



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:16 AM GMT

PDB ID : 3HFA
Title : Crystal Structure of Mycobacterium Tuberculosis Proteasome open-gate mutant
Authors : Li, D.; Li, H.
Deposited on : 2009-05-11
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

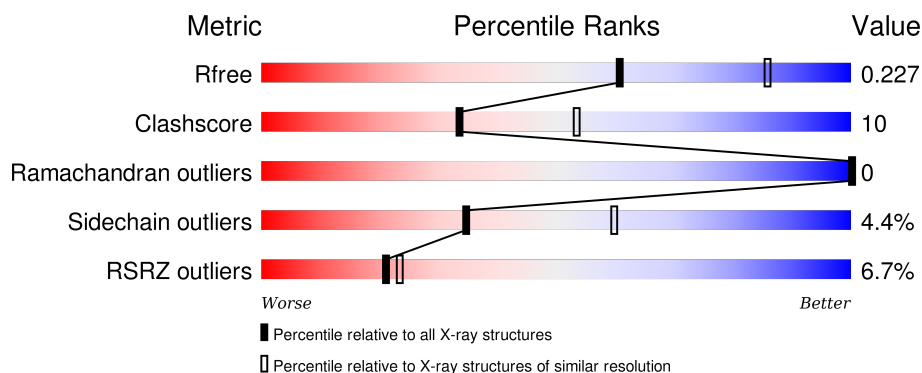
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	2	240	
1	C	240	
1	E	240	
1	G	240	
1	H	240	

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Mol	Chain	Length	Quality of chain
1	J	240	
1	L	240	
1	N	240	
1	P	240	
1	R	240	
1	T	240	
1	V	240	
1	X	240	
1	Z	240	
2	1	240	
2	A	240	
2	B	240	
2	D	240	
2	F	240	
2	I	240	
2	K	240	
2	M	240	
2	O	240	
2	Q	240	
2	S	240	
2	U	240	
2	W	240	
2	Y	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
3	DMF	1	249	-	-	-	X
3	DMF	1	250	-	-	-	X
3	DMF	2	145	-	-	-	X
3	DMF	A	249	-	-	-	X
3	DMF	B	249	-	-	-	X
3	DMF	C	15	-	-	-	X
3	DMF	E	104	-	-	-	X
3	DMF	E	113	-	-	-	X
3	DMF	G	137	-	-	-	X
3	DMF	G	140	-	-	-	X
3	DMF	G	20	-	-	-	X
3	DMF	H	142	-	-	-	X
3	DMF	H	41	-	-	-	X
3	DMF	I	249	-	-	-	X
3	DMF	J	46	-	-	-	X
3	DMF	K	250	-	-	-	X
3	DMF	K	251	-	-	X	X
3	DMF	L	138	-	-	-	X
3	DMF	L	14	-	-	-	X
3	DMF	L	60	-	-	-	X
3	DMF	L	9	-	-	-	X
3	DMF	M	249	-	-	-	X
3	DMF	N	36	-	-	-	X
3	DMF	O	7	-	-	-	X
3	DMF	P	107	-	-	-	X
3	DMF	P	133	-	-	-	X
3	DMF	P	23	-	-	-	X
3	DMF	Q	249	-	-	-	X
3	DMF	Q	251	-	-	X	X
3	DMF	S	249	-	-	-	X
3	DMF	T	62	-	-	-	X
3	DMF	U	249	-	-	-	X
3	DMF	U	250	-	-	-	X
3	DMF	V	117	-	-	-	X
3	DMF	V	121	-	-	-	X
3	DMF	V	135	-	-	-	X
3	DMF	V	136	-	-	-	X
3	DMF	W	249	-	-	-	X
3	DMF	Y	249	-	-	-	X
3	DMF	Y	251	-	-	-	X
3	DMF	Z	105	-	-	-	X
3	DMF	Z	122	-	-	-	X
3	DMF	Z	27	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	DMF	Z	50	-	-	-	X
3	DMF	Z	69	-	-	X	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 48523 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 (Beta subunit) PrcB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	223	Total	C	N	O	S	0	0	0
			1642	1029	283	325	5			
1	C	222	Total	C	N	O	S	0	0	0
			1638	1027	282	324	5			
1	E	223	Total	C	N	O	S	0	0	0
			1642	1029	283	325	5			
1	G	222	Total	C	N	O	S	0	0	0
			1638	1027	282	324	5			
1	J	222	Total	C	N	O	S	0	0	0
			1638	1027	282	324	5			
1	L	222	Total	C	N	O	S	0	0	0
			1638	1027	282	324	5			
1	N	223	Total	C	N	O	S	0	0	0
			1642	1029	283	325	5			
1	P	222	Total	C	N	O	S	0	0	0
			1638	1027	282	324	5			
1	R	222	Total	C	N	O	S	0	0	0
			1638	1027	282	324	5			
1	T	223	Total	C	N	O	S	0	0	0
			1642	1029	283	325	5			
1	V	226	Total	C	N	O	S	0	0	0
			1662	1040	286	331	5			
1	X	224	Total	C	N	O	S	0	0	0
			1647	1032	284	326	5			
1	Z	224	Total	C	N	O	S	0	0	0
			1647	1032	284	326	5			
1	2	222	Total	C	N	O	S	0	0	0
			1638	1027	282	324	5			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	535	HIS	-	EXPRESSION TAG	UNP O33245

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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	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
E	539	HIS	-	EXPRESSION TAG	UNP O33245
E	540	HIS	-	EXPRESSION TAG	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
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	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	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	535	HIS	-	EXPRESSION TAG	UNP O33245

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Chain	Residue	Modelled	Actual	Comment	Reference
P	536	HIS	-	EXPRESSION TAG	UNP O33245
P	537	HIS	-	EXPRESSION TAG	UNP O33245
P	538	HIS	-	EXPRESSION TAG	UNP O33245
P	539	HIS	-	EXPRESSION TAG	UNP O33245
P	540	HIS	-	EXPRESSION TAG	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	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	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	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	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	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
2	539	HIS	-	EXPRESSION TAG	UNP O33245
2	540	HIS	-	EXPRESSION TAG	UNP O33245

- Molecule 2 is a protein called Proteasome (Alpha subunit) PrcA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	213	Total	C	N	O	S	0	0	0
			1645	1030	301	310	4			
2	A	214	Total	C	N	O	S	0	0	0
			1653	1036	302	311	4			
2	B	214	Total	C	N	O	S	0	0	0
			1653	1036	302	311	4			
2	F	214	Total	C	N	O	S	0	0	0
			1653	1036	302	311	4			
2	I	214	Total	C	N	O	S	0	0	0
			1653	1036	302	311	4			
2	K	214	Total	C	N	O	S	0	0	0
			1653	1036	302	311	4			
2	M	214	Total	C	N	O	S	0	0	0
			1653	1036	302	311	4			
2	O	214	Total	C	N	O	S	0	0	0
			1653	1036	302	311	4			
2	Q	214	Total	C	N	O	S	0	0	0
			1653	1036	302	311	4			
2	S	214	Total	C	N	O	S	0	0	0
			1653	1036	302	311	4			
2	U	214	Total	C	N	O	S	0	0	0
			1653	1036	302	311	4			
2	W	214	Total	C	N	O	S	0	0	0
			1653	1036	302	311	4			
2	Y	214	Total	C	N	O	S	0	0	0
			1653	1036	302	311	4			
2	1	214	Total	C	N	O	S	0	0	0
			1653	1036	302	311	4			

There are 14 discrepancies between the modelled and reference sequences:

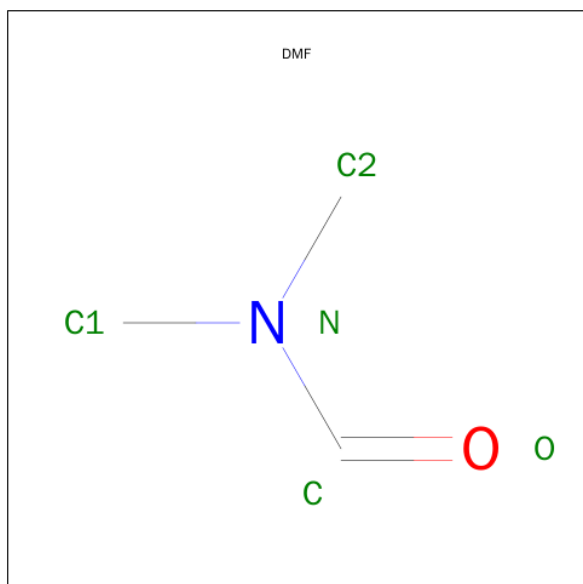
Chain	Residue	Modelled	Actual	Comment	Reference
D	9	MET	-	INITIATING METHIONINE	UNP O33244
A	9	MET	-	INITIATING METHIONINE	UNP O33244
B	9	MET	-	INITIATING METHIONINE	UNP O33244
F	9	MET	-	INITIATING METHIONINE	UNP O33244
I	9	MET	-	INITIATING METHIONINE	UNP O33244
K	9	MET	-	INITIATING METHIONINE	UNP O33244
M	9	MET	-	INITIATING METHIONINE	UNP O33244
O	9	MET	-	INITIATING METHIONINE	UNP O33244
Q	9	MET	-	INSERTION	UNP O33244
S	9	MET	-	INITIATING METHIONINE	UNP O33244
U	9	MET	-	INITIATING METHIONINE	UNP O33244
W	9	MET	-	INITIATING METHIONINE	UNP O33244

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Chain	Residue	Modelled	Actual	Comment	Reference
Y	9	MET	-	INITIATING METHIONINE	UNP O33244
1	9	MET	-	INITIATING METHIONINE	UNP O33244

- Molecule 3 is DIMETHYLFORMAMIDE (three-letter code: DMF) (formula: C_3H_7NO).



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	C	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	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	J	1	Total	C	N	O	0	0
			5	3	1	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	L	1	Total 5	C 3	N 1	O 1	0	0
3	L	1	Total 5	C 3	N 1	O 1	0	0
3	L	1	Total 5	C 3	N 1	O 1	0	0
3	L	1	Total 5	C 3	N 1	O 1	0	0
3	N	1	Total 5	C 3	N 1	O 1	0	0
3	N	1	Total 5	C 3	N 1	O 1	0	0
3	P	1	Total 5	C 3	N 1	O 1	0	0
3	P	1	Total 5	C 3	N 1	O 1	0	0
3	P	1	Total 5	C 3	N 1	O 1	0	0
3	T	1	Total 5	C 3	N 1	O 1	0	0
3	T	1	Total 5	C 3	N 1	O 1	0	0
3	V	1	Total 5	C 3	N 1	O 1	0	0
3	V	1	Total 5	C 3	N 1	O 1	0	0
3	V	1	Total 5	C 3	N 1	O 1	0	0
3	V	1	Total 5	C 3	N 1	O 1	0	0
3	V	1	Total 5	C 3	N 1	O 1	0	0
3	V	1	Total 5	C 3	N 1	O 1	0	0
3	X	1	Total 5	C 3	N 1	O 1	0	0
3	Z	1	Total 5	C 3	N 1	O 1	0	0
3	Z	1	Total 5	C 3	N 1	O 1	0	0
3	Z	1	Total 5	C 3	N 1	O 1	0	0
3	Z	1	Total 5	C 3	N 1	O 1	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	Z	1	Total 5	C 3	N 1	O 1	0	0
3	2	1	Total 5	C 3	N 1	O 1	0	0
3	2	1	Total 5	C 3	N 1	O 1	0	0
3	D	1	Total 5	C 3	N 1	O 1	0	0
3	A	1	Total 5	C 3	N 1	O 1	0	0
3	A	1	Total 5	C 3	N 1	O 1	0	0
3	B	1	Total 5	C 3	N 1	O 1	0	0
3	F	1	Total 5	C 3	N 1	O 1	0	0
3	I	1	Total 5	C 3	N 1	O 1	0	0
3	K	1	Total 5	C 3	N 1	O 1	0	0
3	K	1	Total 5	C 3	N 1	O 1	0	0
3	K	1	Total 5	C 3	N 1	O 1	0	0
3	M	1	Total 5	C 3	N 1	O 1	0	0
3	O	1	Total 5	C 3	N 1	O 1	0	0
3	Q	1	Total 5	C 3	N 1	O 1	0	0
3	Q	1	Total 5	C 3	N 1	O 1	0	0
3	Q	1	Total 5	C 3	N 1	O 1	0	0
3	S	1	Total 5	C 3	N 1	O 1	0	0
3	S	1	Total 5	C 3	N 1	O 1	0	0
3	U	1	Total 5	C 3	N 1	O 1	0	0
3	U	1	Total 5	C 3	N 1	O 1	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	W	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	Y	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	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		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	92	Total	O	0	0
			92	92		
4	C	43	Total	O	0	0
			43	43		
4	E	121	Total	O	0	0
			121	121		
4	G	113	Total	O	0	0
			113	113		
4	J	73	Total	O	0	0
			73	73		
4	L	109	Total	O	0	0
			109	109		
4	N	124	Total	O	0	0
			124	124		
4	P	93	Total	O	0	0
			93	93		
4	R	89	Total	O	0	0
			89	89		
4	T	96	Total	O	0	0
			96	96		
4	V	133	Total	O	0	0
			133	133		
4	X	101	Total	O	0	0
			101	101		
4	Z	91	Total	O	0	0
			91	91		

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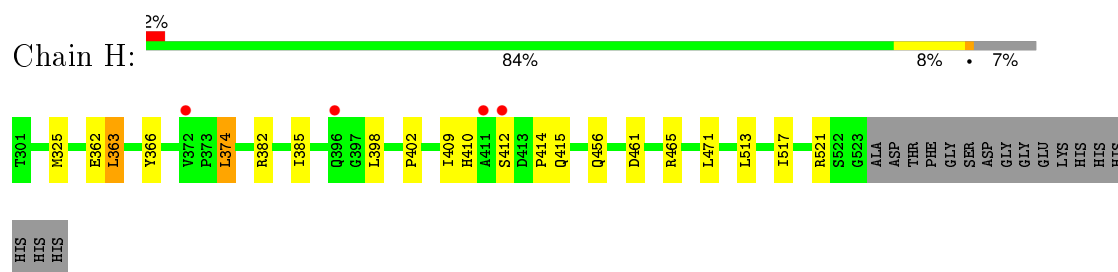
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	2	107	Total 107	O 107	0	0
4	D	49	Total 49	O 49	0	0
4	A	31	Total 31	O 31	0	0
4	B	37	Total 37	O 37	0	0
4	F	60	Total 60	O 60	0	0
4	I	35	Total 35	O 35	0	0
4	K	52	Total 52	O 52	0	0
4	M	74	Total 74	O 74	0	0
4	O	34	Total 34	O 34	0	0
4	Q	71	Total 71	O 71	0	0
4	S	35	Total 35	O 35	0	0
4	U	48	Total 48	O 48	0	0
4	W	62	Total 62	O 62	0	0
4	Y	68	Total 68	O 68	0	0
4	1	68	Total 68	O 68	0	0

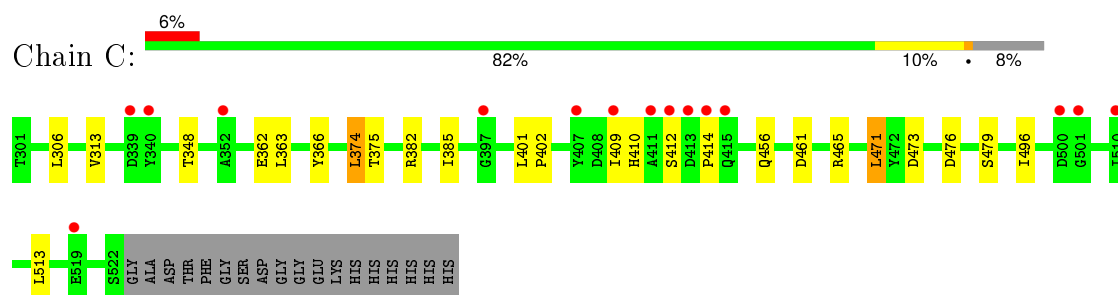
3 Residue-property plots [i](#)

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

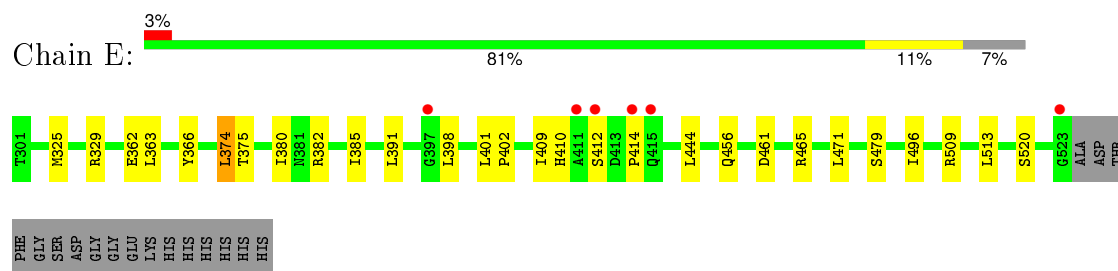
- Molecule 1: Proteasome (Beta subunit) PrcB



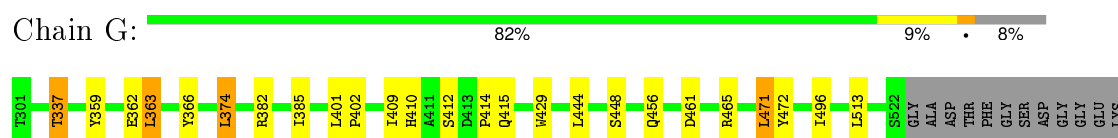
- Molecule 1: Proteasome (Beta subunit) PrcB



- Molecule 1: Proteasome (Beta subunit) PrcB




- Molecule 1: Proteasome (Beta subunit) PrcB



HIS
HIS
HIS
HIS
HIS


• Molecule 1: Proteasome (Beta subunit) PrcB

Chain J: 



GLY
GLU
LYS
HIS
HIS
HIS
HIS
HIS


• Molecule 1: Proteasome (Beta subunit) PrcB

Chain L: 



GLY
GLU
LYS
HIS
HIS
HIS
HIS
HIS


• Molecule 1: Proteasome (Beta subunit) PrcB

Chain N: 



HIS
HIS
HIS


• Molecule 1: Proteasome (Beta subunit) PrcB

Chain P: 



HIS

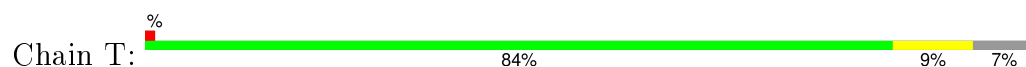
• Molecule 1: Proteasome (Beta subunit) PrcB

Chain R: 



ASP
GLY
GLY
GLY
LYS
HIS
HIS
HIS
HIS
HIS

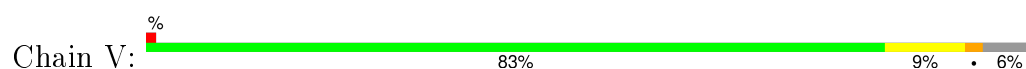
• Molecule 1: Proteasome (Beta subunit) PrcB



T301 L306 E362 E363 Y366 L374 A377 R382 I385 L391 N395 L398 I409 H410 A411 S412 D413 P414 Q415 L444 Q456 D461 R465 L471 Y472 L513 G523 ALA ASP THR PHE GLY SER ASP GLY LYS HIS HIS HIS HIS

HIS
HIS

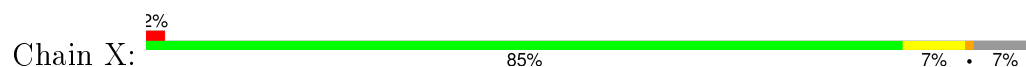
• Molecule 1: Proteasome (Beta subunit) PrcB



T301 R319 I345 R357 E362 E363 Y366 L374 T375 R382 I385 R388 I409 H410 A411 S412 D413 P414 W429 K451 Q456 D461 R465 L471 I496 L513 A514 R515 G523 T526 PHE GLY SER ASP GLY LYS HIS HIS HIS HIS HIS

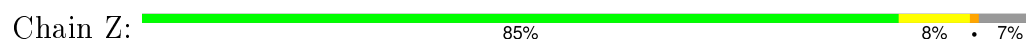
HIS
HIS
HIS
HIS

• Molecule 1: Proteasome (Beta subunit) PrcB



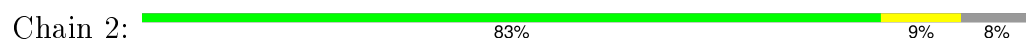
T301 N325 E362 E363 Y366 L374 R382 I385 L391 Q396 I409 H410 A411 S412 D413 P414 Q415 K451 Q456 D461 R465 L471 D509 L513 G523 A524 ASP THR PHE GLY SER ASP GLY GLY GLY LYS HIS HIS HIS HIS HIS HIS

• Molecule 1: Proteasome (Beta subunit) PrcB



T301 R357 E362 E363 Y366 L374 R382 I385 L398 I409 H410 A411 S412 D413 P414 W429 L444 Q456 D461 R465 L471 I496 R509 L513 A524 ASP THR PHE GLY SER ASP GLY GLY GLY LYS HIS HIS HIS HIS HIS HIS

