4.21	4 (66%)	3,10,12	1.33	1 (33%)
2	YCP	g	5	2	6,8,9	1.48	2 (33%)	7,9,11	1.22	0
2	YCP	f	5	2	6,8,9	1.36	2 (33%)	7,9,11	1.23	1 (14%)
2	MP8	q	7	2	6,8,9	4.10	3 (50%)	3,10,12	1.61	1 (33%)
2	WFP	r	2	3,2	12,13,14	1.23	0	12,17,19	1.44	3 (25%)
2	YCP	c	5	2	6,8,9	1.41	1 (16%)	7,9,11	1.57	2 (28%)
2	ALO	h	3	2	5,6,7	0.84	0	5,7,9	0.80	0
2	ALO	g	3	2	5,6,7	0.70	0	5,7,9	1.19	0
2	MP8	u	7	2	6,8,9	4.16	4 (66%)	3,10,12	1.31	0
2	WFP	o	2	3,2	12,13,14	1.09	0	12,17,19	1.53	3 (25%)
2	WFP	t	2	3,2	12,13,14	1.12	0	12,17,19	1.32	1 (8%)
2	MP8	v	7	2	6,8,9	4.27	4 (66%)	3,10,12	1.16	0
2	ALO	u	3	2	5,6,7	0.67	0	5,7,9	0.86	0
2	ALO	e	3	2	5,6,7	0.81	0	5,7,9	0.76	0
2	YCP	q	5	2	6,8,9	1.39	1 (16%)	7,9,11	1.21	1 (14%)
2	WFP	k	2	3,2	12,13,14	1.05	0	12,17,19	1.67	3 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	WFP	p	2	3,2	12,13,14	1.10	0	12,17,19	1.80	5 (41%)
2	MP8	w	7	2	6,8,9	4.05	3 (50%)	3,10,12	1.20	0
2	ALO	r	3	2	5,6,7	0.74	0	5,7,9	0.63	0
2	MP8	y	7	2	6,8,9	3.96	4 (66%)	3,10,12	1.02	0
2	ALO	2	3	2	5,6,7	0.65	0	5,7,9	0.67	0
2	WFP	x	2	3,2	12,13,14	1.05	0	12,17,19	1.53	2 (16%)
2	WFP	q	2	3,2	12,13,14	0.96	0	12,17,19	1.72	3 (25%)
2	ALO	p	3	2	5,6,7	0.77	0	5,7,9	0.77	0
2	YCP	3	5	2	6,8,9	1.31	1 (16%)	7,9,11	2.22	3 (42%)

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	MP8	i	7	2	-	0/0/11/13	0/1/1/1
2	YCP	o	5	2	-	0/1/10/12	1/1/1/1
2	WFP	3	2	3,2	-	0/5/6/8	0/1/1/1
2	WFP	s	2	3,2	-	0/5/6/8	0/1/1/1
2	ALO	t	3	2	-	1/5/6/8	-
2	WFP	i	2	3,2	-	0/5/6/8	0/1/1/1
2	ALO	o	3	2	-	1/5/6/8	-
2	ALO	w	3	2	-	1/5/6/8	-
2	MP8	h	7	2	-	0/0/11/13	0/1/1/1
2	ALO	q	3	2	-	1/5/6/8	-
2	WFP	l	2	3,2	-	0/5/6/8	0/1/1/1
2	MP8	c	7	2	-	0/0/11/13	0/1/1/1
2	YCP	2	5	2	-	0/1/10/12	0/1/1/1
2	ALO	i	3	2	-	1/5/6/8	-
2	YCP	t	5	2	-	0/1/10/12	1/1/1/1
2	WFP	4	2	3,2	-	0/5/6/8	0/1/1/1
2	MP8	3	7	2	-	0/0/11/13	0/1/1/1
2	MP8	o	7	2	-	0/0/11/13	0/1/1/1
2	YCP	p	5	2	-	1/1/10/12	1/1/1/1
2	ALO	s	3	2	-	1/5/6/8	-
2	WFP	g	2	3,2	-	0/5/6/8	0/1/1/1
2	WFP	n	2	3,2	-	0/5/6/8	0/1/1/1
2	ALO	f	3	2	-	1/5/6/8	-
2	ALO	l	3	2	-	1/5/6/8	-
2	ALO	v	3	2	-	1/5/6/8	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	WFP	j	2	3,2	-	0/5/6/8	0/1/1/1
2	YCP	y	5	2	-	1/1/10/12	1/1/1/1
2	MP8	l	7	2	-	0/0/11/13	0/1/1/1
2	MP8	x	7	2	-	0/0/11/13	0/1/1/1
2	MP8	t	7	2	-	0/0/11/13	0/1/1/1
2	WFP	0	2	3,2	-	0/5/6/8	0/1/1/1
2	ALO	l	3	2	-	1/5/6/8	-
2	MP8	r	7	2	-	0/0/11/13	0/1/1/1
2	YCP	s	5	2	-	0/1/10/12	1/1/1/1
2	YCP	z	5	2	-	1/1/10/12	1/1/1/1
2	MP8	f	7	2	-	0/0/11/13	0/1/1/1
2	ALO	y	3	2	-	1/5/6/8	-
2	MP8	p	7	2	-	0/0/11/13	0/1/1/1
2	YCP	w	5	2	-	0/1/10/12	0/1/1/1
2	MP8	m	7	2	-	0/0/11/13	0/1/1/1
2	YCP	k	5	2	-	0/1/10/12	1/1/1/1
2	WFP	z	2	3,2	-	0/5/6/8	0/1/1/1
2	YCP	4	5	2	-	0/1/10/12	1/1/1/1
2	WFP	y	2	3,2	-	0/5/6/8	0/1/1/1
2	ALO	j	3	2	-	1/5/6/8	-
2	WFP	e	2	3,2	-	0/5/6/8	0/1/1/1
2	YCP	i	5	2	-	0/1/10/12	1/1/1/1
2	YCP	0	5	2	-	0/1/10/12	1/1/1/1
2	MP8	s	7	2	-	0/0/11/13	0/1/1/1
2	MP8	z	7	2	-	0/0/11/13	0/1/1/1
2	YCP	r	5	2	-	0/1/10/12	1/1/1/1
2	ALO	c	3	2	-	1/5/6/8	-
2	WFP	c	2	3,2	-	0/5/6/8	0/1/1/1
2	ALO	0	3	2	-	1/5/6/8	-
2	YCP	l	5	2	-	0/1/10/12	1/1/1/1
2	MP8	4	7	2	-	0/0/11/13	0/1/1/1
2	WFP	1	2	3,2	-	0/5/6/8	0/1/1/1
2	YCP	n	5	2	-	1/1/10/12	1/1/1/1
2	WFP	m	2	3,2	-	0/5/6/8	0/1/1/1
2	YCP	j	5	2	-	1/1/10/12	1/1/1/1
2	YCP	m	5	2	-	1/1/10/12	1/1/1/1
2	YCP	e	5	2	-	0/1/10/12	1/1/1/1
2	YCP	1	5	2	-	0/1/10/12	0/1/1/1
2	WFP	h	2	3,2	-	0/5/6/8	0/1/1/1
2	MP8	j	7	2	-	0/0/11/13	0/1/1/1
2	WFP	2	2	3,2	-	0/5/6/8	0/1/1/1
2	WFP	f	2	3,2	-	0/5/6/8	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MP8	0	7	2	-	0/0/11/13	0/1/1/1
2	ALO	n	3	2	-	1/5/6/8	-
2	YCP	u	5	2	-	1/1/10/12	1/1/1/1
2	MP8	1	7	2	-	0/0/11/13	0/1/1/1
2	WFP	v	2	3,2	-	0/5/6/8	0/1/1/1
2	ALO	k	3	2	-	1/5/6/8	-
2	YCP	v	5	2	-	1/1/10/12	1/1/1/1
2	MP8	g	7	2	-	0/0/11/13	0/1/1/1
2	ALO	3	3	2	-	1/5/6/8	-
2	MP8	n	7	2	-	0/0/11/13	0/1/1/1
2	YCP	x	5	2	-	0/1/10/12	0/1/1/1
2	WFP	w	2	3,2	-	0/5/6/8	0/1/1/1
2	MP8	k	7	2	-	0/0/11/13	0/1/1/1
2	YCP	h	5	2	-	0/1/10/12	1/1/1/1
2	MP8	e	7	2	-	0/0/11/13	0/1/1/1
2	ALO	m	3	2	-	1/5/6/8	-
2	ALO	x	3	2	-	1/5/6/8	-
2	WFP	u	2	3,2	-	0/5/6/8	0/1/1/1
2	ALO	4	3	2	-	1/5/6/8	-
2	ALO	z	3	2	-	1/5/6/8	-
2	MP8	2	7	2	-	0/0/11/13	0/1/1/1
2	YCP	g	5	2	-	0/1/10/12	1/1/1/1
2	YCP	f	5	2	-	0/1/10/12	1/1/1/1
2	MP8	q	7	2	-	0/0/11/13	0/1/1/1
2	WFP	r	2	3,2	-	0/5/6/8	0/1/1/1
2	YCP	c	5	2	-	1/1/10/12	1/1/1/1
2	ALO	h	3	2	-	1/5/6/8	-
2	ALO	g	3	2	-	1/5/6/8	-
2	MP8	u	7	2	-	0/0/11/13	0/1/1/1
2	WFP	o	2	3,2	-	0/5/6/8	0/1/1/1
2	WFP	t	2	3,2	-	0/5/6/8	0/1/1/1
2	MP8	v	7	2	-	0/0/11/13	0/1/1/1
2	ALO	u	3	2	-	1/5/6/8	-
2	ALO	e	3	2	-	1/5/6/8	-
2	YCP	q	5	2	-	0/1/10/12	1/1/1/1
2	WFP	k	2	3,2	-	0/5/6/8	0/1/1/1
2	WFP	p	2	3,2	-	0/5/6/8	0/1/1/1
2	MP8	w	7	2	-	0/0/11/13	0/1/1/1
2	ALO	r	3	2	-	1/5/6/8	-
2	MP8	y	7	2	-	0/0/11/13	0/1/1/1
2	ALO	2	3	2	-	1/5/6/8	-
2	WFP	x	2	3,2	-	0/5/6/8	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	WFP	q	2	3,2	-	0/5/6/8	0/1/1/1
2	ALO	p	3	2	-	1/5/6/8	-
2	YCP	3	5	2	-	0/1/10/12	0/1/1/1

