



Full wwPDB EM Validation Report ⓘ

Mar 16, 2026 – 07:42 PM UTC

PDB ID : 7UMS / pdb_00007ums
EMDB ID : EMD-26608
Title : Structure of the VP5*/VP8* assembly from the human rotavirus strain CDC-9
in complex with antibody 41 - Upright conformation
Authors : Jenni, S.; Zongli, L.; Wang, Y.; Bessey, T.; Salgado, E.N.; Schmidt, A.G.;
Greenberg, H.B.; Jiang, B.; Harrison, S.C.
Deposited on : 2022-04-07
Resolution : 3.50 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

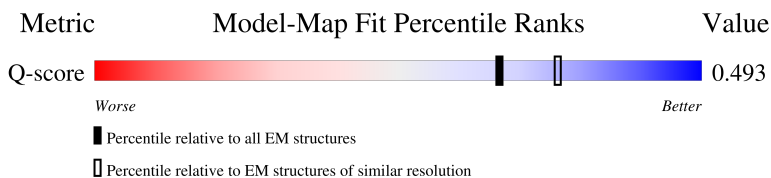
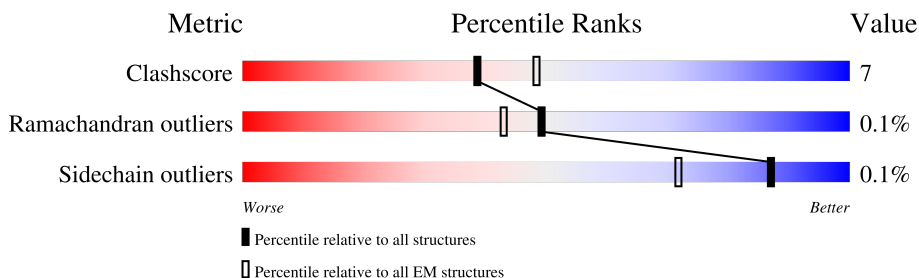
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 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.



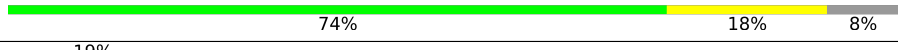
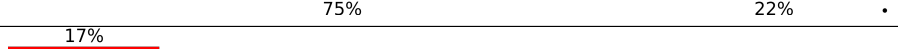
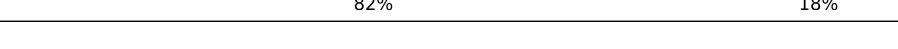
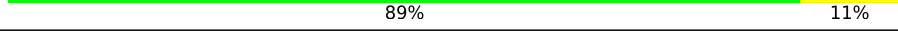
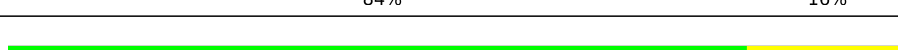
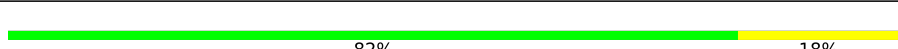
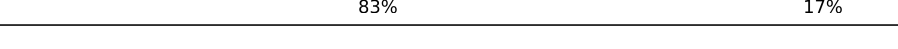
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13950 (3.00 - 4.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	T	230	
1	U	230	
1	V	230	
2	1	529	

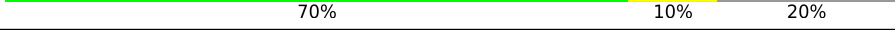
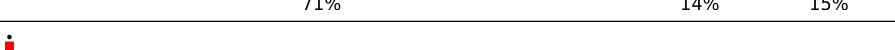
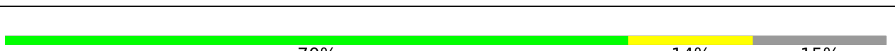
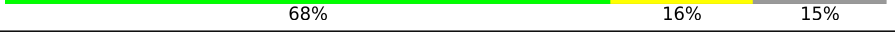
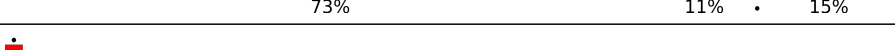
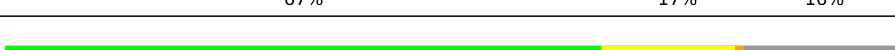
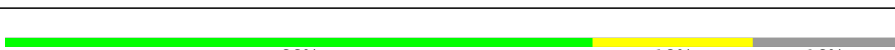
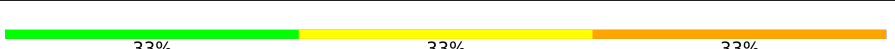
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Mol	Chain	Length	Quality of chain
2	2	529	
2	3	529	
3	4	236	
3	6	236	
4	5	217	
4	7	217	
5	A	397	
5	B	397	
5	C	397	
5	D	397	
5	E	397	
5	F	397	
5	G	397	
5	H	397	
5	I	397	
5	J	397	
5	K	397	
5	L	397	
5	M	397	
5	N	397	
5	O	397	
5	P	397	
5	Q	397	
5	R	397	
6	a	326	