• Molecule 1: Proteasome (Beta subunit) PrcB

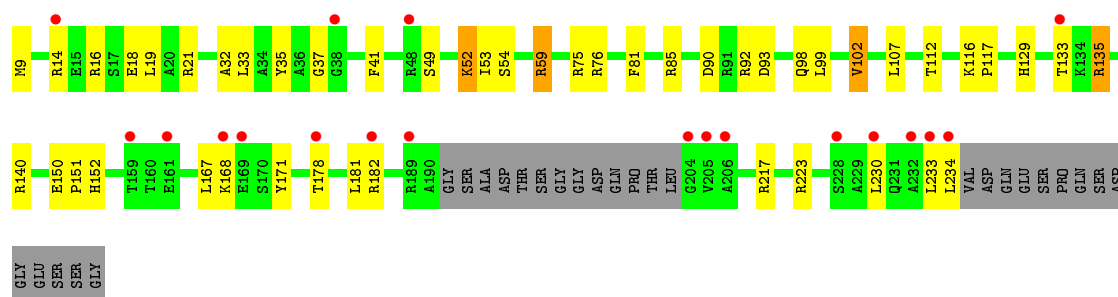


T301 L306 N335 E362 E363 Y366 L374 R382 I385 Q396 S397 L398 L401 P402 I409 H410 A411 S412 D413 P414 Q415 W429 S448 D461 L471 Y472 L513 S522 GLY ALA ASP THR PHE GLY GLY SER ASP GLY LYS HIS HIS HIS HIS HIS

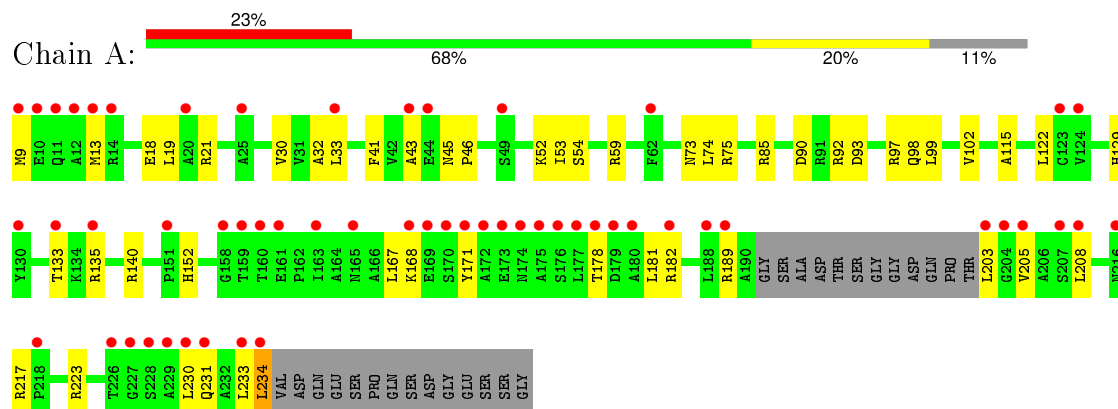
HIS
HIS

• Molecule 2: Proteasome (Alpha subunit) PrcA

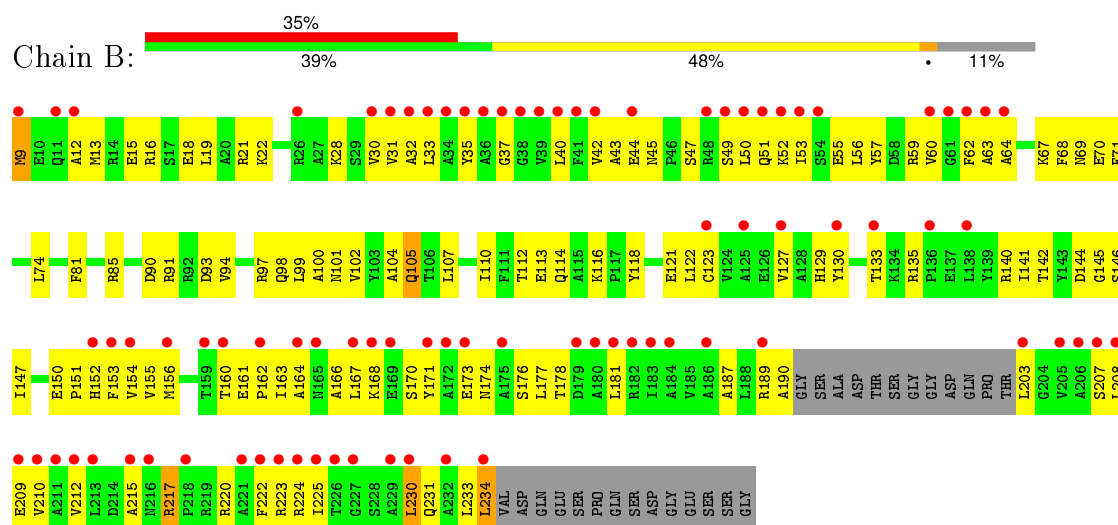




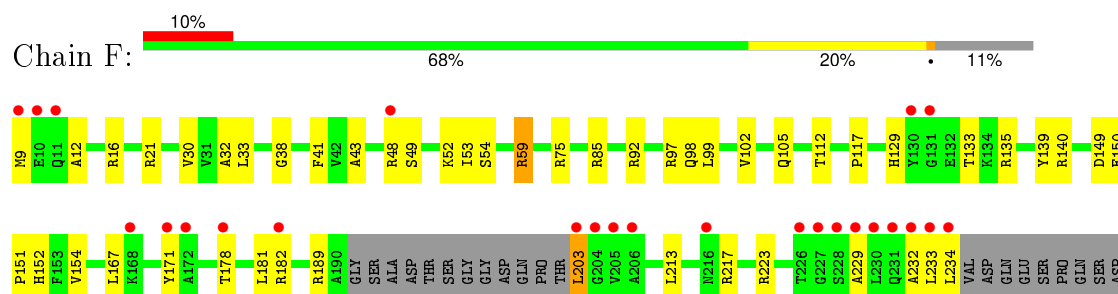
• Molecule 2: Proteasome (Alpha subunit) PrcA



• Molecule 2: Proteasome (Alpha subunit) PrcA

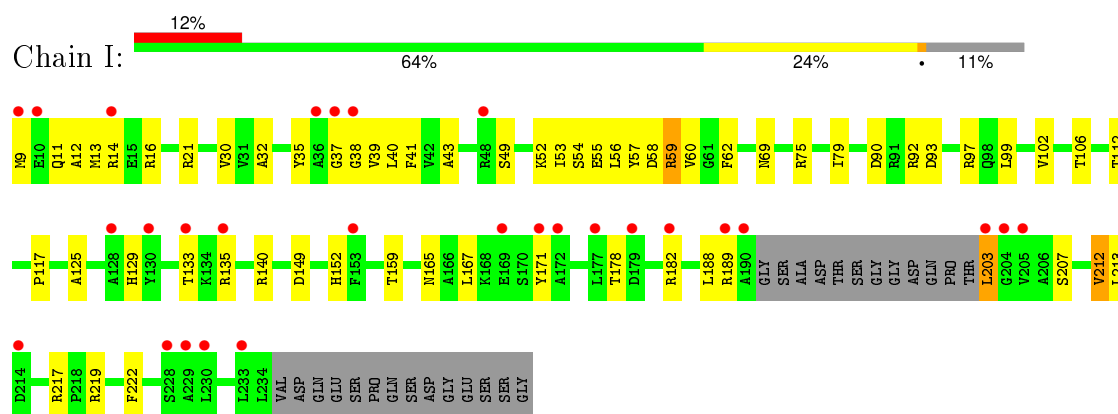


• Molecule 2: Proteasome (Alpha subunit) PrcA

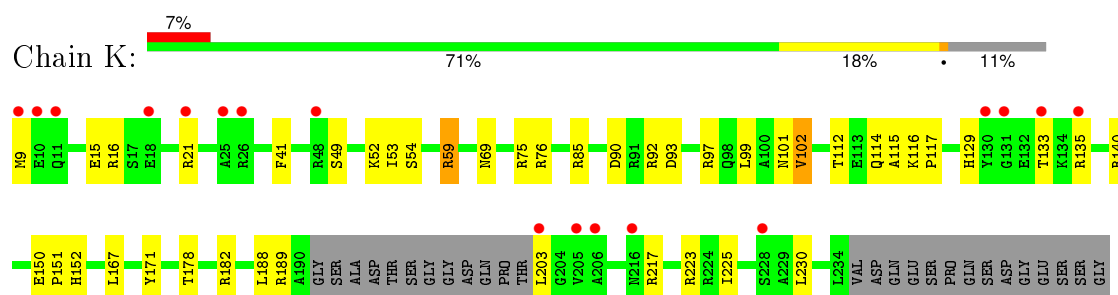


GLY
GLU
SER
SER
GLY

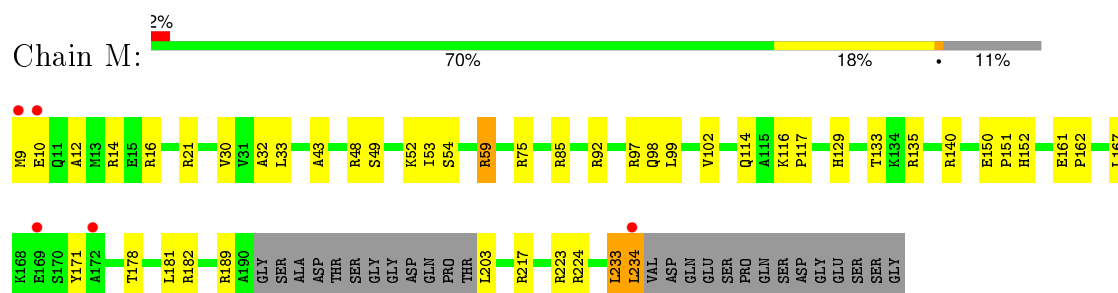
• Molecule 2: Proteasome (Alpha subunit) PrcA



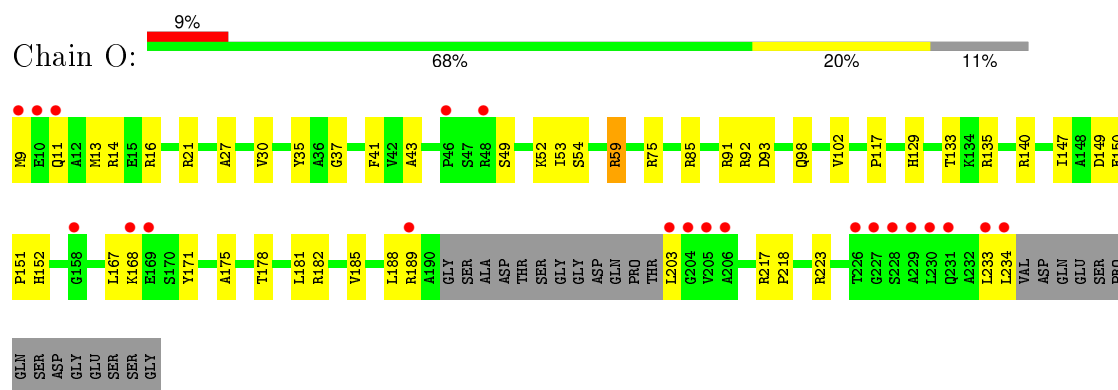
• Molecule 2: Proteasome (Alpha subunit) PrcA



• Molecule 2: Proteasome (Alpha subunit) PrcA

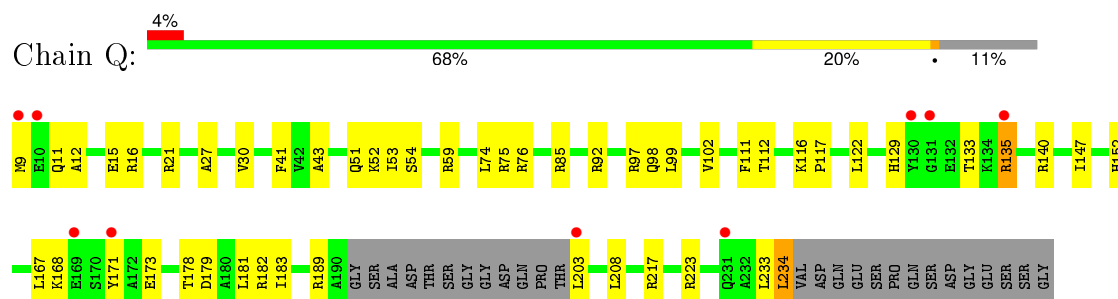


• Molecule 2: Proteasome (Alpha subunit) PrcA

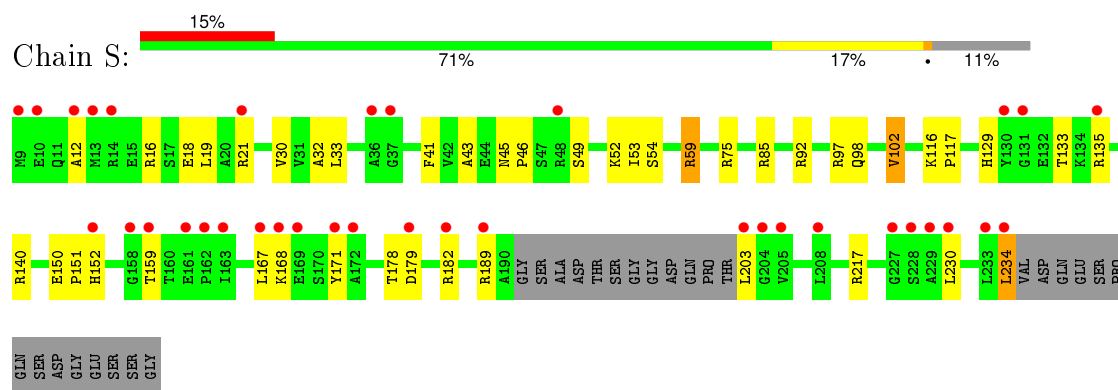


GLN
SER
ASP
GLY
GLU
SER
SER
GLY

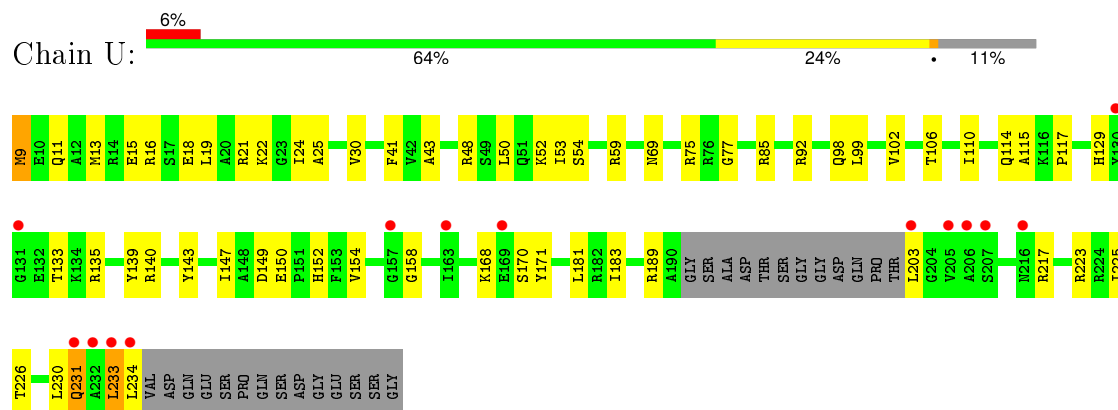
- Molecule 2: Proteasome (Alpha subunit) PrcA



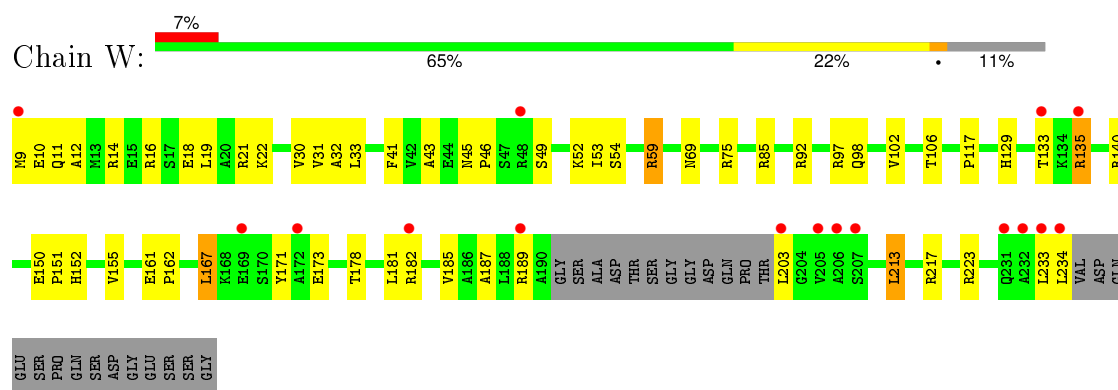
- Molecule 2: Proteasome (Alpha subunit) PrcA



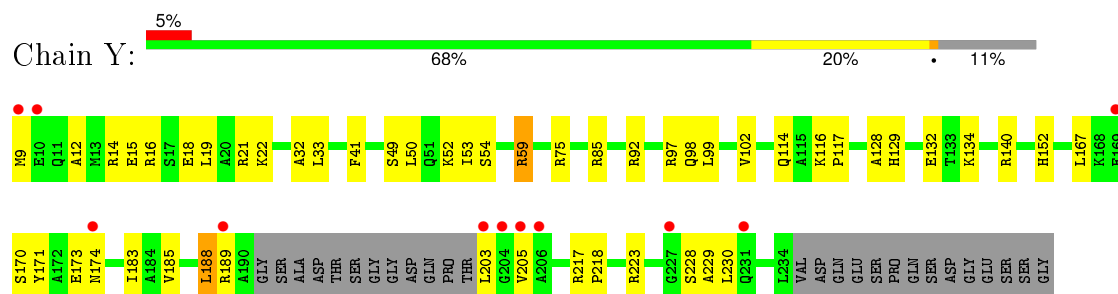
- Molecule 2: Proteasome (Alpha subunit) PrcA



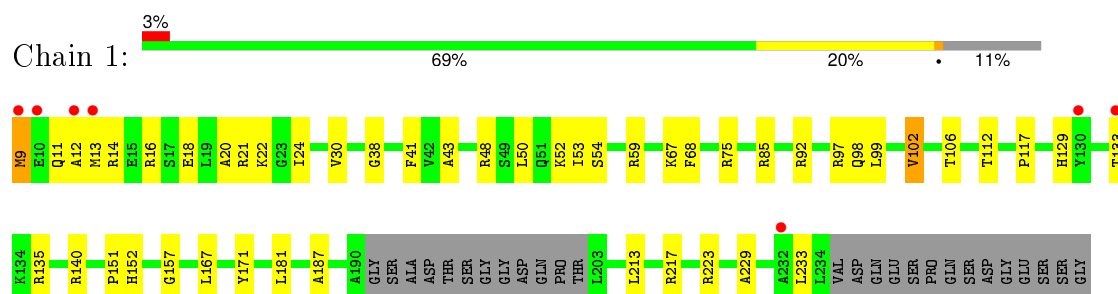
- Molecule 2: Proteasome (Alpha subunit) PrcA



- Molecule 2: Proteasome (Alpha subunit) PrcA



- Molecule 2: Proteasome (Alpha subunit) PrcA



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	119.16Å 221.44Å 137.12Å 90.00° 104.89° 90.00°	Depositor
Resolution (Å)	29.86 – 2.50 29.86 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.8 (29.86-2.50) 97.8 (29.86-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.173 , 0.214 0.198 , 0.227	Depositor DCC
R_{free} test set	11466 reflections (4.99%)	DCC
Wilson B-factor (Å ²)	33.3	Xtriage
Anisotropy	0.164	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 55.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 229841 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	48523	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.15% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 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	2	0.33	0/1662	0.49	0/2254
1	C	0.32	0/1662	0.50	0/2254
1	E	0.34	0/1666	0.49	0/2259
1	G	0.36	0/1662	0.50	0/2254
1	H	0.33	0/1666	0.48	0/2259
1	J	0.31	0/1662	0.48	0/2254
1	L	0.33	0/1662	0.49	0/2254
1	N	0.35	0/1666	0.49	0/2259
1	P	0.32	0/1662	0.49	0/2254
1	R	0.33	0/1662	0.49	0/2254
1	T	0.33	0/1666	0.49	0/2259
1	V	0.36	0/1686	0.50	0/2287
1	X	0.32	0/1671	0.49	0/2266
1	Z	0.32	0/1671	0.50	0/2266
2	1	0.32	0/1677	0.54	0/2264
2	A	0.32	0/1677	0.53	0/2264
2	B	0.32	0/1677	0.56	0/2264
2	D	0.36	0/1669	0.54	0/2253
2	F	0.32	0/1677	0.54	0/2264
2	I	0.32	0/1677	0.55	1/2264 (0.0%)
2	K	0.32	0/1677	0.53	0/2264
2	M	0.32	0/1677	0.54	0/2264
2	O	0.31	0/1677	0.51	0/2264
2	Q	0.33	0/1677	0.57	0/2264
2	S	0.29	0/1677	0.51	0/2264
2	U	0.32	0/1677	0.54	0/2264
2	W	0.33	0/1677	0.55	1/2264 (0.0%)
2	Y	0.33	0/1677	0.55	0/2264
All	All	0.33	0/46796	0.52	2/63318 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	I	213	LEU	N-CA-C	-5.19	96.99	111.00
2	W	213	LEU	N-CA-C	-5.11	97.20	111.00