All (156) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	2	7	MP8	CA-N	-7.80	1.35	1.48
2	z	7	MP8	CA-N	-7.75	1.35	1.48
2	v	7	MP8	CA-N	-7.53	1.35	1.48
2	u	7	MP8	CA-N	-7.50	1.35	1.48
2	q	7	MP8	CA-N	-7.48	1.35	1.48
2	x	7	MP8	CA-N	-7.38	1.35	1.48
2	0	7	MP8	CA-N	-7.36	1.35	1.48
2	s	7	MP8	CA-N	-7.35	1.35	1.48
2	f	7	MP8	CA-N	-7.26	1.36	1.48
2	i	7	MP8	CA-N	-7.24	1.36	1.48
2	t	7	MP8	CA-N	-7.18	1.36	1.48
2	j	7	MP8	CA-N	-7.17	1.36	1.48
2	o	7	MP8	CA-N	-7.17	1.36	1.48
2	w	7	MP8	CA-N	-7.07	1.36	1.48
2	r	7	MP8	CA-N	-7.05	1.36	1.48
2	3	7	MP8	CA-N	-6.99	1.36	1.48
2	m	7	MP8	CA-N	-6.94	1.36	1.48
2	l	7	MP8	CA-N	-6.93	1.36	1.48
2	y	7	MP8	CA-N	-6.93	1.36	1.48
2	c	7	MP8	CA-N	-6.92	1.36	1.48
2	h	7	MP8	CA-N	-6.91	1.36	1.48
2	n	7	MP8	CA-N	-6.89	1.36	1.48
2	1	7	MP8	CA-N	-6.88	1.36	1.48
2	g	7	MP8	CA-N	-6.84	1.36	1.48
2	e	7	MP8	CA-N	-6.75	1.36	1.48
2	4	7	MP8	CA-N	-6.72	1.37	1.48
2	p	7	MP8	CA-N	-6.68	1.37	1.48
2	k	7	MP8	CA-N	-6.65	1.37	1.48
2	u	7	MP8	CD-N	5.96	1.67	1.47
2	e	7	MP8	CD-N	5.94	1.67	1.47
2	r	7	MP8	CD-N	5.93	1.67	1.47
2	1	7	MP8	CD-N	5.93	1.67	1.47
2	g	7	MP8	CD-N	5.92	1.67	1.47
2	h	7	MP8	CD-N	5.90	1.67	1.47
2	v	7	MP8	CD-N	5.89	1.67	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	w	7	MP8	CD-N	5.86	1.67	1.47
2	i	7	MP8	CD-N	5.85	1.67	1.47
2	o	7	MP8	CD-N	5.83	1.67	1.47
2	0	7	MP8	CD-N	5.81	1.67	1.47
2	m	7	MP8	CD-N	5.81	1.67	1.47
2	j	7	MP8	CD-N	5.81	1.66	1.47
2	t	7	MP8	CD-N	5.81	1.66	1.47
2	4	7	MP8	CD-N	5.76	1.66	1.47
2	f	7	MP8	CD-N	5.74	1.66	1.47
2	q	7	MP8	CD-N	5.74	1.66	1.47
2	l	7	MP8	CD-N	5.73	1.66	1.47
2	z	7	MP8	CD-N	5.71	1.66	1.47
2	n	7	MP8	CD-N	5.69	1.66	1.47
2	x	7	MP8	CD-N	5.63	1.66	1.47
2	s	7	MP8	CD-N	5.63	1.66	1.47
2	2	7	MP8	CD-N	5.61	1.66	1.47
2	y	7	MP8	CD-N	5.60	1.66	1.47
2	c	7	MP8	CD-N	5.59	1.66	1.47
2	3	7	MP8	CD-N	5.55	1.66	1.47
2	k	7	MP8	CD-N	5.53	1.66	1.47
2	p	7	MP8	CD-N	5.53	1.66	1.47
2	v	7	MP8	CB-CG	3.19	1.61	1.52
2	w	7	MP8	CB-CG	2.99	1.61	1.52
2	f	7	MP8	CB-CG	2.98	1.61	1.52
2	r	7	MP8	CB-CG	2.95	1.60	1.52
2	2	7	MP8	CB-CG	2.89	1.60	1.52
2	x	7	MP8	CB-CG	2.87	1.60	1.52
2	k	7	MP8	CB-CG	2.86	1.60	1.52
2	e	7	MP8	CB-CG	2.86	1.60	1.52
2	y	7	MP8	CB-CG	2.85	1.60	1.52
2	g	7	MP8	CB-CG	2.84	1.60	1.52
2	l	7	MP8	CB-CG	2.84	1.60	1.52
2	i	7	MP8	CB-CG	2.81	1.60	1.52
2	4	7	MP8	CB-CG	2.81	1.60	1.52
2	o	7	MP8	CB-CG	2.77	1.60	1.52
2	p	7	MP8	CB-CG	2.76	1.60	1.52
2	s	7	MP8	CB-CG	2.74	1.60	1.52
2	z	7	MP8	CB-CG	2.73	1.60	1.52
2	0	5	YCP	CG-CB	-2.71	1.46	1.53
2	n	7	MP8	CB-CG	2.69	1.60	1.52
2	q	7	MP8	CB-CG	2.66	1.60	1.52
2	t	7	MP8	CB-CG	2.66	1.60	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	j	7	MP8	CB-CG	2.63	1.60	1.52
2	m	7	MP8	CB-CG	2.62	1.59	1.52
2	c	7	MP8	CB-CG	2.60	1.59	1.52
2	h	7	MP8	CB-CG	2.59	1.59	1.52
2	s	5	YCP	CG-CB	-2.57	1.46	1.53
2	l	7	MP8	CB-CG	2.57	1.59	1.52
2	o	7	MP8	CB-CG	2.57	1.59	1.52
2	x	5	YCP	CE-N	2.57	1.54	1.47
2	v	5	YCP	CE-N	2.56	1.53	1.47
2	3	7	MP8	CB-CG	2.56	1.59	1.52
2	p	5	YCP	CE-N	2.53	1.53	1.47
2	z	5	YCP	CG-CB	-2.51	1.46	1.53
2	u	7	MP8	CB-CG	2.49	1.59	1.52
2	u	5	YCP	CE-N	2.46	1.53	1.47
2	y	5	YCP	CE-N	2.45	1.53	1.47
2	r	5	YCP	CG-CB	-2.45	1.47	1.53
2	t	5	YCP	CG-CB	-2.45	1.47	1.53
2	i	7	MP8	CB-CA	2.44	1.59	1.54
2	m	5	YCP	CE-N	2.41	1.53	1.47
2	v	7	MP8	CB-CA	2.39	1.59	1.54
2	4	5	YCP	CE-N	2.39	1.53	1.47
2	r	7	MP8	CB-CA	2.39	1.59	1.54
2	i	5	YCP	CG-CB	-2.38	1.47	1.53
2	j	5	YCP	CG-CB	-2.38	1.47	1.53
2	u	2	WFP	CD2-CE2	2.35	1.41	1.37
2	u	5	YCP	CG-CB	-2.35	1.47	1.53
2	h	5	YCP	CG-CB	-2.35	1.47	1.53
2	q	5	YCP	CG-CB	-2.35	1.47	1.53
2	v	5	YCP	CG-CB	-2.34	1.47	1.53
2	c	5	YCP	CG-CB	-2.34	1.47	1.53
2	k	5	YCP	CE-N	2.33	1.53	1.47
2	g	5	YCP	CG-CB	-2.33	1.47	1.53
2	2	5	YCP	CG-CB	-2.33	1.47	1.53
2	k	5	YCP	CG-CB	-2.31	1.47	1.53
2	g	5	YCP	CE-N	2.27	1.53	1.47
2	g	7	MP8	CB-CA	2.27	1.59	1.54
2	y	7	MP8	CB-CA	2.26	1.59	1.54
2	n	5	YCP	CG-CB	-2.26	1.47	1.53
2	p	7	MP8	CB-CA	2.25	1.59	1.54
2	4	5	YCP	CG-CB	-2.25	1.47	1.53
2	h	5	YCP	CE-N	2.24	1.53	1.47
2	f	7	MP8	CB-CA	2.21	1.59	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	3	7	MP8	CB-CA	2.21	1.59	1.54
2	w	5	YCP	CG-CB	-2.20	1.47	1.53
2	0	5	YCP	CE-N	2.19	1.53	1.47
2	n	5	YCP	CE-N	2.19	1.53	1.47
2	t	7	MP8	CB-CA	2.19	1.59	1.54
2	l	5	YCP	CE-N	2.19	1.53	1.47
2	z	2	WFP	CD1-CE1	2.18	1.41	1.37
2	p	5	YCP	CG-CB	-2.16	1.47	1.53
2	m	5	YCP	CG-CB	-2.16	1.47	1.53
2	e	7	MP8	CB-CA	2.16	1.59	1.54
2	s	7	MP8	CB-CA	2.14	1.59	1.54
2	e	5	YCP	CE-N	2.14	1.52	1.47
2	s	5	YCP	CE-N	2.14	1.52	1.47
2	o	5	YCP	CG-CB	-2.14	1.47	1.53
2	l	5	YCP	CG-CB	-2.13	1.47	1.53
2	3	5	YCP	CG-CB	-2.12	1.47	1.53
2	j	5	YCP	CE-N	2.12	1.52	1.47
2	f	5	YCP	CG-CB	-2.11	1.47	1.53
2	j	2	WFP	CD2-CE2	2.11	1.41	1.37
2	c	7	MP8	CB-CA	2.11	1.58	1.54
2	t	5	YCP	CE-N	2.10	1.52	1.47
2	e	5	YCP	CG-CB	-2.10	1.47	1.53
2	0	7	MP8	CB-CA	2.10	1.58	1.54
2	w	5	YCP	CE-N	2.09	1.52	1.47
2	u	7	MP8	CB-CA	2.09	1.58	1.54
2	l	2	WFP	CD2-CE2	2.08	1.41	1.37
2	h	7	MP8	CB-CA	2.07	1.58	1.54
2	w	2	WFP	CD2-CE2	2.07	1.41	1.37
2	y	5	YCP	CG-CB	-2.07	1.48	1.53
2	k	7	MP8	CB-CA	2.07	1.58	1.54
2	2	7	MP8	CB-CA	2.05	1.58	1.54
2	j	7	MP8	CB-CA	2.05	1.58	1.54
2	x	5	YCP	CG-CB	-2.04	1.48	1.53
2	1	5	YCP	CE-N	2.04	1.52	1.47
2	f	5	YCP	CE-N	2.02	1.52	1.47
2	1	7	MP8	CB-CA	2.02	1.58	1.54
2	m	7	MP8	CB-CA	2.01	1.58	1.54