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Mol	Chain	Length	Quality of chain
6	b	326	
6	c	326	
6	d	326	
6	e	326	
6	f	326	
6	g	326	
6	h	326	
6	i	326	
6	j	326	
6	k	326	
6	l	326	
6	m	326	
6	n	326	
6	o	326	
6	p	326	
6	q	326	
6	r	326	
7	S	3	

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 236285 atoms, of which 116945 are hydrogens and 0 are deuteriums.

In the tables below, 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 Outer capsid protein VP8*.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	T	227	Total	C	H	N	O	S	0	0
			3553	1151	1726	309	364	3		
1	U	227	Total	C	H	N	O	S	0	0
			3553	1151	1726	309	364	3		
1	V	25	Total	C	H	N	O		0	0
			401	130	197	31	43			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
T	28	GLY	GLU	conflict	UNP X4YMN0
T	51	ASP	GLY	conflict	UNP X4YMN0
U	28	GLY	GLU	conflict	UNP X4YMN0
U	51	ASP	GLY	conflict	UNP X4YMN0
V	28	GLY	GLU	conflict	UNP X4YMN0
V	51	ASP	GLY	conflict	UNP X4YMN0

- Molecule 2 is a protein called Outer capsid protein VP5*.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	1	499	Total	C	H	N	O	S	0	0
			7914	2514	3951	664	768	17		
2	2	499	Total	C	H	N	O	S	0	0
			7914	2514	3951	664	768	17		
2	3	486	Total	C	H	N	O	S	0	0
			7719	2455	3853	649	745	17		

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
1	250	ASP	ASN	conflict	UNP X4YMN0
1	331	PHE	SER	conflict	UNP X4YMN0
1	364	ILE	MET	conflict	UNP X4YMN0

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Chain	Residue	Modelled	Actual	Comment	Reference
1	378	ARG	LYS	conflict	UNP X4YMN0
1	385	HIS	ASP	conflict	UNP X4YMN0
1	388	LEU	ILE	conflict	UNP X4YMN0
1	499	ASN	ASP	conflict	UNP X4YMN0
1	605	ASN	SER	conflict	UNP X4YMN0
2	250	ASP	ASN	conflict	UNP X4YMN0
2	331	PHE	SER	conflict	UNP X4YMN0
2	364	ILE	MET	conflict	UNP X4YMN0
2	378	ARG	LYS	conflict	UNP X4YMN0
2	385	HIS	ASP	conflict	UNP X4YMN0
2	388	LEU	ILE	conflict	UNP X4YMN0
2	499	ASN	ASP	conflict	UNP X4YMN0
2	605	ASN	SER	conflict	UNP X4YMN0
3	250	ASP	ASN	conflict	UNP X4YMN0
3	331	PHE	SER	conflict	UNP X4YMN0
3	364	ILE	MET	conflict	UNP X4YMN0
3	378	ARG	LYS	conflict	UNP X4YMN0
3	385	HIS	ASP	conflict	UNP X4YMN0
3	388	LEU	ILE	conflict	UNP X4YMN0
3	499	ASN	ASP	conflict	UNP X4YMN0
3	605	ASN	SER	conflict	UNP X4YMN0

- Molecule 3 is a protein called Fab 41 heavy chain.

Mol	Chain	Residues	Atoms						AltConf	Trace
3	4	230	Total	C	H	N	O	S	0	0
			3396	1094	1667	289	339	7		
3	6	230	Total	C	H	N	O	S	0	0
			3396	1094	1667	289	339	7		

- Molecule 4 is a protein called Fab 41 light chain.

