



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 30, 2026 – 01:06 PM UTC

PDB ID : 3H6I / pdb_00003h6i
Title : Crystal Structure of Mycobacterium Tuberculosis Proteasome Modified by inhibitor GL1
Authors : Li, D.; Li, H.; Lin, G.
Deposited on : 2009-04-23
Resolution : 2.43 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:

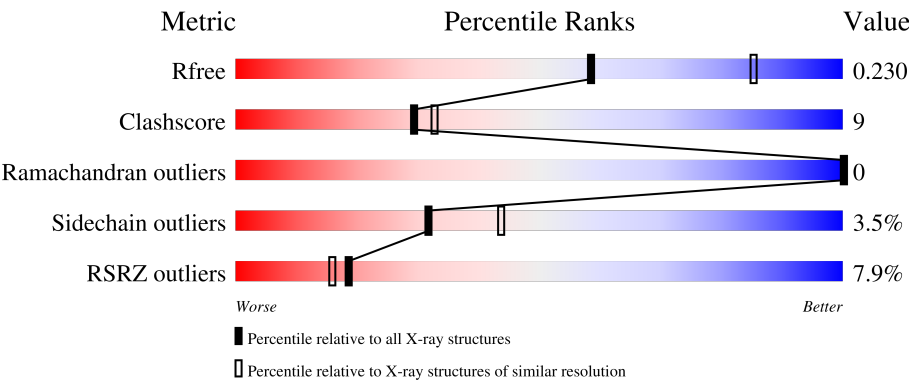
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.43 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	2340 (2.46-2.42)
Clashscore	190562	2400 (2.46-2.42)
Ramachandran outliers	187476	2379 (2.46-2.42)
Sidechain outliers	187428	2379 (2.46-2.42)
RSRZ outliers	180081	2340 (2.46-2.42)

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	1	248	10% 73% 12% • 13%
1	A	248	11% 74% 11% • 13%
1	B	248	15% 72% 13% • 14%
1	D	248	15% 70% 13% • 14%
1	F	248	15% 69% 15% • 12%

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Mol	Chain	Length	Quality of chain
1	I	248	
1	K	248	
1	M	248	
1	O	248	
1	Q	248	
1	S	248	
1	U	248	
1	W	248	
1	Y	248	
2	2	240	
2	C	240	
2	E	240	
2	G	240	
2	H	240	
2	J	240	
2	L	240	
2	N	240	
2	P	240	
2	R	240	
2	T	240	
2	V	240	
2	X	240	
2	Z	240	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	OZT	P	301	-	-	X	-
3	DMF	C	5	-	-	X	-
3	DMF	C	64	-	-	X	-
3	DMF	C	73	-	-	X	-
3	DMF	G	96	-	-	X	-
3	DMF	H	19	-	-	X	-
3	DMF	J	6	-	-	X	-
3	DMF	O	250	-	-	X	-
3	DMF	T	35	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 48540 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 Proteasome (Alpha subunit) PrcA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	S	0	0	0
			1662	1041	304	314	3			
1	B	214	Total	C	N	O	S	0	0	0
			1650	1033	302	312	3			
1	D	213	Total	C	N	O	S	0	0	0
			1644	1030	301	310	3			
1	F	217	Total	C	N	O	S	0	0	0
			1669	1045	305	316	3			
1	I	215	Total	C	N	O	S	0	0	0
			1656	1038	303	312	3			
1	K	217	Total	C	N	O	S	0	0	0
			1670	1047	305	315	3			
1	M	216	Total	C	N	O	S	0	0	0
			1662	1041	304	314	3			
1	O	215	Total	C	N	O	S	0	0	0
			1654	1035	303	313	3			
1	Q	218	Total	C	N	O	S	0	0	0
			1677	1051	306	317	3			
1	S	215	Total	C	N	O	S	0	0	0
			1654	1035	303	313	3			
1	U	215	Total	C	N	O	S	0	0	0
			1654	1035	303	313	3			
1	W	217	Total	C	N	O	S	0	0	0
			1670	1047	305	315	3			
1	Y	213	Total	C	N	O	S	0	0	0
			1644	1030	301	310	3			
1	1	215	Total	C	N	O	S	0	0	0
			1656	1038	303	312	3			

- Molecule 2 is a protein called Proteasome (Beta subunit) PrcB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	215	Total 1593	C 998	N 274	O 317	S 4	0	0	0
2	E	216	Total 1601	C 1004	N 275	O 318	S 4	0	0	0
2	G	216	Total 1601	C 1004	N 275	O 318	S 4	0	0	0
2	H	213	Total 1583	C 992	N 272	O 315	S 4	0	0	0
2	J	216	Total 1601	C 1004	N 275	O 318	S 4	0	0	0
2	L	216	Total 1601	C 1004	N 275	O 318	S 4	0	0	0
2	N	213	Total 1580	C 989	N 272	O 315	S 4	0	0	0
2	P	216	Total 1601	C 1004	N 275	O 318	S 4	0	0	0
2	R	223	Total 1646	C 1031	N 282	O 329	S 4	0	0	0
2	T	216	Total 1601	C 1004	N 275	O 318	S 4	0	0	0
2	V	221	Total 1632	C 1022	N 280	O 326	S 4	0	0	0
2	X	216	Total 1601	C 1004	N 275	O 318	S 4	0	0	0
2	Z	215	Total 1593	C 998	N 274	O 317	S 4	0	0	0
2	2	215	Total 1593	C 998	N 274	O 317	S 4	0	0	0

There are 98 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	301	OZT	-	insertion	UNP O33245
C	535	HIS	-	expression tag	UNP O33245
C	536	HIS	-	expression tag	UNP O33245
C	537	HIS	-	expression tag	UNP O33245
C	538	HIS	-	expression tag	UNP O33245
C	539	HIS	-	expression tag	UNP O33245
C	540	HIS	-	expression tag	UNP O33245
E	301	OZT	-	insertion	UNP O33245
E	535	HIS	-	expression tag	UNP O33245
E	536	HIS	-	expression tag	UNP O33245
E	537	HIS	-	expression tag	UNP O33245
E	538	HIS	-	expression tag	UNP O33245

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Chain	Residue	Modelled	Actual	Comment	Reference
E	539	HIS	-	expression tag	UNP O33245
E	540	HIS	-	expression tag	UNP O33245
G	301	OZT	-	insertion	UNP O33245
G	535	HIS	-	expression tag	UNP O33245
G	536	HIS	-	expression tag	UNP O33245
G	537	HIS	-	expression tag	UNP O33245
G	538	HIS	-	expression tag	UNP O33245
G	539	HIS	-	expression tag	UNP O33245
G	540	HIS	-	expression tag	UNP O33245
H	301	OZT	-	insertion	UNP O33245
H	535	HIS	-	expression tag	UNP O33245
H	536	HIS	-	expression tag	UNP O33245
H	537	HIS	-	expression tag	UNP O33245
H	538	HIS	-	expression tag	UNP O33245
H	539	HIS	-	expression tag	UNP O33245
H	540	HIS	-	expression tag	UNP O33245
J	301	OZT	-	insertion	UNP O33245
J	535	HIS	-	expression tag	UNP O33245
J	536	HIS	-	expression tag	UNP O33245
J	537	HIS	-	expression tag	UNP O33245
J	538	HIS	-	expression tag	UNP O33245
J	539	HIS	-	expression tag	UNP O33245
J	540	HIS	-	expression tag	UNP O33245
L	301	OZT	-	insertion	UNP O33245
L	535	HIS	-	expression tag	UNP O33245
L	536	HIS	-	expression tag	UNP O33245
L	537	HIS	-	expression tag	UNP O33245
L	538	HIS	-	expression tag	UNP O33245
L	539	HIS	-	expression tag	UNP O33245
L	540	HIS	-	expression tag	UNP O33245
N	301	OZT	-	insertion	UNP O33245
N	535	HIS	-	expression tag	UNP O33245
N	536	HIS	-	expression tag	UNP O33245
N	537	HIS	-	expression tag	UNP O33245
N	538	HIS	-	expression tag	UNP O33245
N	539	HIS	-	expression tag	UNP O33245
N	540	HIS	-	expression tag	UNP O33245
P	301	OZT	-	insertion	UNP O33245
P	535	HIS	-	expression tag	UNP O33245
P	536	HIS	-	expression tag	UNP O33245
P	537	HIS	-	expression tag	UNP O33245
P	538	HIS	-	expression tag	UNP O33245

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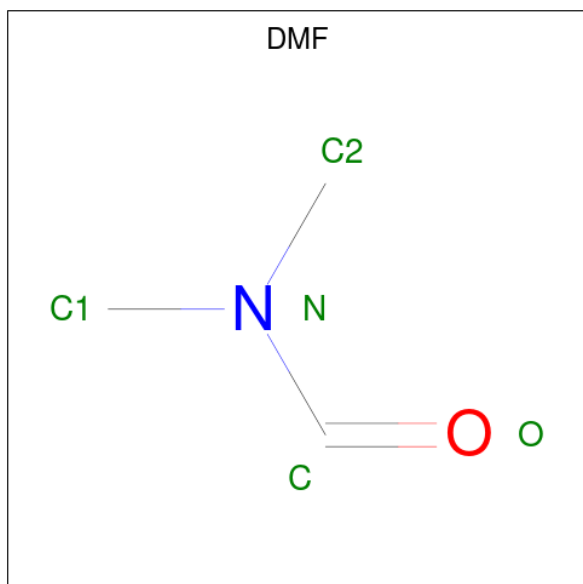
Chain	Residue	Modelled	Actual	Comment	Reference
P	539	HIS	-	expression tag	UNP O33245
P	540	HIS	-	expression tag	UNP O33245
R	301	OZT	-	insertion	UNP O33245
R	535	HIS	-	expression tag	UNP O33245
R	536	HIS	-	expression tag	UNP O33245
R	537	HIS	-	expression tag	UNP O33245
R	538	HIS	-	expression tag	UNP O33245
R	539	HIS	-	expression tag	UNP O33245
R	540	HIS	-	expression tag	UNP O33245
T	301	OZT	-	insertion	UNP O33245
T	535	HIS	-	expression tag	UNP O33245
T	536	HIS	-	expression tag	UNP O33245
T	537	HIS	-	expression tag	UNP O33245
T	538	HIS	-	expression tag	UNP O33245
T	539	HIS	-	expression tag	UNP O33245
T	540	HIS	-	expression tag	UNP O33245
V	301	OZT	-	insertion	UNP O33245
V	535	HIS	-	expression tag	UNP O33245
V	536	HIS	-	expression tag	UNP O33245
V	537	HIS	-	expression tag	UNP O33245
V	538	HIS	-	expression tag	UNP O33245
V	539	HIS	-	expression tag	UNP O33245
V	540	HIS	-	expression tag	UNP O33245
X	301	OZT	-	insertion	UNP O33245
X	535	HIS	-	expression tag	UNP O33245
X	536	HIS	-	expression tag	UNP O33245
X	537	HIS	-	expression tag	UNP O33245
X	538	HIS	-	expression tag	UNP O33245
X	539	HIS	-	expression tag	UNP O33245
X	540	HIS	-	expression tag	UNP O33245
Z	301	OZT	-	insertion	UNP O33245
Z	535	HIS	-	expression tag	UNP O33245
Z	536	HIS	-	expression tag	UNP O33245
Z	537	HIS	-	expression tag	UNP O33245
Z	538	HIS	-	expression tag	UNP O33245
Z	539	HIS	-	expression tag	UNP O33245
Z	540	HIS	-	expression tag	UNP O33245
2	301	OZT	-	insertion	UNP O33245
2	535	HIS	-	expression tag	UNP O33245
2	536	HIS	-	expression tag	UNP O33245
2	537	HIS	-	expression tag	UNP O33245
2	538	HIS	-	expression tag	UNP O33245

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Chain	Residue	Modelled	Actual	Comment	Reference
2	539	HIS	-	expression tag	UNP O33245
2	540	HIS	-	expression tag	UNP O33245

- Molecule 3 is DIMETHYLFORMAMIDE (CCD ID: DMF) (formula: C_3H_7NO).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			5	3	1	1		
3	A	1	Total	C	N	O	0	0
			5	3	1	1		
3	B	1	Total	C	N	O	0	0
			5	3	1	1		
3	B	1	Total	C	N	O	0	0
			5	3	1	1		
3	B	1	Total	C	N	O	0	0
			5	3	1	1		
3	C	1	Total	C	N	O	0	0
			5	3	1	1		
3	C	1	Total	C	N	O	0	0
			5	3	1	1		
3	C	1	Total	C	N	O	0	0
			5	3	1	1		
3	C	1	Total	C	N	O	0	0
			5	3	1	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	N	O	0	0
			5	3	1	1		
3	C	1	Total	C	N	O	0	0
			5	3	1	1		
3	C	1	Total	C	N	O	0	0
			5	3	1	1		
3	D	1	Total	C	N	O	0	0
			5	3	1	1		
3	D	1	Total	C	N	O	0	0
			5	3	1	1		
3	D	1	Total	C	N	O	0	0
			5	3	1	1		
3	D	1	Total	C	N	O	0	0
			5	3	1	1		
3	E	1	Total	C	N	O	0	0
			5	3	1	1		
3	E	1	Total	C	N	O	0	0
			5	3	1	1		
3	E	1	Total	C	N	O	0	0
			5	3	1	1		
3	E	1	Total	C	N	O	0	0
			5	3	1	1		
3	E	1	Total	C	N	O	0	0
			5	3	1	1		
3	E	1	Total	C	N	O	0	0
			5	3	1	1		
3	F	1	Total	C	N	O	0	0
			5	3	1	1		
3	F	1	Total	C	N	O	0	0
			5	3	1	1		
3	G	1	Total	C	N	O	0	0
			5	3	1	1		
3	G	1	Total	C	N	O	0	0
			5	3	1	1		
3	G	1	Total	C	N	O	0	0
			5	3	1	1		
3	G	1	Total	C	N	O	0	0
			5	3	1	1		
3	H	1	Total	C	N	O	0	0
			5	3	1	1		
3	H	1	Total	C	N	O	0	0
			5	3	1	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	H	1	Total	C	N	O	0	0
			5	3	1	1		
3	H	1	Total	C	N	O	0	0
			5	3	1	1		
3	H	1	Total	C	N	O	0	0
			5	3	1	1		
3	H	1	Total	C	N	O	0	0
			5	3	1	1		
3	H	1	Total	C	N	O	0	0
			5	3	1	1		
3	I	1	Total	C	N	O	0	0
			5	3	1	1		
3	I	1	Total	C	N	O	0	0
			5	3	1	1		
3	J	1	Total	C	N	O	0	0
			5	3	1	1		
3	J	1	Total	C	N	O	0	0
			5	3	1	1		
3	J	1	Total	C	N	O	0	0
			5	3	1	1		
3	K	1	Total	C	N	O	0	0
			5	3	1	1		
3	K	1	Total	C	N	O	0	0
			5	3	1	1		
3	L	1	Total	C	N	O	0	0
			5	3	1	1		
3	L	1	Total	C	N	O	0	0
			5	3	1	1		
3	L	1	Total	C	N	O	0	0
			5	3	1	1		
3	M	1	Total	C	N	O	0	0
			5	3	1	1		
3	M	1	Total	C	N	O	0	0
			5	3	1	1		
3	N	1	Total	C	N	O	0	0
			5	3	1	1		
3	N	1	Total	C	N	O	0	0
			5	3	1	1		
3	N	1	Total	C	N	O	0	0
			5	3	1	1		
3	N	1	Total	C	N	O	0	0
			5	3	1	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	N	1	Total	C	N	O	0	0
			5	3	1	1		
3	N	1	Total	C	N	O	0	0
			5	3	1	1		
3	N	1	Total	C	N	O	0	0
			5	3	1	1		
3	O	1	Total	C	N	O	0	0
			5	3	1	1		
3	O	1	Total	C	N	O	0	0
			5	3	1	1		
3	P	1	Total	C	N	O	0	0
			5	3	1	1		
3	P	1	Total	C	N	O	0	0
			5	3	1	1		
3	P	1	Total	C	N	O	0	0
			5	3	1	1		
3	P	1	Total	C	N	O	0	0
			5	3	1	1		
3	P	1	Total	C	N	O	0	0
			5	3	1	1		
3	Q	1	Total	C	N	O	0	0
			5	3	1	1		
3	Q	1	Total	C	N	O	0	0
			5	3	1	1		
3	R	1	Total	C	N	O	0	0
			5	3	1	1		
3	R	1	Total	C	N	O	0	0
			5	3	1	1		
3	R	1	Total	C	N	O	0	0
			5	3	1	1		
3	R	1	Total	C	N	O	0	0
			5	3	1	1		
3	S	1	Total	C	N	O	0	0
			5	3	1	1		
3	S	1	Total	C	N	O	0	0
			5	3	1	1		
3	S	1	Total	C	N	O	0	0
			5	3	1	1		
3	T	1	Total	C	N	O	0	0
			5	3	1	1		
3	T	1	Total	C	N	O	0	0
			5	3	1	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	T	1	Total	C	N	O	0	0
			5	3	1	1		
3	T	1	Total	C	N	O	0	0
			5	3	1	1		
3	T	1	Total	C	N	O	0	0
			5	3	1	1		
3	U	1	Total	C	N	O	0	0
			5	3	1	1		
3	U	1	Total	C	N	O	0	0
			5	3	1	1		
3	U	1	Total	C	N	O	0	0
			5	3	1	1		
3	U	1	Total	C	N	O	0	0
			5	3	1	1		
3	V	1	Total	C	N	O	0	0
			5	3	1	1		
3	V	1	Total	C	N	O	0	0
			5	3	1	1		
3	V	1	Total	C	N	O	0	0
			5	3	1	1		
3	V	1	Total	C	N	O	0	0
			5	3	1	1		
3	W	1	Total	C	N	O	0	0
			5	3	1	1		
3	W	1	Total	C	N	O	0	0
			5	3	1	1		
3	X	1	Total	C	N	O	0	0
			5	3	1	1		
3	X	1	Total	C	N	O	0	0
			5	3	1	1		
3	X	1	Total	C	N	O	0	0
			5	3	1	1		
3	Y	1	Total	C	N	O	0	0
			5	3	1	1		
3	Z	1	Total	C	N	O	0	0
			5	3	1	1		
3	Z	1	Total	C	N	O	0	0
			5	3	1	1		
3	Z	1	Total	C	N	O	0	0
			5	3	1	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	Z	1	Total	C	N	O	0	0
			5	3	1	1		
3	Z	1	Total	C	N	O	0	0
			5	3	1	1		
3	1	1	Total	C	N	O	0	0
			5	3	1	1		
3	1	1	Total	C	N	O	0	0
			5	3	1	1		
3	1	1	Total	C	N	O	0	0
			5	3	1	1		
3	2	1	Total	C	N	O	0	0
			5	3	1	1		
3	2	1	Total	C	N	O	0	0
			5	3	1	1		
3	2	1	Total	C	N	O	0	0
			5	3	1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	43	Total	O	0	0
			43	43		
4	B	45	Total	O	0	0
			45	45		
4	C	142	Total	O	0	0
			142	142		
4	D	43	Total	O	0	0
			43	43		
4	E	124	Total	O	0	0
			124	124		
4	F	34	Total	O	0	0
			34	34		
4	G	136	Total	O	0	0
			136	136		
4	H	122	Total	O	0	0
			122	122		
4	I	37	Total	O	0	0
			37	37		
4	J	128	Total	O	0	0
			128	128		
4	K	45	Total	O	0	0
			45	45		

Continued on next page...

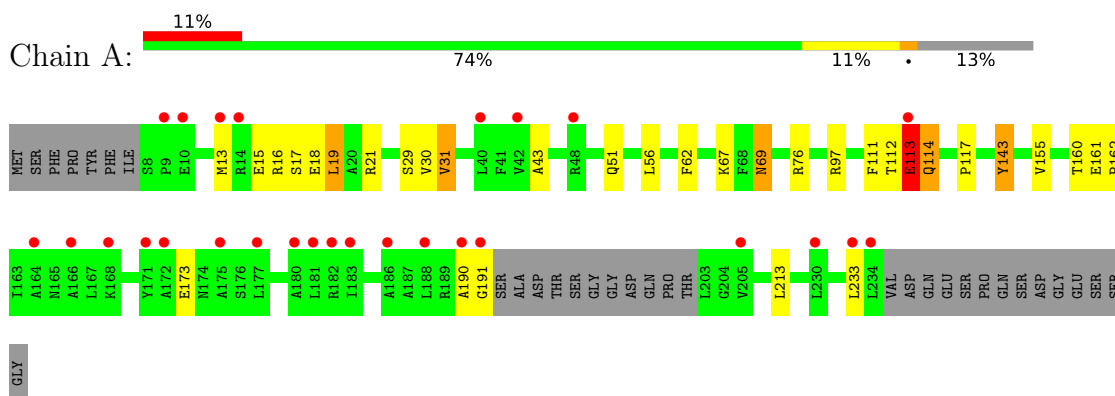
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	138	Total 138	O 138	0	0
4	M	52	Total 52	O 52	0	0
4	N	118	Total 118	O 118	0	0
4	O	51	Total 51	O 51	0	0
4	P	111	Total 111	O 111	0	0
4	Q	41	Total 41	O 41	0	0
4	R	141	Total 141	O 141	0	0
4	S	31	Total 31	O 31	0	0
4	T	112	Total 112	O 112	0	0
4	U	43	Total 43	O 43	0	0
4	V	154	Total 154	O 154	0	0
4	W	31	Total 31	O 31	0	0
4	X	130	Total 130	O 130	0	0
4	Y	41	Total 41	O 41	0	0
4	Z	113	Total 113	O 113	0	0
4	1	42	Total 42	O 42	0	0
4	2	133	Total 133	O 133	0	0

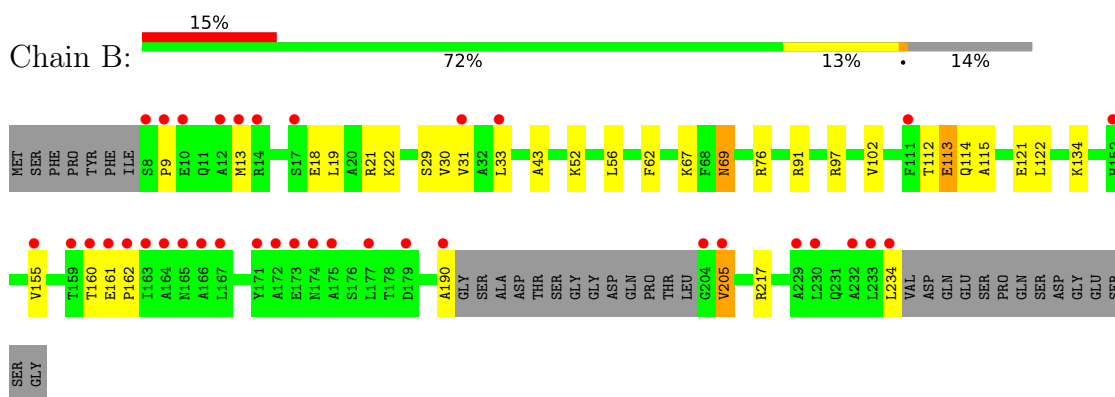
3 Residue-property plots [i](#)

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

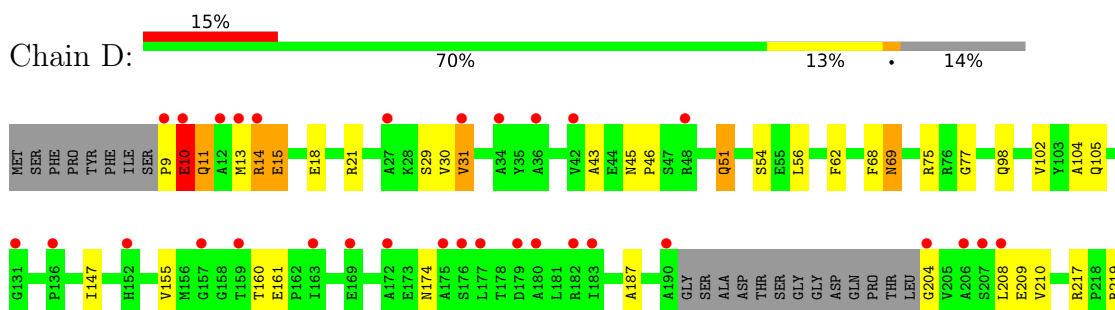
• Molecule 1: Proteasome (Alpha subunit) PrcA

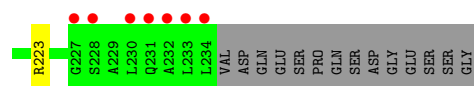


• Molecule 1: Proteasome (Alpha subunit) PrcA

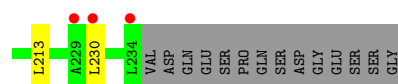
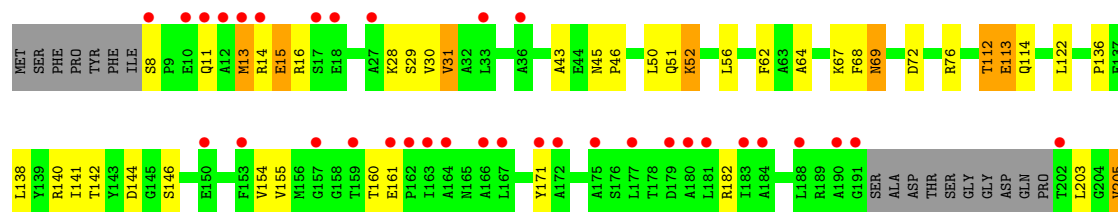