All (133) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	1	5	YCP	CD-CG-CB	4.46	120.58	111.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	3	5	YCP	CG-CB-CA	3.92	116.49	110.93
2	x	5	YCP	CD-CG-CB	3.85	119.33	111.42
2	i	2	WFP	CD2-CE2-CZ	-3.68	119.03	123.50
2	n	2	WFP	CD1-CE1-CZ	-3.58	119.15	123.50
2	l	5	YCP	CG-CB-CA	3.56	115.98	110.93
2	s	2	WFP	CD2-CE2-CZ	-3.52	119.22	123.50
2	z	2	WFP	CD1-CE1-CZ	-3.38	119.39	123.50
2	m	2	WFP	CD1-CE1-CZ	-3.38	119.39	123.50
2	k	2	WFP	CD1-CE1-CZ	-3.36	119.42	123.50
2	i	2	WFP	CE2-CZ-CE1	3.33	120.93	116.08
2	2	2	WFP	CD1-CE1-CZ	-3.24	119.56	123.50
2	u	2	WFP	F2-CE2-CD2	3.22	122.86	118.28
2	n	5	YCP	O-C-CA	-3.22	116.48	124.77
2	z	2	WFP	CE2-CZ-CE1	3.16	120.68	116.08
2	p	2	WFP	CD2-CE2-CZ	-3.16	119.66	123.50
2	u	2	WFP	CD2-CE2-CZ	-3.13	119.70	123.50
2	o	5	YCP	O-C-CA	-3.07	116.87	124.77
2	s	2	WFP	CG-CD2-CE2	3.07	121.43	118.75
2	q	2	WFP	CD1-CE1-CZ	-3.06	119.78	123.50
2	0	2	WFP	CD2-CE2-CZ	-3.04	119.80	123.50
2	x	2	WFP	CD1-CE1-CZ	-3.03	119.81	123.50
2	n	2	WFP	CE2-CZ-CE1	3.03	120.49	116.08
2	i	2	WFP	CD1-CE1-CZ	-3.01	119.85	123.50
2	u	2	WFP	CE2-CZ-CE1	2.96	120.40	116.08
2	h	2	WFP	CD2-CE2-CZ	-2.95	119.91	123.50
2	r	5	YCP	O-C-CA	-2.94	117.22	124.77
2	w	2	WFP	CD2-CE2-CZ	-2.93	119.94	123.50
2	q	2	WFP	CE2-CZ-CE1	2.91	120.31	116.08
2	c	2	WFP	CD2-CE2-CZ	-2.90	119.97	123.50
2	j	5	YCP	O-C-CA	-2.90	117.32	124.77
2	l	2	WFP	CD2-CE2-CZ	-2.85	120.04	123.50
2	4	2	WFP	CD2-CE2-CZ	-2.84	120.05	123.50
2	v	2	WFP	CD2-CE2-CZ	-2.82	120.08	123.50
2	s	5	YCP	O-C-CA	-2.81	117.53	124.77
2	l	2	WFP	CD1-CE1-CZ	-2.81	120.09	123.50
2	o	2	WFP	CD1-CE1-CZ	-2.81	120.09	123.50
2	p	2	WFP	CE2-CZ-CE1	2.80	120.17	116.08
2	y	2	WFP	CE2-CZ-CE1	2.80	120.16	116.08
2	3	2	WFP	CD1-CE1-CZ	-2.80	120.10	123.50
2	0	2	WFP	CE2-CZ-CE1	2.80	120.16	116.08
2	e	2	WFP	CD1-CE1-CZ	-2.80	120.10	123.50
2	y	2	WFP	CD1-CE1-CZ	-2.78	120.12	123.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	m	2	WFP	CE2-CZ-CE1	2.77	120.12	116.08
2	3	5	YCP	CD-CG-CB	2.77	117.12	111.42
2	m	5	YCP	O-C-CA	-2.75	117.69	124.77
2	t	5	YCP	O-C-CA	-2.75	117.70	124.77
2	p	5	YCP	CG-CB-CA	2.72	114.79	110.93
2	l	2	WFP	CD2-CE2-CZ	-2.69	120.23	123.50
2	o	5	YCP	CG-CB-CA	2.69	114.75	110.93
2	y	2	WFP	CD2-CE2-CZ	-2.67	120.25	123.50
2	v	3	ALO	CG2-CB-CA	-2.66	106.11	112.16
2	q	2	WFP	CD2-CE2-CZ	-2.66	120.27	123.50
2	p	2	WFP	F2-CE2-CD2	2.64	122.04	118.28
2	x	5	YCP	CG-CB-CA	2.64	114.67	110.93
2	k	2	WFP	CE2-CZ-CE1	2.64	119.92	116.08
2	n	2	WFP	F1-CE1-CD1	2.64	122.03	118.28
2	z	2	WFP	CD2-CE2-CZ	-2.60	120.34	123.50
2	t	5	YCP	CB-CA-N	2.60	113.88	109.23
2	f	2	WFP	CD2-CE2-CZ	-2.60	120.34	123.50
2	i	5	YCP	O-C-CA	-2.58	118.13	124.77
2	2	2	WFP	CE2-CZ-CE1	2.57	119.83	116.08
2	m	2	WFP	CD2-CE2-CZ	-2.57	120.38	123.50
2	0	5	YCP	O-C-CA	-2.55	118.20	124.77
2	v	2	WFP	CD1-CE1-CZ	-2.55	120.40	123.50
2	q	5	YCP	O-C-CA	-2.55	118.21	124.77
2	v	2	WFP	CE2-CZ-CE1	2.54	119.78	116.08
2	r	2	WFP	CD2-CE2-CZ	-2.54	120.42	123.50
2	e	2	WFP	CD2-CE2-CZ	-2.52	120.44	123.50
2	g	2	WFP	CD1-CE1-CZ	-2.49	120.47	123.50
2	j	2	WFP	CD1-CE1-CZ	-2.49	120.48	123.50
2	j	2	WFP	CD2-CE2-CZ	-2.48	120.49	123.50
2	l	2	WFP	CE2-CZ-CE1	2.47	119.68	116.08
2	c	5	YCP	CG-CB-CA	2.45	114.40	110.93
2	0	2	WFP	CD1-CE1-CZ	-2.42	120.56	123.50
2	3	5	YCP	O-C-CA	-2.42	118.54	124.77
2	s	2	WFP	CE2-CZ-CE1	2.42	119.61	116.08
2	x	2	WFP	CE2-CZ-CE1	2.41	119.60	116.08
2	w	5	YCP	CD-CG-CB	2.40	116.36	111.42
2	j	2	WFP	CE2-CZ-CE1	2.40	119.58	116.08
2	t	2	WFP	CD2-CE2-CZ	-2.40	120.58	123.50
2	g	2	WFP	CD2-CE2-CZ	-2.39	120.59	123.50
2	w	5	YCP	CG-CB-CA	2.39	114.32	110.93
2	w	2	WFP	CE2-CZ-CE1	2.39	119.56	116.08
2	n	2	WFP	CD2-CE2-CZ	-2.38	120.61	123.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	4	2	WFP	CE2-CZ-CE1	2.35	119.50	116.08
2	4	2	WFP	CD1-CE1-CZ	-2.34	120.65	123.50
2	u	2	WFP	CD1-CE1-CZ	-2.34	120.66	123.50
2	s	2	WFP	CD1-CE1-CZ	-2.33	120.67	123.50
2	c	5	YCP	O-C-CA	-2.32	118.80	124.77
2	v	5	YCP	O-C-CA	-2.30	118.84	124.77
2	z	7	MP8	O-C-CA	-2.30	118.86	124.77
2	j	2	WFP	F2-CE2-CD2	2.29	121.54	118.28
2	p	5	YCP	CD-CG-CB	2.29	116.13	111.42
2	l	2	WFP	CD1-CE1-CZ	-2.29	120.72	123.50
2	w	3	ALO	OG1-CB-CA	-2.28	104.19	109.01
2	g	2	WFP	F2-CE2-CZ	2.28	121.52	118.28
2	k	5	YCP	CD-CG-CB	2.28	116.10	111.42
2	f	2	WFP	CD1-CE1-CZ	-2.26	120.75	123.50
2	p	2	WFP	F1-CE1-CD1	2.26	121.49	118.28
2	z	2	WFP	F1-CE1-CD1	2.25	121.48	118.28
2	r	2	WFP	CD1-CE1-CZ	-2.25	120.77	123.50
2	l	5	YCP	CD-CG-CB	2.25	116.03	111.42
2	w	5	YCP	O-C-CA	-2.21	119.08	124.77
2	f	2	WFP	CE2-CZ-CE1	2.21	119.30	116.08
2	l	2	WFP	CE2-CZ-CE1	2.21	119.30	116.08
2	p	2	WFP	CD1-CE1-CZ	-2.20	120.82	123.50
2	f	5	YCP	O-C-CA	-2.20	119.11	124.77
2	w	2	WFP	CD1-CE1-CZ	-2.19	120.84	123.50
2	z	5	YCP	O-C-CA	-2.19	119.14	124.77
2	h	2	WFP	CE2-CZ-CE1	2.18	119.27	116.08
2	3	2	WFP	CE2-CZ-CE1	2.18	119.26	116.08
2	e	2	WFP	CE2-CZ-CE1	2.18	119.26	116.08
2	x	5	YCP	O-C-CA	-2.18	119.17	124.77
2	y	2	WFP	F1-CE1-CD1	2.17	121.36	118.28
2	o	5	YCP	CD-CG-CB	2.16	115.87	111.42
2	k	2	WFP	CD2-CE2-CZ	-2.16	120.88	123.50
2	h	5	YCP	CG-CB-CA	2.15	113.98	110.93
2	e	2	WFP	F1-CE1-CZ	2.15	121.33	118.28
2	2	7	MP8	O-C-CA	-2.13	119.28	124.77
2	3	2	WFP	CD2-CE2-CZ	-2.12	120.92	123.50
2	n	2	WFP	F2-CE2-CD2	2.10	121.26	118.28
2	h	2	WFP	CG-CD2-CE2	2.09	120.57	118.75
2	j	7	MP8	CE-CG-CB	-2.08	108.86	114.00
2	h	7	MP8	O-C-CA	-2.08	119.43	124.77
2	o	2	WFP	F1-CE1-CD1	2.08	121.23	118.28
2	2	2	WFP	CD2-CE2-CZ	-2.07	120.98	123.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	q	7	MP8	O-C-CA	-2.06	119.46	124.77
2	s	3	ALO	OG1-CB-CA	-2.06	104.67	109.01
2	r	2	WFP	CE2-CZ-CE1	2.04	119.06	116.08
2	4	2	WFP	F2-CE2-CD2	2.02	121.15	118.28
2	g	2	WFP	CE2-CZ-CE1	2.02	119.03	116.08
2	o	2	WFP	CE2-CZ-CE1	2.02	119.03	116.08