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	2	1638	0	1630	16	0
1	C	1638	0	1630	23	0
1	E	1642	0	1633	28	0
1	G	1638	0	1630	21	0
1	H	1642	0	1633	15	0
1	J	1638	0	1630	15	0
1	L	1638	0	1630	21	0
1	N	1642	0	1633	19	0
1	P	1638	0	1630	18	0
1	R	1638	0	1630	23	0
1	T	1642	0	1633	19	0
1	V	1662	0	1649	22	0
1	X	1647	0	1638	16	0
1	Z	1647	0	1638	22	0
2	1	1653	0	1656	49	0
2	A	1653	0	1656	42	0
2	B	1653	0	1656	161	0
2	D	1645	0	1645	49	0
2	F	1653	0	1656	40	0
2	I	1653	0	1656	62	0
2	K	1653	0	1656	35	0
2	M	1653	0	1656	38	0
2	O	1653	0	1656	48	0
2	Q	1653	0	1656	39	0
2	S	1653	0	1656	51	0
2	U	1653	0	1656	47	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	W	1653	0	1656	52	0
2	Y	1653	0	1656	46	0
3	1	10	0	14	2	0
3	2	10	0	14	2	0
3	A	10	0	14	1	0
3	B	5	0	7	0	0
3	C	5	0	7	0	0
3	D	5	0	7	0	0
3	E	15	0	21	4	0
3	F	5	0	7	0	0
3	G	15	0	21	4	0
3	H	10	0	14	1	0
3	I	5	0	7	1	0
3	J	5	0	7	0	0
3	K	15	0	21	5	0
3	L	20	0	28	4	0
3	M	5	0	7	0	0
3	N	10	0	14	1	0
3	O	5	0	7	0	0
3	P	15	0	21	3	0
3	Q	15	0	21	5	0
3	S	10	0	14	1	0
3	T	10	0	14	5	0
3	U	10	0	14	1	0
3	V	25	0	35	7	0
3	W	5	0	7	1	0
3	X	5	0	7	0	0
3	Y	15	0	21	1	0
3	Z	25	0	35	7	0
4	1	68	0	0	1	0
4	2	107	0	0	3	0
4	A	31	0	0	0	0
4	B	37	0	0	4	0
4	C	43	0	0	1	0
4	D	49	0	0	5	0
4	E	121	0	0	8	0
4	F	60	0	0	4	0
4	G	113	0	0	2	0
4	H	92	0	0	0	0
4	I	35	0	0	6	0
4	J	73	0	0	2	0
4	K	52	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	L	109	0	0	4	0
4	M	74	0	0	5	0
4	N	124	0	0	6	0
4	O	34	0	0	4	0
4	P	93	0	0	2	0
4	Q	71	0	0	2	0
4	R	89	0	0	2	0
4	S	35	0	0	3	0
4	T	96	0	0	1	0
4	U	48	0	0	2	0
4	V	133	0	0	2	0
4	W	62	0	0	4	0
4	X	101	0	0	1	0
4	Y	68	0	0	3	0
4	Z	91	0	0	3	0
All	All	48523	0	46446	945	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 10.