• Molecule 1: Proteasome (Alpha subunit) PrcA

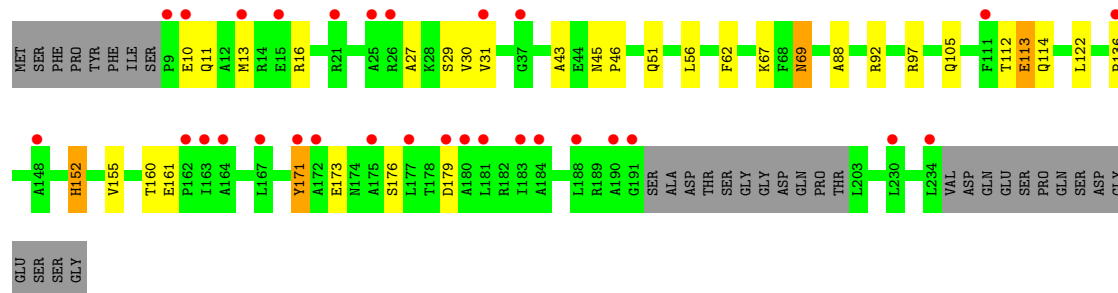




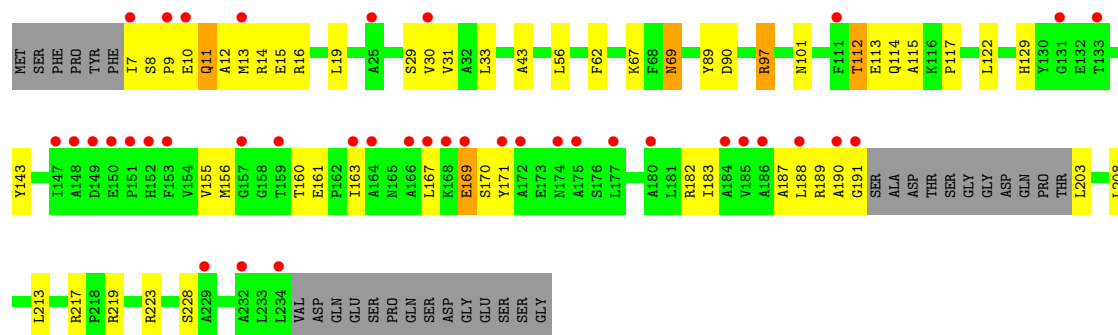
• Molecule 1: Proteasome (Alpha subunit) PrcA



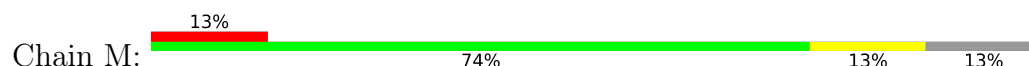
• Molecule 1: Proteasome (Alpha subunit) PrcA

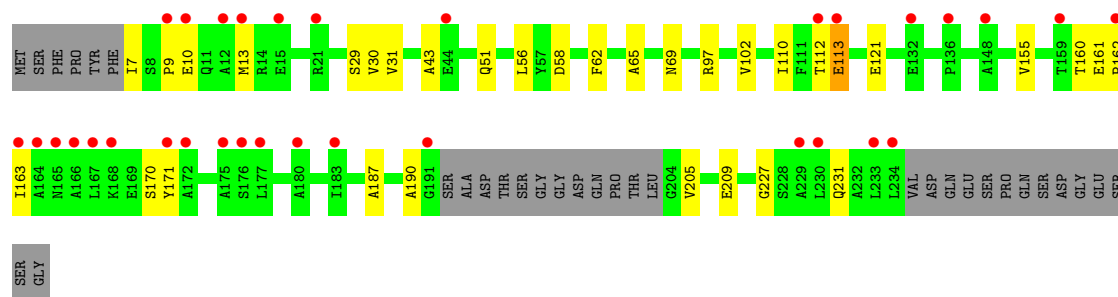


• Molecule 1: Proteasome (Alpha subunit) PrcA

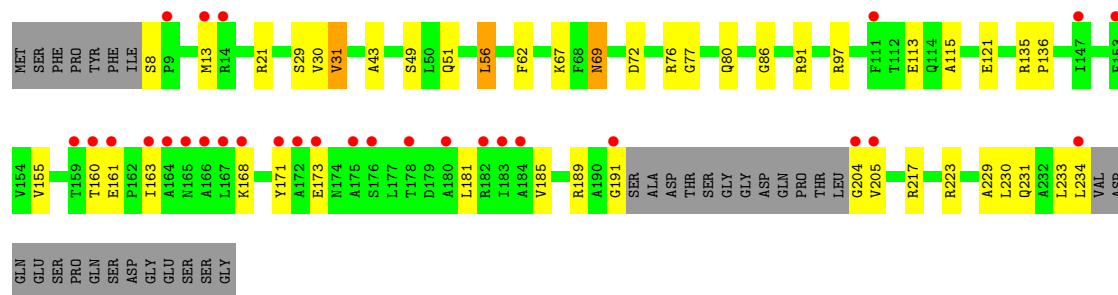


• Molecule 1: Proteasome (Alpha subunit) PrcA

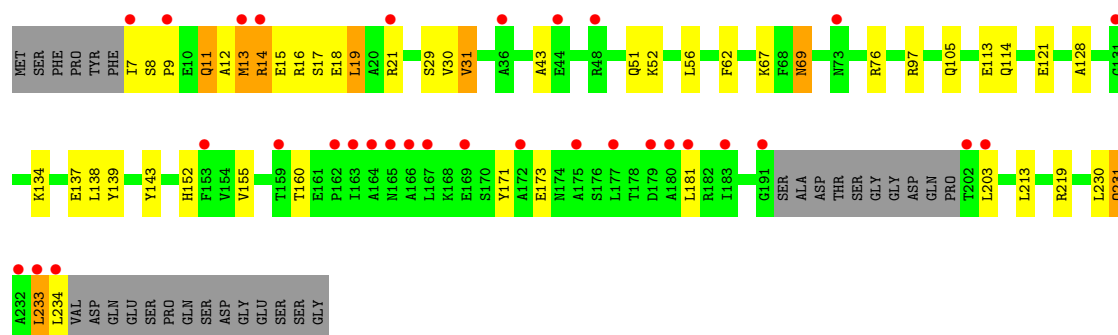




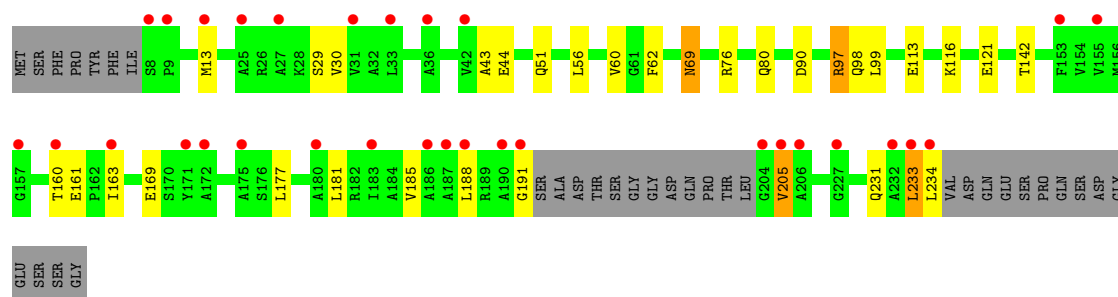
• Molecule 1: Proteasome (Alpha subunit) PrcA



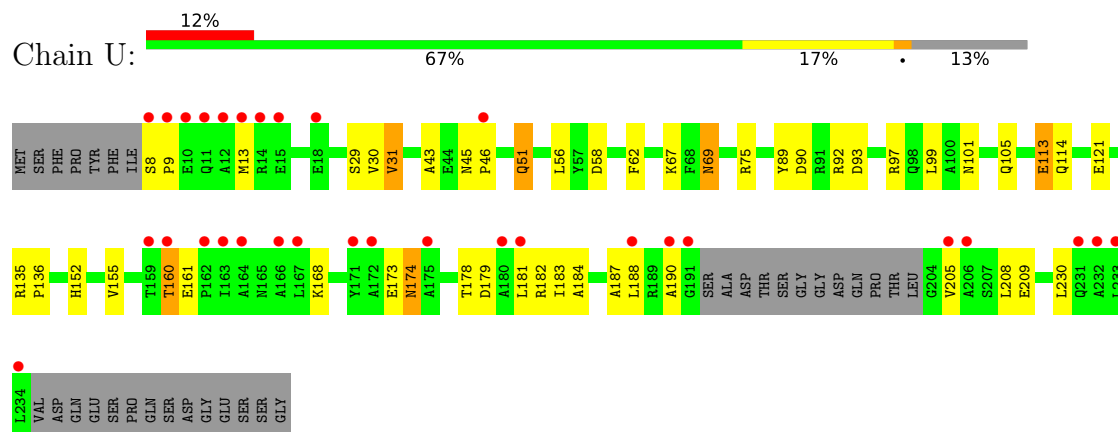
• Molecule 1: Proteasome (Alpha subunit) PrcA



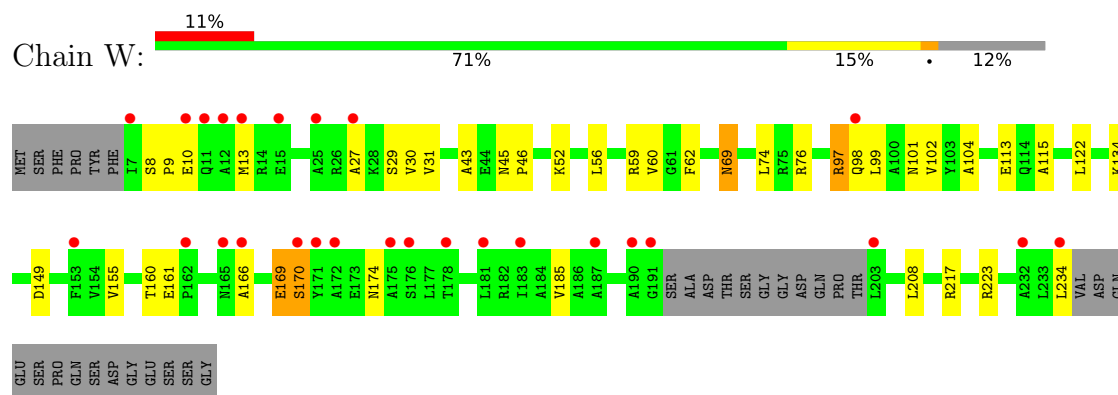
• Molecule 1: Proteasome (Alpha subunit) PrcA



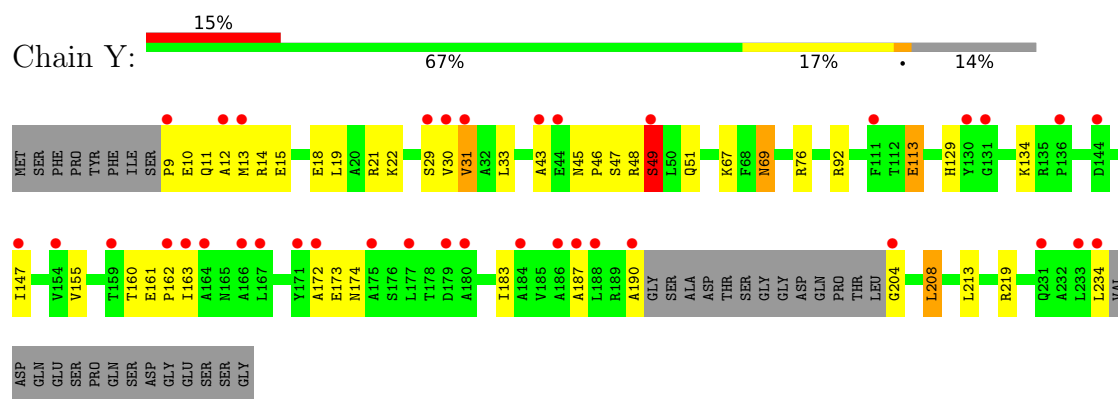
- Molecule 1: Proteasome (Alpha subunit) PrcA



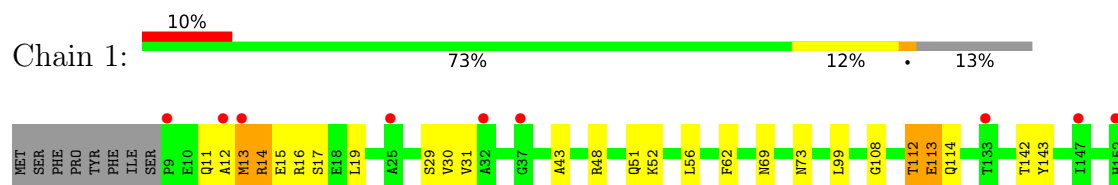
- Molecule 1: Proteasome (Alpha subunit) PrcA

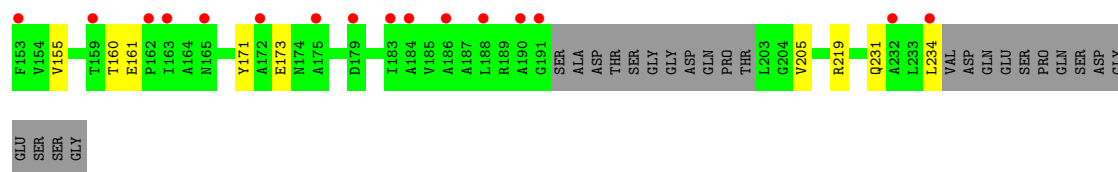


- Molecule 1: Proteasome (Alpha subunit) PrcA

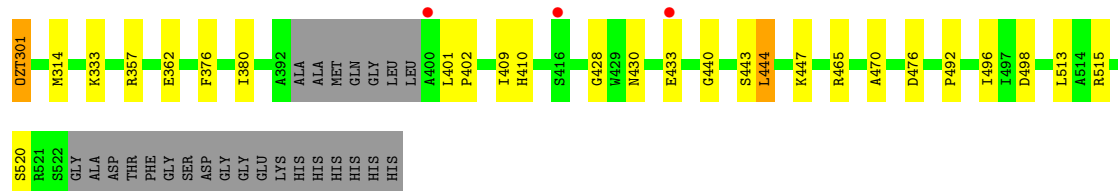
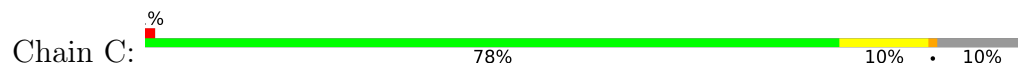


- Molecule 1: Proteasome (Alpha subunit) PrcA

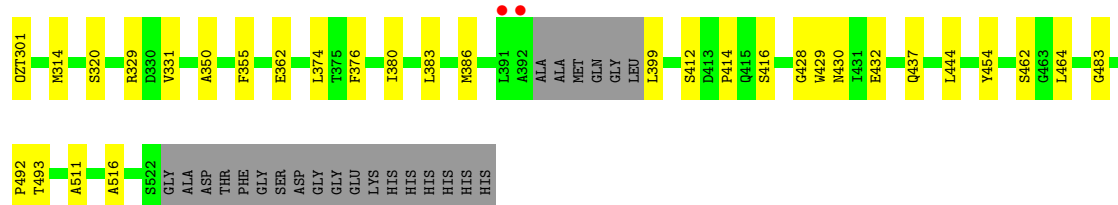
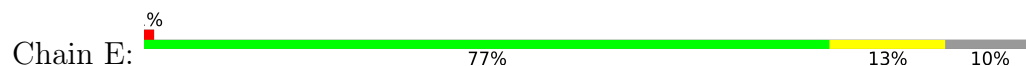




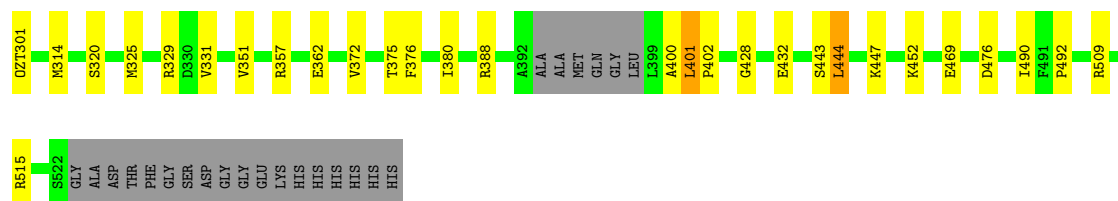
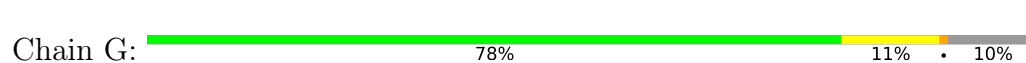
- Molecule 2: Proteasome (Beta subunit) PrcB



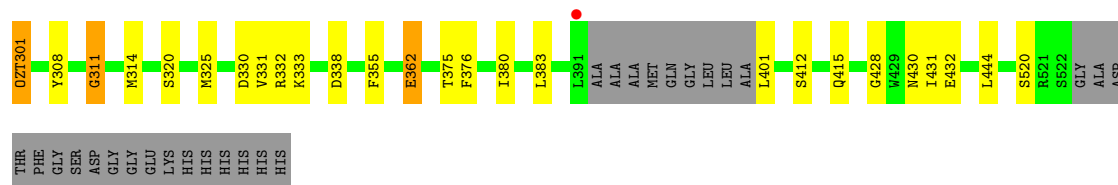
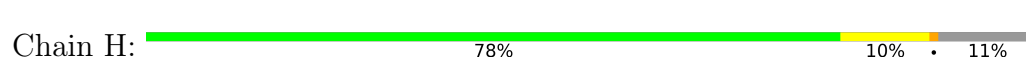
- Molecule 2: Proteasome (Beta subunit) PrcB



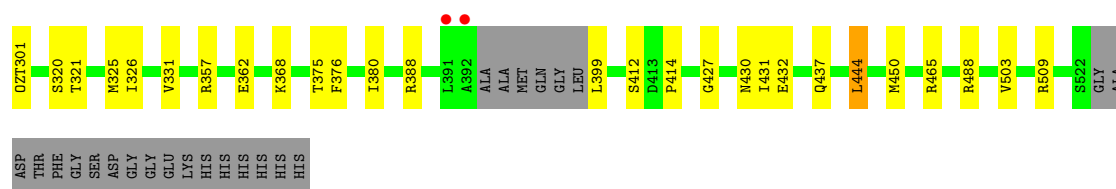
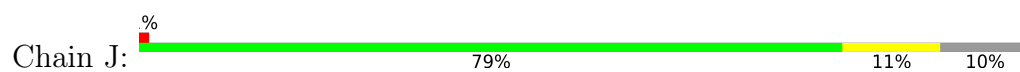
- Molecule 2: Proteasome (Beta subunit) PrcB



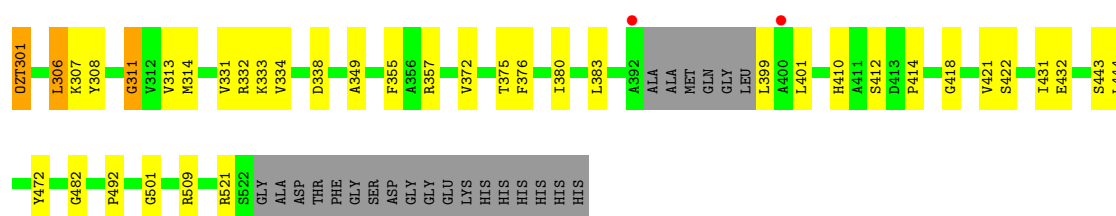
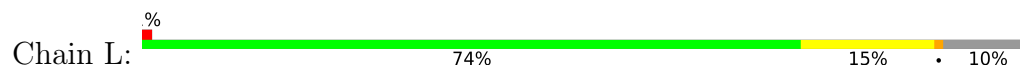
- Molecule 2: Proteasome (Beta subunit) PrcB



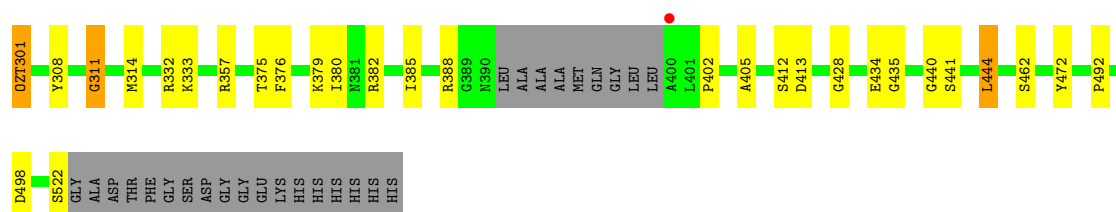
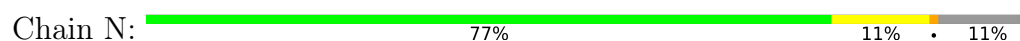
- Molecule 2: Proteasome (Beta subunit) PrcB



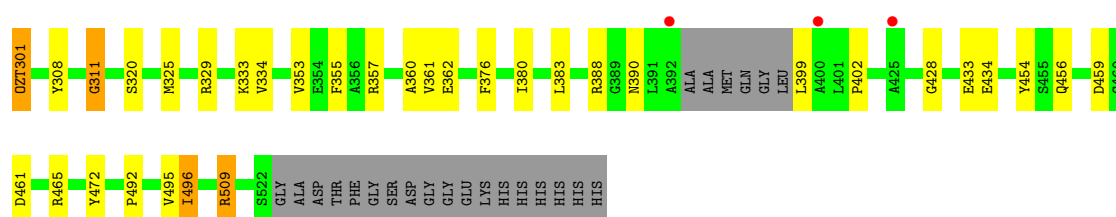
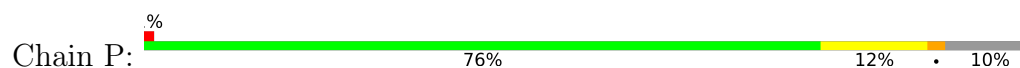
• Molecule 2: Proteasome (Beta subunit) PrcB



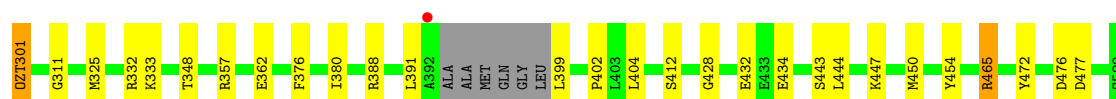
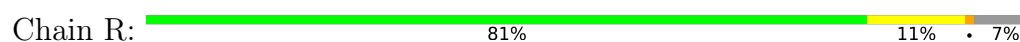
• Molecule 2: Proteasome (Beta subunit) PrcB



• Molecule 2: Proteasome (Beta subunit) PrcB



• Molecule 2: Proteasome (Beta subunit) PrcB



ASP
GLY
GLY
GLY
LYS
HIS
HIS
HIS
HIS
HIS

• Molecule 2: Proteasome (Beta subunit) PrcB

Chain T:  2% 78% 12% 10%

OZT301 M314 S320 D330 V331 R332 K333 V334 D339 F355 A360 V372 P373 L383 R388 L391 A392 ALA MET GLN GLY LEU L399 A400 A401 P402 L403 G428 W429 N430 I431 E432 Q437 L444 K447 K451 D459 S462 P492 E507

S522 GLY ALA ASP THR PHE GLY SER ASP K333 GLY GLY GLU LYS HIS HIS HIS HIS HIS

• Molecule 2: Proteasome (Beta subunit) PrcB

Chain V:  0% 79% 13% 8%

OZT301 M314 S320 K325 I326 R329 D330 V331 R332 K333 A349 E354 F355 K368 F376 I380 L383 R388 L391 A392 ALA MET GLN GLY LEU LEU A400 L401 P402 D413 G428 E434 S443 L444 Y454 L464 R465 V468 P492

I496 S528 SER ASP GLY GLY LYS HIS HIS HIS HIS HIS

• Molecule 2: Proteasome (Beta subunit) PrcB

Chain X:  2% 78% 12% 10%

OZT301 M314 S320 V331 R332 K333 A349 F355 E362 V372 T375 F376 I380 L383 G389 N390 L391 A392 ALA MET GLN GLY LEU L399 A400 D413 G428 E432 E433 Q437 L444 D473 D476 P492 V495 I496 V503

R509 R515 S522 GLY ALA ASP THR PHE GLY SER ASP GLY GLY GLU LYS HIS HIS HIS HIS HIS

• Molecule 2: Proteasome (Beta subunit) PrcB

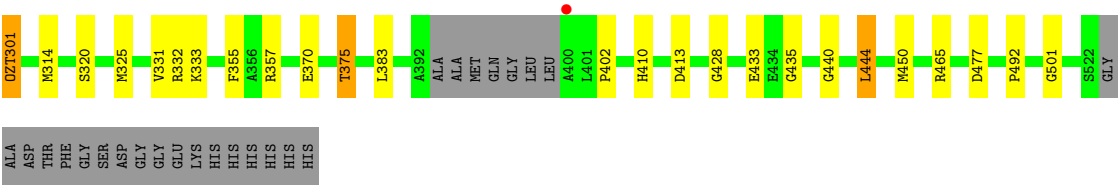
Chain Z:  80% 8% 10%

OZT301 K307 Y308 G311 M314 R332 K333 D338 F355 E362 T375 L383 M386 V387 R388 A392 ALA MET GLN GLY LEU LEU A400 P402 L403 G428 W429 N430 I431 L444 G463 P492 T493 S522 GLY ALA ASP THR PHE GLY SER ASP