There are no chirality outliers.

All (37) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	e	3	ALO	O-C-CA-CB
2	f	3	ALO	O-C-CA-CB
2	g	3	ALO	O-C-CA-CB
2	h	3	ALO	O-C-CA-CB
2	i	3	ALO	O-C-CA-CB
2	j	3	ALO	O-C-CA-CB
2	k	3	ALO	O-C-CA-CB
2	l	3	ALO	O-C-CA-CB
2	m	3	ALO	O-C-CA-CB
2	n	3	ALO	O-C-CA-CB
2	o	3	ALO	O-C-CA-CB
2	p	3	ALO	O-C-CA-CB
2	q	3	ALO	O-C-CA-CB
2	r	3	ALO	O-C-CA-CB
2	s	3	ALO	O-C-CA-CB
2	t	3	ALO	O-C-CA-CB
2	u	3	ALO	O-C-CA-CB
2	v	3	ALO	O-C-CA-CB
2	w	3	ALO	O-C-CA-CB
2	x	3	ALO	O-C-CA-CB
2	y	3	ALO	O-C-CA-CB
2	z	3	ALO	O-C-CA-CB
2	0	3	ALO	O-C-CA-CB
2	1	3	ALO	O-C-CA-CB
2	2	3	ALO	O-C-CA-CB
2	3	3	ALO	O-C-CA-CB
2	4	3	ALO	O-C-CA-CB
2	c	5	YCP	O-C-CA-CB
2	m	5	YCP	O-C-CA-CB
2	p	5	YCP	O-C-CA-CB
2	u	5	YCP	O-C-CA-CB