All (945) 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:A:115:ALA:HB3	2:B:112:THR:CG2	1.66	1.25
2:S:116:LYS:HB2	2:1:13:MET:CE	1.68	1.23
2:B:9:MET:CE	2:B:13:MET:HE3	1.68	1.22
2:S:230:LEU:O	2:S:234:LEU:HD13	1.34	1.22
2:O:9:MET:HE2	2:O:13:MET:CE	1.77	1.15
2:B:64:ALA:HB2	2:B:122:LEU:CD2	1.79	1.12
1:Z:509:ARG:HH12	3:Z:69:DMF:H23	1.15	1.11
2:B:217:ARG:HG2	2:B:217:ARG:HH11	1.16	1.10
2:A:115:ALA:HB3	2:B:112:THR:HG23	1.32	1.08
2:B:64:ALA:CB	2:B:122:LEU:HD23	1.84	1.08
2:O:9:MET:CE	2:O:13:MET:CE	2.33	1.07
2:B:9:MET:CE	2:B:13:MET:CE	2.33	1.05
2:D:16:ARG:HG2	2:K:9:MET:HE1	1.37	1.05
2:B:64:ALA:HB2	2:B:122:LEU:HD23	1.36	1.03
2:O:9:MET:CE	2:O:13:MET:HE3	1.88	1.02
2:B:9:MET:HE1	2:B:13:MET:HE3	1.41	1.02
2:S:116:LYS:HB2	2:1:13:MET:HE2	1.41	1.01
2:I:152:HIS:HB3	2:I:171:TYR:CE2	1.97	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:11:GLN:HA	2:I:11:GLN:OE1	1.57	0.99
2:I:11:GLN:HG3	2:I:14:ARG:NH1	1.79	0.98
2:S:116:LYS:HB2	2:I:13:MET:HE3	1.43	0.98
1:C:306:LEU:HD12	1:C:313:VAL:HG12	1.46	0.97
1:Z:496:ILE:HD12	3:Z:69:DMF:H22	1.46	0.97
2:B:35:TYR:CZ	2:B:177:LEU:HD13	1.99	0.96
2:B:181:LEU:HD23	2:B:181:LEU:O	1.66	0.95
2:B:217:ARG:HG3	2:B:223:ARG:HH21	1.29	0.95
2:B:35:TYR:OH	2:B:177:LEU:HD13	1.67	0.94
1:C:306:LEU:CD1	1:C:313:VAL:CG1	2.46	0.93
2:B:35:TYR:CZ	2:B:177:LEU:CD1	2.52	0.93
2:I:11:GLN:CG	2:I:14:ARG:NH1	2.32	0.93
2:W:181:LEU:O	2:W:181:LEU:HD23	1.69	0.92
2:O:9:MET:HE2	2:O:13:MET:HE2	1.48	0.92
2:O:9:MET:HE2	2:O:13:MET:HE3	1.42	0.92
2:B:31:VAL:CG1	2:B:33:LEU:HD11	2.02	0.89
2:B:181:LEU:HD23	2:B:181:LEU:C	1.93	0.88
2:I:11:GLN:HG3	2:I:14:ARG:HH12	1.37	0.88
2:A:115:ALA:CB	2:B:112:THR:HG23	2.04	0.87
2:B:155:VAL:HG11	2:B:164:ALA:HB2	1.54	0.87
2:B:189:ARG:HH21	2:B:203:LEU:HD22	1.40	0.87
2:D:16:ARG:HG2	2:K:9:MET:CE	2.03	0.87
2:D:35:TYR:CE1	2:D:37:GLY:HA3	2.09	0.87
1:C:306:LEU:HD11	1:C:313:VAL:CG1	2.04	0.87
2:B:217:ARG:HG2	2:B:217:ARG:NH1	1.87	0.86
2:B:9:MET:HE3	2:B:13:MET:HE3	1.57	0.86
2:I:35:TYR:CD2	2:I:38:GLY:O	2.30	0.85
1:C:306:LEU:CD1	1:C:313:VAL:HG12	2.05	0.85
2:D:85:ARG:NH2	3:Q:251:DMF:H23	1.92	0.85
2:M:59:ARG:HD2	2:M:129:HIS:HA	1.58	0.85
2:I:203:LEU:N	2:I:203:LEU:HD12	1.90	0.85
2:O:59:ARG:HD2	2:O:129:HIS:HA	1.58	0.85
2:B:64:ALA:CB	2:B:122:LEU:CD2	2.49	0.84
1:Z:509:ARG:NH1	3:Z:69:DMF:H23	1.92	0.84
2:F:59:ARG:HD2	2:F:129:HIS:HA	1.60	0.84
2:I:59:ARG:HD2	2:I:129:HIS:HA	1.58	0.84
2:A:9:MET:CE	2:A:9:MET:HA	2.07	0.84
2:K:59:ARG:HD2	2:K:129:HIS:HA	1.60	0.83
2:Q:59:ARG:HD2	2:Q:129:HIS:HA	1.61	0.83
1:H:398:LEU:HD11	1:P:391:LEU:HD21	1.60	0.83
2:D:14:ARG:HD3	4:D:1624:HOH:O	1.78	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:S:19:LEU:HD13	2:1:9:MET:CE	2.08	0.83
2:S:59:ARG:HD2	2:S:129:HIS:HA	1.60	0.82
2:Q:178:THR:O	2:Q:182:ARG:HG2	1.80	0.82
2:D:59:ARG:HD2	2:D:129:HIS:HA	1.60	0.82
2:B:49:SER:OG	2:I:97:ARG:HD2	1.78	0.82
2:S:230:LEU:CD1	2:S:234:LEU:HD11	2.10	0.82
2:U:59:ARG:HD2	2:U:129:HIS:HA	1.61	0.81
2:S:19:LEU:HD13	2:1:9:MET:HE3	1.63	0.81
2:1:59:ARG:HD2	2:1:129:HIS:HA	1.59	0.81
2:B:33:LEU:HG	2:B:153:PHE:HB2	1.62	0.81
2:A:59:ARG:HD2	2:A:129:HIS:HA	1.59	0.81
2:W:59:ARG:HD2	2:W:129:HIS:HA	1.62	0.81
2:B:30:VAL:HG12	2:B:43:ALA:CB	2.12	0.80
2:O:85:ARG:NH1	2:O:98:GLN:NE2	2.30	0.80
2:A:115:ALA:CB	2:B:112:THR:CG2	2.56	0.80
1:C:306:LEU:HD12	1:C:313:VAL:CG1	2.08	0.80
2:B:209:GLU:HB3	4:B:1703:HOH:O	1.80	0.79
2:O:35:TYR:CE1	2:O:37:GLY:HA3	2.17	0.79
2:W:189:ARG:HH12	2:W:203:LEU:N	1.81	0.78
1:T:395:MET:HE2	4:2:1676:HOH:O	1.84	0.78
1:T:472:TYR:CZ	3:T:62:DMF:H13	2.18	0.77
2:B:49:SER:HG	2:I:97:ARG:HD2	1.48	0.77
2:B:33:LEU:HG	2:B:153:PHE:CB	2.14	0.77
2:O:9:MET:HE1	2:O:13:MET:HE3	1.66	0.76
2:B:166:ALA:HB3	2:B:187:ALA:CB	2.16	0.76
2:S:230:LEU:HD12	2:S:234:LEU:HD11	1.66	0.76
2:W:181:LEU:C	2:W:181:LEU:HD23	2.06	0.76
2:B:19:LEU:O	2:B:19:LEU:HD23	1.85	0.76
2:W:167:LEU:HD13	2:W:187:ALA:HB2	1.68	0.76
3:H:142:DMF:HC	3:L:9:DMF:O	1.85	0.76
2:W:167:LEU:HD13	2:W:187:ALA:CB	2.16	0.75
2:B:47:SER:HB2	2:I:149:ASP:OD2	1.87	0.75
2:B:166:ALA:HB3	2:B:187:ALA:HB2	1.69	0.75
2:B:189:ARG:NH2	2:B:203:LEU:HD22	2.01	0.75
2:S:12:ALA:O	2:S:16:ARG:HG3	1.86	0.75
2:I:178:THR:HG22	2:I:182:ARG:HH21	1.52	0.74
2:Q:76:ARG:NH1	3:Q:251:DMF:H22	2.02	0.74
2:B:30:VAL:HG12	2:B:43:ALA:HB1	1.67	0.74
1:G:337:THR:HG21	1:G:359:TYR:HD2	1.52	0.74
2:D:99:LEU:O	2:D:102:VAL:HG12	1.88	0.74
1:H:362:GLU:OE2	1:H:382:ARG:HD3	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:18:GLU:CD	2:A:21:ARG:HH12	1.90	0.73
2:B:31:VAL:HG12	2:B:33:LEU:CD1	2.18	0.73
2:I:99:LEU:O	2:I:102:VAL:HG12	1.89	0.73
2:U:231:GLN:OE1	2:U:231:GLN:HA	1.86	0.73
2:B:31:VAL:CG1	2:B:33:LEU:CD1	2.67	0.72
2:B:231:GLN:O	2:B:234:LEU:HG	1.88	0.72
2:W:189:ARG:NH1	2:W:203:LEU:N	2.38	0.72
2:I:159:THR:HG23	2:I:159:THR:O	1.88	0.72
2:B:181:LEU:CD2	2:B:181:LEU:C	2.57	0.71
2:Y:203:LEU:HD12	2:Y:203:LEU:N	2.06	0.71
1:N:414:PRO:HD2	4:N:1177:HOH:O	1.89	0.71
2:U:85:ARG:NH1	2:U:98:GLN:NE2	2.38	0.71
2:W:152:HIS:HB3	2:W:171:TYR:CE2	2.25	0.71
2:A:205:VAL:HG21	2:A:231:GLN:OE1	1.89	0.71
2:B:59:ARG:HD2	2:B:127:VAL:HG13	1.71	0.70
2:O:9:MET:HE1	2:O:13:MET:CE	2.22	0.70
2:W:135:ARG:NH2	2:W:173:GLU:OE2	2.23	0.70
2:W:106:THR:HG21	3:W:249:DMF:HC	1.71	0.70
2:B:32:ALA:C	2:B:33:LEU:HD12	2.11	0.70
2:Q:152:HIS:HB3	2:Q:171:TYR:CE2	2.26	0.70
2:B:9:MET:HE2	2:B:13:MET:CE	2.20	0.69
2:I:11:GLN:HG2	2:I:14:ARG:NH1	2.06	0.69
2:1:18:GLU:HG3	2:1:22:LYS:HE3	1.74	0.69
2:A:115:ALA:HB3	2:B:112:THR:HG22	1.71	0.69
2:O:21:ARG:HB3	2:O:21:ARG:HH11	1.57	0.69
2:B:166:ALA:CB	2:B:187:ALA:HA	2.23	0.69
2:O:11:GLN:HG3	2:O:14:ARG:CZ	2.23	0.69
2:Q:97:ARG:HD2	2:Y:49:SER:HB2	1.75	0.69
2:K:21:ARG:HB3	2:K:21:ARG:HH11	1.58	0.69
2:O:21:ARG:HB3	2:O:21:ARG:NH1	2.09	0.68
2:S:21:ARG:HH11	2:S:21:ARG:HB3	1.58	0.68
2:D:21:ARG:NH1	2:D:21:ARG:HB3	2.09	0.68
2:S:21:ARG:NH1	2:S:21:ARG:HB3	2.09	0.68
2:D:21:ARG:HH11	2:D:21:ARG:HB3	1.58	0.67
2:W:30:VAL:HG13	2:W:43:ALA:HB2	1.76	0.67
1:T:362:GLU:OE2	1:T:382:ARG:HD3	1.94	0.67
2:I:189:ARG:NH1	2:I:203:LEU:N	2.42	0.67
2:Q:12:ALA:O	2:Q:16:ARG:HG3	1.95	0.67
2:K:21:ARG:HB3	2:K:21:ARG:NH1	2.09	0.67
2:F:21:ARG:HB3	2:F:21:ARG:NH1	2.10	0.67
2:S:116:LYS:CB	2:1:13:MET:HE2	2.23	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:92:ARG:HD3	2:M:129:HIS:CE1	2.30	0.67
2:B:166:ALA:O	2:B:170:SER:HB3	1.93	0.66
1:Z:362:GLU:OE2	1:Z:382:ARG:HD3	1.95	0.66
2:B:164:ALA:O	2:B:168:LYS:HB2	1.95	0.66
2:S:116:LYS:CB	2:I:13:MET:HE3	2.23	0.66
2:F:21:ARG:HH11	2:F:21:ARG:HB3	1.61	0.66
2:S:117:PRO:HD2	2:I:9:MET:HE1	1.77	0.66
2:Y:16:ARG:NH2	2:Y:114:GLN:O	2.23	0.66
2:W:178:THR:HG22	2:W:182:ARG:HH21	1.60	0.66
2:B:189:ARG:HA	2:B:189:ARG:NE	2.10	0.66
2:B:225:ILE:O	2:B:225:ILE:HG13	1.94	0.66
2:M:21:ARG:NH1	2:M:21:ARG:HB3	2.11	0.66
2:F:178:THR:HG22	2:F:182:ARG:HH21	1.61	0.66
2:I:21:ARG:HH11	2:I:21:ARG:HB3	1.61	0.66
2:I:92:ARG:HD3	2:I:129:HIS:CE1	2.31	0.66
2:M:21:ARG:HH11	2:M:21:ARG:HB3	1.60	0.66
2:I:21:ARG:HB3	2:I:21:ARG:NH1	2.11	0.66
2:K:178:THR:HG22	2:K:182:ARG:HH21	1.61	0.66
1:C:306:LEU:HD11	1:C:313:VAL:HG11	1.76	0.65
2:A:178:THR:HG22	2:A:182:ARG:HH21	1.61	0.65
2:U:9:MET:HE1	2:U:13:MET:CE	2.26	0.65
2:O:178:THR:HG22	2:O:182:ARG:HH21	1.61	0.65
2:W:19:LEU:HD23	2:W:19:LEU:C	2.16	0.65
2:U:92:ARG:HD3	2:U:129:HIS:CE1	2.32	0.65
2:K:15:GLU:OE1	2:M:9:MET:N	2.30	0.65
1:X:362:GLU:OE2	1:X:382:ARG:HD3	1.96	0.65
2:A:90:ASP:HB3	2:A:93:ASP:OD2	1.95	0.65
2:W:11:GLN:OE1	2:W:14:ARG:NH1	2.30	0.65
2:Q:189:ARG:HH12	2:Q:203:LEU:N	1.95	0.65
2:Q:21:ARG:HB3	2:Q:21:ARG:NH1	2.12	0.65
2:B:9:MET:HE3	2:B:13:MET:CE	2.18	0.65
2:B:15:GLU:CD	2:I:9:MET:N	2.50	0.65
1:N:362:GLU:OE2	1:N:382:ARG:HD3	1.96	0.65
2:Y:128:ALA:HB1	2:Y:132:GLU:HG3	1.78	0.65
2:D:85:ARG:HH21	3:Q:251:DMF:H23	1.60	0.64
2:B:163:ILE:O	2:B:187:ALA:HB1	1.97	0.64
2:B:121:GLU:OE2	2:B:140:ARG:NH1	2.30	0.64
2:B:33:LEU:HD21	2:B:167:LEU:HD13	1.78	0.64
2:Q:21:ARG:HH11	2:Q:21:ARG:HB3	1.62	0.64
2:B:31:VAL:HG11	2:B:33:LEU:HD11	1.77	0.64
1:R:362:GLU:OE2	1:R:382:ARG:HD3	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:35:TYR:OH	2:B:177:LEU:CD1	2.40	0.64
2:F:92:ARG:HD3	2:F:129:HIS:CE1	2.32	0.64
2:1:229:ALA:O	2:1:233:LEU:HD13	1.97	0.64
2:M:178:THR:HG22	2:M:182:ARG:HH21	1.63	0.64
2:U:9:MET:CE	2:U:13:MET:CE	2.76	0.64
2:I:106:THR:HG21	3:I:249:DMF:HC	1.79	0.64
2:1:181:LEU:HD23	2:1:233:LEU:HB3	1.80	0.64
1:G:362:GLU:OE2	1:G:382:ARG:HD3	1.97	0.64
1:X:391:LEU:HD21	1:Z:398:LEU:HD11	1.80	0.64
1:E:362:GLU:OE2	1:E:382:ARG:HD3	1.98	0.64
1:T:391:LEU:HD21	1:2:398:LEU:HD11	1.79	0.64
2:D:90:ASP:HB3	2:D:93:ASP:OD2	1.98	0.64
2:Y:152:HIS:HB3	2:Y:171:TYR:CE2	2.33	0.64
1:T:382:ARG:NH2	1:T:385:ILE:HD13	2.13	0.63
2:D:178:THR:HG22	2:D:182:ARG:HH21	1.61	0.63
2:F:97:ARG:HD2	2:M:49:SER:HB2	1.80	0.63
2:I:165:ASN:ND2	4:I:1026:HOH:O	2.30	0.63
2:B:35:TYR:CZ	2:B:177:LEU:HD12	2.31	0.63
1:H:398:LEU:HD11	1:P:391:LEU:CD2	2.29	0.63
2:O:85:ARG:HH12	2:O:98:GLN:NE2	1.96	0.63
2:O:11:GLN:NE2	4:O:2134:HOH:O	2.30	0.63
2:M:161:GLU:HB3	2:M:162:PRO:HD3	1.79	0.63
1:V:362:GLU:OE2	1:V:382:ARG:HD3	1.98	0.63
2:1:92:ARG:HD3	2:1:129:HIS:CE1	2.33	0.63
2:B:19:LEU:HD23	2:B:19:LEU:C	2.18	0.63
1:C:362:GLU:OE2	1:C:382:ARG:HD3	1.99	0.63
2:B:47:SER:CB	4:I:1752:HOH:O	2.47	0.63
2:Y:189:ARG:NH1	2:Y:203:LEU:HD12	2.14	0.63
1:R:382:ARG:NH2	1:R:385:ILE:HD13	2.13	0.63
1:E:382:ARG:NH2	1:E:385:ILE:HD13	2.13	0.63
2:M:85:ARG:HH12	2:M:98:GLN:NE2	1.96	0.63
2:S:159:THR:HA	4:S:1892:HOH:O	1.98	0.63
2:K:92:ARG:HD3	2:K:129:HIS:CE1	2.34	0.63
2:O:151:PRO:HD2	4:O:249:HOH:O	1.97	0.63
2:B:123:CYS:SG	2:B:154:VAL:HG21	2.39	0.63
2:B:217:ARG:CG	2:B:223:ARG:HH21	2.08	0.63
2:D:92:ARG:HD3	2:D:129:HIS:CE1	2.34	0.63
2:S:116:LYS:CB	2:1:13:MET:CE	2.62	0.62
2:D:16:ARG:CG	2:K:9:MET:HE1	2.22	0.62
2:I:152:HIS:CB	2:I:171:TYR:CE2	2.80	0.62
2:Q:85:ARG:HH12	2:Q:98:GLN:NE2	1.96	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:366:TYR:CD2	1:V:374:LEU:HD13	2.35	0.62
2:S:85:ARG:HH12	2:S:98:GLN:NE2	1.96	0.62
2:A:92:ARG:HD3	2:A:129:HIS:CE1	2.34	0.62
2:W:85:ARG:HH12	2:W:98:GLN:NE2	1.98	0.62
2:W:18:GLU:OE2	2:W:21:ARG:NH1	2.30	0.62
2:Q:152:HIS:HB3	2:Q:171:TYR:CZ	2.35	0.62
1:P:362:GLU:OE2	1:P:382:ARG:HD3	1.98	0.62
2:W:181:LEU:C	2:W:181:LEU:CD2	2.68	0.62
2:O:149:ASP:OD2	2:U:48:ARG:HG2	1.99	0.62
2:D:81:PHE:HE2	2:D:98:GLN:HE21	1.47	0.62
2:S:92:ARG:HD3	2:S:129:HIS:CE1	2.35	0.62
2:W:92:ARG:HD3	2:W:129:HIS:CE1	2.35	0.62
2:B:152:HIS:HB3	2:B:171:TYR:CZ	2.35	0.62
2:A:203:LEU:HD23	2:A:208:LEU:HD21	1.80	0.62
1:G:429:TRP:CH2	3:G:140:DMF:H13	2.35	0.62
1:2:362:GLU:OE2	1:2:382:ARG:HD3	1.99	0.62
2:U:11:GLN:O	2:U:15:GLU:HG3	2.00	0.62
2:S:19:LEU:HD13	2:1:9:MET:HE1	1.80	0.61
2:W:12:ALA:O	2:W:16:ARG:HG3	1.99	0.61
2:Y:59:ARG:HD2	2:Y:129:HIS:HA	1.81	0.61
2:Y:85:ARG:HH12	2:Y:98:GLN:NE2	1.98	0.61
2:B:217:ARG:HD3	2:B:220:ARG:O	2.00	0.61
2:I:203:LEU:N	2:I:203:LEU:CD1	2.63	0.61
2:O:92:ARG:HD3	2:O:129:HIS:CE1	2.35	0.61
2:A:205:VAL:CG2	2:A:231:GLN:OE1	2.47	0.61
2:1:12:ALA:O	2:1:16:ARG:HG3	2.01	0.61
1:L:362:GLU:OE2	1:L:382:ARG:HD3	2.01	0.61
2:1:18:GLU:OE2	2:1:21:ARG:NH2	2.30	0.61
1:X:456:GLN:HE22	1:X:465:ARG:NH2	1.99	0.61
2:S:234:LEU:N	2:S:234:LEU:CD1	2.63	0.61
2:B:166:ALA:HB3	2:B:187:ALA:CA	2.30	0.61
2:U:85:ARG:NH1	2:U:98:GLN:HE22	1.97	0.61
2:Y:152:HIS:HB3	2:Y:171:TYR:CZ	2.35	0.61
2:D:35:TYR:CE1	2:D:37:GLY:CA	2.81	0.61
1:H:382:ARG:NH2	1:H:385:ILE:HD13	2.15	0.61
1:Z:509:ARG:HH12	3:Z:69:DMF:C2	2.04	0.61
1:C:382:ARG:NH2	1:C:385:ILE:HD13	2.16	0.61
1:L:444:LEU:HB2	4:L:290:HOH:O	2.01	0.61
2:B:47:SER:HB3	4:I:1752:HOH:O	2.00	0.60
2:B:166:ALA:HB3	2:B:187:ALA:HA	1.82	0.60
2:Q:92:ARG:HD3	2:Q:129:HIS:CE1	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:391:LEU:HD21	1:N:398:LEU:HD11	1.83	0.60
2:B:68:PHE:HA	2:B:71:PHE:CE2	2.36	0.60
2:S:117:PRO:CD	2:1:9:MET:CE	2.78	0.60
2:B:127:VAL:CG2	2:B:215:ALA:HB2	2.30	0.60
2:B:127:VAL:HG21	2:B:215:ALA:HB2	1.83	0.60
2:D:135:ARG:NH1	4:D:2073:HOH:O	2.32	0.60
1:R:366:TYR:CD2	1:R:374:LEU:HD13	2.36	0.60
1:V:382:ARG:NH2	1:V:385:ILE:HD13	2.16	0.60
2:A:85:ARG:HH12	2:A:98:GLN:NE2	2.00	0.60
1:C:456:GLN:HE22	1:C:465:ARG:NH2	2.00	0.60
2:B:152:HIS:HB3	2:B:171:TYR:CE2	2.37	0.60
2:U:25:ALA:O	2:U:158:GLY:HA2	2.02	0.60
2:K:90:ASP:HB3	2:K:93:ASP:OD2	2.02	0.60
3:V:16:DMF:HC	4:2:554:HOH:O	2.01	0.60
2:A:9:MET:HE1	2:A:9:MET:HA	1.84	0.59
2:Q:203:LEU:HD13	2:Q:208:LEU:HD21	1.84	0.59
1:N:382:ARG:NH2	1:N:385:ILE:HD13	2.17	0.59
1:J:456:GLN:HE22	1:J:465:ARG:NH2	1.99	0.59
2:U:15:GLU:O	2:U:18:GLU:HG2	2.03	0.59
2:S:49:SER:HB2	2:1:97:ARG:HD2	1.83	0.59
1:V:429:TRP:CZ2	3:V:121:DMF:H22	2.36	0.59
2:W:49:SER:HB2	2:Y:97:ARG:HD2	1.83	0.59
1:J:362:GLU:OE2	1:J:382:ARG:HD3	2.01	0.59
2:U:77:GLY:HA3	3:U:249:DMF:O	2.01	0.59
1:E:380:ILE:HB	3:E:104:DMF:H13	1.85	0.59
1:H:366:TYR:CD2	1:H:374:LEU:HD13	2.38	0.59
2:B:35:TYR:CE1	2:B:37:GLY:HA3	2.37	0.59
1:X:382:ARG:NH2	1:X:385:ILE:HD13	2.17	0.59
1:G:448:SER:HB3	1:2:448:SER:HB3	1.83	0.59
2:B:49:SER:OG	2:I:97:ARG:CD	2.51	0.59
2:I:59:ARG:NH2	2:I:217:ARG:O	2.35	0.59
2:S:230:LEU:HD12	2:S:234:LEU:CD1	2.33	0.59
1:P:396:GLN:NE2	4:P:1778:HOH:O	2.35	0.59
2:M:181:LEU:HD23	2:M:233:LEU:HB3	1.84	0.59
2:B:40:LEU:HA	2:B:212:VAL:HG12	1.84	0.59
2:B:85:ARG:HB3	2:B:93:ASP:OD2	2.02	0.59
2:B:33:LEU:HD21	2:B:167:LEU:HD22	1.85	0.58
1:X:366:TYR:CD2	1:X:374:LEU:HD13	2.38	0.58
2:Y:205:VAL:HG13	2:Y:230:LEU:HD23	1.84	0.58
2:B:64:ALA:HB1	2:B:122:LEU:HD23	1.81	0.58
2:I:11:GLN:HG3	2:I:14:ARG:CZ	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:382:ARG:NH2	1:J:385:ILE:HD13	2.17	0.58
2:B:18:GLU:HA	2:B:21:ARG:HG2	1.85	0.58
1:P:456:GLN:HE22	1:P:465:ARG:NH2	2.00	0.58
2:S:230:LEU:HD11	2:S:234:LEU:HD11	1.86	0.58
2:B:97:ARG:O	2:B:101:ASN:HB2	2.03	0.58
1:P:382:ARG:NH2	1:P:385:ILE:HD13	2.19	0.58
1:H:456:GLN:HE22	1:H:465:ARG:NH2	2.01	0.58
1:T:456:GLN:HE22	1:T:465:ARG:NH2	2.01	0.58
2:W:18:GLU:O	2:W:22:LYS:HG3	2.04	0.58
2:F:85:ARG:HH12	2:F:98:GLN:NE2	2.02	0.58
2:U:9:MET:HG3	2:1:16:ARG:HG2	1.86	0.58
1:Z:456:GLN:HE22	1:Z:465:ARG:NH2	2.01	0.58
2:O:9:MET:CE	2:O:13:MET:HE2	2.18	0.58
1:V:366:TYR:CE2	1:V:374:LEU:HD13	2.38	0.58
3:Y:249:DMF:H21	4:Y:944:HOH:O	2.03	0.58
2:I:178:THR:CG2	2:I:182:ARG:HH21	2.17	0.58
2:B:28:LYS:HE3	2:B:44:GLU:HB3	1.85	0.58
1:G:337:THR:HG21	1:G:359:TYR:CD2	2.37	0.57
1:E:456:GLN:HE22	1:E:465:ARG:NH2	2.01	0.57
2:F:189:ARG:HH12	2:F:203:LEU:N	2.02	0.57
2:B:45:ASN:ND2	2:B:50:LEU:O	2.36	0.57
1:Z:382:ARG:NH2	1:Z:385:ILE:HD13	2.19	0.57
2:M:189:ARG:NH1	4:M:1093:HOH:O	2.33	0.57
2:W:161:GLU:HB3	2:W:162:PRO:HD3	1.86	0.57
2:M:10:GLU:O	2:M:14:ARG:HB2	2.05	0.57
2:A:9:MET:HE2	2:A:9:MET:HA	1.87	0.57
1:N:456:GLN:HE22	1:N:465:ARG:NH2	2.02	0.57
2:S:59:ARG:NH2	2:S:217:ARG:O	2.36	0.57
2:B:121:GLU:HG2	2:B:156:MET:HG2	1.85	0.57