GLY
GLY
GLU
LYS
HIS
HIS
HIS
HIS
HIS

• Molecule 2: Proteasome (Beta subunit) PrcB

Chain 2:  79% 9% 10%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	171.62Å 117.67Å 196.34Å 90.00° 113.65° 90.00°	Depositor
Resolution (Å)	29.75 – 2.43 29.75 – 2.44	Depositor EDS
% Data completeness (in resolution range)	97.8 (29.75-2.43) 97.9 (29.75-2.44)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.82 (at 2.45Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.204 , 0.232 0.203 , 0.230	Depositor DCC
R_{free} test set	13191 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	30.5	Xtriage
Anisotropy	0.198	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 43.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.003 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	48540	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.73% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	1	0.83	0/1681	1.14	7/2270 (0.3%)
1	A	0.82	0/1687	1.15	7/2279 (0.3%)
1	B	0.83	0/1675	1.15	6/2263 (0.3%)
1	D	0.83	0/1669	1.12	6/2254 (0.3%)
1	F	0.85	0/1694	1.13	8/2289 (0.3%)
1	I	0.81	0/1681	1.14	7/2270 (0.3%)
1	K	0.83	0/1695	1.11	6/2290 (0.3%)
1	M	0.84	0/1687	1.18	9/2279 (0.4%)
1	O	0.85	0/1679	1.15	6/2268 (0.3%)
1	Q	0.85	0/1702	1.11	8/2300 (0.3%)
1	S	0.78	0/1679	1.13	5/2268 (0.2%)
1	U	0.84	0/1679	1.15	11/2268 (0.5%)
1	W	0.84	0/1695	1.15	8/2290 (0.3%)
1	Y	0.81	0/1669	1.18	9/2254 (0.4%)
2	2	1.14	3/1607 (0.2%)	1.11	5/2178 (0.2%)
2	C	1.12	2/1607 (0.1%)	1.13	1/2178 (0.0%)
2	E	1.15	4/1615 (0.2%)	1.15	2/2189 (0.1%)
2	G	1.08	5/1615 (0.3%)	1.13	3/2189 (0.1%)
2	H	1.07	1/1597 (0.1%)	1.14	3/2164 (0.1%)
2	J	1.11	3/1615 (0.2%)	1.10	2/2189 (0.1%)
2	L	1.13	1/1615 (0.1%)	1.14	4/2189 (0.2%)
2	N	1.08	1/1594 (0.1%)	1.15	6/2160 (0.3%)
2	P	1.04	3/1615 (0.2%)	1.10	6/2189 (0.3%)
2	R	1.09	1/1661 (0.1%)	1.15	6/2251 (0.3%)
2	T	1.05	1/1615 (0.1%)	1.16	3/2189 (0.1%)
2	V	1.16	2/1647 (0.1%)	1.10	6/2232 (0.3%)
2	X	1.12	4/1615 (0.2%)	1.11	3/2189 (0.1%)
2	Z	1.06	1/1607 (0.1%)	1.11	3/2178 (0.1%)
All	All	0.97	32/46197 (0.1%)	1.14	156/62506 (0.2%)

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	2	314	MET	SD-CE	-13.72	1.45	1.79
2	Z	314	MET	SD-CE	-12.28	1.48	1.79
2	C	314	MET	SD-CE	-11.65	1.50	1.79
2	X	314	MET	SD-CE	-11.36	1.51	1.79
2	V	314	MET	SD-CE	-10.91	1.52	1.79
2	E	314	MET	SD-CE	-9.83	1.54	1.79
2	G	314	MET	SD-CE	-8.94	1.57	1.79
2	P	496	ILE	C-N	-8.83	1.21	1.33
2	H	314	MET	SD-CE	-7.98	1.59	1.79
2	J	503	VAL	CA-CB	7.83	1.66	1.54
2	T	372	VAL	CA-CB	7.65	1.64	1.54
2	X	495	VAL	C-N	-6.50	1.25	1.33
2	X	503	VAL	CA-CB	6.46	1.64	1.54
2	P	495	VAL	C-N	-6.33	1.25	1.33
2	2	325	MET	SD-CE	6.32	1.95	1.79
2	V	326	ILE	C-O	-6.31	1.17	1.24
2	E	516	ALA	CA-CB	-6.13	1.43	1.53
2	X	372	VAL	CA-CB	6.08	1.62	1.54
2	E	511	ALA	CA-CB	5.98	1.62	1.53
2	2	450	MET	SD-CE	5.97	1.94	1.79
2	J	450	MET	SD-CE	5.92	1.94	1.79
2	C	470	ALA	CA-C	5.83	1.60	1.52
2	N	462	SER	CA-C	5.83	1.60	1.52
2	G	325	MET	SD-CE	5.82	1.94	1.79
2	P	353	VAL	CA-CB	5.79	1.61	1.54
2	J	326	ILE	C-O	-5.57	1.17	1.23
2	G	469	GLU	CA-C	5.36	1.59	1.52
2	R	450	MET	SD-CE	5.34	1.92	1.79
2	E	350	ALA	CA-CB	-5.27	1.44	1.53
2	L	372	VAL	CA-CB	5.24	1.60	1.54
2	G	372	VAL	CA-CB	5.10	1.60	1.54
2	G	400	ALA	CA-C	5.05	1.58	1.52

All (156) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	173	GLU	N-CA-C	9.52	121.66	111.28
1	M	205	VAL	N-CA-C	8.98	120.86	110.62
1	A	113	GLU	N-CA-C	8.69	120.83	111.36
1	B	113	GLU	N-CA-C	8.56	123.01	112.23
1	O	113	GLU	N-CA-C	8.43	122.06	111.69
1	B	205	VAL	N-CA-C	8.31	119.18	111.45
1	S	113	GLU	N-CA-C	8.17	121.37	111.40

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Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	W	113	GLU	N-CA-C	7.99	123.16	113.41
1	Q	114	GLN	N-CA-C	7.95	120.19	110.41
1	I	113	GLU	N-CA-C	7.69	121.92	112.23
1	D	10	GLU	N-CA-C	-7.56	104.03	113.55
1	M	171	TYR	N-CA-C	7.51	120.74	109.25
1	S	205	VAL	N-CA-C	7.34	118.11	110.62
2	P	496	ILE	O-C-N	-7.33	115.28	123.20
2	R	311	GLY	N-CA-C	7.28	121.07	110.80
1	B	160	THR	N-CA-C	7.20	119.21	111.36
1	D	160	THR	N-CA-C	7.08	119.08	111.36
1	O	160	THR	N-CA-C	7.02	119.01	111.36
1	M	160	THR	N-CA-C	6.96	118.95	111.36
1	Q	160	THR	N-CA-C	6.92	118.90	111.36
1	I	160	THR	N-CA-C	6.87	118.85	111.36
1	Y	160	THR	N-CA-C	6.82	118.79	111.36
2	L	311	GLY	N-CA-C	6.82	121.73	110.55
2	N	375	THR	N-CA-C	-6.78	102.07	110.41
1	U	160	THR	N-CA-C	6.78	118.75	111.36
1	S	160	THR	N-CA-C	6.75	118.72	111.36
1	D	208	LEU	N-CA-C	6.74	119.68	109.23
1	W	160	THR	N-CA-C	6.73	118.70	111.36
1	M	113	GLU	N-CA-C	6.72	121.77	113.50
1	1	173	GLU	N-CA-C	6.71	118.59	111.28
1	W	208	LEU	N-CA-C	6.70	119.90	109.52
1	1	113	GLU	N-CA-C	6.58	118.53	111.36
1	1	160	THR	N-CA-C	6.53	118.48	111.36
1	I	171	TYR	N-CA-C	6.49	120.08	109.76
1	A	160	THR	N-CA-C	6.48	118.43	111.36
1	K	160	THR	N-CA-C	6.47	118.42	111.36
1	W	115	ALA	N-CA-C	6.39	117.90	111.07
1	Y	208	LEU	N-CA-C	6.36	119.89	109.72
1	Y	49	SER	N-CA-C	6.30	122.88	114.12
1	F	160	THR	N-CA-C	6.28	118.20	111.36
1	W	170	SER	N-CA-C	6.23	119.05	111.82
2	P	325	MET	N-CA-C	6.22	119.64	109.76
2	H	311	GLY	N-CA-C	6.06	120.49	110.55
2	N	413	ASP	CA-C-N	6.05	125.77	119.05
2	N	413	ASP	C-N-CA	6.05	125.77	119.05
2	P	402	PRO	N-CA-C	6.01	120.71	111.57
1	D	29	SER	N-CA-C	6.00	119.83	110.17
1	1	29	SER	N-CA-C	6.00	119.82	110.17
2	R	443	SER	N-CA-C	5.98	118.58	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	R	402	PRO	N-CA-C	5.96	120.86	111.38
1	Y	51	GLN	N-CA-C	5.94	119.07	110.28
2	H	375	THR	N-CA-C	-5.93	103.11	110.41
1	Y	47	SER	N-CA-C	5.93	118.93	110.10
1	W	29	SER	N-CA-C	5.92	119.69	110.17
1	F	203	LEU	N-CA-C	5.89	119.77	111.52
1	M	29	SER	N-CA-C	5.89	119.65	110.17
2	X	413	ASP	CA-C-N	5.88	125.75	119.28
2	X	413	ASP	C-N-CA	5.88	125.75	119.28
2	P	311	GLY	N-CA-C	5.87	120.17	110.55
1	F	113	GLU	N-CA-C	5.84	121.13	113.30
1	Y	113	GLU	N-CA-C	5.84	118.53	111.40
1	A	29	SER	N-CA-C	5.83	119.55	110.17
1	K	208	LEU	N-CA-C	5.82	118.95	109.24
1	M	170	SER	N-CA-C	5.77	120.18	113.20
1	O	56	LEU	N-CA-C	-5.72	105.39	112.72
1	U	208	LEU	N-CA-C	5.72	118.54	109.50
2	2	501	GLY	N-CA-C	5.72	119.02	111.70
1	A	51	GLN	N-CA-C	5.70	119.20	110.20
1	1	51	GLN	N-CA-C	5.69	119.20	110.20
1	U	29	SER	N-CA-C	5.69	119.33	110.17
1	I	152	HIS	N-CA-C	5.68	119.98	112.72
2	2	413	ASP	CA-C-N	5.67	126.05	119.47
2	2	413	ASP	C-N-CA	5.67	126.05	119.47
1	B	29	SER	N-CA-C	5.65	119.26	110.17
2	C	443	SER	N-CA-C	5.65	117.44	111.28
2	N	435	GLY	N-CA-C	5.64	120.39	113.79
2	P	454	TYR	N-CA-C	5.60	118.15	111.71
2	R	391	LEU	N-CA-C	-5.59	101.83	109.71
1	S	51	GLN	N-CA-C	5.56	118.98	110.20
1	I	29	SER	N-CA-C	5.53	119.07	110.17
2	Z	311	GLY	N-CA-C	5.52	119.61	110.55
2	G	443	SER	N-CA-C	5.50	117.27	111.28
1	U	121	GLU	N-CA-C	-5.50	98.96	108.69
1	D	51	GLN	N-CA-C	5.48	118.85	110.20
1	O	29	SER	N-CA-C	5.47	118.97	110.17
1	K	33	LEU	N-CA-C	5.45	117.67	109.23
2	N	311	GLY	N-CA-C	5.45	119.48	110.55
2	R	412	SER	N-CA-C	5.44	117.21	111.28
2	V	413	ASP	CA-C-N	5.44	125.26	119.28
2	V	413	ASP	C-N-CA	5.44	125.26	119.28
1	K	29	SER	N-CA-C	5.43	118.91	110.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	S	29	SER	N-CA-C	5.41	118.87	110.17
1	F	29	SER	N-CA-C	5.40	118.87	110.17
2	Z	375	THR	N-CA-C	-5.40	103.57	110.53
1	O	51	GLN	N-CA-C	5.36	118.68	110.20
1	B	121	GLU	N-CA-C	-5.36	99.21	108.69
2	E	462	SER	N-CA-C	-5.35	105.61	111.82
1	Y	29	SER	N-CA-C	5.35	118.78	110.17
2	G	401	LEU	N-CA-C	5.34	117.68	110.31
1	Q	29	SER	N-CA-C	5.34	118.77	110.17
2	J	325	MET	N-CA-C	5.34	118.25	109.76
2	V	443	SER	N-CA-C	5.33	117.08	111.28
2	T	401	LEU	CA-C-N	5.32	125.22	119.85
2	T	401	LEU	C-N-CA	5.32	125.22	119.85
2	V	320	SER	N-CA-C	-5.30	101.01	109.23
2	V	454	TYR	N-CA-C	5.28	117.78	111.71
1	W	169	GLU	N-CA-C	5.27	116.83	111.14
1	K	143	TYR	N-CA-C	5.26	118.49	111.75
2	Z	402	PRO	N-CA-C	5.26	119.57	111.57
1	U	58	ASP	N-CA-C	5.25	116.69	111.07
1	Q	121	GLU	N-CA-C	-5.24	99.41	108.69
1	Y	33	LEU	N-CA-C	5.24	117.35	109.23
1	M	51	GLN	N-CA-C	5.23	118.47	110.20
1	U	51	GLN	N-CA-C	5.22	118.45	110.20
1	U	89	TYR	O-C-N	5.22	124.69	120.83
2	H	325	MET	N-CA-C	5.19	117.74	109.96
2	N	402	PRO	N-CA-C	5.19	119.45	111.57
1	U	113	GLU	N-CA-C	5.18	119.73	113.41
1	Q	51	GLN	N-CA-C	5.17	118.12	110.48
1	Q	171	TYR	N-CA-C	5.16	117.66	109.96
1	1	171	TYR	N-CA-C	5.16	117.17	108.96
2	L	501	GLY	N-CA-C	5.15	118.29	111.70
2	V	402	PRO	N-CA-C	5.15	119.39	111.57
1	M	227	GLY	N-CA-C	5.14	120.70	111.34
2	L	443	SER	N-CA-C	5.14	116.88	111.28
2	E	454	TYR	N-CA-C	5.13	117.61	111.71
1	U	209	GLU	N-CA-C	-5.13	99.92	108.34
1	O	171	TYR	N-CA-C	5.13	116.54	109.15
1	W	174	ASN	N-CA-C	5.12	118.80	112.24
2	J	375	THR	N-CA-C	-5.12	103.93	110.53
1	Y	213	LEU	N-CA-C	-5.11	97.09	107.41
1	I	173	GLU	N-CA-C	5.11	116.93	111.36
1	1	143	TYR	N-CA-C	5.11	118.29	111.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	112	THR	N-CA-C	5.11	116.92	111.36
1	I	51	GLN	N-CA-C	5.10	118.26	110.20
2	T	403	LEU	N-CA-C	-5.10	100.10	108.41
2	G	375	THR	N-CA-C	-5.09	103.96	110.53
2	L	375	THR	N-CA-C	-5.09	104.14	110.41
1	A	114	GLN	N-CA-C	5.09	117.97	110.59
1	M	58	ASP	N-CA-C	5.08	116.51	111.07
1	D	68	PHE	N-CA-C	5.07	117.21	111.02
1	Q	213	LEU	N-CA-C	-5.07	97.17	107.41
1	B	33	LEU	N-CA-C	5.07	117.08	109.23
1	Q	233	LEU	N-CA-C	5.06	116.80	111.28
2	2	375	THR	N-CA-C	-5.06	104.00	110.53
2	X	375	THR	N-CA-C	-5.06	103.85	110.53
1	A	143	TYR	N-CA-C	5.05	118.22	111.75
1	U	8	SER	CA-C-N	5.05	125.08	119.32
1	U	8	SER	C-N-CA	5.05	125.08	119.32
2	2	402	PRO	N-CA-C	5.04	119.24	111.57
1	F	68	PHE	N-CA-C	5.04	117.17	111.02
1	F	171	TYR	N-CA-C	5.03	116.95	109.25
1	F	213	LEU	N-CA-C	-5.03	97.25	107.41
1	K	213	LEU	N-CA-C	-5.03	97.25	107.41
2	P	459	ASP	N-CA-C	5.02	114.96	107.88
2	R	454	TYR	N-CA-C	5.01	117.47	111.71

There are no chirality outliers.

There are no planarity outliers.

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	1	1656	0	1658	34	0
1	A	1662	0	1662	30	0
1	B	1650	0	1648	32	0
1	D	1644	0	1644	33	0
1	F	1669	0	1669	38	0
1	I	1656	0	1658	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	K	1670	0	1673	53	0
1	M	1662	0	1662	16	0
1	O	1654	0	1651	44	0
1	Q	1677	0	1680	39	0
1	S	1654	0	1651	20	0
1	U	1654	0	1651	38	0
1	W	1670	0	1673	27	0
1	Y	1644	0	1644	49	0
2	2	1593	0	1577	19	0
2	C	1593	0	1577	32	0
2	E	1601	0	1588	19	0
2	G	1601	0	1588	24	0
2	H	1583	0	1567	18	0
2	J	1601	0	1588	24	0
2	L	1601	0	1588	30	0
2	N	1580	0	1561	21	0
2	P	1601	0	1588	24	0
2	R	1646	0	1624	23	0
2	T	1601	0	1588	26	0
2	V	1632	0	1608	17	0
2	X	1601	0	1588	17	0
2	Z	1593	0	1577	19	0
3	1	15	0	21	3	0
3	2	15	0	21	1	0
3	A	10	0	14	2	0
3	B	15	0	21	2	0
3	C	40	0	56	20	0
3	D	20	0	28	2	0
3	E	30	0	42	5	0
3	F	10	0	14	1	0
3	G	20	0	28	9	0
3	H	35	0	49	8	0
3	I	10	0	14	0	0
3	J	15	0	21	5	0
3	K	10	0	14	2	0
3	L	15	0	21	5	0
3	M	10	0	14	1	0
3	N	35	0	49	7	0
3	O	10	0	14	6	0
3	P	25	0	35	2	0
3	Q	10	0	14	0	0
3	R	20	0	28	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	S	15	0	21	1	0
3	T	25	0	35	9	0
3	U	20	0	28	4	0
3	V	20	0	28	1	0
3	W	10	0	14	1	0
3	X	15	0	21	1	0
3	Y	5	0	7	0	0
3	Z	30	0	42	10	0
4	1	42	0	0	9	0
4	2	133	0	0	13	0
4	A	43	0	0	6	0
4	B	45	0	0	13	0
4	C	142	0	0	12	0
4	D	43	0	0	15	0
4	E	124	0	0	8	0
4	F	34	0	0	6	0
4	G	136	0	0	18	0
4	H	122	0	0	13	0
4	I	37	0	0	9	0
4	J	128	0	0	16	0
4	K	45	0	0	17	0
4	L	138	0	0	7	0
4	M	52	0	0	6	0
4	N	118	0	0	11	0
4	O	51	0	0	17	0
4	P	111	0	0	10	0
4	Q	41	0	0	4	0
4	R	141	0	0	14	0
4	S	31	0	0	5	0
4	T	112	0	0	9	0
4	U	43	0	0	13	0
4	V	154	0	0	9	0
4	W	31	0	0	5	0
4	X	130	0	0	5	0
4	Y	41	0	0	21	0
4	Z	113	0	0	18	0
All	All	48540	0	46145	791	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 9.