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Mol	Chain	Res	Type	Atoms
2	y	5	YCP	O-C-CA-CB
2	z	5	YCP	O-C-CA-CB
2	j	5	YCP	O-C-CA-CB
2	n	5	YCP	O-C-CA-CB
2	v	5	YCP	O-C-CA-CB
2	c	3	ALO	O-C-CA-CB

All (23) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	t	5	YCP	CA-CB-CD-CE-CG-N
2	o	5	YCP	CA-CB-CD-CE-CG-N
2	z	5	YCP	CA-CB-CD-CE-CG-N
2	p	5	YCP	CA-CB-CD-CE-CG-N
2	k	5	YCP	CA-CB-CD-CE-CG-N
2	c	5	YCP	CA-CB-CD-CE-CG-N
2	h	5	YCP	CA-CB-CD-CE-CG-N
2	v	5	YCP	CA-CB-CD-CE-CG-N
2	j	5	YCP	CA-CB-CD-CE-CG-N
2	4	5	YCP	CA-CB-CD-CE-CG-N
2	n	5	YCP	CA-CB-CD-CE-CG-N
2	0	5	YCP	CA-CB-CD-CE-CG-N
2	u	5	YCP	CA-CB-CD-CE-CG-N
2	i	5	YCP	CA-CB-CD-CE-CG-N
2	f	5	YCP	CA-CB-CD-CE-CG-N
2	g	5	YCP	CA-CB-CD-CE-CG-N
2	l	5	YCP	CA-CB-CD-CE-CG-N
2	m	5	YCP	CA-CB-CD-CE-CG-N
2	e	5	YCP	CA-CB-CD-CE-CG-N
2	q	5	YCP	CA-CB-CD-CE-CG-N
2	r	5	YCP	CA-CB-CD-CE-CG-N
2	y	5	YCP	CA-CB-CD-CE-CG-N
2	s	5	YCP	CA-CB-CD-CE-CG-N

33 monomers are involved in 40 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	i	7	MP8	1	0
2	o	5	YCP	1	0
2	l	2	WFP	1	0
2	c	7	MP8	1	0
2	4	2	WFP	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	3	7	MP8	1	0
2	o	7	MP8	1	0
2	n	2	WFP	1	0
2	f	3	ALO	1	0
2	l	7	MP8	2	0
2	t	7	MP8	1	0
2	r	7	MP8	2	0
2	s	5	YCP	1	0
2	f	7	MP8	2	0
2	y	3	ALO	1	0
2	p	7	MP8	1	0
2	s	7	MP8	1	0
2	f	2	WFP	1	0
2	l	7	MP8	1	0
2	g	7	MP8	1	0
2	3	3	ALO	1	0
2	n	7	MP8	1	0
2	k	7	MP8	2	0
2	4	3	ALO	2	0
2	2	7	MP8	2	0
2	q	7	MP8	1	0
2	h	3	ALO	1	0
2	u	7	MP8	1	0
2	e	3	ALO	1	0
2	p	2	WFP	1	0
2	w	7	MP8	2	0
2	y	7	MP8	1	0
2	2	3	ALO	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

28 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	OCA	g	101	2	7,8,9	1.90	2 (28%)	6,7,9	0.79	0
3	OCA	2	101	2	7,8,9	1.90	2 (28%)	6,7,9	0.82	0
3	OCA	p	101	2	7,8,9	1.91	2 (28%)	6,7,9	0.86	0
3	OCA	i	101	2	7,8,9	1.85	2 (28%)	6,7,9	0.82	0
3	OCA	u	101	2	7,8,9	1.97	2 (28%)	6,7,9	0.85	0
3	OCA	4	101	2	7,8,9	1.83	2 (28%)	6,7,9	0.97	0
3	OCA	s	101	2	7,8,9	2.05	2 (28%)	6,7,9	0.73	0
3	OCA	x	101	2	7,8,9	1.81	2 (28%)	6,7,9	1.02	0
3	OCA	q	101	2	7,8,9	1.89	2 (28%)	6,7,9	0.87	0
3	OCA	t	101	2	7,8,9	1.92	2 (28%)	6,7,9	0.98	0
3	OCA	k	101	2	7,8,9	1.96	2 (28%)	6,7,9	0.81	0
3	OCA	l	101	2	7,8,9	1.98	2 (28%)	6,7,9	0.75	0
3	OCA	0	101	2	7,8,9	1.85	2 (28%)	6,7,9	0.97	0
3	OCA	r	101	2	7,8,9	1.89	2 (28%)	6,7,9	0.87	0
3	OCA	n	101	2	7,8,9	1.84	2 (28%)	6,7,9	0.87	0
3	OCA	w	101	2	7,8,9	1.87	2 (28%)	6,7,9	0.99	0
3	OCA	v	101	2	7,8,9	1.85	2 (28%)	6,7,9	0.82	0
3	OCA	m	101	2	7,8,9	1.91	2 (28%)	6,7,9	0.86	0
3	OCA	c	101	2	7,8,9	1.91	2 (28%)	6,7,9	0.79	0
3	OCA	z	101	2	7,8,9	1.87	2 (28%)	6,7,9	0.87	0
3	OCA	1	101	2	7,8,9	1.85	2 (28%)	6,7,9	0.98	0
3	OCA	y	101	2	7,8,9	1.90	2 (28%)	6,7,9	0.78	0
3	OCA	j	101	2	7,8,9	1.82	2 (28%)	6,7,9	0.90	0
3	OCA	3	101	2	7,8,9	1.91	2 (28%)	6,7,9	0.75	0
3	OCA	o	101	2	7,8,9	1.86	2 (28%)	6,7,9	0.95	0
3	OCA	f	101	2	7,8,9	1.84	2 (28%)	6,7,9	0.85	0
3	OCA	h	101	2	7,8,9	1.93	2 (28%)	6,7,9	0.82	0
3	OCA	e	101	2	7,8,9	1.87	2 (28%)	6,7,9	0.83	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
3	OCA	g	101	2	-	2/6/6/7	-
3	OCA	2	101	2	-	2/6/6/7	-
3	OCA	p	101	2	-	1/6/6/7	-
3	OCA	i	101	2	-	1/6/6/7	-
3	OCA	u	101	2	-	1/6/6/7	-
3	OCA	4	101	2	-	2/6/6/7	-
3	OCA	s	101	2	-	2/6/6/7	-
3	OCA	x	101	2	-	1/6/6/7	-
3	OCA	q	101	2	-	2/6/6/7	-
3	OCA	t	101	2	-	2/6/6/7	-
3	OCA	k	101	2	-	1/6/6/7	-
3	OCA	l	101	2	-	2/6/6/7	-
3	OCA	0	101	2	-	1/6/6/7	-
3	OCA	r	101	2	-	2/6/6/7	-
3	OCA	n	101	2	-	2/6/6/7	-
3	OCA	w	101	2	-	3/6/6/7	-
3	OCA	v	101	2	-	1/6/6/7	-
3	OCA	m	101	2	-	1/6/6/7	-
3	OCA	c	101	2	-	1/6/6/7	-
3	OCA	z	101	2	-	1/6/6/7	-
3	OCA	1	101	2	-	0/6/6/7	-
3	OCA	y	101	2	-	1/6/6/7	-
3	OCA	j	101	2	-	2/6/6/7	-
3	OCA	3	101	2	-	2/6/6/7	-
3	OCA	o	101	2	-	3/6/6/7	-
3	OCA	f	101	2	-	2/6/6/7	-
3	OCA	h	101	2	-	2/6/6/7	-
3	OCA	e	101	2	-	2/6/6/7	-