2:B:142:THR:CG2	2:B:146:SER:HB2	2.34	0.57
2:S:117:PRO:HD3	2:1:9:MET:CE	2.34	0.57
1:N:366:TYR:CD2	1:N:374:LEU:HD13	2.39	0.57
1:P:366:TYR:CD2	1:P:374:LEU:HD13	2.39	0.57
2:S:234:LEU:HD12	2:S:234:LEU:N	2.18	0.57
2:D:85:ARG:NH2	3:Q:251:DMF:C2	2.65	0.57
1:R:456:GLN:HE22	1:R:465:ARG:NH2	2.03	0.57
2:1:85:ARG:HH12	2:1:98:GLN:NE2	2.03	0.57
2:B:178:THR:OG1	2:B:233:LEU:HD11	2.03	0.57
2:I:189:ARG:HH12	2:I:203:LEU:N	2.02	0.57
1:L:382:ARG:NH2	1:L:385:ILE:HD13	2.19	0.57
1:L:456:GLN:HE22	1:L:465:ARG:NH2	2.01	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:56:LEU:HD13	2:I:99:LEU:HD13	1.86	0.56
1:V:375:THR:HG23	2:O:93:ASP:OD1	2.05	0.56
2:K:85:ARG:NH2	3:K:251:DMF:C1	2.68	0.56
1:G:429:TRP:CZ2	3:G:140:DMF:H22	2.39	0.56
2:B:30:VAL:HG12	2:B:43:ALA:HB2	1.85	0.56
2:O:85:ARG:NH1	2:O:98:GLN:HE22	2.02	0.56
4:U:1698:HOH:O	2:1:67:LYS:HE3	2.06	0.56
2:F:48:ARG:HD2	4:F:1871:HOH:O	2.05	0.56
1:G:456:GLN:HE22	1:G:465:ARG:NH2	2.02	0.56
2:D:151:PRO:HD2	4:D:1854:HOH:O	2.05	0.56
2:M:181:LEU:HD21	2:M:234:LEU:CD1	2.35	0.56
2:K:85:ARG:NH2	3:K:251:DMF:H11	2.21	0.56
2:B:18:GLU:OE2	2:B:21:ARG:HD3	2.06	0.56
1:E:465:ARG:NH2	4:E:1535:HOH:O	2.38	0.56
2:D:18:GLU:OE2	2:D:21:ARG:NH1	2.39	0.56
2:B:33:LEU:HG	2:B:153:PHE:HB3	1.87	0.56
1:G:382:ARG:NH2	1:G:385:ILE:HD13	2.20	0.56
1:V:523:GLY:O	1:V:526:THR:HG22	2.05	0.56
2:B:35:TYR:CE1	2:B:177:LEU:HD13	2.39	0.56
1:V:429:TRP:CH2	3:V:121:DMF:H13	2.41	0.56
1:L:366:TYR:CD2	1:L:374:LEU:HD13	2.41	0.56
2:U:139:TYR:CD1	2:U:149:ASP:HB3	2.41	0.56
1:G:366:TYR:CD2	1:G:374:LEU:HD13	2.41	0.56
2:B:173:GLU:HG2	2:B:174:ASN:ND2	2.20	0.56
2:B:74:LEU:HD13	2:B:122:LEU:HD11	1.88	0.56
2:Y:18:GLU:OE1	2:Y:21:ARG:NH1	2.30	0.56
2:F:49:SER:HB2	2:W:97:ARG:HD2	1.88	0.56
1:J:366:TYR:CD2	1:J:374:LEU:HD13	2.41	0.56
2:Q:181:LEU:HD23	2:Q:233:LEU:HB3	1.87	0.55
2:I:152:HIS:CG	2:I:171:TYR:HE2	2.24	0.55
1:2:382:ARG:NH2	1:2:385:ILE:HD13	2.21	0.55
1:G:465:ARG:NH1	4:G:2051:HOH:O	2.39	0.55
1:G:444:LEU:HD21	1:X:325:MET:SD	2.47	0.55
2:I:90:ASP:HB3	2:I:93:ASP:OD2	2.06	0.55
2:S:85:ARG:NH1	2:S:98:GLN:NE2	2.54	0.55
1:N:500:ASP:HB2	4:N:1447:HOH:O	2.06	0.55
1:Z:496:ILE:HD12	3:Z:69:DMF:C2	2.29	0.55
2:O:35:TYR:CE1	2:O:37:GLY:CA	2.90	0.55
1:V:456:GLN:HE22	1:V:465:ARG:NH2	2.04	0.55
2:A:33:LEU:HD12	2:A:33:LEU:N	2.22	0.55
1:Z:366:TYR:CD2	1:Z:374:LEU:HD13	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:1:167:LEU:HG	2:1:187:ALA:CB	2.37	0.55
2:D:16:ARG:CG	2:K:9:MET:CE	2.82	0.54
2:1:11:GLN:OE1	2:1:14:ARG:NE	2.40	0.54
2:B:68:PHE:HA	2:B:71:PHE:CZ	2.42	0.54
1:E:380:ILE:HB	3:E:104:DMF:C1	2.36	0.54
2:B:55:GLU:HB3	2:B:222:PHE:CG	2.42	0.54
1:E:366:TYR:CD2	1:E:374:LEU:HD13	2.42	0.54
1:C:366:TYR:CD2	1:C:374:LEU:HD13	2.42	0.54
2:B:62:PHE:CZ	2:B:122:LEU:HD22	2.43	0.54
2:Y:189:ARG:HH12	2:Y:203:LEU:N	2.05	0.54
2:M:85:ARG:NH1	2:M:98:GLN:NE2	2.55	0.54
1:E:465:ARG:NH1	4:E:1397:HOH:O	2.39	0.54
2:F:151:PRO:HD2	4:F:729:HOH:O	2.07	0.54
2:B:31:VAL:HG11	2:B:167:LEU:HD11	1.89	0.54
1:R:366:TYR:CE2	1:R:374:LEU:HD13	2.42	0.54
2:Q:11:GLN:HA	2:Q:11:GLN:NE2	2.22	0.54
2:B:32:ALA:HB3	2:B:154:VAL:CG2	2.38	0.54
2:O:85:ARG:NH1	2:O:98:GLN:HE21	2.02	0.54
1:V:319:ARG:HH22	3:V:135:DMF:H11	1.72	0.54
2:W:16:ARG:HB3	2:W:117:PRO:HG3	1.88	0.54
4:N:1268:HOH:O	1:V:451:LYS:HE3	2.08	0.54
2:I:49:SER:HB2	2:S:97:ARG:HD2	1.90	0.54
2:B:112:THR:HG22	2:B:113:GLU:HG3	1.90	0.54
1:E:325:MET:SD	1:R:444:LEU:HD21	2.48	0.54
2:D:116:LYS:HG3	2:D:117:PRO:HD2	1.90	0.53
2:Q:112:THR:HG21	2:Y:116:LYS:HE3	1.90	0.53
2:W:9:MET:HG2	2:W:9:MET:O	2.09	0.53
2:1:11:GLN:OE1	2:1:11:GLN:CA	2.40	0.53
2:Q:16:ARG:HB3	2:Q:117:PRO:HG3	1.91	0.53
2:W:18:GLU:HG3	2:W:22:LYS:HE3	1.88	0.53
1:2:366:TYR:CD2	1:2:374:LEU:HD13	2.43	0.53
2:A:74:LEU:HD13	2:A:122:LEU:HD11	1.89	0.53
2:Y:12:ALA:O	2:Y:16:ARG:HG3	2.09	0.53
2:Y:170:SER:OG	2:Y:183:ILE:HG23	2.08	0.53
1:E:366:TYR:CE2	1:E:374:LEU:HD13	2.43	0.53
1:T:377:ALA:HB1	3:T:134:DMF:H13	1.91	0.53
2:A:152:HIS:HB3	2:A:171:TYR:CE2	2.44	0.53
2:Y:97:ARG:NH1	4:Y:849:HOH:O	2.42	0.53
2:Q:11:GLN:HE21	2:Q:11:GLN:HA	1.73	0.53
1:L:483:GLY:HA2	4:L:789:HOH:O	2.08	0.53
2:W:152:HIS:HB3	2:W:171:TYR:CZ	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Q:85:ARG:NH1	2:Q:98:GLN:NE2	2.56	0.53
2:B:63:ALA:HB3	2:B:123:CYS:HB3	1.89	0.53
1:T:366:TYR:CD2	1:T:374:LEU:HD13	2.44	0.53
2:W:11:GLN:OE1	2:W:14:ARG:CZ	2.57	0.52
2:U:18:GLU:O	2:U:22:LYS:HG3	2.09	0.52
2:D:76:ARG:HD3	4:D:252:HOH:O	2.09	0.52
2:S:19:LEU:HD23	2:S:19:LEU:C	2.29	0.52
2:W:85:ARG:NH1	2:W:98:GLN:NE2	2.58	0.52
2:I:40:LEU:HA	2:I:212:VAL:HG12	1.91	0.52
2:S:117:PRO:CD	2:1:9:MET:HE2	2.38	0.52
2:I:60:VAL:HG11	2:I:99:LEU:HD12	1.91	0.52
2:Q:217:ARG:HH11	2:Q:223:ARG:HD3	1.74	0.52
2:S:116:LYS:HG3	2:1:9:MET:HE1	1.91	0.52
2:I:178:THR:HG22	2:I:182:ARG:NH2	2.21	0.52
2:M:10:GLU:O	2:M:14:ARG:CB	2.58	0.52
2:U:217:ARG:HH11	2:U:223:ARG:HD3	1.75	0.52
2:U:150:GLU:HG3	2:U:154:VAL:HG22	1.91	0.52
2:A:19:LEU:C	2:A:19:LEU:HD23	2.30	0.52
2:B:155:VAL:HG11	2:B:164:ALA:CB	2.33	0.52
2:Y:14:ARG:O	2:Y:18:GLU:HG2	2.09	0.52
2:K:189:ARG:HH12	2:K:203:LEU:N	2.08	0.52
1:E:329:ARG:HD2	1:R:434:GLU:OE1	2.09	0.52
2:1:20:ALA:O	2:1:24:ILE:HG13	2.09	0.52
1:R:465:ARG:NH1	4:R:2061:HOH:O	2.43	0.52
2:B:162:PRO:HB2	2:B:190:ALA:HB1	1.92	0.52
2:W:217:ARG:HH11	2:W:223:ARG:HD3	1.75	0.52
2:F:181:LEU:HD23	2:F:233:LEU:HB3	1.91	0.52
1:J:515:ARG:HB3	4:J:959:HOH:O	2.09	0.52
2:F:9:MET:O	2:F:9:MET:HG2	2.09	0.52
2:B:33:LEU:HD12	2:B:33:LEU:N	2.25	0.52
2:I:203:LEU:HB3	2:I:207:SER:OG	2.09	0.52
3:G:137:DMF:H12	2:W:85:ARG:NH2	2.25	0.52
1:2:409:ILE:HG13	1:2:410:HIS:ND1	2.24	0.52
2:B:47:SER:HB2	4:I:1752:HOH:O	2.10	0.51
1:J:456:GLN:HE22	1:J:465:ARG:HH21	1.58	0.51
2:D:217:ARG:HH11	2:D:223:ARG:HD3	1.75	0.51
2:S:32:ALA:C	2:S:33:LEU:HD12	2.29	0.51
2:I:35:TYR:CE1	2:I:37:GLY:HA3	2.46	0.51
2:I:97:ARG:HB2	4:I:990:HOH:O	2.09	0.51
2:U:85:ARG:HH12	2:U:98:GLN:NE2	2.07	0.51
2:W:85:ARG:NH1	4:W:659:HOH:O	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:456:GLN:HE22	1:C:465:ARG:HH21	1.58	0.51
2:S:117:PRO:CD	2:1:9:MET:HE1	2.38	0.51
2:A:85:ARG:NH1	2:A:98:GLN:NE2	2.58	0.51
2:F:139:TYR:CE1	2:F:149:ASP:HB3	2.46	0.51
2:B:31:VAL:HG12	2:B:33:LEU:HD12	1.93	0.51
2:K:85:ARG:HH22	3:K:251:DMF:C1	2.23	0.51
2:F:9:MET:HE1	2:M:116:LYS:HG3	1.92	0.51
2:D:112:THR:HG21	2:Q:116:LYS:HE3	1.91	0.51
2:K:16:ARG:NH2	2:K:114:GLN:O	2.31	0.51
2:U:110:ILE:O	2:U:114:GLN:HB2	2.11	0.51
2:I:159:THR:O	2:I:159:THR:CG2	2.56	0.51
2:1:16:ARG:HB3	2:1:117:PRO:HG3	1.93	0.51
2:W:19:LEU:O	2:W:19:LEU:HD23	2.10	0.51
1:T:409:ILE:HG13	1:T:410:HIS:ND1	2.26	0.51
1:P:409:ILE:HG13	1:P:410:HIS:ND1	2.26	0.51
2:A:217:ARG:HH11	2:A:223:ARG:HD3	1.76	0.51
2:O:147:ILE:HG12	2:U:50:LEU:HD11	1.91	0.51
2:I:58:ASP:OD1	2:I:219:ARG:NH1	2.44	0.51
2:Y:92:ARG:HD3	2:Y:129:HIS:CE1	2.46	0.51
2:O:152:HIS:HB3	2:O:171:TYR:CE2	2.46	0.51
2:Y:217:ARG:HH11	2:Y:223:ARG:HD3	1.76	0.51
1:N:412:SER:O	1:N:414:PRO:HD3	2.11	0.51
1:P:465:ARG:HD2	4:P:1128:HOH:O	2.10	0.51
1:G:472:TYR:CZ	3:G:20:DMF:H22	2.46	0.51
2:B:21:ARG:HG3	2:B:22:LYS:N	2.26	0.51
2:M:217:ARG:HH11	2:M:223:ARG:HD3	1.75	0.51
1:X:456:GLN:HE22	1:X:465:ARG:HH21	1.59	0.51
2:K:115:ALA:HB3	4:K:1445:HOH:O	2.11	0.51
2:F:217:ARG:HH11	2:F:223:ARG:HD3	1.76	0.51
2:S:117:PRO:HD3	2:1:9:MET:HE2	1.93	0.50
1:T:472:TYR:CE2	3:T:62:DMF:H13	2.45	0.50
1:Z:409:ILE:HG13	1:Z:410:HIS:ND1	2.26	0.50
2:O:217:ARG:HH11	2:O:223:ARG:HD3	1.76	0.50
1:R:409:ILE:HG13	1:R:410:HIS:ND1	2.26	0.50
2:S:18:GLU:OE2	2:S:21:ARG:NH1	2.44	0.50
2:B:12:ALA:O	2:B:16:ARG:HG3	2.11	0.50
1:G:415:GLN:NE2	4:G:784:HOH:O	2.38	0.50
2:U:16:ARG:NH2	2:U:114:GLN:O	2.30	0.50
2:F:105:GLN:HB2	4:F:258:HOH:O	2.11	0.50
1:Z:366:TYR:CE2	1:Z:374:LEU:HD13	2.46	0.50
2:A:189:ARG:HH12	2:A:203:LEU:N	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Y:85:ARG:NH1	2:Y:98:GLN:NE2	2.59	0.50
1:H:366:TYR:CE2	1:H:374:LEU:HD13	2.45	0.50
2:F:189:ARG:NH1	2:F:203:LEU:N	2.59	0.50
2:K:49:SER:HB2	2:M:97:ARG:HD2	1.93	0.50
2:1:217:ARG:HH11	2:1:223:ARG:HD3	1.75	0.50
1:N:409:ILE:HG13	1:N:410:HIS:ND1	2.26	0.50
2:U:152:HIS:HB3	2:U:171:TYR:CE2	2.46	0.50
1:L:409:ILE:HG13	1:L:410:HIS:ND1	2.27	0.50
2:D:181:LEU:HD23	2:D:233:LEU:HB3	1.94	0.50
2:M:152:HIS:HB3	2:M:171:TYR:CE2	2.46	0.50
1:L:339:ASP:HB2	3:L:138:DMF:H12	1.94	0.50
1:G:409:ILE:HG13	1:G:410:HIS:ND1	2.27	0.50
1:J:412:SER:O	1:J:414:PRO:HD3	2.12	0.50
2:S:152:HIS:HB3	2:S:171:TYR:CE2	2.47	0.50
2:D:152:HIS:HB3	2:D:171:TYR:CE2	2.47	0.50
2:I:55:GLU:HB2	2:I:222:PHE:CG	2.47	0.50
2:B:32:ALA:HB3	2:B:154:VAL:HG23	1.92	0.49
2:1:85:ARG:NH1	2:1:98:GLN:NE2	2.60	0.49
2:F:16:ARG:HB3	2:F:117:PRO:HG3	1.94	0.49
1:E:391:LEU:HD21	1:L:398:LEU:HD11	1.93	0.49
2:K:116:LYS:HG2	2:K:117:PRO:HD2	1.93	0.49
2:F:85:ARG:NH1	2:F:98:GLN:NE2	2.60	0.49
1:Z:456:GLN:HE22	1:Z:465:ARG:HH21	1.60	0.49
2:K:217:ARG:HH11	2:K:223:ARG:HD3	1.76	0.49
1:P:456:GLN:HE22	1:P:465:ARG:HH21	1.59	0.49
2:F:150:GLU:HG3	2:F:154:VAL:HG22	1.93	0.49
2:S:189:ARG:HH12	2:S:203:LEU:N	2.11	0.49
2:F:97:ARG:NH1	4:F:866:HOH:O	2.45	0.49
1:X:409:ILE:HG13	1:X:410:HIS:ND1	2.27	0.49
1:R:412:SER:O	1:R:414:PRO:HD3	2.12	0.49
1:T:456:GLN:HE22	1:T:465:ARG:HH21	1.59	0.49
2:I:11:GLN:CG	2:I:14:ARG:HH12	2.09	0.49
2:I:35:TYR:CE2	2:I:38:GLY:O	2.66	0.49
2:Y:189:ARG:NH1	2:Y:203:LEU:N	2.59	0.49
1:J:391:LEU:HD21	1:T:398:LEU:HD11	1.95	0.49
1:C:409:ILE:HG13	1:C:410:HIS:ND1	2.27	0.49
2:K:152:HIS:HB3	2:K:171:TYR:CE2	2.47	0.49
2:I:182:ARG:NH2	4:I:719:HOH:O	2.45	0.49
2:U:18:GLU:HA	2:U:21:ARG:NH2	2.27	0.49
1:N:456:GLN:HE22	1:N:465:ARG:HH21	1.61	0.49
1:L:456:GLN:HE22	1:L:465:ARG:HH21	1.60	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Z:412:SER:O	1:Z:414:PRO:HD3	2.13	0.49
1:E:520:SER:HB3	4:E:1913:HOH:O	2.12	0.49
2:B:35:TYR:CE1	2:B:177:LEU:CD1	2.96	0.49
2:F:149:ASP:O	2:F:149:ASP:OD1	2.30	0.49
2:S:178:THR:HG23	2:S:179:ASP:N	2.27	0.49
1:Z:444:LEU:HB2	4:Z:692:HOH:O	2.12	0.49
1:C:348:THR:HA	4:C:1685:HOH:O	2.13	0.49
2:B:189:ARG:HE	2:B:189:ARG:HA	1.75	0.49
1:X:366:TYR:CE2	1:X:374:LEU:HD13	2.47	0.49
2:I:11:GLN:CG	2:I:14:ARG:CZ	2.90	0.49
2:U:9:MET:HE2	2:U:13:MET:CE	2.42	0.49
1:V:345:ILE:O	3:V:136:DMF:H21	2.12	0.49
2:1:152:HIS:HB3	2:1:171:TYR:CE2	2.48	0.49
1:E:412:SER:O	1:E:414:PRO:HD3	2.13	0.49
1:Z:429:TRP:CH2	3:Z:105:DMF:H13	2.48	0.49
2:A:97:ARG:HD2	2:O:49:SER:HB2	1.95	0.49
2:M:85:ARG:HB3	4:M:781:HOH:O	2.12	0.48
1:R:456:GLN:HE22	1:R:465:ARG:HH21	1.61	0.48
1:J:366:TYR:CE2	1:J:374:LEU:HD13	2.48	0.48
2:1:102:VAL:HG13	3:1:249:DMF:H12	1.94	0.48
1:V:409:ILE:HG13	1:V:410:HIS:ND1	2.28	0.48
2:Q:9:MET:N	2:Y:15:GLU:OE1	2.46	0.48
1:N:366:TYR:CE2	1:N:374:LEU:HD13	2.48	0.48
2:F:9:MET:CE	2:M:116:LYS:HA	2.43	0.48
1:J:409:ILE:HG13	1:J:410:HIS:ND1	2.28	0.48
2:I:16:ARG:HB3	2:I:117:PRO:HG3	1.96	0.48
2:B:67:LYS:HE3	2:B:69:ASN:OD1	2.14	0.48
2:F:152:HIS:HB3	2:F:171:TYR:CE2	2.47	0.48
2:B:166:ALA:CB	2:B:187:ALA:CA	2.90	0.48
2:Q:217:ARG:NH1	2:Q:223:ARG:HD3	2.29	0.48
1:E:409:ILE:HG13	1:E:410:HIS:ND1	2.28	0.48
1:G:412:SER:O	1:G:414:PRO:HD3	2.13	0.48
1:P:366:TYR:CE2	1:P:374:LEU:HD13	2.49	0.48
2:U:16:ARG:HB3	2:U:117:PRO:HG2	1.95	0.48
2:B:110:ILE:O	2:B:114:GLN:HB2	2.13	0.48
1:H:409:ILE:HG13	1:H:410:HIS:ND1	2.28	0.48
2:O:27:ALA:HB1	4:O:1192:HOH:O	2.14	0.48
2:D:19:LEU:C	2:D:19:LEU:HD23	2.34	0.48
1:E:444:LEU:HD21	1:L:325:MET:SD	2.54	0.48
2:U:19:LEU:HD23	2:U:19:LEU:C	2.33	0.48
2:B:161:GLU:N	2:B:162:PRO:HD2	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:181:LEU:HD23	2:O:233:LEU:HB3	1.95	0.48
1:C:412:SER:O	1:C:414:PRO:HD3	2.13	0.48
2:D:35:TYR:CZ	2:D:37:GLY:HA3	2.47	0.48
1:N:465:ARG:NH1	4:N:611:HOH:O	2.47	0.48
1:L:366:TYR:CE2	1:L:374:LEU:HD13	2.49	0.48
1:2:366:TYR:CE2	1:2:374:LEU:HD13	2.48	0.48
2:B:94:VAL:HA	2:B:98:GLN:OE1	2.12	0.48
2:Q:74:LEU:HD13	2:Q:122:LEU:HD11	1.95	0.48
2:Y:185:VAL:O	2:Y:189:ARG:HG3	2.14	0.48
2:U:16:ARG:HB3	2:U:117:PRO:CG	2.44	0.48
1:V:357:ARG:HA	3:V:117:DMF:HC	1.95	0.48
2:B:49:SER:OG	2:I:97:ARG:HG3	2.13	0.48
1:E:456:GLN:HE22	1:E:465:ARG:HH21	1.60	0.48
2:I:12:ALA:O	2:I:16:ARG:HG3	2.14	0.48
2:S:41:PHE:HB3	2:S:53:ILE:HD13	1.96	0.48
1:P:412:SER:O	1:P:414:PRO:HD3	2.14	0.48
2:B:64:ALA:HB2	2:B:122:LEU:HD22	1.83	0.47
2:W:18:GLU:CD	2:W:21:ARG:NH1	2.67	0.47
2:B:129:HIS:HD2	4:B:1916:HOH:O	1.97	0.47
2:Y:203:LEU:CD1	2:Y:203:LEU:N	2.77	0.47
1:V:412:SER:O	1:V:414:PRO:HD3	2.14	0.47
1:L:412:SER:O	1:L:414:PRO:HD3	2.12	0.47
1:C:375:THR:HG21	2:I:92:ARG:HB3	1.96	0.47
2:U:9:MET:HE2	2:U:13:MET:HE2	1.95	0.47
2:D:217:ARG:NH1	2:D:223:ARG:HD3	2.30	0.47
2:Q:51:GLN:HB2	4:Q:255:HOH:O	2.14	0.47
1:T:412:SER:O	1:T:414:PRO:HD3	2.14	0.47
1:2:412:SER:HB2	4:2:543:HOH:O	2.14	0.47
2:I:152:HIS:CG	2:I:171:TYR:CE2	3.03	0.47
2:A:205:VAL:HG22	2:A:230:LEU:HD23	1.97	0.47
2:O:21:ARG:NH2	4:O:680:HOH:O	2.47	0.47
2:Y:97:ARG:CZ	4:Y:849:HOH:O	2.61	0.47
2:B:60:VAL:HG11	2:B:99:LEU:HD12	1.95	0.47
1:H:456:GLN:HE22	1:H:465:ARG:HH21	1.61	0.47
2:F:12:ALA:O	2:F:16:ARG:HG3	2.14	0.47
1:E:398:LEU:HD11	1:R:391:LEU:HD21	1.95	0.47
2:U:233:LEU:CD1	2:U:233:LEU:N	2.77	0.47
2:B:141:ILE:N	2:B:141:ILE:HD12	2.30	0.47
1:H:325:MET:SD	1:P:444:LEU:HD21	2.55	0.47
2:Y:19:LEU:HD23	2:Y:19:LEU:C	2.35	0.47
2:W:135:ARG:NE	4:W:2032:HOH:O	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Y:18:GLU:O	2:Y:22:LYS:HG3	2.13	0.47
2:Q:11:GLN:OE1	2:Q:15:GLU:HG3	2.14	0.47
1:2:412:SER:O	1:2:414:PRO:HD3	2.15	0.47
2:M:114:GLN:NE2	4:M:1233:HOH:O	2.48	0.47
2:A:18:GLU:CD	2:A:21:ARG:NH1	2.64	0.47
2:B:42:VAL:HG23	2:B:210:VAL:HG22	1.96	0.47
1:T:366:TYR:CE2	1:T:374:LEU:HD13	2.49	0.47
2:A:217:ARG:NH1	2:A:223:ARG:HD3	2.30	0.47
3:E:113:DMF:H22	1:R:429:TRP:CZ2	2.50	0.47
2:M:217:ARG:NH1	2:M:223:ARG:HD3	2.29	0.47
1:H:412:SER:O	1:H:414:PRO:HD3	2.15	0.47
2:A:73:ASN:OD1	2:B:105:GLN:NE2	2.48	0.47
2:I:56:LEU:HG	2:I:62:PHE:HB2	1.96	0.46
2:A:85:ARG:NH2	3:A:249:DMF:H22	2.30	0.46
2:F:32:ALA:C	2:F:33:LEU:HD12	2.36	0.46
2:B:212:VAL:HG23	2:B:223:ARG:HG3	1.97	0.46
2:D:18:GLU:CD	2:D:21:ARG:NH1	2.68	0.46
2:B:210:VAL:HG12	2:B:225:ILE:HG12	1.97	0.46
1:J:456:GLN:NE2	1:J:465:ARG:HH21	2.13	0.46
2:W:217:ARG:NH1	2:W:223:ARG:HD3	2.29	0.46
3:P:133:DMF:H22	2:O:91:ARG:NH1	2.31	0.46
1:X:412:SER:O	1:X:414:PRO:HD3	2.15	0.46
2:W:161:GLU:CB	2:W:162:PRO:HD3	2.45	0.46
1:T:377:ALA:CB	3:T:134:DMF:H13	2.45	0.46
2:1:217:ARG:NH1	2:1:223:ARG:HD3	2.30	0.46
2:1:106:THR:HG21	3:1:249:DMF:HC	1.97	0.46
2:Q:179:ASP:O	2:Q:183:ILE:HG13	2.15	0.46
2:B:74:LEU:CD1	2:B:122:LEU:HD11	2.46	0.46
1:P:456:GLN:NE2	1:P:465:ARG:HH21	2.14	0.46
2:Q:181:LEU:HD21	2:Q:234:LEU:HD13	1.96	0.46
2:U:217:ARG:NH1	2:U:223:ARG:HD3	2.29	0.46
2:O:217:ARG:NH1	2:O:223:ARG:HD3	2.31	0.46
1:E:520:SER:CB	4:E:1913:HOH:O	2.64	0.46
2:B:55:GLU:HB3	2:B:222:PHE:CB	2.45	0.46
2:M:12:ALA:O	2:M:16:ARG:HG3	2.15	0.46