All (791) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:V:400:ALA:HA	4:V:2033:HOH:O	1.27	1.32
2:C:430:ASN:HB2	4:C:2190:HOH:O	1.33	1.26
2:C:447:LYS:CD	3:C:5:DMF:H12	1.66	1.26
2:V:332:ARG:HB3	4:V:2317:HOH:O	1.32	1.25
1:U:160:THR:HG22	4:U:2365:HOH:O	1.34	1.24
2:C:447:LYS:HD2	3:C:5:DMF:C1	1.64	1.24
1:S:121:GLU:HG3	4:S:2023:HOH:O	1.34	1.23
1:A:162:PRO:HB2	4:A:2159:HOH:O	1.36	1.21
1:Y:18:GLU:OE1	1:Y:22:LYS:HE3	1.43	1.19
2:G:402:PRO:HD3	4:G:2237:HOH:O	1.41	1.18
2:X:332:ARG:HD2	4:X:2151:HOH:O	1.44	1.16
2:C:447:LYS:CD	3:C:5:DMF:C1	2.22	1.15
1:1:11:GLN:HA	1:1:14:ARG:HG2	1.21	1.12
1:W:134:LYS:HE2	4:W:2357:HOH:O	1.49	1.11
2:C:447:LYS:HD3	3:C:5:DMF:H12	1.13	1.10
1:S:76:ARG:HD3	4:S:2370:HOH:O	1.52	1.09
2:C:447:LYS:HD2	3:C:5:DMF:H11	1.29	1.08
1:1:11:GLN:HG3	1:1:14:ARG:NE	1.68	1.08
1:O:229:ALA:O	1:O:233:LEU:HD23	1.55	1.06
2:H:332:ARG:HD2	4:H:2352:HOH:O	1.55	1.06
1:1:11:GLN:HG3	1:1:14:ARG:HE	0.93	1.06
2:L:306:LEU:CD1	2:L:313:VAL:CG1	2.33	1.05
2:R:357:ARG:HD2	4:Z:2095:HOH:O	1.56	1.05
2:E:428:GLY:HA2	4:E:2161:HOH:O	1.57	1.04
2:T:388:ARG:HG2	4:T:2215:HOH:O	1.57	1.04
1:D:104:ALA:HB2	4:D:2360:HOH:O	1.58	1.04
2:L:306:LEU:HD12	2:L:313:VAL:CG1	1.89	1.03
1:B:52:LYS:HD3	4:B:2129:HOH:O	1.58	1.01
1:U:51:GLN:HB2	4:U:2105:HOH:O	1.59	1.01
2:G:351:VAL:HG13	4:G:2353:HOH:O	1.59	1.01
1:M:231:GLN:HG3	4:M:2387:HOH:O	1.58	1.01
1:Q:14:ARG:HH11	1:Q:14:ARG:HB3	1.24	1.00
2:L:306:LEU:HD12	2:L:313:VAL:HG13	1.40	1.00
2:2:465:ARG:HD3	4:2:2340:HOH:O	1.61	1.00
2:L:306:LEU:CD1	2:L:313:VAL:HG13	1.91	1.00
3:Z:91:DMF:HC	4:Z:1148:HOH:O	1.59	1.00
2:X:432:GLU:HB3	4:X:2235:HOH:O	1.62	1.00
3:U:252:DMF:H13	4:2:2315:HOH:O	1.62	0.99
1:F:28:LYS:HE2	4:F:2298:HOH:O	1.60	0.99
1:U:9:PRO:HD2	1:1:15:GLU:OE1	1.63	0.99
2:C:496:ILE:HD13	3:C:64:DMF:HC	1.43	0.99
1:O:76:ARG:HD3	4:O:2287:HOH:O	1.61	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:386:MET:SD	4:E:2398:HOH:O	2.21	0.97
1:D:54:SER:HB3	4:D:2426:HOH:O	1.63	0.97
3:E:55:DMF:HC	4:K:2288:HOH:O	1.65	0.96
1:Y:18:GLU:CD	1:Y:21:ARG:HH21	1.73	0.95
2:L:314:MET:HE3	2:L:334:VAL:HG13	1.46	0.95
1:Q:14:ARG:HB3	1:Q:14:ARG:NH1	1.81	0.95
2:J:388:ARG:HG2	4:J:2405:HOH:O	1.67	0.94
2:G:509:ARG:HG3	4:G:2167:HOH:O	1.69	0.92
2:P:433:GLU:HG3	4:P:2308:HOH:O	1.67	0.92
1:Q:97:ARG:HD2	1:Y:49:SER:OG	1.66	0.92
1:O:21:ARG:HD3	4:O:2361:HOH:O	1.69	0.92
2:X:428:GLY:HA2	4:X:2092:HOH:O	1.71	0.91
2:T:399:LEU:HA	4:T:2342:HOH:O	1.71	0.90
3:L:60:DMF:H12	4:N:2035:HOH:O	1.71	0.90
1:D:147:ILE:HB	4:D:2360:HOH:O	1.70	0.90
1:D:11:GLN:HG2	1:D:14:ARG:NH2	1.87	0.89
2:J:509:ARG:HG2	4:J:2068:HOH:O	1.71	0.89
1:I:92:ARG:HB2	4:I:2214:HOH:O	1.73	0.88
1:A:76:ARG:HD3	4:A:2166:HOH:O	1.72	0.88
1:1:11:GLN:HA	1:1:14:ARG:CG	2.04	0.88
2:R:428:GLY:HA2	4:R:2208:HOH:O	1.72	0.88
2:Z:493:THR:HB	4:Z:2402:HOH:O	1.72	0.88
1:1:11:GLN:CA	1:1:14:ARG:HG2	2.03	0.88
2:R:447:LYS:HD3	4:R:2181:HOH:O	1.72	0.87
1:K:217:ARG:HD3	4:K:2228:HOH:O	1.73	0.87
2:X:433:GLU:HG3	4:Z:2252:HOH:O	1.73	0.87
2:C:444:LEU:HB2	3:C:73:DMF:H22	1.56	0.87
3:T:94:DMF:H23	4:T:2165:HOH:O	1.75	0.85
2:J:444:LEU:HD13	4:J:2376:HOH:O	1.77	0.85
2:2:465:ARG:CD	4:2:2340:HOH:O	2.21	0.85
3:G:20:DMF:H11	4:G:2328:HOH:O	1.77	0.85
3:N:2:DMF:HC	4:N:134:HOH:O	1.77	0.84
1:A:19:LEU:C	1:A:19:LEU:HD23	2.02	0.84
2:R:348:THR:HG23	4:R:2216:HOH:O	1.77	0.84
2:H:431:ILE:H	3:H:53:DMF:HC	1.43	0.84
2:H:430:ASN:HB2	4:H:2304:HOH:O	1.76	0.84
2:J:430:ASN:HB2	4:J:2411:HOH:O	1.77	0.84
1:F:76:ARG:HD3	4:F:2225:HOH:O	1.78	0.83
3:G:20:DMF:C1	4:G:2328:HOH:O	2.25	0.83
1:1:11:GLN:CG	1:1:14:ARG:HE	1.87	0.83
2:G:509:ARG:CG	4:G:2167:HOH:O	2.23	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:J:82:DMF:H12	4:J:2332:HOH:O	1.78	0.82
2:R:362:GLU:HB2	4:R:2079:HOH:O	1.78	0.82
1:1:11:GLN:O	1:1:14:ARG:HG3	1.79	0.82
2:C:428:GLY:HA2	4:C:2031:HOH:O	1.79	0.82
1:O:97:ARG:HD3	4:O:2218:HOH:O	1.78	0.82
2:L:306:LEU:HD11	2:L:313:VAL:CG1	2.09	0.81
1:1:12:ALA:O	1:1:16:ARG:HG3	1.80	0.81
1:D:217:ARG:NH2	1:D:223:ARG:HG3	1.95	0.81
1:Y:21:ARG:CG	4:Y:2251:HOH:O	2.29	0.81
3:V:16:DMF:H13	4:2:2372:HOH:O	1.81	0.81
1:K:217:ARG:CD	4:K:2228:HOH:O	2.30	0.80
1:O:97:ARG:CD	4:O:2218:HOH:O	2.28	0.80
2:P:334:VAL:HB	4:P:2373:HOH:O	1.81	0.80
1:U:155:VAL:HG12	4:U:2365:HOH:O	1.82	0.80
2:J:431:ILE:HD13	4:J:2185:HOH:O	1.82	0.80
1:S:60:VAL:HG11	1:S:99:LEU:HD12	1.64	0.80
1:Y:18:GLU:OE1	1:Y:22:LYS:CE	2.29	0.80
1:K:101:ASN:HB2	4:K:1174:HOH:O	1.80	0.79
1:Y:183:ILE:HD13	4:Y:2086:HOH:O	1.81	0.79
2:P:433:GLU:CG	4:P:2308:HOH:O	2.23	0.79
1:K:15:GLU:HG2	1:M:10:GLU:HG2	1.62	0.79
1:Q:19:LEU:C	1:Q:19:LEU:HD23	2.06	0.79
2:2:410:HIS:HE1	4:2:2267:HOH:O	1.65	0.79
2:L:308:TYR:CZ	2:L:311:GLY:HA3	2.18	0.79
1:U:168:LYS:CD	4:U:2067:HOH:O	2.31	0.79
2:T:428:GLY:HA2	4:T:2055:HOH:O	1.83	0.78
2:Z:430:ASN:HB2	4:Z:2171:HOH:O	1.82	0.78
2:2:301:OZT:H27	2:2:333:LYS:HE2	1.65	0.78
2:V:301:OZT:H27	2:V:333:LYS:HE2	1.66	0.78
1:K:12:ALA:O	1:K:16:ARG:HG3	1.83	0.78
2:T:373:PRO:HG2	3:T:35:DMF:HC	1.66	0.78
2:T:314:MET:HE3	2:T:334:VAL:HG13	1.66	0.77
1:O:168:LYS:CD	4:O:2321:HOH:O	2.31	0.77
2:L:410:HIS:CE1	4:L:2102:HOH:O	2.37	0.77
2:P:301:OZT:C7	2:P:333:LYS:HE2	2.14	0.77
1:Y:183:ILE:CD1	4:Y:2086:HOH:O	2.31	0.77
2:G:401:LEU:HA	4:G:2237:HOH:O	1.83	0.77
2:R:332:ARG:HG3	4:R:2085:HOH:O	1.83	0.77
1:O:163:ILE:HG12	1:O:191:GLY:HA3	1.67	0.76
1:B:30:VAL:HG22	4:B:2129:HOH:O	1.86	0.76
2:2:428:GLY:HA2	4:2:2057:HOH:O	1.85	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:228:SER:HA	4:K:2191:HOH:O	1.87	0.75
2:T:373:PRO:HG2	3:T:35:DMF:C	2.17	0.75
1:Y:10:GLU:HA	1:Y:10:GLU:OE2	1.86	0.75
3:Z:69:DMF:C1	4:Z:2391:HOH:O	2.34	0.74
3:U:251:DMF:C1	4:V:2355:HOH:O	2.35	0.74
2:T:332:ARG:HH11	3:T:94:DMF:H22	1.51	0.74
1:O:168:LYS:HD2	4:O:2321:HOH:O	1.87	0.74
1:Q:19:LEU:HD23	1:Q:19:LEU:O	1.87	0.74
2:G:432:GLU:HG2	4:G:2096:HOH:O	1.86	0.74
1:Q:7:ILE:HD12	1:Q:11:GLN:HB3	1.70	0.74
1:1:11:GLN:HB3	4:1:2326:HOH:O	1.88	0.74
2:R:447:LYS:CD	4:R:2181:HOH:O	2.32	0.74
1:Y:12:ALA:O	1:Y:15:GLU:HB2	1.87	0.74
1:Y:21:ARG:HG2	4:Y:2251:HOH:O	1.86	0.73
3:G:96:DMF:HC	4:G:2327:HOH:O	1.88	0.73
2:H:401:LEU:HB3	4:H:2201:HOH:O	1.87	0.73
3:U:251:DMF:H13	4:V:2355:HOH:O	1.89	0.73
1:K:129:HIS:HE1	4:K:2227:HOH:O	1.72	0.73
1:S:205:VAL:HG21	1:S:231:GLN:HG2	1.71	0.73
1:K:15:GLU:HB3	1:M:9:PRO:HG2	1.70	0.72
1:U:152:HIS:NE2	1:U:173:GLU:OE2	2.22	0.72
1:D:104:ALA:CB	4:D:2360:HOH:O	2.27	0.72
1:B:19:LEU:C	1:B:19:LEU:HD23	2.15	0.71
3:H:31:DMF:H23	4:L:2384:HOH:O	1.89	0.71
1:U:105:GLN:NE2	4:U:925:HOH:O	2.23	0.71
2:E:355:PHE:CE2	4:E:2398:HOH:O	2.44	0.71
1:F:140:ARG:NH1	1:F:142:THR:HG22	2.05	0.71
1:Q:230:LEU:O	1:Q:234:LEU:HD12	1.91	0.71
1:S:44:GLU:HG3	1:S:188:LEU:HD12	1.72	0.71
1:I:97:ARG:HG2	4:I:2313:HOH:O	1.90	0.71
1:O:181:LEU:O	1:O:185:VAL:HG23	1.90	0.71
2:G:432:GLU:CG	4:G:2096:HOH:O	2.38	0.71
1:M:121:GLU:HG3	4:M:2114:HOH:O	1.91	0.70
3:G:96:DMF:C	4:G:2327:HOH:O	2.40	0.70
1:D:9:PRO:HD2	4:D:2330:HOH:O	1.92	0.70
1:K:15:GLU:HG2	1:M:10:GLU:CG	2.21	0.70
1:F:140:ARG:HH12	1:F:142:THR:HG22	1.57	0.70
2:J:321:THR:O	3:J:6:DMF:H21	1.91	0.70
1:D:11:GLN:HG2	1:D:14:ARG:CZ	2.21	0.70
2:E:376:PHE:CE2	2:E:380:ILE:HD11	2.26	0.70
1:1:11:GLN:O	1:1:15:GLU:HG3	1.92	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:444:LEU:HD13	4:G:2406:HOH:O	1.93	0.69
1:I:92:ARG:CG	4:I:2214:HOH:O	2.41	0.69
1:Q:137:GLU:C	1:Q:138:LEU:HD12	2.17	0.69
1:Y:172:ALA:CB	4:Y:2086:HOH:O	2.41	0.69
2:N:388:ARG:NH1	4:N:2048:HOH:O	2.24	0.69
1:Y:19:LEU:C	1:Y:19:LEU:HD23	2.18	0.69
1:K:217:ARG:HH21	1:K:223:ARG:HG3	1.58	0.69
2:C:357:ARG:HD2	4:C:2061:HOH:O	1.92	0.69
2:C:496:ILE:HD13	3:C:64:DMF:C	2.20	0.69
1:K:191:GLY:C	4:K:2150:HOH:O	2.35	0.69
1:O:21:ARG:CD	4:O:2361:HOH:O	2.32	0.69
4:A:2243:HOH:O	1:B:112:THR:HG21	1.92	0.68
1:D:187:ALA:HB2	4:D:2331:HOH:O	1.93	0.68
2:J:321:THR:H	3:J:6:DMF:H22	1.56	0.68
1:U:168:LYS:HD2	4:U:2067:HOH:O	1.93	0.68
1:D:10:GLU:HA	1:D:10:GLU:OE1	1.94	0.68
1:Q:219:ARG:NH2	3:R:89:DMF:H22	2.08	0.68
1:Y:92:ARG:HG3	4:Y:2307:HOH:O	1.94	0.68
1:Y:172:ALA:HB3	4:Y:2086:HOH:O	1.94	0.67
1:I:73:ASN:CB	4:I:2207:HOH:O	2.41	0.67
1:F:13:MET:HE2	1:F:16:ARG:HD3	1.74	0.67
2:N:444:LEU:H	3:N:97:DMF:H12	1.59	0.67
1:F:182:ARG:NH2	4:F:2248:HOH:O	2.26	0.67
1:W:60:VAL:HG11	1:W:99:LEU:HD12	1.77	0.67
3:H:19:DMF:H22	4:L:1904:HOH:O	1.94	0.67
3:Z:69:DMF:H13	4:Z:2391:HOH:O	1.94	0.67
2:J:357:ARG:HD3	4:J:2224:HOH:O	1.95	0.67
1:K:169:GLU:OE1	1:K:169:GLU:C	2.38	0.67
3:A:250:DMF:C1	4:H:2272:HOH:O	2.42	0.66
2:E:355:PHE:CZ	4:E:2398:HOH:O	2.47	0.66
2:N:308:TYR:CZ	2:N:311:GLY:HA3	2.30	0.66
1:A:111:PHE:HE1	1:A:143:TYR:CD2	2.14	0.66
3:Z:69:DMF:H12	4:Z:2391:HOH:O	1.96	0.66
2:C:498:ASP:OD2	3:C:64:DMF:C1	2.43	0.66
1:K:182:ARG:HG3	4:K:2028:HOH:O	1.96	0.66
1:F:140:ARG:NH1	1:F:142:THR:CG2	2.59	0.66
1:K:163:ILE:HG23	1:K:187:ALA:C	2.21	0.65
1:B:97:ARG:NH2	4:B:2348:HOH:O	2.28	0.65
1:U:168:LYS:HD3	4:U:2067:HOH:O	1.94	0.65
1:A:16:ARG:NH2	1:A:114:GLN:O	2.30	0.65
1:O:163:ILE:CG1	1:O:191:GLY:HA3	2.25	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:301:OZT:H27	2:C:333:LYS:HE2	1.78	0.65
2:P:301:OZT:H27	2:P:333:LYS:HE2	1.78	0.65
2:T:430:ASN:HB2	4:T:1606:HOH:O	1.97	0.65
1:Y:18:GLU:OE2	1:Y:21:ARG:NH2	2.30	0.65
1:A:18:GLU:OE1	1:A:21:ARG:NH2	2.29	0.65
1:D:147:ILE:CB	4:D:2360:HOH:O	2.37	0.65
2:C:444:LEU:HD22	3:C:73:DMF:H13	1.79	0.65
1:B:19:LEU:HD23	1:B:19:LEU:O	1.97	0.65
1:O:205:VAL:HG21	1:O:231:GLN:HG2	1.79	0.65
1:B:43:ALA:HB1	4:B:2129:HOH:O	1.97	0.64
1:A:113:GLU:N	1:A:113:GLU:OE1	2.30	0.64
3:H:19:DMF:H23	4:H:2377:HOH:O	1.95	0.64
1:U:31:VAL:HB	1:U:188:LEU:HD11	1.79	0.64
2:R:432:GLU:OE2	2:R:434:GLU:HB2	1.96	0.64
3:S:250:DMF:HC	2:T:360:ALA:HB1	1.78	0.64
1:B:114:GLN:NE2	4:B:2199:HOH:O	2.30	0.64
1:D:18:GLU:OE1	1:D:21:ARG:NH2	2.30	0.64
2:G:452:LYS:HB3	3:G:39:DMF:H13	1.79	0.64
1:Y:18:GLU:OE2	1:Y:21:ARG:NE	2.30	0.64
1:1:19:LEU:C	1:1:19:LEU:HD23	2.23	0.64
1:Q:14:ARG:NH1	1:Q:14:ARG:CB	2.58	0.64
2:2:332:ARG:NH2	4:2:1264:HOH:O	2.30	0.63
2:C:433:GLU:HG3	4:C:2311:HOH:O	1.98	0.63
2:G:388:ARG:NH1	4:G:2024:HOH:O	2.31	0.63
2:G:444:LEU:HB3	4:G:1218:HOH:O	1.98	0.63
2:2:357:ARG:NE	4:2:2213:HOH:O	2.31	0.63
1:K:11:GLN:HE21	1:K:14:ARG:NH1	1.96	0.63
1:Q:128:ALA:HB2	1:Q:134:LYS:HB3	1.81	0.63
1:U:92:ARG:HD2	2:2:375:THR:HG21	1.80	0.63
2:V:401:LEU:HB2	4:V:2153:HOH:O	1.99	0.63
2:N:332:ARG:NH2	4:N:2264:HOH:O	2.28	0.63
1:Y:147:ILE:HD13	4:Y:2303:HOH:O	1.98	0.63
1:K:163:ILE:HG23	1:K:187:ALA:O	1.98	0.63
1:B:162:PRO:HB2	1:B:190:ALA:HB1	1.81	0.63
1:A:111:PHE:HE1	1:A:143:TYR:HD2	1.46	0.62
1:D:204:GLY:N	4:D:2400:HOH:O	2.32	0.62
2:N:357:ARG:HD3	4:N:2035:HOH:O	1.97	0.62
1:A:15:GLU:HB3	1:B:9:PRO:HG2	1.81	0.62
3:H:19:DMF:C2	4:H:2377:HOH:O	2.47	0.62
1:K:228:SER:HB3	4:K:2191:HOH:O	1.99	0.62
2:Z:428:GLY:HA2	4:Z:2109:HOH:O	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:401:LEU:HB2	4:G:2255:HOH:O	1.99	0.62
2:L:306:LEU:HD12	2:L:313:VAL:HG12	1.77	0.62
1:M:102:VAL:HG12	3:M:249:DMF:H22	1.81	0.62
2:H:301:OZT:H27	2:H:333:LYS:HE2	1.80	0.62
1:Y:19:LEU:HD23	1:Y:19:LEU:O	2.00	0.62
1:Y:204:GLY:N	4:Y:1615:HOH:O	2.33	0.62
2:X:515:ARG:HD2	4:X:2081:HOH:O	2.00	0.62
1:F:8:SER:N	4:F:2029:HOH:O	2.33	0.62
1:F:140:ARG:HH12	1:F:142:THR:CG2	2.13	0.62
1:O:76:ARG:O	1:O:80:GLN:HG3	1.99	0.62
2:C:515:ARG:HD2	4:C:2187:HOH:O	2.00	0.62
1:D:219:ARG:NH2	3:D:252:DMF:HC	2.15	0.62
2:H:308:TYR:CZ	2:H:311:GLY:HA3	2.35	0.62
1:I:11:GLN:O	1:I:14:ARG:CG	2.47	0.62
2:N:444:LEU:HB2	3:N:97:DMF:H12	1.82	0.61
2:V:428:GLY:HA2	4:V:2197:HOH:O	1.99	0.61
1:W:76:ARG:NH2	3:W:249:DMF:H22	2.15	0.61
2:2:301:OZT:C7	2:2:333:LYS:HE2	2.29	0.61
1:Y:18:GLU:OE1	1:Y:21:ARG:NH2	2.30	0.61
1:I:73:ASN:CG	4:I:2207:HOH:O	2.42	0.61
1:A:112:THR:CG2	1:O:115:ALA:HB3	2.30	0.61
1:D:210:VAL:N	4:D:2263:HOH:O	2.22	0.61
1:W:97:ARG:NH1	4:W:2056:HOH:O	2.30	0.61
1:Y:21:ARG:HG3	4:Y:2251:HOH:O	1.96	0.61
1:F:136:PRO:HG3	4:F:2210:HOH:O	1.99	0.61
2:V:325:MET:SD	2:2:444:LEU:HD21	2.41	0.61
1:Q:13:MET:HE2	1:Q:13:MET:HA	1.83	0.61
1:I:92:ARG:CB	4:I:2214:HOH:O	2.36	0.61
1:U:179:ASP:O	1:U:183:ILE:HG13	2.01	0.61
2:Z:483:GLY:HA3	3:Z:27:DMF:O	2.01	0.61
1:B:134:LYS:NZ	4:B:2220:HOH:O	2.33	0.60
2:G:444:LEU:CD1	4:G:2406:HOH:O	2.48	0.60
1:M:209:GLU:HA	4:M:2429:HOH:O	2.00	0.60
2:G:428:GLY:HA2	4:G:2024:HOH:O	2.02	0.60
1:I:27:ALA:HB1	4:I:2062:HOH:O	2.00	0.60
1:W:166:ALA:O	1:W:169:GLU:HG3	2.01	0.60
2:J:509:ARG:CG	4:J:2068:HOH:O	2.41	0.60
1:I:16:ARG:NH2	1:I:114:GLN:O	2.30	0.60
1:K:182:ARG:CG	4:K:2028:HOH:O	2.49	0.60
2:Z:429:TRP:HZ3	2:Z:431:ILE:HG13	1.66	0.60
1:A:191:GLY:N	4:A:2159:HOH:O	2.35	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:141:ILE:HD12	1:F:141:ILE:N	2.17	0.60
1:O:233:LEU:N	1:O:233:LEU:HD22	2.16	0.60
1:D:147:ILE:CG2	4:D:2360:HOH:O	2.50	0.59
1:F:30:VAL:HG22	1:F:52:LYS:HZ3	1.67	0.59
2:L:331:VAL:HG13	2:L:349:ALA:HB2	1.84	0.59
1:B:52:LYS:CD	4:B:2129:HOH:O	2.32	0.59
1:A:111:PHE:CE1	1:A:143:TYR:HD2	2.20	0.59
3:A:250:DMF:H11	4:H:2272:HOH:O	2.02	0.59
2:C:444:LEU:H	3:C:73:DMF:H22	1.67	0.59
1:D:77:GLY:HA3	3:D:251:DMF:H13	1.83	0.59
1:O:121:GLU:HG3	4:O:2044:HOH:O	2.03	0.59
1:D:30:VAL:HG13	1:D:43:ALA:HB2	1.85	0.59
2:X:515:ARG:NH2	4:X:2341:HOH:O	2.28	0.59
2:H:428:GLY:HA2	4:H:2132:HOH:O	2.02	0.59
1:I:92:ARG:HG3	4:I:2214:HOH:O	2.01	0.59
2:Z:388:ARG:NH1	4:Z:2109:HOH:O	2.35	0.59
2:G:357:ARG:HA	3:G:1:DMF:HC	1.85	0.59
1:U:31:VAL:CG2	1:U:188:LEU:HD11	2.33	0.59
1:Y:147:ILE:CD1	4:Y:2303:HOH:O	2.49	0.59
1:D:11:GLN:HG2	1:D:14:ARG:HH22	1.68	0.58
2:H:332:ARG:HB2	4:H:2352:HOH:O	2.02	0.58
1:Y:129:HIS:HE1	4:Y:2152:HOH:O	1.86	0.58
2:J:321:THR:N	3:J:6:DMF:H22	2.18	0.58
2:N:332:ARG:HG3	4:N:2380:HOH:O	2.02	0.58
2:G:476:ASP:OD1	3:G:20:DMF:H12	2.03	0.58
3:K:250:DMF:HC	3:L:4:DMF:H21	1.85	0.58
2:L:332:ARG:HB2	4:L:2140:HOH:O	2.02	0.58
1:Q:105:GLN:NE2	4:Q:2244:HOH:O	2.36	0.58
1:S:97:ARG:HG2	1:S:98:GLN:N	2.19	0.58
2:C:465:ARG:HD3	4:C:1567:HOH:O	2.03	0.58
1:Y:92:ARG:HG2	4:Y:2152:HOH:O	2.02	0.58
2:2:357:ARG:HD3	4:2:2213:HOH:O	2.03	0.58
1:W:52:LYS:NZ	4:W:2259:HOH:O	2.33	0.58
2:C:409:ILE:HG22	4:C:2174:HOH:O	2.02	0.58
1:O:91:ARG:HH12	3:O:250:DMF:H13	1.69	0.58
4:A:2428:HOH:O	3:H:41:DMF:C2	2.52	0.58
1:K:189:ARG:HD3	1:K:203:LEU:HD12	1.85	0.58
2:L:306:LEU:HD11	2:L:313:VAL:HG13	1.72	0.58
2:H:401:LEU:N	4:H:1585:HOH:O	2.37	0.58
2:L:306:LEU:HD11	2:L:313:VAL:HG11	1.86	0.58
1:K:129:HIS:CE1	4:K:2227:HOH:O	2.52	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:76:ARG:O	1:O:80:GLN:CG	2.52	0.58
1:S:56:LEU:HD13	1:S:99:LEU:HD13	1.86	0.58
1:W:169:GLU:HG3	1:W:170:SER:H	1.69	0.58
1:Q:19:LEU:C	1:Q:19:LEU:CD2	2.77	0.57
2:T:332:ARG:NH1	3:T:94:DMF:H22	2.18	0.57
2:2:410:HIS:CE1	4:2:2267:HOH:O	2.45	0.57
1:O:77:GLY:HA3	3:O:249:DMF:C	2.34	0.57
2:2:357:ARG:CD	4:2:2213:HOH:O	2.52	0.57
1:Q:76:ARG:HD3	4:Q:527:HOH:O	2.04	0.57
2:V:464:LEU:O	2:V:468:VAL:HG23	2.03	0.57
1:F:31:VAL:HG22	1:F:155:VAL:HG22	1.86	0.57
1:B:217:ARG:NH1	4:B:2322:HOH:O	2.37	0.57
1:K:167:LEU:O	1:K:171:TYR:N	2.37	0.57
2:L:307:LYS:HE3	2:L:418:GLY:O	2.04	0.57
1:Y:129:HIS:CE1	4:Y:2152:HOH:O	2.56	0.57
2:J:412:SER:O	2:J:414:PRO:HD3	2.05	0.57
1:I:30:VAL:HG13	1:I:43:ALA:HB2	1.86	0.57
1:Q:230:LEU:HG	1:Q:234:LEU:HD11	1.85	0.57
2:R:465:ARG:HD3	4:R:1483:HOH:O	2.05	0.57
1:U:184:ALA:O	1:U:188:LEU:HD13	2.05	0.57
1:O:229:ALA:O	1:O:233:LEU:CD2	2.44	0.56
1:O:233:LEU:N	1:O:233:LEU:CD2	2.68	0.56
1:Q:67:LYS:HE3	4:Q:389:HOH:O	2.04	0.56
1:Q:69:ASN:HD22	1:Q:69:ASN:H	1.52	0.56
2:P:428:GLY:HA2	4:P:1465:HOH:O	2.05	0.56
1:B:18:GLU:OE1	1:B:21:ARG:NE	2.36	0.56
1:M:97:ARG:NH1	4:M:2407:HOH:O	2.38	0.56
2:T:451:LYS:HE2	2:X:473:ASP:OD1	2.05	0.56
1:K:228:SER:CB	4:K:2191:HOH:O	2.53	0.56
2:R:357:ARG:CD	4:Z:2095:HOH:O	2.30	0.56
2:C:465:ARG:HG3	2:C:513:LEU:HD22	1.87	0.56
2:J:321:THR:H	3:J:6:DMF:C2	2.18	0.56
1:F:30:VAL:CG2	1:F:52:LYS:NZ	2.69	0.56
1:M:110:ILE:HG22	4:M:2182:HOH:O	2.05	0.56
1:A:19:LEU:C	1:A:19:LEU:CD2	2.73	0.55
1:W:60:VAL:HG11	1:W:99:LEU:CD1	2.35	0.55
1:B:52:LYS:NZ	4:B:2129:HOH:O	2.35	0.55
1:I:12:ALA:N	4:I:2326:HOH:O	2.38	0.55
2:P:461:ASP:OD1	2:P:509:ARG:NH2	2.39	0.55
1:U:31:VAL:CB	1:U:188:LEU:HD11	2.36	0.55
3:C:17:DMF:H22	4:C:1355:HOH:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:97:ARG:HD2	4:O:2218:HOH:O	1.99	0.55
1:A:112:THR:HG22	1:O:115:ALA:HB3	1.87	0.55
1:Q:8:SER:HB2	1:Q:9:PRO:CD	2.37	0.55
1:F:69:ASN:H	1:F:69:ASN:HD22	1.55	0.55
1:K:16:ARG:NH2	1:K:114:GLN:O	2.40	0.55
1:B:91:ARG:HD2	4:B:2302:HOH:O	2.07	0.55
1:K:228:SER:CA	4:K:2191:HOH:O	2.50	0.55
1:Q:18:GLU:OE1	1:Q:21:ARG:NH2	2.30	0.55
2:T:432:GLU:HG3	2:T:437:GLN:HB2	1.89	0.55
3:1:251:DMF:HC	3:2:78:DMF:O	2.06	0.55
1:A:19:LEU:HD23	1:A:19:LEU:O	2.07	0.54
1:D:98:GLN:O	1:D:102:VAL:HG23	2.07	0.54
1:U:181:LEU:HD12	1:U:181:LEU:O	2.08	0.54
2:H:401:LEU:CB	4:H:2201:HOH:O	2.52	0.54
1:Y:92:ARG:CG	4:Y:2152:HOH:O	2.56	0.54
1:Y:92:ARG:CG	4:Y:2307:HOH:O	2.52	0.54
1:I:73:ASN:HB2	4:I:2207:HOH:O	2.04	0.54
1:F:46:PRO:HG3	4:F:2298:HOH:O	2.07	0.54
2:L:357:ARG:NH1	4:L:2273:HOH:O	2.30	0.54
2:E:416:SER:HA	4:E:2094:HOH:O	2.06	0.54
1:F:52:LYS:HE2	1:F:64:ALA:O	2.08	0.54
1:U:114:GLN:NE2	4:U:2039:HOH:O	2.41	0.54
2:E:429:TRP:CE2	3:E:70:DMF:H12	2.43	0.54
2:V:388:ARG:NH1	4:V:2197:HOH:O	2.41	0.54
1:I:219:ARG:CZ	3:1:251:DMF:H23	2.38	0.54
2:C:444:LEU:N	3:C:73:DMF:H22	2.22	0.53
2:E:320:SER:HB3	2:E:331:VAL:HG21	1.90	0.53
1:Q:181:LEU:HD23	1:Q:233:LEU:HB3	1.90	0.53
2:C:444:LEU:HB2	3:C:73:DMF:C2	2.33	0.53
2:C:498:ASP:OD2	3:C:64:DMF:H12	2.08	0.53
2:V:301:OZT:C7	2:V:333:LYS:HE2	2.36	0.53
1:Y:92:ARG:HB2	4:Y:2307:HOH:O	2.08	0.53
1:F:16:ARG:NH2	1:F:114:GLN:O	2.32	0.53
1:O:91:ARG:NH1	3:O:250:DMF:H13	2.23	0.53
1:I:30:VAL:HG13	1:I:43:ALA:HB2	1.90	0.53
2:R:332:ARG:NH2	4:R:1931:HOH:O	2.42	0.53
2:C:444:LEU:CB	3:C:73:DMF:H22	2.33	0.53
1:O:21:ARG:NE	4:O:2361:HOH:O	2.39	0.53
1:S:30:VAL:HG13	1:S:43:ALA:HB2	1.90	0.53
2:T:451:LYS:NZ	2:X:476:ASP:OD2	2.42	0.53
2:X:320:SER:HB3	2:X:331:VAL:HG21	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:162:PRO:HB2	1:M:190:ALA:O	2.10	0.52
1:Y:234:LEU:HA	4:Y:2005:HOH:O	2.08	0.52
2:Z:493:THR:CB	4:Z:2402:HOH:O	2.42	0.52
1:D:9:PRO:CD	4:D:2330:HOH:O	2.52	0.52
1:U:97:ARG:NH2	2:2:370:GLU:O	2.43	0.52
1:I:69:ASN:H	1:I:69:ASN:HD22	1.58	0.52
1:O:136:PRO:HG3	4:O:2149:HOH:O	2.09	0.52
2:X:496:ILE:HD13	3:X:67:DMF:H13	1.92	0.52
2:J:357:ARG:CD	4:J:2224:HOH:O	2.56	0.52
1:S:163:ILE:HD11	1:S:191:GLY:HA3	1.92	0.52
1:U:178:THR:HG22	1:U:182:ARG:CZ	2.39	0.52
1:A:16:ARG:HB3	1:A:117:PRO:HG3	1.92	0.52
1:Y:69:ASN:HD22	1:Y:69:ASN:H	1.58	0.52
1:F:15:GLU:HB2	1:W:9:PRO:HG2	1.92	0.52
1:Q:52:LYS:NZ	4:Q:2084:HOH:O	2.27	0.52
1:B:67:LYS:HG2	1:B:69:ASN:HD21	1.76	0.51
2:L:357:ARG:HD3	4:L:2131:HOH:O	2.09	0.51
1:I:52:LYS:NZ	4:I:2042:HOH:O	2.41	0.51
1:F:205:VAL:HG22	1:F:230:LEU:HD23	1.91	0.51
1:I:11:GLN:N	4:I:2112:HOH:O	2.43	0.51
1:Q:231:GLN:OE1	1:Q:231:GLN:HA	2.09	0.51
1:F:50:LEU:HD12	1:W:149:ASP:HB2	1.92	0.51
2:E:355:PHE:CD2	4:E:2398:HOH:O	2.60	0.51
1:K:190:ALA:CB	4:K:2004:HOH:O	2.58	0.51
1:Q:231:GLN:OE1	1:Q:231:GLN:CA	2.58	0.51
1:Y:30:VAL:HG13	1:Y:43:ALA:HB2	1.92	0.51
1:B:21:ARG:NH2	1:B:22:LYS:CG	2.74	0.51
1:Q:8:SER:HB2	1:Q:9:PRO:HD2	1.91	0.51
2:J:368:LYS:NZ	4:J:2318:HOH:O	2.39	0.51
3:O:250:DMF:H12	2:P:361:VAL:HA	1.92	0.51
1:S:177:LEU:HG	1:S:233:LEU:HD11	1.93	0.51
1:W:31:VAL:HG22	1:W:155:VAL:HG22	1.93	0.51
1:Y:19:LEU:C	1:Y:19:LEU:CD2	2.82	0.51
1:Y:31:VAL:HG22	1:Y:155:VAL:HG22	1.92	0.51
3:C:43:DMF:HC	1:I:88:ALA:HA	1.92	0.51
1:O:217:ARG:NH2	1:O:223:ARG:HG3	2.26	0.51
2:Z:308:TYR:CZ	2:Z:311:GLY:HA3	2.46	0.51
1:B:21:ARG:NH2	1:B:22:LYS:HG2	2.26	0.51
1:K:19:LEU:HD23	1:K:19:LEU:C	2.36	0.51
1:K:69:ASN:H	1:K:69:ASN:HD22	1.58	0.51
2:J:431:ILE:CD1	4:J:2185:HOH:O	2.50	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:U:2295:HOH:O	1:I:48:ARG:NH2	2.44	0.50
2:L:355:PHE:HE1	2:L:383:LEU:HD11	1.76	0.50
1:U:75:ARG:HD3	4:U:2058:HOH:O	2.09	0.50
1:Y:76:ARG:HD3	4:Y:1321:HOH:O	2.10	0.50
3:Z:91:DMF:C	4:Z:1148:HOH:O	2.37	0.50
1:K:169:GLU:OE1	1:K:169:GLU:O	2.30	0.50
1:K:170:SER:O	1:K:183:ILE:HD13	2.11	0.50
1:K:217:ARG:NH2	1:K:223:ARG:HG3	2.24	0.50
1:W:185:VAL:HG21	1:W:234:LEU:HD11	1.93	0.50
1:A:69:ASN:H	1:A:69:ASN:HD22	1.60	0.50
3:C:64:DMF:C	4:C:2195:HOH:O	2.60	0.50
2:P:355:PHE:HE2	4:P:2412:HOH:O	1.95	0.50
2:C:496:ILE:CD1	3:C:64:DMF:HC	2.28	0.50
2:H:362:GLU:HG3	4:H:1702:HOH:O	2.12	0.50
2:R:376:PHE:CE2	2:R:380:ILE:HD11	2.46	0.50
1:S:69:ASN:HD22	1:S:69:ASN:H	1.59	0.50
1:S:163:ILE:CG1	1:S:191:GLY:HA3	2.42	0.50
1:Y:10:GLU:O	1:Y:13:MET:HB3	2.11	0.50
2:C:410:HIS:CE1	4:C:2344:HOH:O	2.64	0.50
1:A:111:PHE:CE1	1:A:143:TYR:CD2	2.96	0.50
1:B:115:ALA:HB3	1:I:112:THR:HG23	1.93	0.50
2:L:472:TYR:CE1	3:L:9:DMF:C1	2.95	0.50
1:Q:31:VAL:HG22	1:Q:155:VAL:HG22	1.94	0.50
1:U:101:ASN:HB3	3:U:250:DMF:HC	1.94	0.50
1:F:15:GLU:OE2	1:W:8:SER:OG	2.30	0.50
1:K:97:ARG:NH1	1:K:97:ARG:HG2	2.27	0.50
2:R:301:OZT:H27	2:R:333:LYS:HE2	1.94	0.49
1:B:22:LYS:NZ	1:I:10:GLU:OE2	2.38	0.49
2:G:476:ASP:O	2:V:329:ARG:NH2	2.44	0.49
2:R:432:GLU:HB2	4:R:2012:HOH:O	2.12	0.49
2:T:373:PRO:HG2	3:T:35:DMF:C1	2.42	0.49
2:L:472:TYR:CE1	3:L:9:DMF:H13	2.48	0.49
2:N:301:OZT:H17	2:N:333:LYS:NZ	2.27	0.49
2:P:329:ARG:HD2	2:V:434:GLU:OE2	2.12	0.49
2:P:390:ASN:CG	4:P:2412:HOH:O	2.55	0.49
1:A:30:VAL:HG13	1:A:43:ALA:HB2	1.95	0.49
1:B:102:VAL:HG12	3:B:249:DMF:H12	1.93	0.49
2:J:432:GLU:HG2	2:J:437:GLN:HB2	1.95	0.49
2:T:459:ASP:OD1	2:T:462:SER:OG	2.26	0.49
2:Z:332:ARG:NH2	4:Z:2163:HOH:O	2.45	0.49
1:D:56:LEU:HG	1:D:62:PHE:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Y:173:GLU:O	1:Y:174:ASN:HB2	2.13	0.49
1:K:8:SER:OG	1:K:11:GLN:HB2	2.12	0.49
1:S:142:THR:HA	4:S:2023:HOH:O	2.12	0.49
2:2:320:SER:HB3	2:2:331:VAL:HG21	1.94	0.49
1:A:97:ARG:HD3	1:O:49:SER:HB2	1.95	0.49
2:H:320:SER:HB3	2:H:331:VAL:HG21	1.94	0.49
1:O:76:ARG:NH1	4:O:2287:HOH:O	2.30	0.49
2:P:376:PHE:CE2	2:P:380:ILE:HD11	2.48	0.49
1:Y:204:GLY:O	1:Y:208:LEU:HG	2.12	0.49
1:F:11:GLN:O	1:F:15:GLU:HG3	2.13	0.49
1:U:173:GLU:C	1:U:174:ASN:OD1	2.56	0.49
2:N:472:TYR:CZ	3:N:36:DMF:H22	2.48	0.48
4:O:2280:HOH:O	2:P:357:ARG:NH1	2.34	0.48
2:R:301:OZT:H17	2:R:333:LYS:NZ	2.28	0.48
2:T:430:ASN:HA	3:T:92:DMF:HC	1.94	0.48
1:I:176:SER:OG	1:I:179:ASP:HB2	2.13	0.48
2:R:388:ARG:NH1	4:R:2208:HOH:O	2.47	0.48
1:U:136:PRO:HG3	4:U:2145:HOH:O	2.13	0.48
1:U:51:GLN:CB	4:U:2105:HOH:O	2.38	0.48
1:Y:92:ARG:CB	4:Y:2307:HOH:O	2.61	0.48
1:B:19:LEU:C	1:B:19:LEU:CD2	2.83	0.48
2:Z:429:TRP:CZ3	2:Z:431:ILE:HG13	2.48	0.48
1:F:30:VAL:HG22	1:F:52:LYS:NZ	2.28	0.48
1:S:163:ILE:CG1	1:S:191:GLY:CA	2.92	0.48
2:L:376:PHE:CE2	2:L:380:ILE:HD11	2.48	0.48
1:I:205:VAL:HG21	1:I:231:GLN:HG2	1.95	0.48
1:D:10:GLU:HG2	1:Q:15:GLU:HG3	1.95	0.48
1:I:152:HIS:HB3	1:I:171:TYR:CZ	2.49	0.48
2:G:376:PHE:CE2	2:G:380:ILE:HD11	2.49	0.47
1:F:52:LYS:HE2	1:F:52:LYS:HB3	1.56	0.47
2:P:388:ARG:NH1	4:P:1465:HOH:O	2.46	0.47
2:P:456:GLN:OE1	2:P:465:ARG:NH1	2.46	0.47
1:A:15:GLU:HB3	1:B:9:PRO:CG	2.45	0.47
1:B:56:LEU:HG	1:B:62:PHE:HB2	1.97	0.47
2:L:308:TYR:CE1	2:L:311:GLY:HA3	2.48	0.47
2:H:301:OZT:H17	2:H:333:LYS:NZ	2.30	0.47
2:T:320:SER:HB3	2:T:331:VAL:HG21	1.96	0.47
1:W:76:ARG:HD2	4:W:2090:HOH:O	2.14	0.47
2:2:355:PHE:HE1	2:2:383:LEU:HD11	1.79	0.47
1:I:152:HIS:CD2	1:I:171:TYR:HE2	2.32	0.47
2:L:306:LEU:CD1	2:L:313:VAL:HG12	2.34	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:U:9:PRO:CD	1:1:15:GLU:OE1	2.48	0.47
2:Z:386:MET:HE2	4:Z:2339:HOH:O	2.13	0.47
2:Z:429:TRP:CE2	3:Z:71:DMF:H22	2.50	0.47
3:1:250:DMF:HC	4:1:1918:HOH:O	2.14	0.47
2:C:444:LEU:HD22	3:C:73:DMF:C1	2.45	0.47
1:K:11:GLN:O	1:K:12:ALA:C	2.56	0.47
1:K:190:ALA:HB3	4:K:2004:HOH:O	2.14	0.47
2:Z:301:OZT:H27	2:Z:333:LYS:HE2	1.96	0.47
1:1:13:MET:HE2	1:1:13:MET:HB3	1.69	0.47
4:B:2199:HOH:O	1:I:112:THR:HG21	2.14	0.47
1:K:97:ARG:HG2	1:K:97:ARG:HH11	1.80	0.47
2:N:301:OZT:H27	2:N:333:LYS:HE2	1.97	0.47
1:O:168:LYS:HD3	4:O:2321:HOH:O	2.08	0.47
2:R:332:ARG:CG	4:R:2085:HOH:O	2.51	0.47
1:W:30:VAL:HG13	1:W:43:ALA:HB2	1.97	0.47
1:I:105:GLN:NE2	4:I:2193:HOH:O	2.47	0.46
1:K:170:SER:OG	1:K:183:ILE:HG23	2.15	0.46
1:O:135:ARG:HD3	4:O:1427:HOH:O	2.15	0.46
1:Q:13:MET:HE3	1:Q:13:MET:HB2	1.72	0.46
2:Z:338:ASP:OD1	2:Z:338:ASP:C	2.58	0.46
1:B:69:ASN:HD22	1:B:69:ASN:H	1.64	0.46
2:2:477:ASP:OD2	4:2:1889:HOH:O	2.21	0.46
1:F:67:LYS:HG2	1:F:69:ASN:HD21	1.80	0.46
1:Y:219:ARG:NH2	3:Z:91:DMF:H22	2.30	0.46
3:N:2:DMF:H23	4:P:2424:HOH:O	2.14	0.46
2:E:430:ASN:HA	3:E:95:DMF:HC	1.96	0.46
1:F:30:VAL:CG2	1:F:52:LYS:HZ1	2.28	0.46
2:P:320:SER:CB	4:P:2168:HOH:O	2.63	0.46
2:E:483:GLY:HA3	3:E:32:DMF:H23	1.97	0.46
2:J:465:ARG:NH2	4:J:2247:HOH:O	2.41	0.46
2:L:355:PHE:CE1	2:L:383:LEU:HD11	2.51	0.46
1:W:98:GLN:O	1:W:102:VAL:HG23	2.16	0.46
1:Y:18:GLU:OE2	1:Y:21:ARG:CZ	2.64	0.46
1:M:31:VAL:HG22	1:M:155:VAL:HG22	1.98	0.46
2:E:429:TRP:CZ2	3:E:70:DMF:H12	2.50	0.46
1:F:138:LEU:HD23	1:F:154:VAL:HG23	1.98	0.46
3:H:19:DMF:H12	2:L:521:ARG:HB3	1.98	0.46
1:K:217:ARG:HD2	4:K:2228:HOH:O	2.05	0.46
1:O:31:VAL:HG22	1:O:155:VAL:HG22	1.96	0.46
1:D:69:ASN:H	1:D:69:ASN:HD22	1.64	0.45
2:H:432:GLU:HB2	4:H:2119:HOH:O	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:30:VAL:HG13	1:M:43:ALA:HB2	1.98	0.45
2:P:301:OZT:H27	2:P:333:LYS:CE	2.47	0.45
1:S:56:LEU:HG	1:S:62:PHE:HB2	1.97	0.45
1:Y:163:ILE:HG23	1:Y:187:ALA:O	2.17	0.45
1:K:19:LEU:HD22	1:K:117:PRO:HG2	1.98	0.45
1:U:69:ASN:H	1:U:69:ASN:HD22	1.65	0.45
1:U:205:VAL:HG13	1:U:230:LEU:HD23	1.98	0.45
2:H:355:PHE:HE1	2:H:383:LEU:HD11	1.82	0.45
1:I:152:HIS:CD2	1:I:171:TYR:CE2	3.05	0.45
1:O:67:LYS:HG2	1:O:69:ASN:HD21	1.82	0.45
1:O:69:ASN:HD22	1:O:69:ASN:H	1.63	0.45
1:Q:12:ALA:O	1:Q:16:ARG:HG3	2.16	0.45
1:W:169:GLU:HG3	1:W:170:SER:N	2.31	0.45
1:I:108:GLY:O	1:I:112:THR:HG23	2.16	0.45
1:F:56:LEU:HG	1:F:62:PHE:HB2	1.99	0.45
2:N:428:GLY:HA2	4:N:2048:HOH:O	2.16	0.45
1:W:10:GLU:OE1	1:W:10:GLU:HA	2.16	0.45
1:W:74:LEU:HD13	1:W:122:LEU:HD21	1.99	0.45
3:Z:27:DMF:C2	4:Z:2076:HOH:O	2.64	0.45
2:H:431:ILE:H	3:H:53:DMF:C	2.22	0.45
1:Y:9:PRO:HD2	1:Y:11:GLN:HB3	1.99	0.45
1:D:9:PRO:N	4:D:2330:HOH:O	2.49	0.45
1:F:11:GLN:O	1:F:15:GLU:CG	2.64	0.45
1:I:31:VAL:HG22	1:I:155:VAL:HG22	1.98	0.45
2:J:444:LEU:CD1	4:J:2376:HOH:O	2.49	0.45
2:X:355:PHE:HE1	2:X:383:LEU:HD11	1.82	0.45
2:L:412:SER:O	2:L:414:PRO:HD3	2.17	0.45
2:N:314:MET:HE3	2:N:405:ALA:HB2	1.99	0.45
1:I:67:LYS:HG2	1:I:69:ASN:HD21	1.81	0.45
1:O:30:VAL:HG13	1:O:43:ALA:HB2	1.97	0.45
2:V:355:PHE:HE1	2:V:383:LEU:HD11	1.82	0.45
2:J:320:SER:HB3	2:J:331:VAL:HG21	1.99	0.44
2:J:427:GLY:HA3	4:J:2405:HOH:O	2.16	0.44
2:N:434:GLU:HB2	4:N:2211:HOH:O	2.17	0.44
1:U:178:THR:CG2	1:U:182:ARG:CZ	2.95	0.44
1:W:56:LEU:HG	1:W:62:PHE:HB2	1.99	0.44
1:A:112:THR:HB	1:A:113:GLU:OE1	2.17	0.44
2:J:376:PHE:CE2	2:J:380:ILE:HD11	2.52	0.44
2:Z:307:LYS:HE2	2:Z:307:LYS:HB3	1.76	0.44
1:A:56:LEU:HG	1:A:62:PHE:HB2	2.00	0.44
1:D:147:ILE:HG21	4:D:2360:HOH:O	2.15	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:9:PRO:O	1:K:10:GLU:C	2.60	0.44
1:M:65:ALA:HB1	4:M:2177:HOH:O	2.17	0.44
2:P:472:TYR:CE1	3:P:23:DMF:H22	2.53	0.44
2:R:301:OZT:O6	4:R:2282:HOH:O	2.21	0.44
2:V:376:PHE:CE2	2:V:380:ILE:HD11	2.53	0.44
1:F:51:GLN:HB2	3:F:250:DMF:H21	1.98	0.44
1:1:11:GLN:C	1:1:14:ARG:HG2	2.42	0.44
2:J:488:ARG:NH2	2:R:476:ASP:OD1	2.49	0.44
2:P:308:TYR:CZ	2:P:311:GLY:HA3	2.53	0.44
1:Q:137:GLU:OE2	1:Q:139:TYR:OH	2.30	0.44
2:X:376:PHE:CE2	2:X:380:ILE:HD11	2.52	0.44
1:A:112:THR:HG23	1:O:115:ALA:HB3	1.99	0.44
1:W:69:ASN:H	1:W:69:ASN:HD22	1.63	0.44
1:1:11:GLN:HG3	1:1:14:ARG:CD	2.46	0.44
2:C:444:LEU:HB3	4:C:281:HOH:O	2.16	0.44
1:F:45:ASN:HA	1:F:46:PRO:HD2	1.85	0.44
1:B:31:VAL:HG22	1:B:155:VAL:HG22	1.99	0.44
1:D:105:GLN:NE2	4:D:1127:HOH:O	2.50	0.44
2:G:320:SER:HB3	2:G:331:VAL:HG21	1.98	0.44
1:M:56:LEU:HG	1:M:62:PHE:HB2	2.00	0.44
1:1:56:LEU:HG	1:1:62:PHE:HB2	2.00	0.44
2:J:465:ARG:NE	4:J:2247:HOH:O	2.33	0.44
1:O:8:SER:N	4:O:1604:HOH:O	2.50	0.44
3:O:250:DMF:H23	2:P:360:ALA:HB3	2.00	0.44
1:Q:152:HIS:NE2	1:Q:173:GLU:OE1	2.51	0.44
2:Z:314:MET:HE1	2:Z:403:LEU:O	2.18	0.44
2:L:472:TYR:CZ	3:L:9:DMF:H11	2.53	0.43
1:B:30:VAL:HG13	1:B:43:ALA:HB2	2.00	0.43
2:G:447:LYS:HZ1	3:G:96:DMF:H23	1.83	0.43
1:K:30:VAL:HG13	1:K:43:ALA:HB2	2.00	0.43
2:N:382:ARG:NH2	2:N:385:ILE:HD13	2.34	0.43
2:T:447:LYS:NZ	4:T:2133:HOH:O	2.50	0.43
2:C:476:ASP:O	2:E:329:ARG:NH2	2.50	0.43
2:E:355:PHE:CE1	4:E:2398:HOH:O	2.70	0.43
2:E:355:PHE:HE1	2:E:383:LEU:HD11	1.83	0.43
2:E:374:LEU:HD21	1:K:89:TYR:HD1	1.83	0.43
1:M:163:ILE:HG23	1:M:187:ALA:O	2.18	0.43
1:S:90:ASP:HB2	4:S:901:HOH:O	2.17	0.43
2:T:355:PHE:HE1	2:T:383:LEU:HD11	1.82	0.43
2:T:373:PRO:HG2	3:T:35:DMF:N	2.33	0.43
1:B:52:LYS:CE	4:B:2129:HOH:O	2.62	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:72:ASP:O	1:F:76:ARG:HG3	2.19	0.43
1:Q:8:SER:O	1:Q:11:GLN:HB2	2.18	0.43
4:D:725:HOH:O	1:K:112:THR:HG21	2.19	0.43
2:P:301:OZT:H17	2:P:333:LYS:HE2	1.99	0.43
1:I:11:GLN:CB	4:I:2326:HOH:O	2.58	0.43
2:J:357:ARG:NE	4:J:2224:HOH:O	2.51	0.43
2:L:422:SER:OG	2:L:432:GLU:OE2	2.30	0.43
2:N:441:SER:O	3:N:97:DMF:H13	2.19	0.43
1:U:30:VAL:HG13	1:U:43:ALA:HB2	2.00	0.43
3:B:250:DMF:H13	4:C:1685:HOH:O	2.19	0.43
2:P:496:ILE:HD13	3:P:61:DMF:H13	2.01	0.43
2:X:432:GLU:HG3	2:X:437:GLN:HB2	2.00	0.43
1:A:31:VAL:HG22	1:A:155:VAL:HG22	2.01	0.43
2:C:401:LEU:HA	2:C:402:PRO:HD3	1.89	0.43
1:F:30:VAL:HG13	1:F:43:ALA:HB2	1.99	0.43
1:I:56:LEU:HG	1:I:62:PHE:HB2	2.01	0.43
1:O:56:LEU:HG	1:O:62:PHE:HB2	2.01	0.43
1:Q:30:VAL:HG13	1:Q:43:ALA:HB2	2.00	0.43
2:X:301:OZT:H27	2:X:333:LYS:HE2	2.01	0.43
1:D:51:GLN:HG2	1:D:209:GLU:OE2	2.18	0.43
2:H:444:LEU:HD23	2:H:444:LEU:C	2.44	0.43
1:I:16:ARG:NH2	1:I:114:GLN:O	2.42	0.43
2:N:379:LYS:NZ	4:N:2250:HOH:O	2.51	0.43
2:C:376:PHE:CE2	2:C:380:ILE:HD11	2.54	0.43
1:O:72:ASP:O	1:O:76:ARG:HG3	2.19	0.43
2:T:399:LEU:CA	4:T:2342:HOH:O	2.47	0.43
1:Y:45:ASN:HA	1:Y:46:PRO:HD2	1.79	0.43
1:Y:134:LYS:N	4:Y:2396:HOH:O	2.47	0.43
1:D:98:GLN:HE21	1:D:98:GLN:HB3	1.69	0.42
1:W:97:ARG:O	1:W:101:ASN:HB2	2.19	0.42
2:G:401:LEU:HA	2:G:402:PRO:HD3	1.88	0.42
1:W:27:ALA:HB1	4:W:2259:HOH:O	2.18	0.42
1:D:15:GLU:HG2	1:K:9:PRO:HD2	2.01	0.42
1:F:144:ASP:CG	1:F:146:SER:HG	2.27	0.42
1:K:97:ARG:HH11	1:K:97:ARG:CG	2.32	0.42
2:T:301:OZT:H27	2:T:333:LYS:HE2	2.01	0.42
1:D:54:SER:OG	1:D:75:ARG:HD2	2.20	0.42
1:K:219:ARG:NH2	3:K:250:DMF:H23	2.35	0.42
1:O:77:GLY:HA3	3:O:249:DMF:HC	2.02	0.42
2:R:325:MET:SD	2:Z:444:LEU:HD21	2.59	0.42
2:X:331:VAL:HG13	2:X:349:ALA:HB2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:163:ILE:HD13	1:K:188:LEU:HA	2.02	0.42
1:U:31:VAL:HG21	1:U:188:LEU:HD11	2.01	0.42
2:Z:355:PHE:HE1	2:Z:383:LEU:HD11	1.85	0.42
2:P:433:GLU:HG2	4:P:2308:HOH:O	2.05	0.42
1:Q:56:LEU:HG	1:Q:62:PHE:HB2	2.02	0.42
2:Z:386:MET:CE	4:Z:2339:HOH:O	2.68	0.42
2:2:435:GLY:N	4:2:2018:HOH:O	2.32	0.42
1:A:17:SER:O	1:A:21:ARG:HB2	2.20	0.42
1:F:69:ASN:HB3	1:W:104:ALA:HB1	2.02	0.42
2:H:376:PHE:CE2	2:H:380:ILE:HD11	2.54	0.42
2:V:465:ARG:HD2	2:V:465:ARG:C	2.45	0.42
1:1:31:VAL:HG22	1:1:155:VAL:HG22	2.02	0.42
1:K:31:VAL:HG22	1:K:155:VAL:HG22	2.01	0.42
1:Q:13:MET:HE2	1:Q:13:MET:CA	2.50	0.42
2:R:432:GLU:OE2	2:R:434:GLU:N	2.49	0.42
2:R:477:ASP:OD2	4:R:1890:HOH:O	2.21	0.42
2:T:459:ASP:OD1	2:T:459:ASP:C	2.63	0.42
1:U:67:LYS:HG2	1:U:69:ASN:HD21	1.85	0.42
1:A:15:GLU:HA	1:A:15:GLU:OE1	2.20	0.41
2:L:482:GLY:HA3	4:L:2221:HOH:O	2.20	0.41
2:N:498:ASP:HB2	4:N:543:HOH:O	2.19	0.41
2:V:354:GLU:OE1	4:V:2432:HOH:O	2.22	0.41
1:A:67:LYS:HG2	1:A:69:ASN:HD21	1.86	0.41
1:Q:138:LEU:HD12	1:Q:138:LEU:N	2.35	0.41
1:U:90:ASP:OD1	1:U:93:ASP:N	2.53	0.41
1:U:135:ARG:HD3	4:U:1432:HOH:O	2.20	0.41
1:B:21:ARG:NH2	1:B:22:LYS:HG3	2.35	0.41
1:D:31:VAL:HG22	1:D:155:VAL:HG22	2.01	0.41
1:I:45:ASN:HA	1:I:46:PRO:HD2	1.99	0.41
4:R:1034:HOH:O	3:Z:79:DMF:H22	2.19	0.41
1:F:122:LEU:O	1:F:140:ARG:HA	2.20	0.41
2:G:447:LYS:NZ	3:G:96:DMF:H23	2.35	0.41
1:K:156:MET:HE3	1:K:156:MET:HB2	1.96	0.41
2:N:376:PHE:CE2	2:N:380:ILE:HD11	2.55	0.41
2:P:355:PHE:HE1	2:P:383:LEU:HD11	1.84	0.41
2:T:339:ASP:HB2	3:T:35:DMF:H22	2.03	0.41
1:U:31:VAL:HG22	1:U:155:VAL:HG22	2.02	0.41
2:C:301:OZT:O	2:C:440:GLY:HA3	2.21	0.41
1:I:136:PRO:HG3	4:I:2245:HOH:O	2.21	0.41
1:K:11:GLN:O	1:K:14:ARG:N	2.49	0.41
1:K:115:ALA:HB3	1:M:112:THR:HG23	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:T:391:LEU:HB3	4:T:1428:HOH:O	2.20	0.41
2:T:507:GLU:HG3	4:T:1165:HOH:O	2.20	0.41
1:B:67:LYS:HG2	1:B:69:ASN:ND2	2.35	0.41
1:K:56:LEU:HG	1:K:62:PHE:HB2	2.03	0.41
2:N:357:ARG:CD	4:N:2035:HOH:O	2.63	0.41
1:U:56:LEU:HG	1:U:62:PHE:HB2	2.03	0.41
2:V:368:LYS:NZ	4:V:2049:HOH:O	2.52	0.41
2:X:433:GLU:CG	4:Z:2252:HOH:O	2.47	0.41
1:Y:11:GLN:O	1:Y:14:ARG:HB3	2.20	0.41
1:B:76:ARG:HD3	4:B:2144:HOH:O	2.19	0.41
1:F:142:THR:OG1	1:F:146:SER:HB2	2.21	0.41
2:G:515:ARG:NE	4:G:2281:HOH:O	2.52	0.41
2:X:509:ARG:HA	2:X:509:ARG:HD2	1.91	0.41
1:Y:67:LYS:HG2	1:Y:69:ASN:HD21	1.85	0.41
1:A:190:ALA:C	4:A:2159:HOH:O	2.63	0.41
2:N:301:OZT:O	2:N:440:GLY:HA3	2.21	0.41
2:N:472:TYR:CE1	3:N:36:DMF:H22	2.55	0.41
1:O:76:ARG:O	1:O:80:GLN:HG2	2.21	0.41
1:O:230:LEU:HD12	1:O:230:LEU:HA	1.81	0.41
1:Q:231:GLN:OE1	1:Q:231:GLN:O	2.38	0.41
1:S:181:LEU:O	1:S:185:VAL:HG23	2.20	0.41
1:Y:162:PRO:HB2	1:Y:190:ALA:O	2.21	0.41
1:1:142:THR:HA	4:1:2089:HOH:O	2.20	0.41
2:2:301:OZT:O	2:2:440:GLY:HA3	2.21	0.41
1:D:45:ASN:HA	1:D:46:PRO:HD2	1.99	0.41
2:E:493:THR:CB	4:E:2369:HOH:O	2.69	0.41
1:S:80:GLN:HB2	4:S:2270:HOH:O	2.21	0.41
2:V:331:VAL:HG13	2:V:349:ALA:HB2	2.02	0.41
1:A:213:LEU:HD23	1:A:213:LEU:HA	1.99	0.40
1:O:86:GLY:O	4:O:2280:HOH:O	2.22	0.40
1:Q:56:LEU:HD23	1:Q:56:LEU:HA	1.95	0.40
1:S:44:GLU:HG3	1:S:188:LEU:CD1	2.45	0.40
2:E:412:SER:O	2:E:414:PRO:HD3	2.21	0.40
1:K:67:LYS:HG2	1:K:69:ASN:HD21	1.86	0.40
2:L:301:OZT:C7	2:L:333:LYS:NZ	2.84	0.40
1:W:217:ARG:NH2	1:W:223:ARG:HG3	2.36	0.40
2:E:432:GLU:HG3	2:E:437:GLN:HB2	2.04	0.40
1:K:90:ASP:HB2	4:K:750:HOH:O	2.21	0.40
1:Q:17:SER:OG	1:Q:143:TYR:OH	2.31	0.40
1:S:116:LYS:HD2	1:1:13:MET:SD	2.60	0.40
1:U:45:ASN:HA	1:U:46:PRO:HD2	2.00	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Y:18:GLU:CD	1:Y:21:ARG:NH2	2.58	0.40
1:Y:48:ARG:H	1:Y:48:ARG:HG2	1.62	0.40
2:G:329:ARG:O	2:G:490:ILE:HG21	2.21	0.40
1:O:189:ARG:HH22	1:O:204:GLY:N	2.18	0.40
2:L:421:VAL:HG22	2:L:431:ILE:HG12	2.03	0.40
2:R:472:TYR:CE1	3:R:68:DMF:H12	2.56	0.40
1:U:181:LEU:HD12	1:U:181:LEU:C	2.46	0.40
1:U:187:ALA:O	1:U:190:ALA:HB3	2.21	0.40
1:W:45:ASN:HA	1:W:46:PRO:HD2	2.00	0.40
1:W:59:ARG:HA	1:W:59:ARG:HD2	1.98	0.40
1:1:11:GLN:HA	1:1:14:ARG:NE	2.36	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	1	211/248 (85%)	206 (98%)	5 (2%)	0	100	100
1	A	212/248 (86%)	207 (98%)	5 (2%)	0	100	100
1	B	210/248 (85%)	206 (98%)	4 (2%)	0	100	100
1	D	209/248 (84%)	205 (98%)	4 (2%)	0	100	100
1	F	213/248 (86%)	209 (98%)	4 (2%)	0	100	100
1	I	211/248 (85%)	207 (98%)	4 (2%)	0	100	100
1	K	213/248 (86%)	208 (98%)	5 (2%)	0	100	100
1	M	212/248 (86%)	206 (97%)	6 (3%)	0	100	100
1	O	211/248 (85%)	208 (99%)	3 (1%)	0	100	100
1	Q	214/248 (86%)	207 (97%)	7 (3%)	0	100	100
1	S	211/248 (85%)	208 (99%)	3 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	U	211/248 (85%)	208 (99%)	3 (1%)	0	100	100
1	W	213/248 (86%)	209 (98%)	4 (2%)	0	100	100
1	Y	209/248 (84%)	205 (98%)	4 (2%)	0	100	100
2	2	211/240 (88%)	210 (100%)	1 (0%)	0	100	100
2	C	211/240 (88%)	210 (100%)	1 (0%)	0	100	100
2	E	212/240 (88%)	211 (100%)	1 (0%)	0	100	100
2	G	212/240 (88%)	211 (100%)	1 (0%)	0	100	100
2	H	209/240 (87%)	208 (100%)	1 (0%)	0	100	100
2	J	212/240 (88%)	211 (100%)	1 (0%)	0	100	100
2	L	212/240 (88%)	211 (100%)	1 (0%)	0	100	100
2	N	209/240 (87%)	209 (100%)	0	0	100	100
2	P	212/240 (88%)	211 (100%)	1 (0%)	0	100	100
2	R	219/240 (91%)	217 (99%)	2 (1%)	0	100	100
2	T	212/240 (88%)	211 (100%)	1 (0%)	0	100	100
2	V	217/240 (90%)	214 (99%)	3 (1%)	0	100	100
2	X	212/240 (88%)	210 (99%)	2 (1%)	0	100	100
2	Z	211/240 (88%)	210 (100%)	1 (0%)	0	100	100
All	All	5931/6832 (87%)	5853 (99%)	78 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	1	164/192 (85%)	155 (94%)	9 (6%)	19	26
1	A	165/192 (86%)	158 (96%)	7 (4%)	26	37
1	B	164/192 (85%)	157 (96%)	7 (4%)	26	36
1	D	163/192 (85%)	154 (94%)	9 (6%)	19	26