All (56) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	s	101	OCA	C3-C2	-3.94	1.35	1.52
3	u	101	OCA	C3-C2	-3.75	1.36	1.52
3	l	101	OCA	C3-C2	-3.74	1.36	1.52
3	k	101	OCA	C5-C4	-3.68	1.33	1.51
3	h	101	OCA	C5-C4	-3.66	1.33	1.51
3	p	101	OCA	C5-C4	-3.65	1.33	1.51
3	s	101	OCA	C5-C4	-3.65	1.33	1.51
3	t	101	OCA	C5-C4	-3.63	1.33	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	l	101	OCA	C5-C4	-3.62	1.33	1.51
3	k	101	OCA	C3-C2	-3.62	1.37	1.52
3	3	101	OCA	C3-C2	-3.60	1.37	1.52
3	m	101	OCA	C3-C2	-3.60	1.37	1.52
3	o	101	OCA	C3-C2	-3.59	1.37	1.52
3	c	101	OCA	C5-C4	-3.59	1.33	1.51
3	r	101	OCA	C5-C4	-3.59	1.33	1.51
3	2	101	OCA	C5-C4	-3.56	1.34	1.51
3	g	101	OCA	C3-C2	-3.56	1.37	1.52
3	q	101	OCA	C3-C2	-3.55	1.37	1.52
3	y	101	OCA	C5-C4	-3.55	1.34	1.51
3	1	101	OCA	C3-C2	-3.55	1.37	1.52
3	u	101	OCA	C5-C4	-3.55	1.34	1.51
3	y	101	OCA	C3-C2	-3.52	1.37	1.52
3	w	101	OCA	C5-C4	-3.52	1.34	1.51
3	t	101	OCA	C3-C2	-3.52	1.37	1.52
3	c	101	OCA	C3-C2	-3.51	1.37	1.52
3	f	101	OCA	C5-C4	-3.50	1.34	1.51
3	h	101	OCA	C3-C2	-3.50	1.37	1.52
3	q	101	OCA	C5-C4	-3.49	1.34	1.51
3	m	101	OCA	C5-C4	-3.49	1.34	1.51
3	g	101	OCA	C5-C4	-3.48	1.34	1.51
3	v	101	OCA	C3-C2	-3.48	1.37	1.52
3	0	101	OCA	C3-C2	-3.47	1.37	1.52
3	e	101	OCA	C5-C4	-3.47	1.34	1.51
3	i	101	OCA	C3-C2	-3.46	1.37	1.52
3	3	101	OCA	C5-C4	-3.46	1.34	1.51
3	z	101	OCA	C5-C4	-3.46	1.34	1.51
3	2	101	OCA	C3-C2	-3.46	1.37	1.52
3	e	101	OCA	C3-C2	-3.44	1.37	1.52
3	r	101	OCA	C3-C2	-3.43	1.37	1.52
3	z	101	OCA	C3-C2	-3.43	1.37	1.52
3	j	101	OCA	C5-C4	-3.43	1.34	1.51
3	n	101	OCA	C5-C4	-3.42	1.34	1.51
3	w	101	OCA	C3-C2	-3.41	1.37	1.52
3	n	101	OCA	C3-C2	-3.40	1.37	1.52
3	p	101	OCA	C3-C2	-3.40	1.37	1.52
3	4	101	OCA	C3-C2	-3.40	1.38	1.52
3	x	101	OCA	C5-C4	-3.38	1.35	1.51
3	i	101	OCA	C5-C4	-3.37	1.35	1.51
3	0	101	OCA	C5-C4	-3.36	1.35	1.51
3	4	101	OCA	C5-C4	-3.35	1.35	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	v	101	OCA	C5-C4	-3.32	1.35	1.51
3	f	101	OCA	C3-C2	-3.30	1.38	1.52
3	o	101	OCA	C5-C4	-3.30	1.35	1.51
3	1	101	OCA	C5-C4	-3.29	1.35	1.51
3	j	101	OCA	C3-C2	-3.29	1.38	1.52
3	x	101	OCA	C3-C2	-3.28	1.38	1.52

There are no bond angle outliers.

There are no chirality outliers.

All (45) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	o	101	OCA	O1-C1-C2-C3
3	x	101	OCA	C4-C5-C6-C7
3	i	101	OCA	C4-C5-C6-C7
3	u	101	OCA	C4-C5-C6-C7
3	s	101	OCA	C4-C5-C6-C7
3	f	101	OCA	C4-C5-C6-C7
3	t	101	OCA	C4-C5-C6-C7
3	g	101	OCA	C4-C5-C6-C7
3	w	101	OCA	C4-C5-C6-C7
3	n	101	OCA	C4-C5-C6-C7
3	0	101	OCA	C4-C5-C6-C7
3	z	101	OCA	C4-C5-C6-C7
3	c	101	OCA	C4-C5-C6-C7
3	4	101	OCA	C4-C5-C6-C7
3	3	101	OCA	C4-C5-C6-C7
3	o	101	OCA	C2-C3-C4-C5
3	o	101	OCA	C4-C5-C6-C7
3	h	101	OCA	C4-C5-C6-C7
3	p	101	OCA	C2-C3-C4-C5
3	q	101	OCA	C2-C3-C4-C5
3	y	101	OCA	C4-C5-C6-C7
3	e	101	OCA	C4-C5-C6-C7
3	2	101	OCA	C4-C5-C6-C7
3	m	101	OCA	C4-C5-C6-C7
3	v	101	OCA	C5-C6-C7-C8
3	s	101	OCA	C5-C6-C7-C8
3	j	101	OCA	C5-C6-C7-C8
3	l	101	OCA	C5-C6-C7-C8
3	e	101	OCA	C5-C6-C7-C8

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Mol	Chain	Res	Type	Atoms
3	2	101	OCA	C5-C6-C7-C8
3	k	101	OCA	C4-C5-C6-C7
3	j	101	OCA	C4-C5-C6-C7
3	f	101	OCA	C5-C6-C7-C8
3	w	101	OCA	C2-C3-C4-C5
3	3	101	OCA	C5-C6-C7-C8
3	r	101	OCA	C4-C5-C6-C7
3	r	101	OCA	C2-C3-C4-C5
3	g	101	OCA	C5-C6-C7-C8
3	n	101	OCA	C5-C6-C7-C8
3	w	101	OCA	C5-C6-C7-C8
3	h	101	OCA	C5-C6-C7-C8
3	l	101	OCA	C4-C5-C6-C7
3	q	101	OCA	C5-C6-C7-C8
3	t	101	OCA	C2-C3-C4-C5
3	4	101	OCA	C2-C3-C4-C5

There are no ring outliers.