2:K:76:ARG:HD3	4:K:255:HOH:O	2.15	0.46
2:Q:27:ALA:HB1	4:Q:5:HOH:O	2.15	0.46
2:Y:217:ARG:NH1	2:Y:223:ARG:HD3	2.30	0.46
1:X:456:GLN:NE2	1:X:465:ARG:HH21	2.12	0.46
1:G:456:GLN:HE22	1:G:465:ARG:HH21	1.63	0.46
2:O:16:ARG:HB3	2:O:117:PRO:HG3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:366:TYR:CE2	1:G:374:LEU:HD13	2.50	0.46
2:Q:135:ARG:NH2	2:Q:173:GLU:OE1	2.49	0.46
2:I:39:VAL:HG21	2:I:125:ALA:HB1	1.98	0.46
1:C:456:GLN:NE2	1:C:465:ARG:HH21	2.14	0.46
1:C:401:LEU:HA	1:C:402:PRO:HD3	1.85	0.46
2:K:97:ARG:NH1	4:K:1647:HOH:O	2.49	0.46
2:I:99:LEU:HA	2:I:102:VAL:HG12	1.98	0.45
2:F:9:MET:HE1	2:M:117:PRO:HD3	1.97	0.45
1:E:444:LEU:HB2	4:E:601:HOH:O	2.15	0.45
2:B:90:ASP:OD1	2:B:90:ASP:O	2.34	0.45
2:S:30:VAL:HG13	2:S:43:ALA:HB2	1.98	0.45
1:L:482:GLY:HA3	4:L:582:HOH:O	2.15	0.45
1:C:479:SER:HB2	1:E:479:SER:HB2	1.98	0.45
2:S:85:ARG:NH1	4:S:1099:HOH:O	2.44	0.45
1:J:515:ARG:NH1	4:J:1880:HOH:O	2.48	0.45
2:B:16:ARG:NH2	2:B:114:GLN:O	2.30	0.45
2:W:31:VAL:HG12	2:W:155:VAL:HG22	1.97	0.45
2:Y:9:MET:SD	2:Y:9:MET:O	2.74	0.45
2:D:9:MET:N	2:Q:15:GLU:OE1	2.50	0.45
2:F:54:SER:CB	2:F:75:ARG:HD2	2.47	0.45
2:I:152:HIS:HB3	2:I:171:TYR:CZ	2.48	0.45
2:Q:97:ARG:CD	2:Y:49:SER:HB2	2.46	0.45
2:K:15:GLU:HB3	2:M:9:MET:N	2.31	0.45
2:W:97:ARG:NH1	4:W:1712:HOH:O	2.50	0.45
2:F:217:ARG:NH1	2:F:223:ARG:HD3	2.30	0.45
2:B:9:MET:HE1	2:B:13:MET:CE	2.23	0.45
2:D:35:TYR:CD1	2:D:37:GLY:N	2.80	0.45
2:Y:189:ARG:HH11	2:Y:203:LEU:HD12	1.81	0.45
2:F:112:THR:HG21	2:M:116:LYS:HE3	1.99	0.45
1:L:416:SER:HA	4:L:2014:HOH:O	2.16	0.45
2:W:181:LEU:HD13	2:W:233:LEU:HB3	1.98	0.45
1:T:456:GLN:NE2	1:T:465:ARG:HH21	2.14	0.45
2:K:217:ARG:NH1	2:K:223:ARG:HD3	2.31	0.45
2:B:141:ILE:HG13	2:B:147:ILE:CD1	2.46	0.45
2:Q:54:SER:CB	2:Q:75:ARG:HD2	2.46	0.45
2:O:185:VAL:O	2:O:189:ARG:HG2	2.16	0.45
1:N:473:ASP:OD1	1:N:521:ARG:NH1	2.37	0.45
2:B:155:VAL:HG22	2:B:160:THR:HB	1.98	0.45
2:Y:16:ARG:HB3	2:Y:117:PRO:HG3	1.99	0.45
2:M:151:PRO:HD2	4:M:1408:HOH:O	2.17	0.45
1:E:401:LEU:HA	1:E:402:PRO:HD3	1.84	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:30:VAL:HG13	2:O:43:ALA:HB2	1.99	0.45
2:B:57:TYR:CZ	2:B:91:ARG:NH1	2.84	0.45
2:B:9:MET:CE	2:B:13:MET:HE1	2.40	0.45
1:Z:456:GLN:NE2	1:Z:465:ARG:HH21	2.15	0.45
1:L:456:GLN:NE2	1:L:465:ARG:HH21	2.15	0.45
2:D:150:GLU:HA	2:D:151:PRO:HD3	1.82	0.45
2:Q:11:GLN:CA	2:Q:11:GLN:NE2	2.80	0.45
2:A:41:PHE:HB3	2:A:53:ILE:HD13	1.99	0.45
2:A:230:LEU:O	2:A:231:GLN:C	2.55	0.45
1:C:366:TYR:CE2	1:C:374:LEU:HD13	2.52	0.45
2:S:182:ARG:NH1	4:S:1950:HOH:O	2.49	0.45
2:A:45:ASN:HA	2:A:46:PRO:HD3	1.83	0.45
1:P:471:LEU:HA	1:P:471:LEU:HD12	1.87	0.45
2:F:139:TYR:CD1	2:F:149:ASP:HB3	2.53	0.44
2:B:99:LEU:HA	2:B:99:LEU:HD23	1.86	0.44
2:A:181:LEU:HD23	2:A:233:LEU:HB3	2.00	0.44
2:B:116:LYS:HB2	2:I:13:MET:HE3	1.98	0.44
2:F:229:ALA:O	2:F:232:ALA:HB3	2.17	0.44
2:B:32:ALA:HA	2:B:40:LEU:O	2.17	0.44
1:E:375:THR:HG21	2:K:92:ARG:HB3	2.00	0.44
2:O:189:ARG:CD	2:O:203:LEU:HD12	2.47	0.44
1:C:473:ASP:OD2	1:R:452:LYS:NZ	2.37	0.44
2:1:30:VAL:HG13	2:1:43:ALA:HB2	2.00	0.44
1:H:456:GLN:NE2	1:H:465:ARG:HH21	2.15	0.44
1:X:471:LEU:HA	1:X:471:LEU:HD12	1.87	0.44
2:Y:188:LEU:HA	2:Y:188:LEU:HD12	1.89	0.44
2:D:32:ALA:C	2:D:33:LEU:HD12	2.37	0.44
2:W:181:LEU:HD23	2:W:185:VAL:HG23	1.99	0.44
1:V:456:GLN:HE22	1:V:465:ARG:HH21	1.66	0.44
2:A:32:ALA:C	2:A:33:LEU:HD12	2.38	0.44
1:L:379:LYS:HZ3	3:L:138:DMF:HC	1.81	0.44
2:M:150:GLU:HA	2:M:151:PRO:HD3	1.84	0.44
1:R:509:ARG:O	1:R:513:LEU:HD13	2.18	0.44
1:J:471:LEU:HD12	1:J:471:LEU:HA	1.84	0.44
1:X:396:GLN:HB3	1:X:396:GLN:HE21	1.60	0.44
2:B:52:LYS:O	2:B:53:ILE:HD13	2.18	0.44
1:V:429:TRP:CE2	3:V:121:DMF:H22	2.53	0.44
1:N:456:GLN:NE2	1:N:465:ARG:HH21	2.16	0.44
2:Y:173:GLU:O	2:Y:174:ASN:HB2	2.17	0.44
1:N:429:TRP:CH2	3:N:21:DMF:H13	2.52	0.44
1:2:472:TYR:CE1	3:2:99:DMF:H13	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:107:LEU:HA	2:D:107:LEU:HD23	1.88	0.44
2:S:102:VAL:HG13	3:S:249:DMF:H12	2.00	0.44
2:A:30:VAL:HG13	2:A:43:ALA:HB2	1.99	0.44
2:W:150:GLU:HA	2:W:151:PRO:HD3	1.83	0.44
2:S:168:LYS:HE3	2:S:168:LYS:HB2	1.78	0.44
2:A:168:LYS:HB2	2:A:168:LYS:HE3	1.79	0.44
1:H:363:LEU:HD12	1:H:363:LEU:HA	1.88	0.44
2:U:189:ARG:HG2	2:U:203:LEU:HD12	1.98	0.44
1:C:306:LEU:HD12	1:C:306:LEU:O	2.18	0.44
2:O:85:ARG:HH12	2:O:98:GLN:HE21	1.63	0.44
2:W:189:ARG:NH2	4:W:1576:HOH:O	2.50	0.44
2:B:18:GLU:CD	2:B:21:ARG:HD3	2.38	0.44
1:L:379:LYS:NZ	3:L:138:DMF:HC	2.32	0.44
2:K:116:LYS:CG	2:K:117:PRO:HD2	2.47	0.44
2:W:41:PHE:HB3	2:W:53:ILE:HD13	1.99	0.44
2:D:168:LYS:HB2	2:D:168:LYS:HE3	1.77	0.44
1:R:432:GLU:HB2	4:R:1661:HOH:O	2.17	0.44
2:S:116:LYS:HE3	2:1:112:THR:HG21	2.00	0.44
1:N:388:ARG:NE	4:N:1663:HOH:O	2.47	0.44
2:O:168:LYS:HB2	2:O:168:LYS:HE3	1.78	0.44
2:D:116:LYS:HE3	2:K:112:THR:HG21	2.00	0.43
2:U:147:ILE:HG21	2:1:68:PHE:CD2	2.53	0.43
1:R:366:TYR:OH	2:D:93:ASP:HB3	2.19	0.43
1:G:456:GLN:NE2	1:G:465:ARG:HH21	2.16	0.43
2:F:30:VAL:HG13	2:F:43:ALA:HB2	2.00	0.43
1:E:456:GLN:NE2	1:E:465:ARG:HH21	2.16	0.43
2:F:9:MET:HE1	2:M:117:PRO:CD	2.48	0.43
2:B:56:LEU:HD13	2:B:99:LEU:HD13	1.99	0.43
2:U:181:LEU:CD2	2:U:233:LEU:HB3	2.49	0.43
2:Y:228:SER:O	2:Y:229:ALA:C	2.56	0.43
2:I:41:PHE:HB3	2:I:53:ILE:HD13	2.00	0.43
2:A:9:MET:O	2:A:13:MET:HG2	2.19	0.43
1:L:410:HIS:HD2	4:M:1361:HOH:O	2.00	0.43
2:1:54:SER:CB	2:1:75:ARG:HD2	2.48	0.43
2:M:32:ALA:C	2:M:33:LEU:HD12	2.38	0.43
2:O:41:PHE:HB3	2:O:53:ILE:HD13	2.00	0.43
1:G:363:LEU:HA	1:G:363:LEU:HD12	1.88	0.43
1:R:456:GLN:NE2	1:R:465:ARG:HH21	2.17	0.43
2:A:54:SER:CB	2:A:75:ARG:HD2	2.48	0.43
2:D:230:LEU:O	2:D:234:LEU:HD13	2.19	0.43
2:S:54:SER:CB	2:S:75:ARG:HD2	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:471:LEU:HA	1:C:471:LEU:HD12	1.83	0.43
1:N:375:THR:HG21	2:F:92:ARG:HB3	2.00	0.43
2:U:54:SER:CB	2:U:75:ARG:HD2	2.49	0.43
1:T:444:LEU:HD21	1:2:325:MET:SD	2.58	0.43
2:U:41:PHE:HB3	2:U:53:ILE:HD13	2.01	0.43
2:B:59:ARG:HD2	2:B:127:VAL:CG1	2.45	0.43
2:O:189:ARG:HD2	2:O:203:LEU:CD1	2.48	0.43
2:Y:41:PHE:HB3	2:Y:53:ILE:HD13	2.01	0.43
2:1:151:PRO:HD2	4:1:393:HOH:O	2.19	0.43
1:2:306:LEU:HD12	1:2:306:LEU:O	2.19	0.43
2:B:64:ALA:HB1	2:B:122:LEU:CD2	2.43	0.43
2:U:147:ILE:HG12	2:1:50:LEU:HD11	2.01	0.43
1:2:401:LEU:HA	1:2:402:PRO:HD3	1.86	0.43
1:J:306:LEU:C	1:J:306:LEU:HD12	2.40	0.43
1:G:471:LEU:HA	1:G:471:LEU:HD12	1.91	0.43
2:B:207:SER:O	2:B:208:LEU:HG	2.19	0.43
2:1:41:PHE:HB3	2:1:53:ILE:HD13	2.01	0.43
2:B:217:ARG:HG3	2:B:223:ARG:NH2	2.14	0.43
2:B:40:LEU:CD1	2:B:212:VAL:HG13	2.48	0.43
2:S:33:LEU:N	2:S:33:LEU:HD12	2.34	0.43
1:P:377:ALA:HB1	3:P:107:DMF:C	2.49	0.43
1:V:363:LEU:HD12	1:V:363:LEU:HA	1.87	0.43
2:U:225:ILE:HG22	2:U:230:LEU:HA	2.01	0.42
1:E:509:ARG:NH2	4:E:1830:HOH:O	2.48	0.42
2:O:13:MET:HG3	2:U:19:LEU:HD11	2.01	0.42
2:D:35:TYR:CE1	2:D:37:GLY:N	2.88	0.42
2:D:85:ARG:HH22	3:Q:251:DMF:C2	2.31	0.42
2:U:9:MET:CE	2:U:13:MET:HE3	2.47	0.42
1:E:456:GLN:HG3	4:E:2203:HOH:O	2.19	0.42
2:D:33:LEU:HD12	2:D:33:LEU:N	2.33	0.42
2:D:129:HIS:HE1	4:D:1211:HOH:O	2.01	0.42
1:J:368:LYS:HB3	2:I:79:ILE:HD13	2.00	0.42
2:D:54:SER:CB	2:D:75:ARG:HD2	2.49	0.42
2:K:54:SER:CB	2:K:75:ARG:HD2	2.49	0.42
2:D:41:PHE:HB3	2:D:53:ILE:HD13	2.01	0.42
1:L:375:THR:HG21	2:M:92:ARG:HB3	2.01	0.42
2:Y:128:ALA:HB2	2:Y:134:LYS:HB3	2.01	0.42
2:Y:32:ALA:C	2:Y:33:LEU:HD12	2.39	0.42
2:K:225:ILE:HG22	2:K:230:LEU:HA	2.01	0.42
2:Y:18:GLU:CD	2:Y:21:ARG:HH12	2.20	0.42
2:B:224:ARG:HG3	2:B:224:ARG:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:381:ASN:HB2	3:P:107:DMF:H12	2.02	0.42
1:E:496:ILE:HD12	4:E:1830:HOH:O	2.20	0.42
2:Q:41:PHE:HB3	2:Q:53:ILE:HD13	1.99	0.42
2:B:150:GLU:HA	2:B:151:PRO:HD3	1.82	0.42
1:V:471:LEU:HD12	1:V:471:LEU:HA	1.88	0.42
1:G:401:LEU:HA	1:G:402:PRO:HD3	1.85	0.42
2:B:59:ARG:NH2	2:B:130:TYR:HB2	2.34	0.42
2:B:135:ARG:NH2	2:B:173:GLU:OE2	2.53	0.42
2:W:9:MET:N	2:W:9:MET:SD	2.93	0.42
2:B:161:GLU:OE1	2:B:161:GLU:HA	2.19	0.42
2:U:168:LYS:HB2	2:U:168:LYS:HE3	1.78	0.42
2:B:62:PHE:CE1	2:B:122:LEU:HD22	2.55	0.42
1:E:380:ILE:HG21	3:E:104:DMF:H12	2.01	0.42
2:B:110:ILE:O	2:B:114:GLN:CB	2.68	0.42
2:O:54:SER:CB	2:O:75:ARG:HD2	2.50	0.42
1:R:306:LEU:C	1:R:306:LEU:HD12	2.39	0.42
1:V:388:ARG:NE	4:V:1388:HOH:O	2.40	0.42
2:B:40:LEU:HD12	2:B:212:VAL:HG13	2.01	0.42
2:M:33:LEU:HD12	2:M:33:LEU:N	2.35	0.42
2:S:45:ASN:HA	2:S:46:PRO:HD3	1.82	0.42
2:M:30:VAL:HG13	2:M:43:ALA:HB2	2.01	0.42
2:M:54:SER:CB	2:M:75:ARG:HD2	2.49	0.42
2:U:170:SER:OG	2:U:183:ILE:HG23	2.20	0.42
2:Y:54:SER:CB	2:Y:75:ARG:HD2	2.49	0.42
2:A:115:ALA:HB3	2:B:112:THR:HG21	1.83	0.42
1:2:306:LEU:HD12	1:2:306:LEU:C	2.40	0.42
2:F:150:GLU:HA	2:F:151:PRO:HD3	1.83	0.41
1:V:515:ARG:HD2	4:V:2036:HOH:O	2.21	0.41
2:M:53:ILE:O	2:M:224:ARG:NH2	2.50	0.41
2:K:102:VAL:HG13	3:K:249:DMF:H12	2.02	0.41
2:W:167:LEU:HD12	2:W:167:LEU:HA	1.80	0.41
2:B:230:LEU:CD2	2:B:230:LEU:C	2.88	0.41
1:R:306:LEU:O	1:R:306:LEU:HD12	2.20	0.41
2:W:54:SER:CB	2:W:75:ARG:HD2	2.50	0.41
2:K:150:GLU:HA	2:K:151:PRO:HD3	1.83	0.41
2:W:32:ALA:C	2:W:33:LEU:HD12	2.40	0.41
2:W:19:LEU:C	2:W:19:LEU:CD2	2.86	0.41
2:O:150:GLU:HA	2:O:151:PRO:HD3	1.83	0.41
2:U:106:THR:O	2:U:110:ILE:HG13	2.20	0.41
2:O:217:ARG:HA	2:O:218:PRO:HD3	1.88	0.41
2:W:45:ASN:HA	2:W:46:PRO:HD3	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:41:PHE:HB3	2:K:53:ILE:HD13	2.01	0.41
2:I:30:VAL:HG13	2:I:43:ALA:HB2	2.02	0.41
1:Z:357:ARG:HD3	4:Z:1337:HOH:O	2.19	0.41
1:2:429:TRP:CH2	3:2:145:DMF:H23	2.55	0.41
2:B:19:LEU:CD2	2:B:19:LEU:C	2.86	0.41
1:X:391:LEU:CD2	1:Z:398:LEU:HD11	2.48	0.41
2:1:167:LEU:HG	2:1:187:ALA:HB1	2.01	0.41
2:B:225:ILE:HG13	2:B:230:LEU:HG	2.02	0.41
2:S:150:GLU:HA	2:S:151:PRO:HD3	1.84	0.41
2:Q:16:ARG:NE	2:Q:111:PHE:O	2.53	0.41
2:U:30:VAL:HG13	2:U:43:ALA:HB2	2.03	0.41
1:N:306:LEU:C	1:N:306:LEU:HD12	2.41	0.41
2:W:213:LEU:HD23	2:W:213:LEU:HA	1.93	0.41
1:Z:429:TRP:CZ2	3:Z:105:DMF:H22	2.56	0.41
2:D:49:SER:HB2	2:K:97:ARG:HD2	2.02	0.41
2:Q:168:LYS:HE3	2:Q:168:LYS:HB2	1.80	0.41
1:2:396:GLN:HB3	1:2:396:GLN:HE21	1.60	0.41
2:U:9:MET:CE	2:U:13:MET:HE2	2.50	0.41
2:D:116:LYS:CG	2:D:117:PRO:HD2	2.50	0.41
2:A:19:LEU:O	2:A:19:LEU:HD23	2.19	0.41
2:F:41:PHE:HB3	2:F:53:ILE:HD13	2.03	0.41
2:O:11:GLN:HA	2:O:14:ARG:HB3	2.03	0.41
1:V:456:GLN:NE2	1:V:465:ARG:HH21	2.18	0.41
2:I:32:ALA:HA	2:I:40:LEU:O	2.21	0.41
2:B:100:ALA:HB1	2:B:147:ILE:HD11	2.02	0.41
2:1:38:GLY:HA3	2:1:213:LEU:O	2.21	0.41
1:H:415:GLN:HA	1:H:415:GLN:OE1	2.21	0.41
1:Z:363:LEU:HD12	1:Z:363:LEU:HA	1.85	0.41
2:O:9:MET:HE1	2:U:115:ALA:O	2.20	0.41
2:B:217:ARG:CG	2:B:217:ARG:HH11	2.03	0.41
2:I:54:SER:CB	2:I:75:ARG:HD2	2.51	0.41
1:C:476:ASP:HB3	1:R:451:LYS:HZ1	1.86	0.41
1:H:517:ILE:O	1:H:521:ARG:HG3	2.21	0.41
1:R:401:LEU:HA	1:R:402:PRO:HD3	1.83	0.41
2:A:234:LEU:HD12	2:A:234:LEU:HA	1.90	0.41
1:P:513:LEU:HA	1:P:513:LEU:HD12	1.90	0.41
2:B:70:GLU:HG2	2:B:118:TYR:CE1	2.56	0.41
2:B:53:ILE:HB	4:B:1703:HOH:O	2.21	0.40
2:1:24:ILE:HG22	2:1:157:GLY:HA2	2.03	0.40
2:Y:217:ARG:HA	2:Y:218:PRO:HD3	1.89	0.40
2:Q:147:ILE:HG12	2:Y:50:LEU:HD11	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:52:LYS:HB3	2:D:52:LYS:HE2	1.90	0.40
2:B:133:THR:HG22	2:B:133:THR:O	2.21	0.40
2:B:217:ARG:HA	4:B:751:HOH:O	2.20	0.40
2:O:35:TYR:HB2	2:O:175:ALA:O	2.21	0.40
1:T:472:TYR:OH	3:T:62:DMF:H13	2.22	0.40
2:F:9:MET:HE1	2:M:116:LYS:HA	2.02	0.40
2:U:226:THR:HG22	4:U:252:HOH:O	2.21	0.40
2:B:47:SER:HB2	2:I:149:ASP:CG	2.39	0.40
2:Y:129:HIS:O	2:Y:132:GLU:HG2	2.21	0.40
2:W:16:ARG:NH1	2:Y:9:MET:HG3	2.36	0.40
2:Y:9:MET:SD	2:Y:9:MET:C	2.99	0.40
2:M:189:ARG:HH12	2:M:203:LEU:N	2.20	0.40
2:Q:30:VAL:HG13	2:Q:43:ALA:HB2	2.02	0.40
1:X:465:ARG:NH1	4:X:1509:HOH:O	2.54	0.40
1:Z:465:ARG:NH1	4:Z:1776:HOH:O	2.53	0.40
2:F:38:GLY:HA3	2:F:213:LEU:O	2.21	0.40
1:R:363:LEU:HA	1:R:363:LEU:HD12	1.88	0.40
2:B:81:PHE:CE2	2:B:85:ARG:HG3	2.57	0.40
2:K:101:ASN:HB3	3:K:251:DMF:C2	2.51	0.40
2:B:116:LYS:HE3	2:I:112:THR:HG21	2.03	0.40
4:T:1393:HOH:O	1:X:451:LYS:HE3	2.20	0.40
1:N:332:ARG:HB3	4:N:1560:HOH:O	2.20	0.40
2:B:104:ALA:HB1	2:B:145:GLY:O	2.21	0.40
2:U:24:ILE:CD1	2:U:143:TYR:HD2	2.34	0.40
1:T:306:LEU:HD12	1:T:306:LEU:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	2	220/240 (92%)	218 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	220/240 (92%)	220 (100%)	0	0	100	100
1	E	221/240 (92%)	219 (99%)	2 (1%)	0	100	100
1	G	220/240 (92%)	218 (99%)	2 (1%)	0	100	100
1	H	221/240 (92%)	219 (99%)	2 (1%)	0	100	100
1	J	220/240 (92%)	219 (100%)	1 (0%)	0	100	100
1	L	220/240 (92%)	218 (99%)	2 (1%)	0	100	100
1	N	221/240 (92%)	218 (99%)	3 (1%)	0	100	100
1	P	220/240 (92%)	218 (99%)	2 (1%)	0	100	100
1	R	220/240 (92%)	218 (99%)	2 (1%)	0	100	100
1	T	221/240 (92%)	219 (99%)	2 (1%)	0	100	100
1	V	224/240 (93%)	222 (99%)	2 (1%)	0	100	100
1	X	222/240 (92%)	220 (99%)	2 (1%)	0	100	100
1	Z	222/240 (92%)	221 (100%)	1 (0%)	0	100	100
2	1	210/240 (88%)	206 (98%)	4 (2%)	0	100	100
2	A	210/240 (88%)	206 (98%)	4 (2%)	0	100	100
2	B	210/240 (88%)	201 (96%)	9 (4%)	0	100	100
2	D	209/240 (87%)	205 (98%)	4 (2%)	0	100	100
2	F	210/240 (88%)	206 (98%)	4 (2%)	0	100	100
2	I	210/240 (88%)	206 (98%)	4 (2%)	0	100	100
2	K	210/240 (88%)	206 (98%)	4 (2%)	0	100	100
2	M	210/240 (88%)	206 (98%)	4 (2%)	0	100	100
2	O	210/240 (88%)	205 (98%)	5 (2%)	0	100	100
2	Q	210/240 (88%)	206 (98%)	4 (2%)	0	100	100
2	S	210/240 (88%)	204 (97%)	6 (3%)	0	100	100
2	U	210/240 (88%)	205 (98%)	5 (2%)	0	100	100
2	W	210/240 (88%)	207 (99%)	3 (1%)	0	100	100
2	Y	210/240 (88%)	205 (98%)	5 (2%)	0	100	100
All	All	6031/6720 (90%)	5941 (98%)	90 (2%)	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	2	165/178 (93%)	160 (97%)	5 (3%)	48	76
1	C	165/178 (93%)	159 (96%)	6 (4%)	42	69
1	E	165/178 (93%)	160 (97%)	5 (3%)	48	76
1	G	165/178 (93%)	158 (96%)	7 (4%)	36	62
1	H	165/178 (93%)	159 (96%)	6 (4%)	42	69
1	J	165/178 (93%)	159 (96%)	6 (4%)	42	69
1	L	165/178 (93%)	160 (97%)	5 (3%)	48	76
1	N	165/178 (93%)	161 (98%)	4 (2%)	57	82
1	P	165/178 (93%)	160 (97%)	5 (3%)	48	76
1	R	165/178 (93%)	160 (97%)	5 (3%)	48	76
1	T	165/178 (93%)	160 (97%)	5 (3%)	48	76
1	V	167/178 (94%)	160 (96%)	7 (4%)	36	62
1	X	165/178 (93%)	160 (97%)	5 (3%)	48	76
1	Z	165/178 (93%)	160 (97%)	5 (3%)	48	76
2	1	164/184 (89%)	156 (95%)	8 (5%)	31	55
2	A	164/184 (89%)	156 (95%)	8 (5%)	31	55
2	B	164/184 (89%)	154 (94%)	10 (6%)	23	42
2	D	163/184 (89%)	156 (96%)	7 (4%)	35	61
2	F	164/184 (89%)	154 (94%)	10 (6%)	23	42
2	I	164/184 (89%)	153 (93%)	11 (7%)	20	37
2	K	164/184 (89%)	154 (94%)	10 (6%)	23	42
2	M	164/184 (89%)	153 (93%)	11 (7%)	20	37
2	O	164/184 (89%)	155 (94%)	9 (6%)	27	48
2	Q	164/184 (89%)	156 (95%)	8 (5%)	31	55
2	S	164/184 (89%)	156 (95%)	8 (5%)	31	55
2	U	164/184 (89%)	153 (93%)	11 (7%)	20	37