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	166/192 (86%)	156 (94%)	10 (6%)	17	23
1	I	164/192 (85%)	159 (97%)	5 (3%)	36	48
1	K	166/192 (86%)	156 (94%)	10 (6%)	17	23
1	M	165/192 (86%)	160 (97%)	5 (3%)	36	48
1	O	164/192 (85%)	158 (96%)	6 (4%)	30	41
1	Q	167/192 (87%)	158 (95%)	9 (5%)	20	27
1	S	164/192 (85%)	157 (96%)	7 (4%)	26	36
1	U	164/192 (85%)	157 (96%)	7 (4%)	26	36
1	W	166/192 (86%)	162 (98%)	4 (2%)	43	56
1	Y	163/192 (85%)	158 (97%)	5 (3%)	35	47
2	2	160/177 (90%)	157 (98%)	3 (2%)	50	63
2	C	160/177 (90%)	156 (98%)	4 (2%)	42	54
2	E	161/177 (91%)	156 (97%)	5 (3%)	35	47
2	G	161/177 (91%)	158 (98%)	3 (2%)	50	63
2	H	160/177 (90%)	154 (96%)	6 (4%)	29	41
2	J	161/177 (91%)	158 (98%)	3 (2%)	50	63
2	L	161/177 (91%)	154 (96%)	7 (4%)	26	36
2	N	159/177 (90%)	155 (98%)	4 (2%)	42	54
2	P	161/177 (91%)	156 (97%)	5 (3%)	35	47
2	R	165/177 (93%)	161 (98%)	4 (2%)	43	56
2	T	161/177 (91%)	158 (98%)	3 (2%)	50	63
2	V	163/177 (92%)	159 (98%)	4 (2%)	42	54
2	X	161/177 (91%)	157 (98%)	4 (2%)	42	54
2	Z	160/177 (90%)	156 (98%)	4 (2%)	42	54
All	All	4559/5166 (88%)	4400 (96%)	159 (4%)	32	44