16 monomers are involved in 24 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	2	101	OCA	1	0
3	p	101	OCA	2	0
3	4	101	OCA	2	0
3	s	101	OCA	1	0
3	x	101	OCA	1	0
3	q	101	OCA	1	0
3	t	101	OCA	2	0
3	k	101	OCA	2	0
3	l	101	OCA	1	0
3	r	101	OCA	1	0
3	y	101	OCA	1	0
3	j	101	OCA	1	0
3	3	101	OCA	3	0
3	f	101	OCA	1	0
3	h	101	OCA	2	0
3	e	101	OCA	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	175/217 (80%)	-0.12	2 (1%) 78 76	29, 39, 60, 74	0
1	B	173/217 (79%)	-0.19	3 (1%) 69 67	30, 38, 60, 72	0
1	C	173/217 (79%)	-0.14	2 (1%) 76 75	32, 42, 60, 80	0
1	D	173/217 (79%)	-0.11	1 (0%) 85 85	32, 43, 63, 76	0
1	E	174/217 (80%)	-0.13	3 (1%) 69 67	32, 41, 59, 79	0
1	F	171/217 (78%)	-0.26	2 (1%) 76 75	31, 37, 53, 76	0
1	G	173/217 (79%)	-0.18	0 100 100	30, 39, 57, 69	0
1	H	174/217 (80%)	-0.07	3 (1%) 69 67	32, 42, 59, 70	0
1	I	171/217 (78%)	-0.00	2 (1%) 76 75	34, 43, 62, 76	0
1	J	172/217 (79%)	-0.11	1 (0%) 85 85	33, 42, 60, 76	0
1	K	174/217 (80%)	-0.06	4 (2%) 61 58	32, 39, 61, 72	0
1	L	174/217 (80%)	-0.28	2 (1%) 78 76	29, 37, 57, 79	0
1	M	171/217 (78%)	-0.28	2 (1%) 76 75	27, 37, 56, 76	0
1	N	173/217 (79%)	-0.31	3 (1%) 69 67	31, 38, 55, 68	0
1	O	172/217 (79%)	-0.13	1 (0%) 85 85	30, 41, 64, 90	0
1	P	173/217 (79%)	-0.25	1 (0%) 85 85	27, 37, 59, 75	0
1	Q	173/217 (79%)	-0.26	2 (1%) 76 75	30, 39, 58, 70	0
1	R	175/217 (80%)	-0.18	1 (0%) 85 85	32, 40, 59, 69	0
1	S	173/217 (79%)	-0.01	3 (1%) 69 67	32, 42, 65, 77	0
1	T	171/217 (78%)	-0.03	3 (1%) 67 65	35, 46, 61, 72	0
1	U	174/217 (80%)	-0.07	4 (2%) 61 58	30, 42, 63, 79	0
1	V	171/217 (78%)	-0.32	2 (1%) 76 75	28, 36, 53, 69	0
1	W	173/217 (79%)	-0.17	3 (1%) 69 67	30, 38, 56, 69	0
1	X	174/217 (80%)	-0.12	3 (1%) 69 67	31, 40, 59, 76	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	Y	172/217 (79%)	-0.18	2 (1%) 76 75	28, 39, 56, 80	0
1	Z	172/217 (79%)	-0.29	2 (1%) 76 75	28, 37, 56, 66	0
1	a	173/217 (79%)	-0.29	2 (1%) 76 75	28, 38, 55, 72	0
1	b	172/217 (79%)	-0.33	0 100 100	30, 36, 49, 70	0
2	0	2/6 (33%)	-0.38	0 100 100	41, 41, 41, 45	0
2	1	2/6 (33%)	0.03	0 100 100	40, 40, 40, 42	0
2	2	2/6 (33%)	0.11	0 100 100	41, 41, 41, 45	0
2	3	2/6 (33%)	0.10	0 100 100	37, 37, 37, 41	0
2	4	2/6 (33%)	-0.28	0 100 100	39, 39, 39, 41	0
2	c	2/6 (33%)	0.44	0 100 100	45, 45, 45, 50	0
2	e	2/6 (33%)	0.03	0 100 100	45, 45, 45, 48	0
2	f	2/6 (33%)	0.34	0 100 100	51, 51, 51, 52	0
2	g	2/6 (33%)	-0.09	0 100 100	47, 47, 47, 47	0
2	h	2/6 (33%)	-0.43	0 100 100	44, 44, 44, 47	0
2	i	2/6 (33%)	-0.32	0 100 100	42, 42, 42, 42	0
2	j	2/6 (33%)	-0.01	0 100 100	46, 46, 46, 49	0
2	k	2/6 (33%)	-0.20	0 100 100	42, 42, 42, 43	0
2	l	2/6 (33%)	0.13	0 100 100	44, 44, 44, 49	0
2	m	2/6 (33%)	-0.25	0 100 100	44, 44, 44, 44	0
2	n	2/6 (33%)	0.41	0 100 100	46, 46, 46, 56	0
2	o	2/6 (33%)	-0.11	0 100 100	39, 39, 39, 44	0
2	p	2/6 (33%)	0.25	0 100 100	46, 46, 46, 48	0
2	q	2/6 (33%)	-0.07	0 100 100	43, 43, 43, 44	0
2	r	2/6 (33%)	-0.06	0 100 100	44, 44, 44, 49	0
2	s	2/6 (33%)	-0.21	0 100 100	46, 46, 46, 51	0
2	t	2/6 (33%)	-0.51	0 100 100	40, 40, 40, 42	0
2	u	2/6 (33%)	-0.19	0 100 100	43, 43, 43, 48	0
2	v	2/6 (33%)	0.32	0 100 100	45, 45, 45, 48	0
2	w	2/6 (33%)	0.49	0 100 100	51, 51, 51, 56	0
2	x	2/6 (33%)	-0.13	0 100 100	47, 47, 47, 51	0
2	y	2/6 (33%)	0.00	0 100 100	43, 43, 43, 43	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
2	z	2/6 (33%)	0.06	0 100 100	45, 45, 45, 46	0
All	All	4895/6244 (78%)	-0.17	59 (1%) 76 75	27, 40, 59, 90	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	W	136	GLY	3.9
1	A	197	GLU	3.6
1	S	133	GLY	3.3
1	X	133	GLY	3.3
1	X	155	LYS	3.2
1	C	133	GLY	3.1
1	T	132	SER	3.1
1	V	197	GLU	3.0
1	J	198	ASN	3.0
1	O	24	ILE	3.0
1	E	136	GLY	2.9
1	W	133	GLY	2.8
1	V	198	ASN	2.8
1	K	136	GLY	2.8
1	D	198	ASN	2.8
1	Z	132	SER	2.7
1	A	147	ARG	2.7
1	L	133	GLY	2.6
1	P	136	GLY	2.6
1	Z	184	GLU	2.6
1	T	24	ILE	2.5
1	S	147	ARG	2.5
1	B	136	GLY	2.4
1	F	23	ASP	2.4
1	K	194	GLN	2.4
1	K	198	ASN	2.4
1	H	136	GLY	2.4
1	U	136	GLY	2.4
1	E	184	GLU	2.3
1	S	43	GLU	2.3
1	M	132	SER	2.3
1	Y	23	ASP	2.3
1	H	198	ASN	2.3
1	F	165	ASP	2.3
1	a	23	ASP	2.3
1	B	30	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	a	132	SER	2.2
1	C	172	GLU	2.2
1	X	135	LEU	2.2
1	N	133	GLY	2.2
1	I	132	SER	2.2
1	U	184	GLU	2.2
1	M	194	GLN	2.2
1	L	43	GLU	2.1
1	B	46	ASN	2.1
1	K	132	SER	2.1
1	W	30	LYS	2.1
1	Q	130	LEU	2.1
1	N	136	GLY	2.1
1	U	23	ASP	2.1
1	Y	26	SER	2.1
1	R	130	LEU	2.1
1	T	191	LEU	2.1
1	Q	24	ILE	2.1
1	H	135	LEU	2.0
1	I	54	PHE	2.0
1	E	131	ILE	2.0
1	U	24	ILE	2.0
1	N	27	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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	ALO	u	3	7/8	0.86	0.11	44,46,50,56	0
2	ALO	x	3	7/8	0.87	0.12	42,51,54,56	0
2	ALO	4	3	7/8	0.87	0.12	35,37,42,43	0
2	ALO	3	3	7/8	0.89	0.11	36,42,44,48	0
2	ALO	n	3	7/8	0.90	0.09	41,52,53,53	0
2	ALO	s	3	7/8	0.90	0.09	45,50,53,57	0
2	ALO	c	3	7/8	0.90	0.10	39,45,47,51	0
2	YCP	j	5	8/9	0.90	0.13	47,48,53,53	0
2	YCP	x	5	8/9	0.90	0.12	47,50,51,52	0
2	ALO	y	3	7/8	0.91	0.10	37,43,46,46	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	YCP	c	5	8/9	0.91	0.12	39,46,49,52	0
2	YCP	g	5	8/9	0.91	0.12	41,45,50,51	0
2	ALO	l	3	7/8	0.91	0.12	39,41,44,46	0
2	ALO	q	3	7/8	0.91	0.10	31,42,44,44	0
2	YCP	y	5	8/9	0.91	0.12	41,45,46,48	0
2	YCP	3	5	8/9	0.91	0.11	24,35,39,42	0
2	MP8	k	7	8/9	0.91	0.13	39,43,48,51	0
2	MP8	x	7	8/9	0.91	0.12	46,49,52,54	0
2	ALO	w	3	7/8	0.92	0.09	46,53,56,56	0
2	YCP	i	5	8/9	0.92	0.12	40,41,47,48	0
2	ALO	j	3	7/8	0.92	0.10	40,44,49,54	0
2	ALO	k	3	7/8	0.92	0.09	38,42,44,47	0
2	WFP	c	2	13/14	0.92	0.11	38,46,50,50	0
2	YCP	l	5	8/9	0.92	0.11	39,43,47,48	0
2	ALO	f	3	7/8	0.92	0.11	46,46,50,53	0
2	ALO	g	3	7/8	0.92	0.10	37,44,49,52	0
2	MP8	s	7	8/9	0.92	0.09	45,50,51,51	0
2	ALO	i	3	7/8	0.92	0.11	41,44,46,46	0
2	YCP	r	5	8/9	0.93	0.11	46,49,54,54	0
2	YCP	u	5	8/9	0.93	0.12	42,46,48,50	0
2	ALO	0	3	7/8	0.93	0.08	38,41,44,45	0
2	YCP	e	5	8/9	0.93	0.10	45,49,51,53	0
2	ALO	r	3	7/8	0.93	0.08	39,45,51,51	0
2	ALO	p	3	7/8	0.93	0.09	43,45,50,53	0
2	MP8	f	7	8/9	0.93	0.09	43,46,48,49	0
2	ALO	e	3	7/8	0.93	0.10	46,47,48,51	0
2	MP8	l	7	8/9	0.93	0.10	34,42,43,49	0
2	MP8	r	7	8/9	0.93	0.10	41,44,49,49	0
2	YCP	k	5	8/9	0.93	0.10	38,43,45,45	0
2	YCP	o	5	8/9	0.93	0.10	39,40,43,43	0
2	ALO	z	3	7/8	0.94	0.09	37,43,46,52	0
2	YCP	p	5	8/9	0.94	0.09	41,44,47,47	0
2	ALO	o	3	7/8	0.94	0.09	41,42,45,52	0
2	YCP	t	5	8/9	0.94	0.10	36,40,41,42	0
2	WFP	f	2	13/14	0.94	0.09	44,48,50,52	0
2	YCP	v	5	8/9	0.94	0.08	40,43,44,47	0
2	WFP	g	2	13/14	0.94	0.08	36,41,44,45	0
2	WFP	h	2	13/14	0.94	0.08	41,45,47,48	0
2	YCP	z	5	8/9	0.94	0.12	39,44,47,50	0
2	YCP	0	5	8/9	0.94	0.11	37,43,45,47	0
2	WFP	n	2	13/14	0.94	0.10	43,46,51,55	0
2	YCP	2	5	8/9	0.94	0.09	36,38,42,42	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	WFP	t	2	13/14	0.94	0.09	37,41,47,49	0
2	YCP	4	5	8/9	0.94	0.10	37,42,43,44	0
2	MP8	c	7	8/9	0.94	0.10	39,43,47,48	0
2	ALO	v	3	7/8	0.94	0.09	37,44,47,48	0
2	WFP	x	2	13/14	0.94	0.08	45,48,53,55	0
2	WFP	0	2	13/14	0.94	0.08	35,43,49,49	0
2	WFP	1	2	13/14	0.94	0.09	36,40,43,43	0
2	YCP	l	5	8/9	0.94	0.09	34,44,48,49	0
2	MP8	w	7	8/9	0.94	0.10	46,50,53,56	0
2	YCP	n	5	8/9	0.94	0.09	46,51,55,57	0
2	MP8	z	7	8/9	0.94	0.11	39,40,44,44	0
2	WFP	k	2	13/14	0.95	0.08	36,41,45,48	0
2	WFP	l	2	13/14	0.95	0.08	39,44,48,51	0
2	WFP	m	2	13/14	0.95	0.07	45,47,50,50	0
2	WFP	2	2	13/14	0.95	0.09	34,39,45,46	0
2	YCP	f	5	8/9	0.95	0.09	43,50,53,53	0
2	WFP	3	2	13/14	0.95	0.07	34,37,41,43	0
2	YCP	h	5	8/9	0.95	0.10	37,45,48,50	0
2	WFP	e	2	13/14	0.95	0.09	41,43,47,49	0
2	WFP	o	2	13/14	0.95	0.08	34,40,44,45	0
2	WFP	p	2	13/14	0.95	0.08	35,39,47,47	0
2	WFP	r	2	13/14	0.95	0.08	35,40,47,49	0
2	MP8	e	7	8/9	0.95	0.09	46,47,48,48	0
2	YCP	m	5	8/9	0.95	0.08	38,44,46,47	0
2	WFP	i	2	13/14	0.95	0.09	42,43,46,49	0
2	WFP	u	2	13/14	0.95	0.09	38,43,47,48	0
2	MP8	p	7	8/9	0.95	0.12	39,44,47,50	0
2	WFP	w	2	13/14	0.95	0.09	45,50,53,53	0
2	YCP	q	5	8/9	0.95	0.09	39,43,44,46	0
2	MP8	t	7	8/9	0.95	0.08	35,37,43,45	0
2	MP8	v	7	8/9	0.95	0.11	37,40,42,44	0
2	ALO	l	3	7/8	0.95	0.07	41,43,47,53	0
2	YCP	s	5	8/9	0.95	0.09	44,48,53,53	0
2	MP8	y	7	8/9	0.95	0.09	37,39,42,43	0
2	ALO	m	3	7/8	0.95	0.07	39,48,50,55	0
2	MP8	1	7	8/9	0.95	0.08	39,41,43,44	0
2	MP8	3	7	8/9	0.95	0.09	31,37,40,41	0
2	MP8	n	7	8/9	0.96	0.08	37,46,47,47	0
2	MP8	o	7	8/9	0.96	0.10	38,43,45,47	0
2	ALO	t	3	7/8	0.96	0.07	34,38,40,42	0
2	WFP	s	2	13/14	0.96	0.08	39,47,52,55	0
2	YCP	w	5	8/9	0.96	0.09	46,52,54,55	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	WFP	y	2	13/14	0.96	0.07	33,37,43,43	0
2	MP8	u	7	8/9	0.96	0.09	41,43,44,44	0
2	WFP	4	2	13/14	0.96	0.08	33,38,43,47	0
2	WFP	z	2	13/14	0.96	0.08	39,42,45,46	0
2	MP8	g	7	8/9	0.96	0.08	35,41,45,45	0
2	MP8	h	7	8/9	0.96	0.09	37,39,42,43	0
2	MP8	j	7	8/9	0.96	0.09	38,43,47,47	0
2	WFP	v	2	13/14	0.96	0.08	35,40,44,48	0
2	MP8	2	7	8/9	0.96	0.10	36,39,43,43	0
2	WFP	j	2	13/14	0.96	0.07	41,44,47,48	0
2	ALO	2	3	7/8	0.97	0.07	39,41,44,55	0
2	WFP	q	2	13/14	0.97	0.07	39,43,47,51	0
2	MP8	q	7	8/9	0.97	0.08	36,39,42,42	0
2	ALO	h	3	7/8	0.97	0.07	41,44,48,52	0
2	MP8	i	7	8/9	0.97	0.07	36,42,44,44	0
2	MP8	4	7	8/9	0.97	0.08	28,36,38,41	0
2	MP8	m	7	8/9	0.98	0.05	38,42,45,46	0
2	MP8	0	7	8/9	0.98	0.05	36,40,41,44	0