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	W	164/184 (89%)	154 (94%)	10 (6%)	23	42
2	Y	164/184 (89%)	157 (96%)	7 (4%)	35	61
All	All	4607/5068 (91%)	4403 (96%)	204 (4%)	35	60

All (204) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	363	LEU
1	H	374	LEU
1	H	402	PRO
1	H	461	ASP
1	H	471	LEU
1	H	513	LEU
1	C	363	LEU
1	C	374	LEU
1	C	461	ASP
1	C	471	LEU
1	C	496	ILE
1	C	513	LEU
1	E	363	LEU
1	E	374	LEU
1	E	461	ASP
1	E	471	LEU
1	E	513	LEU
1	G	337	THR
1	G	363	LEU
1	G	374	LEU
1	G	461	ASP
1	G	471	LEU
1	G	496	ILE
1	G	513	LEU
1	J	363	LEU
1	J	374	LEU
1	J	444	LEU
1	J	461	ASP
1	J	471	LEU
1	J	513	LEU
1	L	363	LEU
1	L	374	LEU
1	L	461	ASP
1	L	471	LEU

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Mol	Chain	Res	Type
1	L	513	LEU
1	N	363	LEU
1	N	374	LEU
1	N	461	ASP
1	N	513	LEU
1	P	363	LEU
1	P	374	LEU
1	P	461	ASP
1	P	471	LEU
1	P	513	LEU
1	R	363	LEU
1	R	374	LEU
1	R	402	PRO
1	R	461	ASP
1	R	471	LEU
1	T	363	LEU
1	T	374	LEU
1	T	461	ASP
1	T	471	LEU
1	T	513	LEU
1	V	363	LEU
1	V	374	LEU
1	V	461	ASP
1	V	471	LEU
1	V	496	ILE
1	V	513	LEU
1	V	526	THR
1	X	363	LEU
1	X	374	LEU
1	X	461	ASP
1	X	471	LEU
1	X	513	LEU
1	Z	363	LEU
1	Z	374	LEU
1	Z	461	ASP
1	Z	471	LEU
1	Z	513	LEU
1	2	363	LEU
1	2	374	LEU
1	2	461	ASP
1	2	471	LEU
1	2	513	LEU

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Mol	Chain	Res	Type
2	D	52	LYS
2	D	59	ARG
2	D	102	VAL
2	D	133	THR
2	D	135	ARG
2	D	140	ARG
2	D	167	LEU
2	A	52	LYS
2	A	99	LEU
2	A	102	VAL
2	A	133	THR
2	A	135	ARG
2	A	140	ARG
2	A	167	LEU
2	A	234	LEU
2	B	9	MET
2	B	51	GLN
2	B	102	VAL
2	B	105	GLN
2	B	107	LEU
2	B	144	ASP
2	B	176	SER
2	B	217	ARG
2	B	230	LEU
2	B	234	LEU
2	F	52	LYS
2	F	59	ARG
2	F	99	LEU
2	F	102	VAL
2	F	133	THR
2	F	135	ARG
2	F	140	ARG
2	F	167	LEU
2	F	203	LEU
2	F	234	LEU
2	I	52	LYS
2	I	57	TYR
2	I	59	ARG
2	I	69	ASN
2	I	133	THR
2	I	135	ARG
2	I	140	ARG

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Mol	Chain	Res	Type
2	I	167	LEU
2	I	188	LEU
2	I	203	LEU
2	I	212	VAL
2	K	52	LYS
2	K	59	ARG
2	K	69	ASN
2	K	99	LEU
2	K	102	VAL
2	K	133	THR
2	K	135	ARG
2	K	140	ARG
2	K	167	LEU
2	K	188	LEU
2	M	48	ARG
2	M	52	LYS
2	M	59	ARG
2	M	99	LEU
2	M	102	VAL
2	M	133	THR
2	M	135	ARG
2	M	140	ARG
2	M	167	LEU
2	M	233	LEU
2	M	234	LEU
2	O	52	LYS
2	O	59	ARG
2	O	102	VAL
2	O	133	THR
2	O	135	ARG
2	O	140	ARG
2	O	167	LEU
2	O	188	LEU
2	O	234	LEU
2	Q	52	LYS
2	Q	99	LEU
2	Q	102	VAL
2	Q	133	THR
2	Q	135	ARG
2	Q	140	ARG
2	Q	167	LEU
2	Q	234	LEU

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Mol	Chain	Res	Type
2	S	52	LYS
2	S	59	ARG
2	S	102	VAL
2	S	133	THR
2	S	135	ARG
2	S	140	ARG
2	S	167	LEU
2	S	234	LEU
2	U	9	MET
2	U	52	LYS
2	U	69	ASN
2	U	99	LEU
2	U	102	VAL
2	U	133	THR
2	U	135	ARG
2	U	140	ARG
2	U	231	GLN
2	U	233	LEU
2	U	234	LEU
2	W	10	GLU
2	W	52	LYS
2	W	59	ARG
2	W	69	ASN
2	W	102	VAL
2	W	133	THR
2	W	135	ARG
2	W	140	ARG
2	W	167	LEU
2	W	234	LEU
2	Y	52	LYS
2	Y	59	ARG
2	Y	99	LEU
2	Y	102	VAL
2	Y	140	ARG
2	Y	167	LEU
2	Y	188	LEU
2	1	9	MET
2	1	48	ARG
2	1	52	LYS
2	1	99	LEU
2	1	102	VAL
2	1	133	THR

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Mol	Chain	Res	Type
2	1	135	ARG
2	1	140	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (74) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	396	GLN
1	H	456	GLN
1	C	396	GLN
1	C	456	GLN
1	E	396	GLN
1	E	456	GLN
1	G	396	GLN
1	G	456	GLN
1	J	396	GLN
1	J	456	GLN
1	L	396	GLN
1	L	456	GLN
1	N	396	GLN
1	N	456	GLN
1	P	396	GLN
1	P	456	GLN
1	R	396	GLN
1	R	456	GLN
1	T	396	GLN
1	T	456	GLN
1	V	396	GLN
1	V	456	GLN
1	X	396	GLN
1	X	456	GLN
1	Z	396	GLN
1	Z	456	GLN
1	2	396	GLN
1	2	430	ASN
1	2	456	GLN
2	D	69	ASN
2	D	73	ASN
2	D	98	GLN
2	D	129	HIS
2	A	69	ASN
2	A	73	ASN
2	A	98	GLN

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Mol	Chain	Res	Type
2	A	129	HIS
2	B	105	GLN
2	B	174	ASN
2	F	69	ASN
2	F	98	GLN
2	F	129	HIS
2	I	69	ASN
2	I	73	ASN
2	I	105	GLN
2	I	129	HIS
2	K	69	ASN
2	K	129	HIS
2	M	69	ASN
2	M	98	GLN
2	M	114	GLN
2	M	129	HIS
2	O	69	ASN
2	O	98	GLN
2	O	105	GLN
2	O	129	HIS
2	Q	11	GLN
2	Q	69	ASN
2	Q	98	GLN
2	S	69	ASN
2	S	98	GLN
2	S	129	HIS
2	U	69	ASN
2	U	98	GLN
2	U	129	HIS
2	W	69	ASN
2	W	98	GLN
2	W	129	HIS
2	Y	69	ASN
2	Y	98	GLN
2	Y	129	HIS
2	1	69	ASN
2	1	98	GLN
2	1	129	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

58 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	1	249	-	4,4,4	0.36	0	4,4,4	0.24	0
3	DMF	1	250	-	4,4,4	0.35	0	4,4,4	0.44	0
3	DMF	2	145	-	4,4,4	0.30	0	4,4,4	0.39	0
3	DMF	2	99	-	4,4,4	0.33	0	4,4,4	0.42	0
3	DMF	A	249	-	4,4,4	0.37	0	4,4,4	0.46	0
3	DMF	A	8	-	4,4,4	0.24	0	4,4,4	0.25	0
3	DMF	B	249	-	4,4,4	0.29	0	4,4,4	0.41	0
3	DMF	C	15	-	4,4,4	0.31	0	4,4,4	0.37	0
3	DMF	D	249	-	4,4,4	0.30	0	4,4,4	0.44	0
3	DMF	E	104	-	4,4,4	0.27	0	4,4,4	0.41	0
3	DMF	E	113	-	4,4,4	0.35	0	4,4,4	0.45	0
3	DMF	E	28	-	4,4,4	0.34	0	4,4,4	0.36	0
3	DMF	F	249	-	4,4,4	0.31	0	4,4,4	0.39	0
3	DMF	G	137	-	4,4,4	0.34	0	4,4,4	0.42	0
3	DMF	G	140	-	4,4,4	0.34	0	4,4,4	0.46	0
3	DMF	G	20	-	4,4,4	0.34	0	4,4,4	0.22	0
3	DMF	H	142	-	4,4,4	0.36	0	4,4,4	0.46	0
3	DMF	H	41	-	4,4,4	0.30	0	4,4,4	0.43	0
3	DMF	I	249	-	4,4,4	0.30	0	4,4,4	0.40	0
3	DMF	J	46	-	4,4,4	0.35	0	4,4,4	0.47	0
3	DMF	K	249	-	4,4,4	0.29	0	4,4,4	0.28	0
3	DMF	K	250	-	4,4,4	0.36	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	K	251	-	4,4,4	0.33	0	4,4,4	0.24	0
3	DMF	L	138	-	4,4,4	0.34	0	4,4,4	0.27	0
3	DMF	L	14	-	4,4,4	0.34	0	4,4,4	0.31	0
3	DMF	L	60	-	4,4,4	0.32	0	4,4,4	0.41	0
3	DMF	L	9	-	4,4,4	0.35	0	4,4,4	0.43	0
3	DMF	M	249	-	4,4,4	0.36	0	4,4,4	0.34	0
3	DMF	N	21	-	4,4,4	0.30	0	4,4,4	0.40	0
3	DMF	N	36	-	4,4,4	0.34	0	4,4,4	0.38	0
3	DMF	O	7	-	4,4,4	0.35	0	4,4,4	0.36	0
3	DMF	P	107	-	4,4,4	0.27	0	4,4,4	0.38	0
3	DMF	P	133	-	4,4,4	0.33	0	4,4,4	0.44	0
3	DMF	P	23	-	4,4,4	0.31	0	4,4,4	0.38	0
3	DMF	Q	249	-	4,4,4	0.31	0	4,4,4	0.33	0
3	DMF	Q	250	-	4,4,4	0.33	0	4,4,4	0.25	0
3	DMF	Q	251	-	4,4,4	0.30	0	4,4,4	0.40	0
3	DMF	S	249	-	4,4,4	0.30	0	4,4,4	0.28	0
3	DMF	S	250	-	4,4,4	0.34	0	4,4,4	0.33	0
3	DMF	T	134	-	4,4,4	0.33	0	4,4,4	0.44	0
3	DMF	T	62	-	4,4,4	0.32	0	4,4,4	0.34	0
3	DMF	U	249	-	4,4,4	0.32	0	4,4,4	0.31	0
3	DMF	U	250	-	4,4,4	0.30	0	4,4,4	0.29	0
3	DMF	V	117	-	4,4,4	0.37	0	4,4,4	0.40	0
3	DMF	V	121	-	4,4,4	0.34	0	4,4,4	0.52	0
3	DMF	V	135	-	4,4,4	0.38	0	4,4,4	0.42	0
3	DMF	V	136	-	4,4,4	0.38	0	4,4,4	0.44	0
3	DMF	V	16	-	4,4,4	0.34	0	4,4,4	0.37	0
3	DMF	W	249	-	4,4,4	0.28	0	4,4,4	0.19	0
3	DMF	X	141	-	4,4,4	0.29	0	4,4,4	0.33	0
3	DMF	Y	249	-	4,4,4	0.45	0	4,4,4	0.53	0
3	DMF	Y	250	-	4,4,4	0.28	0	4,4,4	0.38	0
3	DMF	Y	251	-	4,4,4	0.38	0	4,4,4	0.43	0
3	DMF	Z	105	-	4,4,4	0.36	0	4,4,4	0.36	0
3	DMF	Z	122	-	4,4,4	0.33	0	4,4,4	0.35	0
3	DMF	Z	27	-	4,4,4	0.32	0	4,4,4	0.41	0
3	DMF	Z	50	-	4,4,4	0.33	0	4,4,4	0.37	0
3	DMF	Z	69	-	4,4,4	0.34	0	4,4,4	0.32	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	1	249	-	-	0/2/2/2	0/0/0/0
3	DMF	1	250	-	-	0/2/2/2	0/0/0/0
3	DMF	2	145	-	-	0/2/2/2	0/0/0/0
3	DMF	2	99	-	-	0/2/2/2	0/0/0/0
3	DMF	A	249	-	-	0/2/2/2	0/0/0/0
3	DMF	A	8	-	-	0/2/2/2	0/0/0/0
3	DMF	B	249	-	-	0/2/2/2	0/0/0/0
3	DMF	C	15	-	-	0/2/2/2	0/0/0/0
3	DMF	D	249	-	-	0/2/2/2	0/0/0/0
3	DMF	E	104	-	-	0/2/2/2	0/0/0/0
3	DMF	E	113	-	-	0/2/2/2	0/0/0/0
3	DMF	E	28	-	-	0/2/2/2	0/0/0/0
3	DMF	F	249	-	-	0/2/2/2	0/0/0/0
3	DMF	G	137	-	-	0/2/2/2	0/0/0/0
3	DMF	G	140	-	-	0/2/2/2	0/0/0/0
3	DMF	G	20	-	-	0/2/2/2	0/0/0/0
3	DMF	H	142	-	-	0/2/2/2	0/0/0/0
3	DMF	H	41	-	-	0/2/2/2	0/0/0/0
3	DMF	I	249	-	-	0/2/2/2	0/0/0/0
3	DMF	J	46	-	-	0/2/2/2	0/0/0/0
3	DMF	K	249	-	-	0/2/2/2	0/0/0/0
3	DMF	K	250	-	-	0/2/2/2	0/0/0/0
3	DMF	K	251	-	-	0/2/2/2	0/0/0/0
3	DMF	L	138	-	-	0/2/2/2	0/0/0/0
3	DMF	L	14	-	-	0/2/2/2	0/0/0/0
3	DMF	L	60	-	-	0/2/2/2	0/0/0/0
3	DMF	L	9	-	-	0/2/2/2	0/0/0/0
3	DMF	M	249	-	-	0/2/2/2	0/0/0/0
3	DMF	N	21	-	-	0/2/2/2	0/0/0/0
3	DMF	N	36	-	-	0/2/2/2	0/0/0/0
3	DMF	O	7	-	-	0/2/2/2	0/0/0/0
3	DMF	P	107	-	-	0/2/2/2	0/0/0/0
3	DMF	P	133	-	-	0/2/2/2	0/0/0/0
3	DMF	P	23	-	-	0/2/2/2	0/0/0/0
3	DMF	Q	249	-	-	0/2/2/2	0/0/0/0
3	DMF	Q	250	-	-	0/2/2/2	0/0/0/0
3	DMF	Q	251	-	-	0/2/2/2	0/0/0/0
3	DMF	S	249	-	-	0/2/2/2	0/0/0/0
3	DMF	S	250	-	-	0/2/2/2	0/0/0/0
3	DMF	T	134	-	-	0/2/2/2	0/0/0/0
3	DMF	T	62	-	-	0/2/2/2	0/0/0/0
3	DMF	U	249	-	-	0/2/2/2	0/0/0/0
3	DMF	U	250	-	-	0/2/2/2	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	DMF	V	117	-	-	0/2/2/2	0/0/0/0
3	DMF	V	121	-	-	0/2/2/2	0/0/0/0
3	DMF	V	135	-	-	0/2/2/2	0/0/0/0
3	DMF	V	136	-	-	0/2/2/2	0/0/0/0
3	DMF	V	16	-	-	0/2/2/2	0/0/0/0
3	DMF	W	249	-	-	0/2/2/2	0/0/0/0
3	DMF	X	141	-	-	0/2/2/2	0/0/0/0
3	DMF	Y	249	-	-	0/2/2/2	0/0/0/0
3	DMF	Y	250	-	-	0/2/2/2	0/0/0/0
3	DMF	Y	251	-	-	0/2/2/2	0/0/0/0
3	DMF	Z	105	-	-	0/2/2/2	0/0/0/0
3	DMF	Z	122	-	-	0/2/2/2	0/0/0/0
3	DMF	Z	27	-	-	0/2/2/2	0/0/0/0
3	DMF	Z	50	-	-	0/2/2/2	0/0/0/0
3	DMF	Z	69	-	-	0/2/2/2	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