All (159) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	MET
1	A	19	LEU
1	A	31	VAL
1	A	69	ASN
1	A	113	GLU

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Mol	Chain	Res	Type
1	A	161	GLU
1	A	233	LEU
1	B	13	MET
1	B	69	ASN
1	B	113	GLU
1	B	122	LEU
1	B	161	GLU
1	B	205	VAL
1	B	234	LEU
2	C	362	GLU
2	C	444	LEU
2	C	492	PRO
2	C	520	SER
1	D	10	GLU
1	D	11	GLN
1	D	13	MET
1	D	14	ARG
1	D	15	GLU
1	D	31	VAL
1	D	69	ASN
1	D	161	GLU
1	D	174	ASN
2	E	362	GLU
2	E	399	LEU
2	E	444	LEU
2	E	464	LEU
2	E	492	PRO
1	F	13	MET
1	F	14	ARG
1	F	15	GLU
1	F	31	VAL
1	F	52	LYS
1	F	69	ASN
1	F	112	THR
1	F	113	GLU
1	F	161	GLU
1	F	205	VAL
2	G	362	GLU
2	G	444	LEU
2	G	492	PRO
2	H	330	ASP
2	H	338	ASP

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Mol	Chain	Res	Type
2	H	362	GLU
2	H	412	SER
2	H	415	GLN
2	H	520	SER
1	I	13	MET
1	I	69	ASN
1	I	113	GLU
1	I	122	LEU
1	I	161	GLU
2	J	362	GLU
2	J	399	LEU
2	J	444	LEU
1	K	7	ILE
1	K	11	GLN
1	K	13	MET
1	K	69	ASN
1	K	97	ARG
1	K	112	THR
1	K	113	GLU
1	K	122	LEU
1	K	161	GLU
1	K	169	GLU
2	L	306	LEU
2	L	338	ASP
2	L	399	LEU
2	L	401	LEU
2	L	444	LEU
2	L	492	PRO
2	L	509	ARG
1	M	7	ILE
1	M	13	MET
1	M	69	ASN
1	M	113	GLU
1	M	161	GLU
2	N	412	SER
2	N	444	LEU
2	N	492	PRO
2	N	522	SER
1	O	13	MET
1	O	31	VAL
1	O	69	ASN
1	O	161	GLU

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Mol	Chain	Res	Type
1	O	173	GLU
1	O	234	LEU
2	P	362	GLU
2	P	399	LEU
2	P	434	GLU
2	P	492	PRO
2	P	509	ARG
1	Q	11	GLN
1	Q	13	MET
1	Q	14	ARG
1	Q	19	LEU
1	Q	31	VAL
1	Q	69	ASN
1	Q	113	GLU
1	Q	203	LEU
1	Q	231	GLN
2	R	399	LEU
2	R	404	LEU
2	R	444	LEU
2	R	465	ARG
1	S	13	MET
1	S	69	ASN
1	S	97	ARG
1	S	161	GLU
1	S	169	GLU
1	S	233	LEU
1	S	234	LEU
2	T	330	ASP
2	T	444	LEU
2	T	492	PRO
1	U	13	MET
1	U	31	VAL
1	U	69	ASN
1	U	99	LEU
1	U	113	GLU
1	U	161	GLU
1	U	174	ASN
2	V	330	ASP
2	V	444	LEU
2	V	492	PRO
2	V	496	ILE
1	W	13	MET

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Mol	Chain	Res	Type
1	W	69	ASN
1	W	97	ARG
1	W	161	GLU
2	X	362	GLU
2	X	399	LEU
2	X	444	LEU
2	X	492	PRO
1	Y	31	VAL
1	Y	49	SER
1	Y	69	ASN
1	Y	113	GLU
1	Y	161	GLU
2	Z	338	ASP
2	Z	362	GLU
2	Z	444	LEU
2	Z	492	PRO
1	1	13	MET
1	1	14	ARG
1	1	17	SER
1	1	69	ASN
1	1	99	LEU
1	1	112	THR
1	1	113	GLU
1	1	161	GLU
1	1	234	LEU
2	2	433	GLU
2	2	444	LEU
2	2	492	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (96) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	ASN
1	A	98	GLN
1	A	165	ASN
1	B	69	ASN
1	B	98	GLN
1	B	101	ASN
1	B	114	GLN
1	B	129	HIS
1	B	165	ASN
1	B	174	ASN

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Mol	Chain	Res	Type
2	C	390	ASN
2	C	430	ASN
1	D	11	GLN
1	D	69	ASN
1	D	98	GLN
1	D	114	GLN
1	D	129	HIS
1	D	165	ASN
2	E	390	ASN
2	E	430	ASN
1	F	51	GLN
1	F	69	ASN
1	F	98	GLN
1	F	114	GLN
1	F	129	HIS
1	F	165	ASN
2	G	390	ASN
2	H	390	ASN
1	I	45	ASN
1	I	69	ASN
1	I	80	GLN
1	I	98	GLN
1	I	129	HIS
1	I	152	HIS
1	I	165	ASN
2	J	390	ASN
1	K	11	GLN
1	K	69	ASN
1	K	98	GLN
1	K	129	HIS
1	K	165	ASN
2	L	390	ASN
2	L	430	ASN
1	M	69	ASN
1	M	98	GLN
1	M	114	GLN
1	M	129	HIS
1	M	165	ASN
1	M	174	ASN
2	N	390	ASN
2	N	415	GLN
1	O	69	ASN

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Mol	Chain	Res	Type
1	O	98	GLN
1	O	114	GLN
1	O	129	HIS
1	O	165	ASN
2	P	390	ASN
2	P	415	GLN
2	P	430	ASN
1	Q	69	ASN
1	Q	98	GLN
1	Q	114	GLN
1	Q	129	HIS
1	Q	165	ASN
2	R	390	ASN
2	R	430	ASN
1	S	69	ASN
1	S	80	GLN
1	S	165	ASN
2	T	390	ASN
2	T	430	ASN
1	U	69	ASN
1	U	80	GLN
1	U	129	HIS
1	U	165	ASN
2	V	430	ASN
1	W	45	ASN
1	W	69	ASN
1	W	98	GLN
1	W	165	ASN
2	X	390	ASN
2	X	415	GLN
1	Y	69	ASN
1	Y	73	ASN
1	Y	80	GLN
1	Y	98	GLN
1	Y	165	ASN
1	Y	231	GLN
2	Z	390	ASN
2	Z	430	ASN
1	1	69	ASN
1	1	80	GLN
1	1	98	GLN
1	1	114	GLN

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Mol	Chain	Res	Type
1	1	165	ASN
2	2	390	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

14 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	OZT	2	301	2	6,9,10	5.15	5 (83%)	9,12,14	5.91	6 (66%)
2	OZT	R	301	2	6,9,10	5.39	4 (66%)	9,12,14	5.84	6 (66%)
2	OZT	C	301	2	6,9,10	5.10	3 (50%)	9,12,14	6.10	6 (66%)
2	OZT	P	301	2	6,9,10	4.30	3 (50%)	9,12,14	5.85	6 (66%)
2	OZT	J	301	2	6,9,10	5.25	3 (50%)	9,12,14	5.99	6 (66%)
2	OZT	T	301	2	6,9,10	4.75	3 (50%)	9,12,14	6.11	6 (66%)
2	OZT	V	301	2	6,9,10	4.66	3 (50%)	9,12,14	6.21	7 (77%)
2	OZT	H	301	2	6,9,10	4.90	3 (50%)	9,12,14	6.14	6 (66%)
2	OZT	X	301	2	6,9,10	5.36	4 (66%)	9,12,14	6.02	7 (77%)
2	OZT	G	301	2	6,9,10	4.49	3 (50%)	9,12,14	5.96	7 (77%)
2	OZT	L	301	2	6,9,10	5.19	3 (50%)	9,12,14	6.03	6 (66%)
2	OZT	N	301	2	6,9,10	4.85	3 (50%)	9,12,14	5.90	6 (66%)
2	OZT	E	301	2	6,9,10	4.58	3 (50%)	9,12,14	5.72	7 (77%)
2	OZT	Z	301	2	6,9,10	4.79	3 (50%)	9,12,14	6.01	6 (66%)

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	OZT	2	301	2	-	1/1/14/16	0/1/1/1
2	OZT	R	301	2	-	1/1/14/16	0/1/1/1
2	OZT	C	301	2	-	1/1/14/16	0/1/1/1
2	OZT	P	301	2	-	1/1/14/16	0/1/1/1
2	OZT	J	301	2	-	1/1/14/16	0/1/1/1
2	OZT	T	301	2	-	1/1/14/16	0/1/1/1
2	OZT	V	301	2	-	0/1/14/16	0/1/1/1
2	OZT	H	301	2	-	1/1/14/16	0/1/1/1
2	OZT	X	301	2	-	0/1/14/16	0/1/1/1
2	OZT	G	301	2	-	1/1/14/16	0/1/1/1
2	OZT	L	301	2	-	0/1/14/16	0/1/1/1
2	OZT	N	301	2	-	1/1/14/16	0/1/1/1
2	OZT	E	301	2	-	1/1/14/16	0/1/1/1
2	OZT	Z	301	2	-	1/1/14/16	0/1/1/1

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	J	301	OZT	O1-C5	11.04	1.51	1.36
2	X	301	OZT	O1-C5	10.77	1.50	1.36
2	L	301	OZT	O1-C5	10.51	1.50	1.36
2	T	301	OZT	O1-C5	10.09	1.49	1.36
2	R	301	OZT	O1-C5	9.72	1.49	1.36
2	H	301	OZT	O1-C5	9.68	1.49	1.36
2	C	301	OZT	O1-C5	9.53	1.49	1.36
2	E	301	OZT	O1-C5	9.41	1.48	1.36
2	2	301	OZT	O1-C5	9.37	1.48	1.36
2	G	301	OZT	O1-C5	9.26	1.48	1.36
2	Z	301	OZT	O1-C5	9.21	1.48	1.36
2	V	301	OZT	O1-C5	8.65	1.47	1.36
2	N	301	OZT	O1-C5	8.54	1.47	1.36
2	P	301	OZT	O1-C5	8.37	1.47	1.36
2	R	301	OZT	C5-N	7.10	1.43	1.33
2	C	301	OZT	C5-N	6.66	1.42	1.33
2	N	301	OZT	C5-N	6.55	1.42	1.33
2	2	301	OZT	C5-N	6.25	1.42	1.33
2	X	301	OZT	C5-N	5.84	1.41	1.33
2	H	301	OZT	C5-N	5.80	1.41	1.33
2	Z	301	OZT	C5-N	5.36	1.40	1.33
2	L	301	OZT	C5-N	5.27	1.40	1.33
2	J	301	OZT	C5-N	5.11	1.40	1.33
2	V	301	OZT	O1-C2	-5.08	1.38	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	V	301	OZT	C5-N	4.92	1.40	1.33
2	P	301	OZT	C5-N	4.81	1.40	1.33
2	N	301	OZT	O1-C2	-4.73	1.39	1.46
2	G	301	OZT	C5-N	4.63	1.39	1.33
2	2	301	OZT	O1-C2	-4.57	1.39	1.46
2	E	301	OZT	C5-N	4.42	1.39	1.33
2	R	301	OZT	O1-C2	-4.39	1.40	1.46
2	Z	301	OZT	O1-C2	-4.37	1.40	1.46
2	L	301	OZT	O1-C2	-4.23	1.40	1.46
2	T	301	OZT	C5-N	4.01	1.39	1.33
2	X	301	OZT	O1-C2	-4.00	1.40	1.46
2	T	301	OZT	O1-C2	-3.73	1.41	1.46
2	P	301	OZT	O1-C2	-3.69	1.41	1.46
2	J	301	OZT	O1-C2	-3.65	1.41	1.46
2	C	301	OZT	O1-C2	-3.64	1.41	1.46
2	E	301	OZT	O1-C2	-3.58	1.41	1.46
2	H	301	OZT	O1-C2	-3.53	1.41	1.46
2	G	301	OZT	O1-C2	-3.48	1.41	1.46
2	R	301	OZT	O6-C5	2.44	1.26	1.21
2	2	301	OZT	O6-C5	2.38	1.26	1.21
2	X	301	OZT	C7-C2	-2.28	1.46	1.51
2	2	301	OZT	C7-C2	-2.24	1.46	1.51