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($\text{\AA}^2$)	Q<0.9
3	OCA	c	101	9/10	0.92	0.12	35,40,46,47	0
3	OCA	i	101	9/10	0.92	0.15	43,44,58,61	0
3	OCA	m	101	9/10	0.92	0.15	42,46,56,59	0
3	OCA	o	101	9/10	0.92	0.12	38,40,49,52	0
3	OCA	w	101	9/10	0.92	0.12	41,48,52,56	0
3	OCA	t	101	9/10	0.93	0.11	38,41,49,54	0
3	OCA	u	101	9/10	0.93	0.13	40,43,51,57	0
3	OCA	j	101	9/10	0.93	0.13	43,45,50,52	0
3	OCA	g	101	9/10	0.94	0.12	38,44,53,58	0
3	OCA	h	101	9/10	0.94	0.11	39,48,53,56	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	OCA	r	101	9/10	0.94	0.15	39,45,53,69	0
3	OCA	2	101	9/10	0.94	0.11	38,43,54,57	0
3	OCA	3	101	9/10	0.94	0.10	40,41,43,49	0
3	OCA	e	101	9/10	0.95	0.11	40,49,55,58	0
3	OCA	p	101	9/10	0.95	0.08	38,45,51,51	0
3	OCA	v	101	9/10	0.95	0.09	38,43,48,49	0
3	OCA	q	101	9/10	0.95	0.13	37,41,50,59	0
3	OCA	x	101	9/10	0.95	0.10	45,46,53,53	0
3	OCA	z	101	9/10	0.95	0.11	37,43,54,56	0
3	OCA	n	101	9/10	0.95	0.10	44,48,55,57	0
3	OCA	s	101	9/10	0.95	0.11	38,46,53,56	0
3	OCA	f	101	9/10	0.96	0.11	43,49,63,66	0
3	OCA	k	101	9/10	0.96	0.11	31,37,52,53	0
3	OCA	l	101	9/10	0.96	0.09	31,40,50,50	0
3	OCA	4	101	9/10	0.96	0.08	31,35,41,44	0
3	OCA	y	101	9/10	0.97	0.08	40,42,45,49	0
3	OCA	0	101	9/10	0.97	0.07	30,34,44,46	0
3	OCA	1	101	9/10	0.97	0.08	37,41,44,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.