32 monomers are involved in 55 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	1	249	DMF	2	0
3	2	145	DMF	1	0
3	2	99	DMF	1	0
3	A	249	DMF	1	0
3	E	104	DMF	3	0
3	E	113	DMF	1	0
3	G	137	DMF	1	0
3	G	140	DMF	2	0
3	G	20	DMF	1	0
3	H	142	DMF	1	0
3	I	249	DMF	1	0
3	K	249	DMF	1	0
3	K	251	DMF	4	0
3	L	138	DMF	3	0
3	L	9	DMF	1	0
3	N	21	DMF	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	P	107	DMF	2	0
3	P	133	DMF	1	0
3	Q	251	DMF	5	0
3	S	249	DMF	1	0
3	T	134	DMF	2	0
3	T	62	DMF	3	0
3	U	249	DMF	1	0
3	V	117	DMF	1	0
3	V	121	DMF	3	0
3	V	135	DMF	1	0
3	V	136	DMF	1	0
3	V	16	DMF	1	0
3	W	249	DMF	1	0
3	Y	249	DMF	1	0
3	Z	105	DMF	2	0
3	Z	69	DMF	5	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	2	222/240 (92%)	-0.30	1 (0%) 91 92	17, 32, 63, 84	0
1	C	222/240 (92%)	0.59	15 (6%) 20 23	26, 43, 72, 89	0
1	E	223/240 (92%)	-0.14	6 (2%) 58 62	19, 32, 67, 134	0
1	G	222/240 (92%)	-0.24	0 100 100	18, 28, 62, 83	0
1	H	223/240 (92%)	0.11	4 (1%) 71 75	22, 36, 65, 86	0
1	J	222/240 (92%)	0.16	10 (4%) 37 42	24, 38, 71, 89	0
1	L	222/240 (92%)	-0.33	5 (2%) 64 67	18, 30, 64, 83	0
1	N	223/240 (92%)	-0.35	1 (0%) 93 93	15, 30, 64, 118	0
1	P	222/240 (92%)	-0.15	4 (1%) 71 75	19, 33, 67, 85	0
1	R	222/240 (92%)	-0.09	3 (1%) 78 80	22, 34, 66, 84	0
1	T	223/240 (92%)	-0.22	2 (0%) 85 88	21, 33, 67, 98	0
1	V	226/240 (94%)	-0.34	2 (0%) 85 88	16, 29, 65, 81	0
1	X	224/240 (93%)	-0.24	5 (2%) 65 69	19, 30, 66, 124	0
1	Z	224/240 (93%)	-0.18	0 100 100	21, 34, 68, 107	0
2	1	214/240 (89%)	0.17	7 (3%) 50 55	25, 48, 89, 127	0
2	A	214/240 (89%)	1.39	56 (26%) 1 1	37, 64, 111, 131	0
2	B	214/240 (89%)	1.88	85 (39%) 0 0	40, 87, 134, 168	0
2	D	213/240 (88%)	0.34	19 (8%) 12 13	24, 49, 98, 124	0
2	F	214/240 (89%)	0.39	25 (11%) 6 6	24, 48, 96, 127	0
2	I	214/240 (89%)	0.86	28 (13%) 5 4	34, 53, 98, 125	0
2	K	214/240 (89%)	0.44	17 (7%) 15 17	25, 49, 92, 126	0
2	M	214/240 (89%)	0.04	5 (2%) 64 67	22, 45, 90, 124	0
2	O	214/240 (89%)	0.49	21 (9%) 10 10	26, 54, 95, 129	0
2	Q	214/240 (89%)	0.09	9 (4%) 40 45	20, 42, 90, 126	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
2	S	214/240 (89%)	0.73	36 (16%) 2 2	28, 53, 98, 126	0
2	U	214/240 (89%)	0.40	14 (6%) 22 25	22, 51, 99, 126	0
2	W	214/240 (89%)	0.30	16 (7%) 17 19	22, 46, 94, 125	0
2	Y	214/240 (89%)	0.25	11 (5%) 32 36	23, 46, 90, 127	0
All	All	6115/6720 (90%)	0.21	407 (6%) 21 23	15, 40, 90, 168	0

All (407) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	U	234	LEU	8.2
2	A	9	MET	8.0
2	A	203	LEU	7.1
2	Y	9	MET	7.0
2	1	9	MET	6.7
2	B	205	VAL	6.6
2	A	233	LEU	6.6
2	B	42	VAL	6.5
2	B	40	LEU	6.4
2	B	179	ASP	6.3
2	B	36	ALA	6.2
2	B	180	ALA	6.2
2	B	41	PHE	6.2
2	A	205	VAL	6.0
2	F	233	LEU	5.9
2	B	38	GLY	5.8
2	U	206	ALA	5.8
2	B	9	MET	5.7
2	S	205	VAL	5.6
2	D	204	GLY	5.5
2	S	203	LEU	5.4
2	I	203	LEU	5.4
2	A	178	THR	5.3
2	O	231	GLN	5.2
2	A	231	GLN	5.1
2	Q	169	GLU	5.1
2	B	211	ALA	5.0
2	I	205	VAL	5.0
2	A	13	MET	5.0
2	B	33	LEU	4.9
2	O	205	VAL	4.9
2	D	205	VAL	4.8

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Mol	Chain	Res	Type	RSRZ
2	Q	9	MET	4.8
2	S	9	MET	4.8
2	I	9	MET	4.8
2	B	31	VAL	4.7
2	D	206	ALA	4.7
2	Y	203	LEU	4.7
2	Y	10	GLU	4.6
2	F	9	MET	4.6
2	B	230	LEU	4.6
2	A	179	ASP	4.4
2	B	35	TYR	4.4
2	U	203	LEU	4.4
2	Y	169	GLU	4.4
2	B	37	GLY	4.4
2	I	230	LEU	4.3
2	1	12	ALA	4.3
2	B	61	GLY	4.2
2	A	234	LEU	4.2
2	B	39	VAL	4.2
2	B	169	GLU	4.2
2	S	234	LEU	4.2
2	A	227	GLY	4.2
2	B	222	PHE	4.2
2	B	44	GLU	4.2
2	A	171	TYR	4.2
2	A	182	ARG	4.2
2	B	34	ALA	4.2
1	J	414	PRO	4.1
2	B	213	LEU	4.1
2	S	162	PRO	4.1
2	D	182	ARG	4.1
2	M	9	MET	4.1
2	F	10	GLU	4.1
2	F	227	GLY	4.1
2	W	232	ALA	4.1
2	B	60	VAL	4.0
2	O	203	LEU	4.0
2	O	204	GLY	4.0
2	U	205	VAL	4.0
2	A	20	ALA	3.9
2	B	182	ARG	3.9
2	B	234	LEU	3.9

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Mol	Chain	Res	Type	RSRZ
1	J	415	GLN	3.9
2	S	228	SER	3.9
2	B	62	PHE	3.8
1	X	523	GLY	3.8
2	B	212	VAL	3.8
2	A	229	ALA	3.8
2	W	182	ARG	3.8
2	S	169	GLU	3.7
1	C	409	ILE	3.7
2	B	127	VAL	3.7
2	A	168	LYS	3.7
2	Q	131	GLY	3.7
2	A	165	ASN	3.7
2	I	135	ARG	3.6
2	B	133	THR	3.6
2	B	64	ALA	3.6
2	O	9	MET	3.6
1	C	501	GLY	3.6
2	B	48	ARG	3.6
2	B	232	ALA	3.6
2	A	169	GLU	3.6
2	B	153	PHE	3.6
2	W	48	ARG	3.6
2	I	182	ARG	3.5
2	K	203	LEU	3.5
2	I	172	ALA	3.5
2	Y	205	VAL	3.5
2	B	208	LEU	3.5
2	B	171	TYR	3.5
2	I	177	LEU	3.5
2	A	161	GLU	3.5
2	K	26	ARG	3.5
2	S	171	TYR	3.5
2	B	32	ALA	3.4
2	B	181	LEU	3.4
2	F	11	GLN	3.4
1	E	523	GLY	3.4
2	A	218	PRO	3.4
2	B	53	ILE	3.4
1	C	340	TYR	3.3
2	B	218	PRO	3.3
2	O	48	ARG	3.3

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Mol	Chain	Res	Type	RSRZ
2	W	205	VAL	3.3
2	F	178	THR	3.3
1	C	415	GLN	3.3
2	B	225	ILE	3.3
2	B	215	ALA	3.3
2	W	203	LEU	3.3
1	J	413	ASP	3.2
1	H	372	VAL	3.2
2	B	210	VAL	3.2
2	F	171	TYR	3.2
2	A	189	ARG	3.2
1	C	519	GLU	3.2
2	B	226	THR	3.2
2	B	130	TYR	3.2
2	I	38	GLY	3.2
2	I	171	TYR	3.2
2	O	227	GLY	3.2
2	Q	130	TYR	3.2
2	F	203	LEU	3.2
2	K	25	ALA	3.2
2	U	169	GLU	3.2
2	D	169	GLU	3.1
2	K	11	GLN	3.1
1	J	412	SER	3.1
2	Q	10	GLU	3.1
2	B	123	CYS	3.1
2	A	11	GLN	3.1
2	A	135	ARG	3.1
2	B	168	LYS	3.1
2	K	205	VAL	3.1
2	D	189	ARG	3.1
2	W	9	MET	3.1
2	B	227	GLY	3.1
2	O	234	LEU	3.1
2	A	170	SER	3.1
2	B	54	SER	3.0
2	O	10	GLU	3.0
1	X	524	ALA	3.0
2	A	12	ALA	3.0
2	B	30	VAL	3.0
2	I	128	ALA	3.0
2	B	223	ARG	3.0

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Mol	Chain	Res	Type	RSRZ
1	R	412	SER	3.0
2	A	207	SER	3.0
1	N	523	GLY	3.0
2	B	184	ALA	3.0
1	C	407	TYR	3.0
2	I	229	ALA	3.0
2	A	124	VAL	2.9
2	B	52	LYS	2.9
2	F	168	LYS	2.9
2	B	172	ALA	2.9
1	2	415	GLN	2.9
2	A	130	TYR	2.9
2	O	11	GLN	2.9
2	F	205	VAL	2.9
2	A	188	LEU	2.9
2	A	159	THR	2.9
2	B	63	ALA	2.9
2	U	207	SER	2.9
1	V	526	THR	2.9
2	B	51	GLN	2.9
2	S	36	ALA	2.9
2	W	172	ALA	2.9
1	V	523	GLY	2.9
1	C	412	SER	2.9
2	O	233	LEU	2.9
2	I	189	ARG	2.9
2	A	180	ALA	2.9
2	F	206	ALA	2.9
2	B	209	GLU	2.9
2	A	204	GLY	2.9
2	I	228	SER	2.8
2	W	169	GLU	2.8
2	A	123	CYS	2.8
2	D	161	GLU	2.8
2	K	216	ASN	2.8
2	D	234	LEU	2.8
2	B	160	THR	2.8
1	H	412	SER	2.8
2	U	216	ASN	2.8
1	R	413	ASP	2.8
2	S	227	GLY	2.8
2	Q	231	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
2	1	10	GLU	2.8
2	S	37	GLY	2.8
2	A	172	ALA	2.8
2	O	206	ALA	2.8
2	O	229	ALA	2.8
2	O	168	LYS	2.8
2	Y	227	GLY	2.8
2	Q	203	LEU	2.8
2	B	173	GLU	2.8
2	O	169	GLU	2.8
2	B	154	VAL	2.8
2	F	226	THR	2.7
2	D	233	LEU	2.7
2	A	175	ALA	2.7
1	J	506	PRO	2.7
1	L	412	SER	2.7
2	B	175	ALA	2.7
2	B	186	ALA	2.7
1	J	522	SER	2.7
2	B	221	ALA	2.7
2	I	37	GLY	2.7
2	B	165	ASN	2.7
2	F	231	GLN	2.7
2	A	216	ASN	2.7
2	O	158	GLY	2.6
2	B	12	ALA	2.6
2	I	133	THR	2.6
1	P	411	ALA	2.6
2	A	43	ALA	2.6
2	K	21	ARG	2.6
2	A	176	SER	2.6
2	A	133	THR	2.6
2	D	168	LYS	2.6
2	W	206	ALA	2.6
2	F	131	GLY	2.6
2	A	160	THR	2.6
2	I	10	GLU	2.6
2	S	135	ARG	2.6
2	S	167	LEU	2.6
2	S	163	ILE	2.6
2	F	229	ALA	2.6
2	A	230	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
2	U	130	TYR	2.5
1	C	411	ALA	2.5
2	Q	135	ARG	2.5
2	S	158	GLY	2.5
2	B	189	ARG	2.5
2	B	203	LEU	2.5
2	B	183	ILE	2.5
2	S	159	THR	2.5
2	B	216	ASN	2.5
2	Y	204	GLY	2.5
1	C	414	PRO	2.5
1	J	519	GLU	2.5
2	K	133	THR	2.5
2	1	130	TYR	2.5
2	M	172	ALA	2.5
2	1	232	ALA	2.5
1	X	415	GLN	2.5
1	J	503	VAL	2.5
2	A	208	LEU	2.5
1	C	510	ILE	2.5
2	F	48	ARG	2.5
2	A	25	ALA	2.5
2	F	230	LEU	2.4
2	A	174	ASN	2.4
2	S	204	GLY	2.4
2	F	130	TYR	2.4
1	E	414	PRO	2.4
2	S	182	ARG	2.4
2	B	152	HIS	2.4
2	B	159	THR	2.4
2	D	232	ALA	2.4
2	B	125	ALA	2.4
2	S	189	ARG	2.4
2	S	21	ARG	2.4
2	I	153	PHE	2.4
1	P	519	GLU	2.4
2	A	44	GLU	2.4
2	S	10	GLU	2.4
2	S	161	GLU	2.4
2	D	133	THR	2.4
1	R	417	ALA	2.4
2	K	9	MET	2.4

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Mol	Chain	Res	Type	RSRZ
1	P	414	PRO	2.4
2	O	46	PRO	2.4
2	I	204	GLY	2.4
2	K	48	ARG	2.4
2	D	230	LEU	2.4
2	S	179	ASP	2.4
1	J	416	SER	2.4
2	B	138	LEU	2.3
2	D	178	THR	2.3
1	L	414	PRO	2.3
2	K	18	GLU	2.3
2	D	159	THR	2.3
2	U	232	ALA	2.3
2	A	10	GLU	2.3
2	Q	171	TYR	2.3
2	S	230	LEU	2.3
2	U	233	LEU	2.3
2	B	26	ARG	2.3
2	W	189	ARG	2.3
2	I	133	THR	2.3
1	T	415	GLN	2.3
2	B	164	ALA	2.3
2	A	14	ARG	2.3
1	C	397	GLY	2.3
1	J	500	ASP	2.3
1	X	411	ALA	2.3
2	I	233	LEU	2.3
1	P	415	GLN	2.3
2	W	231	GLN	2.3
2	A	163	ILE	2.3
2	U	163	ILE	2.3
2	B	224	ARG	2.3
2	B	50	LEU	2.3
1	C	500	ASP	2.3
2	D	48	ARG	2.3
1	E	411	ALA	2.2
2	B	206	ALA	2.2
2	I	190	ALA	2.2
2	B	207	SER	2.2
1	L	413	ASP	2.2
2	I	214	ASP	2.2
2	M	234	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
2	U	157	GLY	2.2
2	K	10	GLU	2.2
2	M	169	GLU	2.2
2	Y	189	ARG	2.2
2	B	229	ALA	2.2
2	F	172	ALA	2.2
2	A	62	PHE	2.2
2	I	169	GLU	2.2
2	S	208	LEU	2.2
1	H	396	GLN	2.2
2	U	231	GLN	2.2
2	Y	231	GLN	2.2
2	S	12	ALA	2.2
2	Y	174	ASN	2.2
2	S	168	LYS	2.2
2	K	130	TYR	2.2
1	L	519	GLU	2.2
2	A	173	GLU	2.2
2	F	232	ALA	2.2
2	K	206	ALA	2.2
2	I	14	ARG	2.2
2	A	177	LEU	2.2
2	B	167	LEU	2.2
2	B	11	GLN	2.2
2	U	131	GLY	2.2
2	M	10	GLU	2.2
2	I	179	ASP	2.2
1	H	411	ALA	2.2
2	I	130	TYR	2.2
2	A	158	GLY	2.2
2	A	151	PRO	2.2
2	S	152	HIS	2.1
2	B	49	SER	2.1
2	S	13	MET	2.1
2	S	48	ARG	2.1
2	B	136	PRO	2.1
1	E	415	GLN	2.1
2	K	228	SER	2.1
2	B	156	MET	2.1
2	W	133	THR	2.1
1	E	397	GLY	2.1
2	D	228	SER	2.1

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Mol	Chain	Res	Type	RSRZ
2	S	130	TYR	2.1
1	C	339	ASP	2.1
1	E	412	SER	2.1
2	D	14	ARG	2.1
2	I	48	ARG	2.1
2	O	189	ARG	2.1
2	S	14	ARG	2.1
2	W	135	ARG	2.1
2	S	233	LEU	2.1
2	W	234	LEU	2.1
1	T	414	PRO	2.1
2	F	228	SER	2.1
2	W	207	SER	2.1
2	S	229	ALA	2.1
2	Y	206	ALA	2.1
2	1	13	MET	2.1
2	D	38	GLY	2.1
2	A	33	LEU	2.1
2	F	204	GLY	2.1
2	S	131	GLY	2.1
2	A	226	THR	2.1
2	K	135	ARG	2.1
2	O	230	LEU	2.1
1	C	413	ASP	2.1
1	X	500	ASP	2.1
2	F	182	ARG	2.0
2	K	131	GLY	2.0
2	W	233	LEU	2.0
1	L	416	SER	2.0
2	O	228	SER	2.0
2	F	216	ASN	2.0
1	C	352	ALA	2.0
2	I	36	ALA	2.0
2	A	228	SER	2.0
2	S	172	ALA	2.0
2	B	162	PRO	2.0
2	A	49	SER	2.0
2	F	234	LEU	2.0
2	O	226	THR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	LLDF	B-factors(Å ²)	Q<0.9
3	DMF	L	9	5/5	0.86	0.33	19.71	53,54,82,89	0
3	DMF	V	135	5/5	0.78	0.34	18.82	35,51,75,87	0
3	DMF	Z	105	5/5	0.93	0.34	18.82	22,57,81,82	0
3	DMF	G	20	5/5	0.88	0.30	17.04	27,40,70,93	0
3	DMF	Z	50	5/5	0.86	0.40	14.11	86,90,126,146	0
3	DMF	L	14	5/5	0.86	0.40	12.23	30,38,70,90	0
3	DMF	T	62	5/5	0.89	0.33	11.60	45,61,83,95	0
3	DMF	P	23	5/5	0.96	0.20	10.41	49,62,83,93	0
3	DMF	C	15	5/5	0.92	0.27	8.23	54,69,74,107	0
3	DMF	G	137	5/5	0.97	0.20	8.23	27,35,51,57	0
3	DMF	Z	27	5/5	0.91	0.25	7.34	56,66,86,97	0
3	DMF	P	133	5/5	0.93	0.23	6.95	62,69,81,113	0
3	DMF	Z	122	5/5	0.94	0.25	6.70	26,28,76,84	0
3	DMF	E	104	5/5	0.95	0.26	6.50	33,42,51,57	0
3	DMF	Y	249	5/5	0.91	0.21	6.07	19,24,64,70	0
3	DMF	S	249	5/5	0.93	0.29	5.68	37,57,66,75	0
3	DMF	I	249	5/5	0.94	0.26	5.39	19,49,53,60	0
3	DMF	L	60	5/5	0.93	0.20	5.21	35,39,50,58	0
3	DMF	N	36	5/5	0.93	0.22	5.11	35,43,74,83	0
3	DMF	Y	251	5/5	0.86	0.27	4.89	42,45,61,63	0
3	DMF	L	138	5/5	0.94	0.37	4.88	40,47,72,82	0
3	DMF	P	107	5/5	0.92	0.31	4.86	39,58,69,70	0
3	DMF	V	136	5/5	0.91	0.24	4.61	36,44,59,63	0
3	DMF	V	117	5/5	0.86	0.24	4.61	46,50,66,72	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	DMF	H	142	5/5	0.90	0.27	4.54	40,62,68,72	0
3	DMF	Q	249	5/5	0.97	0.20	4.38	31,39,42,58	0
3	DMF	W	249	5/5	0.96	0.16	4.30	25,31,39,60	0
3	DMF	U	249	5/5	0.95	0.20	4.14	11,23,39,54	0
3	DMF	K	251	5/5	0.88	0.23	4.11	39,50,64,80	0
3	DMF	2	145	5/5	0.93	0.22	4.11	29,42,52,58	0
3	DMF	G	140	5/5	0.96	0.19	4.10	24,33,57,59	0
3	DMF	1	250	5/5	0.88	0.22	3.87	43,63,73,87	0
3	DMF	K	250	5/5	0.91	0.25	3.84	42,58,73,78	0
3	DMF	E	113	5/5	0.94	0.31	3.69	48,56,75,77	0
3	DMF	B	249	5/5	0.93	0.27	3.28	30,46,52,59	0
3	DMF	M	249	5/5	0.97	0.18	3.12	12,15,43,51	0
3	DMF	J	46	5/5	0.93	0.21	2.90	29,54,61,63	0
3	DMF	U	250	5/5	0.89	0.20	2.84	50,54,62,63	0
3	DMF	Q	251	5/5	0.90	0.24	2.83	26,56,59,84	0
3	DMF	O	7	5/5	0.92	0.21	2.79	38,48,59,62	0
3	DMF	Z	69	5/5	0.83	0.34	2.61	53,80,102,104	0
3	DMF	H	41	5/5	0.93	0.21	2.54	53,56,63,80	0
3	DMF	V	121	5/5	0.96	0.16	2.41	26,27,49,50	0
3	DMF	I	249	5/5	0.98	0.21	2.36	41,43,50,58	0
3	DMF	A	249	5/5	0.94	0.24	2.35	23,35,53,58	0
3	DMF	2	99	5/5	0.94	0.17	1.67	33,35,64,80	0
3	DMF	D	249	5/5	0.96	0.15	1.52	21,22,39,53	0
3	DMF	T	134	5/5	0.96	0.17	1.29	33,56,59,79	0
3	DMF	N	21	5/5	0.95	0.15	0.87	17,53,56,76	0
3	DMF	Y	250	5/5	0.97	0.15	0.85	14,36,48,55	0
3	DMF	S	250	5/5	0.92	0.18	0.68	26,42,81,85	0
3	DMF	A	8	5/5	0.93	0.20	0.56	16,61,76,85	0
3	DMF	Q	250	5/5	0.95	0.14	0.50	23,37,45,58	0
3	DMF	F	249	5/5	0.97	0.13	0.45	26,28,37,42	0
3	DMF	K	249	5/5	0.98	0.12	-0.29	23,32,46,48	0
3	DMF	E	28	5/5	0.89	0.32	-	36,68,72,88	0
3	DMF	V	16	5/5	0.91	0.29	-	36,61,65,73	0
3	DMF	X	141	5/5	0.89	0.36	-	40,55,82,85	0

6.5 Other polymers ⓘ

There are no such residues in this entry.