All (88) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	301	OZT	O1-C5-N	9.36	117.50	109.87
2	T	301	OZT	O1-C5-N	9.26	117.41	109.87
2	N	301	OZT	O1-C5-N	9.20	117.37	109.87
2	V	301	OZT	O1-C2-CA	9.16	112.05	103.72
2	P	301	OZT	O1-C5-N	9.13	117.31	109.87
2	H	301	OZT	O1-C5-N	9.02	117.22	109.87
2	C	301	OZT	O1-C5-N	8.97	117.17	109.87
2	X	301	OZT	O1-C5-N	8.94	117.15	109.87
2	J	301	OZT	O1-C5-N	8.94	117.15	109.87
2	R	301	OZT	CA-N-C5	-8.89	99.02	112.75
2	Z	301	OZT	CA-N-C5	-8.87	99.04	112.75
2	Z	301	OZT	O1-C5-N	8.80	117.04	109.87
2	L	301	OZT	O1-C5-N	8.67	116.93	109.87
2	V	301	OZT	O1-C5-N	8.64	116.91	109.87
2	C	301	OZT	CA-N-C5	-8.48	99.64	112.75
2	X	301	OZT	CA-N-C5	-8.47	99.66	112.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	2	301	OZT	CA-N-C5	-8.40	99.77	112.75
2	L	301	OZT	O1-C2-CA	8.37	111.33	103.72
2	R	301	OZT	O1-C5-N	8.30	116.63	109.87
2	E	301	OZT	O1-C5-N	8.29	116.62	109.87
2	2	301	OZT	O1-C5-N	8.28	116.61	109.87
2	N	301	OZT	CA-N-C5	-8.23	100.03	112.75
2	H	301	OZT	CA-N-C5	-8.13	100.19	112.75
2	G	301	OZT	CA-N-C5	-8.09	100.25	112.75
2	V	301	OZT	CA-N-C5	-8.01	100.37	112.75
2	J	301	OZT	CA-N-C5	-7.97	100.44	112.75
2	T	301	OZT	CA-N-C5	-7.88	100.57	112.75
2	H	301	OZT	O1-C2-CA	7.83	110.84	103.72
2	E	301	OZT	CA-N-C5	-7.80	100.69	112.75
2	L	301	OZT	CA-N-C5	-7.80	100.69	112.75
2	L	301	OZT	C2-O1-C5	-7.73	98.87	110.33
2	2	301	OZT	O1-C2-CA	7.71	110.73	103.72
2	X	301	OZT	O1-C2-CA	7.69	110.71	103.72
2	P	301	OZT	CA-N-C5	-7.60	101.01	112.75
2	P	301	OZT	O1-C2-CA	7.57	110.61	103.72
2	J	301	OZT	C2-O1-C5	-7.56	99.12	110.33
2	H	301	OZT	C2-O1-C5	-7.55	99.13	110.33
2	T	301	OZT	C2-O1-C5	-7.52	99.18	110.33
2	V	301	OZT	C2-O1-C5	-7.52	99.19	110.33
2	X	301	OZT	C2-O1-C5	-7.49	99.23	110.33
2	P	301	OZT	C2-O1-C5	-7.48	99.24	110.33
2	N	301	OZT	O1-C2-CA	7.45	110.49	103.72
2	R	301	OZT	O1-C2-CA	7.32	110.37	103.72
2	C	301	OZT	O1-C2-CA	7.30	110.36	103.72
2	C	301	OZT	C2-O1-C5	-7.29	99.52	110.33
2	J	301	OZT	O1-C2-CA	7.27	110.33	103.72
2	G	301	OZT	C2-O1-C5	-7.24	99.60	110.33
2	T	301	OZT	O6-C5-N	-7.20	120.67	129.19
2	2	301	OZT	C2-O1-C5	-7.16	99.72	110.33
2	E	301	OZT	O1-C2-CA	7.11	110.19	103.72
2	Z	301	OZT	C2-O1-C5	-7.11	99.80	110.33
2	Z	301	OZT	O1-C2-CA	7.06	110.14	103.72
2	J	301	OZT	O6-C5-N	-7.05	120.84	129.19
2	T	301	OZT	O1-C2-CA	7.04	110.12	103.72
2	E	301	OZT	C2-O1-C5	-6.88	100.13	110.33
2	N	301	OZT	C2-O1-C5	-6.85	100.17	110.33
2	R	301	OZT	C2-O1-C5	-6.83	100.21	110.33
2	G	301	OZT	O1-C2-CA	6.73	109.84	103.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	301	OZT	O6-C5-N	-6.65	121.32	129.19
2	L	301	OZT	O6-C5-N	-6.58	121.41	129.19
2	E	301	OZT	O6-C5-N	-6.42	121.60	129.19
2	H	301	OZT	O6-C5-N	-6.30	121.73	129.19
2	P	301	OZT	O6-C5-N	-6.25	121.79	129.19
2	X	301	OZT	O6-C5-N	-6.25	121.80	129.19
2	Z	301	OZT	O6-C5-N	-6.08	122.00	129.19
2	C	301	OZT	O6-C5-N	-6.03	122.05	129.19
2	C	301	OZT	C2-CA-C	-5.93	105.92	114.16
2	V	301	OZT	O6-C5-N	-5.83	122.30	129.19
2	2	301	OZT	C2-CA-C	-5.61	106.36	114.16
2	H	301	OZT	C2-CA-C	-5.57	106.42	114.16
2	N	301	OZT	O6-C5-N	-5.54	122.63	129.19
2	2	301	OZT	O6-C5-N	-5.47	122.72	129.19
2	V	301	OZT	C2-CA-C	-5.37	106.69	114.16
2	R	301	OZT	C2-CA-C	-5.29	106.81	114.16
2	Z	301	OZT	C2-CA-C	-5.19	106.94	114.16
2	T	301	OZT	C2-CA-C	-5.17	106.98	114.16
2	R	301	OZT	O6-C5-N	-5.11	123.14	129.19
2	N	301	OZT	C2-CA-C	-4.91	107.33	114.16
2	E	301	OZT	C2-CA-C	-4.27	108.22	114.16
2	G	301	OZT	C2-CA-C	-4.21	108.30	114.16
2	L	301	OZT	C2-CA-C	-3.96	108.65	114.16
2	J	301	OZT	C2-CA-C	-3.94	108.68	114.16
2	X	301	OZT	C2-CA-C	-3.85	108.80	114.16
2	P	301	OZT	C2-CA-C	-3.38	109.45	114.16
2	G	301	OZT	O-C-CA	-2.33	118.69	124.86
2	E	301	OZT	O-C-CA	-2.31	118.76	124.86
2	X	301	OZT	O-C-CA	-2.22	118.98	124.86
2	V	301	OZT	O-C-CA	-2.12	119.25	124.86

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	301	OZT	O-C-CA-C2
2	E	301	OZT	O-C-CA-C2
2	G	301	OZT	O-C-CA-C2
2	H	301	OZT	O-C-CA-C2
2	J	301	OZT	O-C-CA-C2
2	N	301	OZT	O-C-CA-C2
2	P	301	OZT	O-C-CA-C2

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Mol	Chain	Res	Type	Atoms
2	R	301	OZT	O-C-CA-C2
2	T	301	OZT	O-C-CA-C2
2	Z	301	OZT	O-C-CA-C2
2	2	301	OZT	O-C-CA-C2

There are no ring outliers.

11 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	2	301	OZT	3	0
2	R	301	OZT	3	0
2	C	301	OZT	2	0
2	P	301	OZT	4	0
2	T	301	OZT	1	0
2	V	301	OZT	2	0
2	H	301	OZT	2	0
2	X	301	OZT	1	0
2	L	301	OZT	1	0
2	N	301	OZT	3	0
2	Z	301	OZT	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

102 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	DMF	O	250	-	4,4,4	0.30	0	4,4,4	0.47	0
3	DMF	I	250	-	4,4,4	0.47	0	4,4,4	0.27	0
3	DMF	B	251	-	4,4,4	0.69	0	4,4,4	0.23	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	DMF	R	3	-	4,4,4	0.26	0	4,4,4	0.45	0
3	DMF	C	48	-	4,4,4	0.48	0	4,4,4	0.29	0
3	DMF	H	53	-	4,4,4	0.33	0	4,4,4	0.38	0
3	DMF	H	52	-	4,4,4	0.36	0	4,4,4	0.55	0
3	DMF	N	21	-	4,4,4	0.25	0	4,4,4	0.43	0
3	DMF	H	41	-	4,4,4	0.55	0	4,4,4	0.37	0
3	DMF	P	85	-	4,4,4	0.84	0	4,4,4	0.39	0
3	DMF	R	54	-	4,4,4	0.38	0	4,4,4	0.43	0
3	DMF	G	39	-	4,4,4	0.72	0	4,4,4	0.35	0
3	DMF	W	249	-	4,4,4	0.34	0	4,4,4	0.39	0
3	DMF	2	78	-	4,4,4	0.21	0	4,4,4	0.41	0
3	DMF	Z	71	-	4,4,4	0.48	0	4,4,4	0.35	0
3	DMF	T	92	-	4,4,4	0.85	0	4,4,4	0.37	0
3	DMF	U	249	-	4,4,4	0.53	0	4,4,4	0.24	0
3	DMF	Q	250	-	4,4,4	0.30	0	4,4,4	0.54	0
3	DMF	Z	27	-	4,4,4	0.15	0	4,4,4	0.39	0
3	DMF	D	251	-	4,4,4	0.45	0	4,4,4	0.49	0
3	DMF	X	67	-	4,4,4	0.72	0	4,4,4	0.32	0
3	DMF	E	70	-	4,4,4	0.33	0	4,4,4	0.54	0
3	DMF	H	19	-	4,4,4	0.49	0	4,4,4	0.30	0
3	DMF	Y	249	-	4,4,4	0.66	0	4,4,4	0.46	0
3	DMF	V	90	-	4,4,4	0.25	0	4,4,4	0.36	0
3	DMF	E	83	-	4,4,4	0.29	0	4,4,4	0.75	0
3	DMF	N	99	-	4,4,4	0.92	0	4,4,4	0.47	0
3	DMF	W	250	-	4,4,4	0.72	0	4,4,4	0.32	0
3	DMF	P	61	-	4,4,4	0.75	0	4,4,4	0.30	0
3	DMF	F	249	-	4,4,4	0.54	0	4,4,4	0.45	0
3	DMF	C	28	-	4,4,4	0.55	0	4,4,4	0.30	0
3	DMF	L	4	-	4,4,4	0.43	0	4,4,4	0.30	0
3	DMF	J	82	-	4,4,4	0.52	0	4,4,4	0.34	0
3	DMF	A	249	-	4,4,4	0.57	0	4,4,4	0.36	0
3	DMF	Z	69	-	4,4,4	0.58	0	4,4,4	0.31	0
3	DMF	P	74	-	4,4,4	0.40	0	4,4,4	0.19	0
3	DMF	O	249	-	4,4,4	0.27	0	4,4,4	0.67	0
3	DMF	M	249	-	4,4,4	0.58	0	4,4,4	0.41	0
3	DMF	P	23	-	4,4,4	0.21	0	4,4,4	0.37	0
3	DMF	P	100	-	4,4,4	0.50	0	4,4,4	0.34	0
3	DMF	2	58	-	4,4,4	0.79	0	4,4,4	0.31	0
3	DMF	1	249	-	4,4,4	0.53	0	4,4,4	0.41	0
3	DMF	B	250	-	4,4,4	0.87	0	4,4,4	0.29	0
3	DMF	K	249	-	4,4,4	0.72	0	4,4,4	0.28	0
3	DMF	J	6	-	4,4,4	0.27	0	4,4,4	0.43	0
3	DMF	F	250	-	4,4,4	0.34	0	4,4,4	0.41	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	DMF	T	46	-	4,4,4	0.53	0	4,4,4	0.37	0
3	DMF	E	44	-	4,4,4	0.29	0	4,4,4	0.54	0
3	DMF	S	249	-	4,4,4	0.59	0	4,4,4	0.21	0
3	DMF	R	68	-	4,4,4	0.26	0	4,4,4	0.52	0
3	DMF	C	64	-	4,4,4	0.27	0	4,4,4	0.46	0
3	DMF	D	252	-	4,4,4	0.24	0	4,4,4	0.37	0
3	DMF	C	17	-	4,4,4	0.26	0	4,4,4	0.42	0
3	DMF	H	102	-	4,4,4	0.48	0	4,4,4	0.40	0
3	DMF	A	250	-	4,4,4	0.64	0	4,4,4	0.46	0
3	DMF	L	9	-	4,4,4	0.21	0	4,4,4	0.46	0
3	DMF	U	250	-	4,4,4	0.30	0	4,4,4	0.49	0
3	DMF	C	5	-	4,4,4	0.32	0	4,4,4	0.57	0
3	DMF	G	1	-	4,4,4	0.36	0	4,4,4	0.60	0
3	DMF	N	36	-	4,4,4	0.33	0	4,4,4	0.46	0
3	DMF	V	65	-	4,4,4	0.89	0	4,4,4	0.29	0
3	DMF	C	43	-	4,4,4	0.94	0	4,4,4	0.36	0
3	DMF	T	62	-	4,4,4	0.47	0	4,4,4	0.32	0
3	DMF	N	2	-	4,4,4	0.17	0	4,4,4	0.43	0
3	DMF	I	249	-	4,4,4	0.74	0	4,4,4	0.39	0
3	DMF	X	25	-	4,4,4	0.59	0	4,4,4	0.47	0
3	DMF	E	95	-	4,4,4	1.11	0	4,4,4	0.43	0
3	DMF	D	249	-	4,4,4	0.29	0	4,4,4	0.34	0
3	DMF	J	59	-	4,4,4	0.23	0	4,4,4	0.39	0
3	DMF	N	37	-	4,4,4	0.68	0	4,4,4	0.41	0
3	DMF	Q	249	-	4,4,4	0.44	0	4,4,4	0.36	0
3	DMF	C	15	-	4,4,4	0.39	0	4,4,4	0.43	0
3	DMF	C	73	-	4,4,4	0.27	0	4,4,4	0.36	0
3	DMF	T	35	-	4,4,4	0.34	0	4,4,4	0.39	0
3	DMF	1	250	-	4,4,4	0.36	0	4,4,4	0.28	0
3	DMF	K	250	-	4,4,4	0.58	0	4,4,4	0.36	0
3	DMF	U	252	-	4,4,4	0.85	0	4,4,4	0.37	0
3	DMF	V	86	-	4,4,4	0.28	0	4,4,4	0.49	0
3	DMF	1	251	-	4,4,4	0.47	0	4,4,4	0.63	0
3	DMF	T	94	-	4,4,4	0.53	0	4,4,4	0.37	0
3	DMF	Z	50	-	4,4,4	0.24	0	4,4,4	0.42	0
3	DMF	G	96	-	4,4,4	0.79	0	4,4,4	0.32	0
3	DMF	L	60	-	4,4,4	0.47	0	4,4,4	0.32	0
3	DMF	2	66	-	4,4,4	0.73	0	4,4,4	0.28	0
3	DMF	E	32	-	4,4,4	0.71	0	4,4,4	0.35	0
3	DMF	N	97	-	4,4,4	0.90	0	4,4,4	0.45	0
3	DMF	E	55	-	4,4,4	0.44	0	4,4,4	0.35	0
3	DMF	S	251	-	4,4,4	0.31	0	4,4,4	0.42	0
3	DMF	G	20	-	4,4,4	0.51	0	4,4,4	0.35	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	DMF	N	24	-	4,4,4	0.46	0	4,4,4	0.34	0
3	DMF	M	250	-	4,4,4	0.73	0	4,4,4	0.53	0
3	DMF	V	16	-	4,4,4	0.36	0	4,4,4	0.57	0
3	DMF	H	31	-	4,4,4	0.86	0	4,4,4	0.32	0
3	DMF	R	89	-	4,4,4	0.30	0	4,4,4	0.57	0
3	DMF	Z	91	-	4,4,4	0.31	0	4,4,4	0.59	0
3	DMF	S	250	-	4,4,4	0.29	0	4,4,4	0.43	0
3	DMF	H	14	-	4,4,4	0.58	0	4,4,4	0.32	0
3	DMF	B	249	-	4,4,4	0.19	0	4,4,4	0.42	0
3	DMF	U	251	-	4,4,4	0.88	0	4,4,4	0.29	0
3	DMF	D	250	-	4,4,4	0.68	0	4,4,4	0.33	0
3	DMF	Z	79	-	4,4,4	0.88	0	4,4,4	0.47	0
3	DMF	X	88	-	4,4,4	0.47	0	4,4,4	0.47	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	DMF	O	250	-	-	2/2/2/2	-
3	DMF	I	250	-	-	0/2/2/2	-
3	DMF	B	251	-	-	0/2/2/2	-
3	DMF	R	3	-	-	2/2/2/2	-
3	DMF	C	48	-	-	0/2/2/2	-
3	DMF	H	53	-	-	2/2/2/2	-
3	DMF	H	52	-	-	2/2/2/2	-
3	DMF	N	21	-	-	2/2/2/2	-
3	DMF	H	41	-	-	0/2/2/2	-
3	DMF	P	85	-	-	0/2/2/2	-
3	DMF	R	54	-	-	0/2/2/2	-
3	DMF	G	39	-	-	0/2/2/2	-
3	DMF	W	249	-	-	2/2/2/2	-
3	DMF	2	78	-	-	0/2/2/2	-
3	DMF	Z	71	-	-	0/2/2/2	-
3	DMF	T	92	-	-	0/2/2/2	-
3	DMF	U	249	-	-	0/2/2/2	-
3	DMF	Q	250	-	-	2/2/2/2	-
3	DMF	Z	27	-	-	2/2/2/2	-
3	DMF	D	251	-	-	0/2/2/2	-
3	DMF	X	67	-	-	0/2/2/2	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	DMF	E	70	-	-	2/2/2/2	-
3	DMF	H	19	-	-	0/2/2/2	-
3	DMF	Y	249	-	-	0/2/2/2	-
3	DMF	V	90	-	-	2/2/2/2	-
3	DMF	E	83	-	-	2/2/2/2	-
3	DMF	N	99	-	-	0/2/2/2	-
3	DMF	W	250	-	-	0/2/2/2	-
3	DMF	P	61	-	-	0/2/2/2	-
3	DMF	F	249	-	-	0/2/2/2	-
3	DMF	C	28	-	-	0/2/2/2	-
3	DMF	L	4	-	-	0/2/2/2	-
3	DMF	J	82	-	-	0/2/2/2	-
3	DMF	A	249	-	-	0/2/2/2	-
3	DMF	Z	69	-	-	0/2/2/2	-
3	DMF	P	74	-	-	0/2/2/2	-
3	DMF	O	249	-	-	2/2/2/2	-
3	DMF	M	249	-	-	0/2/2/2	-
3	DMF	P	23	-	-	0/2/2/2	-
3	DMF	P	100	-	-	0/2/2/2	-
3	DMF	2	58	-	-	0/2/2/2	-
3	DMF	1	249	-	-	0/2/2/2	-
3	DMF	B	250	-	-	0/2/2/2	-
3	DMF	K	249	-	-	0/2/2/2	-
3	DMF	J	6	-	-	2/2/2/2	-
3	DMF	F	250	-	-	2/2/2/2	-
3	DMF	T	46	-	-	0/2/2/2	-
3	DMF	E	44	-	-	2/2/2/2	-
3	DMF	S	249	-	-	0/2/2/2	-
3	DMF	R	68	-	-	2/2/2/2	-
3	DMF	C	64	-	-	2/2/2/2	-
3	DMF	D	252	-	-	2/2/2/2	-
3	DMF	C	17	-	-	0/2/2/2	-
3	DMF	H	102	-	-	0/2/2/2	-
3	DMF	A	250	-	-	0/2/2/2	-
3	DMF	L	9	-	-	2/2/2/2	-
3	DMF	U	250	-	-	2/2/2/2	-
3	DMF	C	5	-	-	2/2/2/2	-
3	DMF	G	1	-	-	2/2/2/2	-
3	DMF	N	36	-	-	0/2/2/2	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	DMF	V	65	-	-	0/2/2/2	-
3	DMF	C	43	-	-	0/2/2/2	-
3	DMF	T	62	-	-	0/2/2/2	-
3	DMF	N	2	-	-	2/2/2/2	-
3	DMF	I	249	-	-	0/2/2/2	-
3	DMF	X	25	-	-	0/2/2/2	-
3	DMF	E	95	-	-	0/2/2/2	-
3	DMF	D	249	-	-	2/2/2/2	-
3	DMF	J	59	-	-	2/2/2/2	-
3	DMF	N	37	-	-	0/2/2/2	-
3	DMF	Q	249	-	-	0/2/2/2	-
3	DMF	C	15	-	-	0/2/2/2	-
3	DMF	C	73	-	-	2/2/2/2	-
3	DMF	T	35	-	-	2/2/2/2	-
3	DMF	1	250	-	-	0/2/2/2	-
3	DMF	K	250	-	-	0/2/2/2	-
3	DMF	U	252	-	-	0/2/2/2	-
3	DMF	V	86	-	-	0/2/2/2	-
3	DMF	1	251	-	-	0/2/2/2	-
3	DMF	T	94	-	-	0/2/2/2	-
3	DMF	Z	50	-	-	0/2/2/2	-
3	DMF	G	96	-	-	0/2/2/2	-
3	DMF	L	60	-	-	0/2/2/2	-
3	DMF	2	66	-	-	0/2/2/2	-
3	DMF	E	32	-	-	0/2/2/2	-
3	DMF	N	97	-	-	0/2/2/2	-
3	DMF	E	55	-	-	0/2/2/2	-
3	DMF	S	251	-	-	2/2/2/2	-
3	DMF	G	20	-	-	0/2/2/2	-
3	DMF	N	24	-	-	0/2/2/2	-
3	DMF	M	250	-	-	0/2/2/2	-
3	DMF	V	16	-	-	2/2/2/2	-
3	DMF	H	31	-	-	0/2/2/2	-
3	DMF	R	89	-	-	2/2/2/2	-
3	DMF	Z	91	-	-	2/2/2/2	-
3	DMF	S	250	-	-	2/2/2/2	-
3	DMF	H	14	-	-	0/2/2/2	-
3	DMF	B	249	-	-	2/2/2/2	-
3	DMF	U	251	-	-	0/2/2/2	-
3	DMF	D	250	-	-	0/2/2/2	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	DMF	Z	79	-	-	0/2/2/2	-
3	DMF	X	88	-	-	0/2/2/2	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (66) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	252	DMF	O-C-N-C1
3	D	252	DMF	O-C-N-C2
3	C	73	DMF	O-C-N-C2
3	C	73	DMF	O-C-N-C1
3	R	89	DMF	O-C-N-C1
3	R	89	DMF	O-C-N-C2
3	N	2	DMF	O-C-N-C1
3	E	70	DMF	O-C-N-C1
3	F	250	DMF	O-C-N-C2
3	N	2	DMF	O-C-N-C2
3	R	68	DMF	O-C-N-C1
3	E	70	DMF	O-C-N-C2
3	F	250	DMF	O-C-N-C1
3	V	90	DMF	O-C-N-C1
3	Z	91	DMF	O-C-N-C1
3	R	68	DMF	O-C-N-C2
3	V	16	DMF	O-C-N-C2
3	N	21	DMF	O-C-N-C1
3	V	16	DMF	O-C-N-C1
3	V	90	DMF	O-C-N-C2
3	Z	91	DMF	O-C-N-C2
3	N	21	DMF	O-C-N-C2
3	C	5	DMF	O-C-N-C1
3	U	250	DMF	O-C-N-C1
3	S	250	DMF	O-C-N-C1
3	G	1	DMF	O-C-N-C1
3	L	9	DMF	O-C-N-C1
3	O	249	DMF	O-C-N-C1
3	T	35	DMF	O-C-N-C1
3	B	249	DMF	O-C-N-C1
3	C	64	DMF	O-C-N-C1
3	D	249	DMF	O-C-N-C1

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Mol	Chain	Res	Type	Atoms
3	E	44	DMF	O-C-N-C1
3	E	83	DMF	O-C-N-C1
3	H	52	DMF	O-C-N-C1
3	H	53	DMF	O-C-N-C1
3	J	6	DMF	O-C-N-C1
3	J	59	DMF	O-C-N-C1
3	O	250	DMF	O-C-N-C1
3	Q	250	DMF	O-C-N-C1
3	R	3	DMF	O-C-N-C1
3	S	251	DMF	O-C-N-C1
3	W	249	DMF	O-C-N-C1
3	Z	27	DMF	O-C-N-C1
3	U	250	DMF	O-C-N-C2
3	C	5	DMF	O-C-N-C2
3	Q	250	DMF	O-C-N-C2
3	S	250	DMF	O-C-N-C2
3	Z	27	DMF	O-C-N-C2
3	J	6	DMF	O-C-N-C2
3	L	9	DMF	O-C-N-C2
3	O	250	DMF	O-C-N-C2
3	H	53	DMF	O-C-N-C2
3	T	35	DMF	O-C-N-C2
3	W	249	DMF	O-C-N-C2
3	D	249	DMF	O-C-N-C2
3	O	249	DMF	O-C-N-C2
3	J	59	DMF	O-C-N-C2
3	R	3	DMF	O-C-N-C2
3	C	64	DMF	O-C-N-C2
3	G	1	DMF	O-C-N-C2
3	E	44	DMF	O-C-N-C2
3	E	83	DMF	O-C-N-C2
3	H	52	DMF	O-C-N-C2
3	B	249	DMF	O-C-N-C2
3	S	251	DMF	O-C-N-C2

There are no ring outliers.

57 monomers are involved in 108 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	O	250	DMF	4	0
3	H	53	DMF	2	0
3	H	41	DMF	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	39	DMF	1	0
3	W	249	DMF	1	0
3	2	78	DMF	1	0
3	Z	71	DMF	1	0
3	T	92	DMF	1	0
3	Z	27	DMF	2	0
3	D	251	DMF	1	0
3	X	67	DMF	1	0
3	E	70	DMF	2	0
3	H	19	DMF	4	0
3	P	61	DMF	1	0
3	L	4	DMF	1	0
3	J	82	DMF	1	0
3	Z	69	DMF	3	0
3	O	249	DMF	2	0
3	M	249	DMF	1	0
3	P	23	DMF	1	0
3	B	250	DMF	1	0
3	J	6	DMF	4	0
3	F	250	DMF	1	0
3	R	68	DMF	1	0
3	C	64	DMF	6	0
3	D	252	DMF	1	0
3	C	17	DMF	1	0
3	A	250	DMF	2	0
3	L	9	DMF	3	0
3	U	250	DMF	1	0
3	C	5	DMF	5	0
3	G	1	DMF	1	0
3	N	36	DMF	2	0
3	C	43	DMF	1	0
3	N	2	DMF	2	0
3	E	95	DMF	1	0
3	C	73	DMF	7	0
3	T	35	DMF	5	0
3	1	250	DMF	1	0
3	K	250	DMF	2	0
3	U	252	DMF	1	0
3	1	251	DMF	2	0
3	T	94	DMF	3	0
3	G	96	DMF	4	0
3	L	60	DMF	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	32	DMF	1	0
3	N	97	DMF	3	0
3	E	55	DMF	1	0
3	G	20	DMF	3	0
3	V	16	DMF	1	0
3	H	31	DMF	1	0
3	R	89	DMF	1	0
3	Z	91	DMF	3	0
3	S	250	DMF	1	0
3	B	249	DMF	1	0
3	U	251	DMF	2	0
3	Z	79	DMF	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	1	215/248 (86%)	0.84	25 (11%) 9 7	20, 55, 92, 101	0
1	A	216/248 (87%)	0.87	27 (12%) 8 6	21, 56, 100, 109	0
1	B	214/248 (86%)	0.96	36 (16%) 4 3	22, 56, 98, 108	0
1	D	213/248 (85%)	1.00	38 (17%) 4 2	21, 56, 96, 105	0
1	F	217/248 (87%)	0.96	37 (17%) 4 3	22, 55, 91, 102	0
1	I	215/248 (86%)	0.93	30 (13%) 6 5	23, 56, 98, 107	0
1	K	217/248 (87%)	0.98	39 (17%) 3 2	19, 57, 94, 109	0
1	M	216/248 (87%)	0.90	32 (14%) 5 4	21, 55, 92, 99	0
1	O	215/248 (86%)	0.87	29 (13%) 7 5	21, 56, 92, 102	0
1	Q	218/248 (87%)	0.88	32 (14%) 6 4	21, 55, 90, 101	0
1	S	215/248 (86%)	0.91	31 (14%) 6 4	21, 57, 102, 112	0
1	U	215/248 (86%)	0.89	31 (14%) 6 4	22, 55, 96, 106	0
1	W	217/248 (87%)	0.88	27 (12%) 8 6	22, 56, 93, 99	0
1	Y	213/248 (85%)	1.19	37 (17%) 4 2	23, 57, 99, 111	0
2	2	214/240 (89%)	-0.43	1 (0%) 87 89	10, 20, 43, 73	0
2	C	214/240 (89%)	-0.49	3 (1%) 73 75	8, 20, 46, 76	0
2	E	215/240 (89%)	-0.47	2 (0%) 81 82	10, 20, 45, 76	0
2	G	215/240 (89%)	-0.50	0 100 100	10, 21, 47, 75	0
2	H	212/240 (88%)	-0.46	1 (0%) 87 89	11, 22, 45, 70	0
2	J	215/240 (89%)	-0.42	2 (0%) 81 82	10, 21, 46, 70	0
2	L	215/240 (89%)	-0.45	2 (0%) 81 82	10, 21, 45, 72	0
2	N	212/240 (88%)	-0.50	1 (0%) 87 89	9, 21, 43, 67	0
2	P	215/240 (89%)	-0.39	3 (1%) 73 75	10, 23, 47, 73	0
2	R	222/240 (92%)	-0.56	1 (0%) 87 89	8, 19, 44, 73	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
2	T	215/240 (89%)	-0.40	4 (1%)	66 67	9, 22, 47, 80	0
2	V	220/240 (91%)	-0.54	3 (1%)	73 75	8, 20, 44, 74	0
2	X	215/240 (89%)	-0.47	4 (1%)	66 67	10, 22, 46, 79	0
2	Z	214/240 (89%)	-0.43	1 (0%)	87 89	10, 23, 46, 68	0
All	All	6029/6832 (88%)	0.23	479 (7%)	18 16	8, 34, 89, 112	0

All (479) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	Y	12	ALA	5.6
1	Y	184	ALA	5.2
1	A	172	ALA	5.1
1	B	234	LEU	4.9
1	F	202	THR	4.8
1	M	12	ALA	4.7
1	D	9	PRO	4.6
1	I	9	PRO	4.5
1	1	191	GLY	4.5
1	M	10	GLU	4.5
1	K	148	ALA	4.4
1	S	188	LEU	4.4
1	Y	188	LEU	4.4
1	Y	9	PRO	4.4
1	U	9	PRO	4.4
1	W	7	ILE	4.3
1	K	7	ILE	4.2
1	F	172	ALA	4.2
1	Y	13	MET	4.2
1	Y	190	ALA	4.2
1	Y	234	LEU	4.2
1	Q	7	ILE	4.2
1	F	13	MET	4.2
1	B	12	ALA	4.1
1	Q	191	GLY	4.1
1	B	13	MET	4.1
1	I	191	GLY	4.1
1	B	10	GLU	4.0
1	U	206	ALA	4.0
1	O	172	ALA	3.9
1	Y	172	ALA	3.9
1	B	190	ALA	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	190	ALA	3.9
1	B	205	VAL	3.9
1	F	191	GLY	3.9
1	K	9	PRO	3.9
1	U	166	ALA	3.8
1	B	171	TYR	3.8
1	K	147	ILE	3.8
1	D	234	LEU	3.8
1	S	190	ALA	3.7
1	B	204	GLY	3.7
1	F	12	ALA	3.6
1	Q	14	ARG	3.6
1	Q	166	ALA	3.6
1	B	175	ALA	3.6
1	O	13	MET	3.6
1	F	162	PRO	3.6
1	S	191	GLY	3.5
1	W	175	ALA	3.5
1	S	234	LEU	3.5
1	Y	171	TYR	3.5
1	I	188	LEU	3.5
1	O	191	GLY	3.5
1	S	8	SER	3.5
1	Q	163	ILE	3.5
1	K	13	MET	3.4
1	U	234	LEU	3.4
1	B	8	SER	3.4
1	O	166	ALA	3.4
1	U	190	ALA	3.4
2	V	400	ALA	3.4
1	S	175	ALA	3.3
1	Y	163	ILE	3.3
1	O	234	LEU	3.3
1	K	172	ALA	3.3
1	K	175	ALA	3.3
1	F	11	GLN	3.3
1	Y	233	LEU	3.3
1	U	191	GLY	3.3
1	Y	166	ALA	3.3
2	Z	392	ALA	3.3
1	1	179	ASP	3.2
1	Y	44	GLU	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	164	ALA	3.2
1	D	180	ALA	3.2
1	W	12	ALA	3.2
1	O	165	ASN	3.2
1	D	233	LEU	3.2
1	D	227	GLY	3.2
1	A	180	ALA	3.2
1	D	175	ALA	3.2
1	O	175	ALA	3.2
1	1	9	PRO	3.2
1	I	163	ILE	3.2
1	S	172	ALA	3.2
1	I	13	MET	3.2
1	1	159	THR	3.2
1	D	10	GLU	3.1
1	K	171	TYR	3.1
1	S	171	TYR	3.1
1	S	153	PHE	3.1
1	O	9	PRO	3.1
1	M	163	ILE	3.1
1	K	164	ALA	3.1
1	1	172	ALA	3.1
1	F	179	ASP	3.1
1	U	181	LEU	3.1
2	V	391	LEU	3.1
1	M	191	GLY	3.1
1	S	204	GLY	3.1
1	A	13	MET	3.1
1	U	13	MET	3.1
1	K	234	LEU	3.1
1	A	175	ALA	3.0
1	Y	180	ALA	3.0
1	D	183	ILE	3.0
2	V	392	ALA	3.0
1	B	9	PRO	3.0
1	B	162	PRO	3.0
1	B	14	ARG	3.0
1	B	166	ALA	3.0
1	W	15	GLU	3.0
1	1	190	ALA	3.0
2	X	400	ALA	3.0
1	Q	162	PRO	3.0

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Mol	Chain	Res	Type	RSRZ
1	W	191	GLY	3.0
1	W	13	MET	3.0
1	1	133	THR	3.0
1	F	175	ALA	3.0
1	Y	164	ALA	3.0
1	U	172	ALA	2.9
1	F	163	ILE	2.9
1	Q	44	GLU	2.9
2	X	392	ALA	2.9
1	A	234	LEU	2.9
1	I	177	LEU	2.9
1	1	234	LEU	2.9
1	B	17	SER	2.9
1	F	8	SER	2.9
1	Y	147	ILE	2.9
1	Y	154	VAL	2.9
1	M	175	ALA	2.9
1	Q	175	ALA	2.9
2	E	392	ALA	2.9
1	A	230	LEU	2.9
1	K	167	LEU	2.9
1	Y	29	SER	2.9
1	Y	131	GLY	2.9
1	Y	30	VAL	2.9
1	F	190	ALA	2.9
1	S	206	ALA	2.9
1	A	181	LEU	2.9
1	F	234	LEU	2.9
1	Y	179	ASP	2.9
1	O	153	PHE	2.9
1	I	184	ALA	2.8
1	K	184	ALA	2.8
2	T	392	ALA	2.8
1	U	188	LEU	2.8
1	M	13	MET	2.8
1	F	10	GLU	2.8
1	I	190	ALA	2.8
2	R	392	ALA	2.8
1	I	234	LEU	2.8
1	I	171	TYR	2.8
1	M	165	ASN	2.8
1	A	40	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	M	234	LEU	2.8
1	W	234	LEU	2.8
1	M	44	GLU	2.8
1	O	161	GLU	2.8
1	Y	231	GLN	2.8
1	D	179	ASP	2.8
1	M	136	PRO	2.8
1	U	162	PRO	2.8
1	I	172	ALA	2.8
1	S	232	ALA	2.8
1	Q	169	GLU	2.8
1	B	163	ILE	2.7
1	W	165	ASN	2.7
1	D	207	SER	2.7
1	O	164	ALA	2.7
1	1	175	ALA	2.7
1	S	205	VAL	2.7
1	B	167	LEU	2.7
1	S	27	ALA	2.7
2	L	392	ALA	2.7
2	P	400	ALA	2.7
1	K	149	ASP	2.7
1	F	153	PHE	2.7
1	1	183	ILE	2.7
1	I	162	PRO	2.7
1	U	46	PRO	2.7
1	A	171	TYR	2.7
1	U	8	SER	2.7
1	F	188	LEU	2.7
1	K	188	LEU	2.7
1	Q	181	LEU	2.7
1	Q	234	LEU	2.7
1	B	173	GLU	2.7
1	Y	204	GLY	2.7
1	W	183	ILE	2.7
1	F	14	ARG	2.7
1	F	171	TYR	2.7
1	M	113	GLU	2.7
1	U	175	ALA	2.7
1	Y	175	ALA	2.7
2	J	392	ALA	2.7
1	K	152	HIS	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	131	GLY	2.6
1	1	163	ILE	2.6
1	D	230	LEU	2.6
1	W	176	SER	2.6
1	D	27	ALA	2.6
1	F	166	ALA	2.6
1	A	205	VAL	2.6
1	F	36	ALA	2.6
1	I	164	ALA	2.6
1	I	180	ALA	2.6
1	O	159	THR	2.6
1	W	25	ALA	2.6
1	1	184	ALA	2.6
1	Y	111	PHE	2.6
1	B	177	LEU	2.6
1	B	230	LEU	2.6
1	M	166	ALA	2.6
1	U	232	ALA	2.6
1	B	152	HIS	2.6
1	Y	159	THR	2.6
1	F	183	ILE	2.6
2	T	391	LEU	2.6
1	M	159	THR	2.6
1	Q	164	ALA	2.6
1	Q	180	ALA	2.6
1	S	180	ALA	2.6
1	W	172	ALA	2.6
1	Y	187	ALA	2.6
1	O	168	LYS	2.6
1	U	171	TYR	2.6
1	Q	13	MET	2.5
1	Q	167	LEU	2.5
1	F	164	ALA	2.5
1	K	186	ALA	2.5
1	Q	36	ALA	2.5
1	B	159	THR	2.5
1	S	183	ILE	2.5
1	A	233	LEU	2.5
2	X	391	LEU	2.5
1	B	172	ALA	2.5
1	B	160	THR	2.5
1	O	171	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
1	Y	136	PRO	2.5
1	M	132	GLU	2.5
1	F	181	LEU	2.5
1	M	177	LEU	2.5
1	W	181	LEU	2.5
1	U	205	VAL	2.5
1	D	172	ALA	2.5
1	D	13	MET	2.5
1	A	177	LEU	2.5
1	U	11	GLN	2.5
1	M	229	ALA	2.5
1	1	32	ALA	2.5
2	C	400	ALA	2.5
1	K	159	THR	2.5
1	M	171	TYR	2.4
1	Y	177	LEU	2.4
2	H	391	LEU	2.4
1	I	26	ARG	2.4
1	A	166	ALA	2.4
1	D	232	ALA	2.4
1	W	162	PRO	2.4
1	O	173	GLU	2.4
1	U	15	GLU	2.4
1	Q	183	ILE	2.4
1	1	152	HIS	2.4
1	S	31	VAL	2.4
1	S	42	VAL	2.4
1	D	12	ALA	2.4
1	M	180	ALA	2.4
2	N	400	ALA	2.4
2	P	392	ALA	2.4
1	F	159	THR	2.4
1	U	163	ILE	2.4
1	B	233	LEU	2.4
1	W	203	LEU	2.4
1	Q	232	ALA	2.4
1	W	190	ALA	2.4
1	M	21	ARG	2.4
1	M	167	LEU	2.4
1	U	231	GLN	2.4
1	1	188	LEU	2.4
2	J	391	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	O	184	ALA	2.4
1	Q	172	ALA	2.4
1	I	15	GLU	2.4
1	Q	9	PRO	2.3
1	S	160	THR	2.3
2	C	416	SER	2.3
1	Q	21	ARG	2.3
1	D	42	VAL	2.3
1	A	164	ALA	2.3
1	A	186	ALA	2.3
1	F	27	ALA	2.3
1	K	10	GLU	2.3
1	S	25	ALA	2.3
1	1	25	ALA	2.3
1	S	9	PRO	2.3
1	O	204	GLY	2.3
1	A	183	ILE	2.3
1	D	231	GLN	2.3
1	D	177	LEU	2.3
1	D	208	LEU	2.3
1	O	167	LEU	2.3
1	B	155	VAL	2.3
1	Q	165	ASN	2.3
1	B	232	ALA	2.3
1	D	36	ALA	2.3
1	M	164	ALA	2.3
1	M	172	ALA	2.3
1	O	180	ALA	2.3
1	U	164	ALA	2.3
1	1	162	PRO	2.3
2	L	400	ALA	2.3
1	A	182	ARG	2.3
1	Q	48	ARG	2.3
1	K	191	GLY	2.3
1	Q	131	GLY	2.3
1	I	230	LEU	2.3
1	M	233	LEU	2.3
1	Q	203	LEU	2.3
1	1	147	ILE	2.3
1	B	161	GLU	2.3
1	F	184	ALA	2.3
1	K	190	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	M	162	PRO	2.3
1	O	14	ARG	2.3
1	D	159	THR	2.3
1	M	183	ILE	2.3
1	U	167	LEU	2.3
1	U	233	LEU	2.3
1	Q	179	ASP	2.3
1	M	15	GLU	2.3
1	A	48	ARG	2.3
2	2	400	ALA	2.2
1	B	33	LEU	2.2
1	F	177	LEU	2.2
1	K	177	LEU	2.2
1	O	183	ILE	2.2
1	S	13	MET	2.2
1	Y	167	LEU	2.2
1	A	10	GLU	2.2
1	U	18	GLU	2.2
1	K	153	PHE	2.2
1	Y	162	PRO	2.2
1	B	165	ASN	2.2
1	F	180	ALA	2.2
1	W	166	ALA	2.2
1	W	232	ALA	2.2
1	1	186	ALA	2.2
1	W	98	GLN	2.2
1	D	157	GLY	2.2
1	D	204	GLY	2.2
1	M	168	LYS	2.2
1	F	17	SER	2.2
1	O	147	ILE	2.2
1	W	170	SER	2.2
1	F	18	GLU	2.2
1	I	21	ARG	2.2
1	S	155	VAL	2.2
1	K	229	ALA	2.2
1	1	165	ASN	2.2
2	P	425	ALA	2.2
1	M	230	LEU	2.2
1	S	227	GLY	2.2
2	T	399	LEU	2.2
1	F	161	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	1	153	PHE	2.2
1	A	190	ALA	2.2
1	B	174	ASN	2.2
1	B	229	ALA	2.2
1	I	148	ALA	2.2
1	U	180	ALA	2.2
1	1	12	ALA	2.2
1	K	133	THR	2.2
1	F	33	LEU	2.2
1	I	37	GLY	2.2
1	D	176	SER	2.2
1	K	150	GLU	2.2
1	K	111	PHE	2.2
1	O	205	VAL	2.1
1	F	229	ALA	2.1
1	K	232	ALA	2.1
1	S	186	ALA	2.1
1	U	12	ALA	2.1
1	W	27	ALA	2.1
1	W	187	ALA	2.1
1	Y	186	ALA	2.1
1	F	167	LEU	2.1
1	U	159	THR	2.1
1	U	160	THR	2.1
1	I	183	ILE	2.1
1	W	171	TYR	2.1
1	I	136	PRO	2.1
1	M	9	PRO	2.1
1	A	42	VAL	2.1
1	D	206	ALA	2.1
1	M	148	ALA	2.1
1	A	188	LEU	2.1
1	I	181	LEU	2.1
1	Q	233	LEU	2.1
1	W	178	THR	2.1
2	E	391	LEU	2.1
1	S	163	ILE	2.1
1	F	150	GLU	2.1
1	D	228	SER	2.1
1	I	179	ASP	2.1
1	M	176	SER	2.1
1	B	111	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	Q	153	PHE	2.1
1	D	31	VAL	2.1
1	W	11	GLN	2.1
1	I	175	ALA	2.1
1	M	112	THR	2.1
1	Q	202	THR	2.1
1	I	37	GLY	2.1
1	D	182	ARG	2.1
1	U	14	ARG	2.1
1	D	163	ILE	2.1
1	Y	130	TYR	2.1
1	A	168	LYS	2.1
1	Y	144	ASP	2.1
1	D	152	HIS	2.1
1	I	13	MET	2.1
1	I	111	PHE	2.1
1	W	153	PHE	2.1
1	Y	31	VAL	2.1
1	D	34	ALA	2.1
1	K	25	ALA	2.1
1	K	180	ALA	2.1
1	A	191	GLY	2.1
1	D	14	ARG	2.1
1	O	182	ARG	2.1
1	S	157	GLY	2.1
2	X	389	GLY	2.1
1	K	169	GLU	2.1
1	O	178	THR	2.1
1	U	10	GLU	2.1
2	C	433	GLU	2.1
1	O	176	SER	2.1
1	D	136	PRO	2.1
1	B	31	VAL	2.0
1	I	31	VAL	2.0
1	K	185	VAL	2.0
1	D	48	ARG	2.0
1	I	25	ALA	2.0
1	K	166	ALA	2.0
1	S	36	ALA	2.0
1	I	232	ALA	2.0
2	T	400	ALA	2.0
1	F	230	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	Q	73	ASN	2.0
1	A	113	GLU	2.0
1	D	169	GLU	2.0
1	I	10	GLU	2.0
1	K	157	GLY	2.0
1	K	168	LYS	2.0
1	O	163	ILE	2.0
1	B	179	ASP	2.0
1	Y	49	SER	2.0
1	A	14	ARG	2.0
1	K	30	VAL	2.0
1	O	111	PHE	2.0
1	I	167	LEU	2.0
1	Q	177	LEU	2.0
1	S	33	LEU	2.0
1	S	187	ALA	2.0
1	S	233	LEU	2.0
1	Y	43	ALA	2.0
1	W	10	GLU	2.0
1	F	157	GLY	2.0
1	K	131	GLY	2.0
1	K	174	ASN	2.0
1	K	163	ILE	2.0
1	O	160	THR	2.0
1	Q	159	THR	2.0
1	A	9	PRO	2.0
1	K	151	PRO	2.0

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

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	OZT	R	301	9/10	0.89	0.12	23,29,35,38	0
2	OZT	N	301	9/10	0.92	0.09	23,28,35,38	0
2	OZT	P	301	9/10	0.92	0.10	25,29,34,39	0
2	OZT	L	301	9/10	0.92	0.09	25,29,34,35	0
2	OZT	V	301	9/10	0.92	0.09	23,30,35,36	0
2	OZT	X	301	9/10	0.92	0.09	28,31,35,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	OZT	2	301	9/10	0.92	0.11	23,30,34,35	0
2	OZT	J	301	9/10	0.93	0.09	24,30,34,38	0
2	OZT	E	301	9/10	0.94	0.08	25,31,33,36	0
2	OZT	Z	301	9/10	0.94	0.07	25,28,32,33	0
2	OZT	H	301	9/10	0.94	0.08	24,30,33,33	0
2	OZT	C	301	9/10	0.95	0.08	19,28,34,37	0
2	OZT	G	301	9/10	0.95	0.07	24,30,34,37	0
2	OZT	T	301	9/10	0.96	0.06	26,29,32,34	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	DMF	V	65	5/5	0.55	0.27	60,60,61,61	0
3	DMF	T	35	5/5	0.63	0.32	75,75,76,77	0
3	DMF	Q	250	5/5	0.69	0.28	69,70,70,71	0
3	DMF	W	249	5/5	0.69	0.25	77,77,78,78	0
3	DMF	F	250	5/5	0.70	0.27	68,69,71,71	0
3	DMF	T	94	5/5	0.70	0.34	104,104,104,104	0
3	DMF	P	61	5/5	0.70	0.21	58,59,61,62	0
3	DMF	D	249	5/5	0.70	0.29	83,83,84,84	0
3	DMF	N	37	5/5	0.71	0.19	63,64,65,67	0
3	DMF	D	252	5/5	0.74	0.23	50,50,54,55	0
3	DMF	N	99	5/5	0.74	0.21	50,52,53,54	0
3	DMF	C	43	5/5	0.75	0.22	57,57,59,59	0
3	DMF	S	251	5/5	0.75	0.23	69,70,70,70	0
3	DMF	U	250	5/5	0.76	0.24	71,71,72,72	0
3	DMF	H	52	5/5	0.76	0.21	65,66,66,67	0
3	DMF	G	96	5/5	0.76	0.27	79,80,80,81	0
3	DMF	X	67	5/5	0.76	0.22	58,59,60,60	0
3	DMF	Y	249	5/5	0.77	0.19	58,59,61,62	0
3	DMF	H	31	5/5	0.78	0.31	64,66,68,68	0
3	DMF	N	21	5/5	0.78	0.27	75,76,76,76	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	DMF	R	54	5/5	0.79	0.21	46,48,49,51	0
3	DMF	D	251	5/5	0.79	0.20	51,53,54,56	0
3	DMF	O	250	5/5	0.79	0.25	59,60,60,61	0
3	DMF	U	252	5/5	0.80	0.17	60,60,61,61	0
3	DMF	2	66	5/5	0.80	0.26	69,70,70,71	0
3	DMF	D	250	5/5	0.81	0.21	67,67,67,68	0
3	DMF	B	251	5/5	0.81	0.18	61,64,64,64	0
3	DMF	C	64	5/5	0.82	0.19	60,60,60,61	0
3	DMF	S	250	5/5	0.82	0.18	48,49,51,51	0
3	DMF	Z	69	5/5	0.82	0.21	69,70,71,71	0
3	DMF	1	251	5/5	0.82	0.18	45,48,49,49	0
3	DMF	U	251	5/5	0.82	0.17	45,45,46,48	0
3	DMF	E	32	5/5	0.83	0.23	60,62,63,65	0
3	DMF	E	83	5/5	0.83	0.24	64,65,65,66	0
3	DMF	E	95	5/5	0.83	0.19	46,49,51,55	0
3	DMF	X	88	5/5	0.83	0.18	46,48,50,51	0
3	DMF	B	249	5/5	0.84	0.17	49,52,53,55	0
3	DMF	G	39	5/5	0.84	0.18	55,57,59,60	0
3	DMF	H	102	5/5	0.84	0.21	73,73,74,74	0
3	DMF	X	25	5/5	0.84	0.17	47,49,50,51	0
3	DMF	2	58	5/5	0.84	0.24	60,62,64,64	0
3	DMF	C	48	5/5	0.84	0.21	68,69,70,70	0
3	DMF	N	36	5/5	0.85	0.20	50,53,54,57	0
3	DMF	Z	27	5/5	0.85	0.25	57,62,64,65	0
3	DMF	Z	50	5/5	0.85	0.17	48,49,50,52	0
3	DMF	W	250	5/5	0.85	0.20	61,62,64,64	0
3	DMF	Z	91	5/5	0.85	0.17	54,55,55,56	0
3	DMF	M	250	5/5	0.85	0.12	35,37,37,38	0
3	DMF	N	97	5/5	0.85	0.17	31,36,38,38	0
3	DMF	C	5	5/5	0.85	0.24	77,77,78,78	0
3	DMF	E	70	5/5	0.86	0.24	77,77,78,78	0
3	DMF	K	250	5/5	0.86	0.20	59,61,62,63	0
3	DMF	O	249	5/5	0.86	0.17	42,46,47,49	0
3	DMF	F	249	5/5	0.86	0.19	55,55,56,56	0
3	DMF	N	2	5/5	0.86	0.17	44,45,48,50	0
3	DMF	H	19	5/5	0.86	0.16	52,53,55,58	0
3	DMF	C	28	5/5	0.86	0.18	53,54,56,57	0
3	DMF	R	89	5/5	0.86	0.18	52,53,53,54	0
3	DMF	S	249	5/5	0.86	0.21	63,64,65,66	0
3	DMF	G	20	5/5	0.86	0.19	46,50,51,53	0
3	DMF	T	62	5/5	0.87	0.20	56,58,58,60	0
3	DMF	T	92	5/5	0.87	0.21	41,44,45,46	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	DMF	R	3	5/5	0.87	0.23	63,63,65,65	0
3	DMF	H	53	5/5	0.87	0.20	63,64,65,65	0
3	DMF	E	44	5/5	0.87	0.16	42,44,46,46	0
3	DMF	C	73	5/5	0.87	0.18	53,54,54,55	0
3	DMF	V	16	5/5	0.87	0.16	48,50,51,52	0
3	DMF	L	60	5/5	0.87	0.18	57,57,58,59	0
3	DMF	V	86	5/5	0.87	0.17	44,44,45,48	0
3	DMF	P	100	5/5	0.87	0.17	53,54,54,55	0
3	DMF	G	1	5/5	0.87	0.17	38,42,43,46	0
3	DMF	1	250	5/5	0.88	0.19	60,60,61,61	0
3	DMF	K	249	5/5	0.88	0.12	37,40,41,41	0
3	DMF	V	90	5/5	0.88	0.19	66,66,68,71	0
3	DMF	A	250	5/5	0.88	0.16	52,54,56,59	0
3	DMF	2	78	5/5	0.88	0.16	43,44,46,48	0
3	DMF	R	68	5/5	0.89	0.20	53,56,57,57	0
3	DMF	C	15	5/5	0.89	0.17	51,53,54,56	0
3	DMF	I	249	5/5	0.89	0.14	42,43,44,45	0
3	DMF	J	82	5/5	0.89	0.18	50,52,54,57	0
3	DMF	H	41	5/5	0.89	0.18	62,63,63,64	0
3	DMF	Z	71	5/5	0.89	0.16	52,52,54,56	0
3	DMF	L	9	5/5	0.90	0.19	61,63,64,65	0
3	DMF	C	17	5/5	0.90	0.11	37,38,41,42	0
3	DMF	B	250	5/5	0.90	0.13	39,39,43,45	0
3	DMF	P	23	5/5	0.90	0.21	60,60,61,63	0
3	DMF	I	250	5/5	0.90	0.15	48,51,51,52	0
3	DMF	P	85	5/5	0.90	0.16	52,53,54,54	0
3	DMF	1	249	5/5	0.91	0.15	56,57,58,59	0
3	DMF	J	6	5/5	0.91	0.14	53,54,55,57	0
3	DMF	A	249	5/5	0.91	0.12	54,55,55,56	0
3	DMF	H	14	5/5	0.91	0.16	63,63,64,64	0
3	DMF	Z	79	5/5	0.91	0.14	38,38,39,40	0
3	DMF	N	24	5/5	0.91	0.12	41,42,44,46	0
3	DMF	P	74	5/5	0.92	0.20	72,73,74,74	0
3	DMF	T	46	5/5	0.92	0.12	45,46,47,48	0
3	DMF	U	249	5/5	0.92	0.12	46,47,48,49	0
3	DMF	Q	249	5/5	0.92	0.14	52,52,53,54	0
3	DMF	M	249	5/5	0.93	0.13	43,43,44,46	0
3	DMF	J	59	5/5	0.93	0.14	65,65,66,66	0
3	DMF	E	55	5/5	0.94	0.09	41,41,42,42	0
3	DMF	L	4	5/5	0.95	0.09	47,47,47,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.