Mol	Chain	Residues	Atoms						AltConf	Trace
4	5	217	Total	C	H	N	O	S	0	0
			3202	1019	1562	276	339	6		
4	7	217	Total	C	H	N	O	S	0	0
			3202	1019	1562	276	339	6		

- Molecule 5 is a protein called Intermediate capsid protein VP6.

Mol	Chain	Residues	Atoms						AltConf	Trace
5	A	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0
5	B	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0
5	C	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0
5	D	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0
5	E	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0
5	F	397	Total 6296	C 2012	H 3128	N 552	O 589	S 15	0	0
5	G	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0
5	H	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0
5	I	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0
5	J	397	Total 6296	C 2012	H 3128	N 552	O 589	S 15	0	0
5	K	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0
5	L	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0
5	M	397	Total 6296	C 2012	H 3128	N 552	O 589	S 15	0	0
5	N	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0
5	O	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0
5	P	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0
5	Q	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0
5	R	397	Total 6295	C 2012	H 3127	N 552	O 589	S 15	0	0

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	281	VAL	ILE	conflict	UNP A0A223GHC7
B	281	VAL	ILE	conflict	UNP A0A223GHC7
C	281	VAL	ILE	conflict	UNP A0A223GHC7
D	281	VAL	ILE	conflict	UNP A0A223GHC7

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Chain	Residue	Modelled	Actual	Comment	Reference
E	281	VAL	ILE	conflict	UNP A0A223GHC7
F	281	VAL	ILE	conflict	UNP A0A223GHC7
G	281	VAL	ILE	conflict	UNP A0A223GHC7
H	281	VAL	ILE	conflict	UNP A0A223GHC7
I	281	VAL	ILE	conflict	UNP A0A223GHC7
J	281	VAL	ILE	conflict	UNP A0A223GHC7
K	281	VAL	ILE	conflict	UNP A0A223GHC7
L	281	VAL	ILE	conflict	UNP A0A223GHC7
M	281	VAL	ILE	conflict	UNP A0A223GHC7
N	281	VAL	ILE	conflict	UNP A0A223GHC7
O	281	VAL	ILE	conflict	UNP A0A223GHC7
P	281	VAL	ILE	conflict	UNP A0A223GHC7
Q	281	VAL	ILE	conflict	UNP A0A223GHC7
R	281	VAL	ILE	conflict	UNP A0A223GHC7

- Molecule 6 is a protein called Outer capsid glycoprotein VP7.

Mol	Chain	Residues	Atoms						AltConf	Trace
6	a	271	Total	C	H	N	O	S	0	0
			4279	1374	2111	351	426	17		
6	b	276	Total	C	H	N	O	S	0	0
			4357	1400	2148	358	434	17		
6	c	260	Total	C	H	N	O	S	0	0
			4096	1314	2023	333	409	17		
6	d	276	Total	C	H	N	O	S	0	0
			4357	1400	2148	358	434	17		
6	e	261	Total	C	H	N	O	S	0	0
			4103	1316	2026	334	410	17		
6	f	276	Total	C	H	N	O	S	0	0
			4357	1400	2148	358	434	17		
6	g	276	Total	C	H	N	O	S	0	0
			4357	1400	2148	358	434	17		
6	h	271	Total	C	H	N	O	S	0	0
			4279	1374	2111	351	426	17		
6	i	276	Total	C	H	N	O	S	0	0
			4357	1400	2148	358	434	17		
6	j	274	Total	C	H	N	O	S	0	0
			4326	1391	2134	354	430	17		
6	k	276	Total	C	H	N	O	S	0	0
			4357	1400	2148	358	434	17		
6	l	276	Total	C	H	N	O	S	0	0
			4357	1400	2148	358	434	17		

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Mol	Chain	Residues	Atoms						AltConf	Trace
6	m	275	Total	C	H	N	O	S	0	0
			4340	1395	2140	356	432	17		
6	n	276	Total	C	H	N	O	S	0	0
			4357	1400	2148	358	434	17		
6	o	276	Total	C	H	N	O	S	0	0
			4357	1400	2148	358	434	17		
6	p	275	Total	C	H	N	O	S	0	0
			4340	1395	2140	356	432	17		
6	q	271	Total	C	H	N	O	S	0	0
			4279	1374	2111	351	426	17		
6	r	275	Total	C	H	N	O	S	0	0
			4340	1395	2140	356	432	17		

There are 36 discrepancies between the modelled and reference sequences:

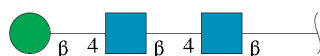
Chain	Residue	Modelled	Actual	Comment	Reference
a	108	ILE	THR	conflict	UNP B1NP55
a	147	SER	ASN	conflict	UNP B1NP55
b	108	ILE	THR	conflict	UNP B1NP55
b	147	SER	ASN	conflict	UNP B1NP55
c	108	ILE	THR	conflict	UNP B1NP55
c	147	SER	ASN	conflict	UNP B1NP55
d	108	ILE	THR	conflict	UNP B1NP55
d	147	SER	ASN	conflict	UNP B1NP55
e	108	ILE	THR	conflict	UNP B1NP55
e	147	SER	ASN	conflict	UNP B1NP55
f	108	ILE	THR	conflict	UNP B1NP55
f	147	SER	ASN	conflict	UNP B1NP55
g	108	ILE	THR	conflict	UNP B1NP55
g	147	SER	ASN	conflict	UNP B1NP55
h	108	ILE	THR	conflict	UNP B1NP55
h	147	SER	ASN	conflict	UNP B1NP55
i	108	ILE	THR	conflict	UNP B1NP55
i	147	SER	ASN	conflict	UNP B1NP55
j	108	ILE	THR	conflict	UNP B1NP55
j	147	SER	ASN	conflict	UNP B1NP55
k	108	ILE	THR	conflict	UNP B1NP55
k	147	SER	ASN	conflict	UNP B1NP55
l	108	ILE	THR	conflict	UNP B1NP55
l	147	SER	ASN	conflict	UNP B1NP55
m	108	ILE	THR	conflict	UNP B1NP55
m	147	SER	ASN	conflict	UNP B1NP55
n	108	ILE	THR	conflict	UNP B1NP55

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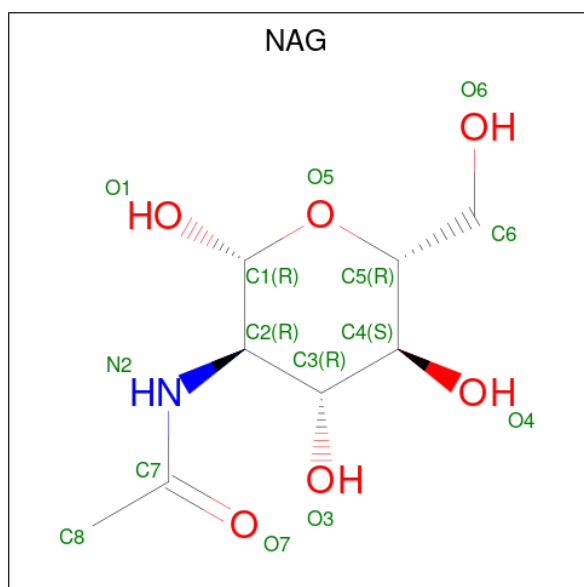
Chain	Residue	Modelled	Actual	Comment	Reference
n	147	SER	ASN	conflict	UNP B1NP55
o	108	ILE	THR	conflict	UNP B1NP55
o	147	SER	ASN	conflict	UNP B1NP55
p	108	ILE	THR	conflict	UNP B1NP55
p	147	SER	ASN	conflict	UNP B1NP55
q	108	ILE	THR	conflict	UNP B1NP55
q	147	SER	ASN	conflict	UNP B1NP55
r	108	ILE	THR	conflict	UNP B1NP55
r	147	SER	ASN	conflict	UNP B1NP55

- Molecule 7 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms					AltConf	Trace
7	S	3	Total	C	H	N	O	0	0
			76	22	37	2	15		

- Molecule 8 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms					AltConf
8	a	1	Total	C	H	N	O	0
			27	8	13	1	5	
8	b	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	b	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	c	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	c	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	d	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	d	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	e	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	e	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	f	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	f	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	g	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	g	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	h	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	h	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	i	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	i	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	j	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	j	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	k	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	k	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	l	1	Total	C	H	N	O	0
			28	8	14	1	5	

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Mol	Chain	Residues	Atoms					AltConf
8	l	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	m	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	m	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	n	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	n	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	o	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	o	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	p	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	p	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	q	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	q	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	r	1	Total	C	H	N	O	0
			28	8	14	1	5	
8	r	1	Total	C	H	N	O	0
			28	8	14	1	5	

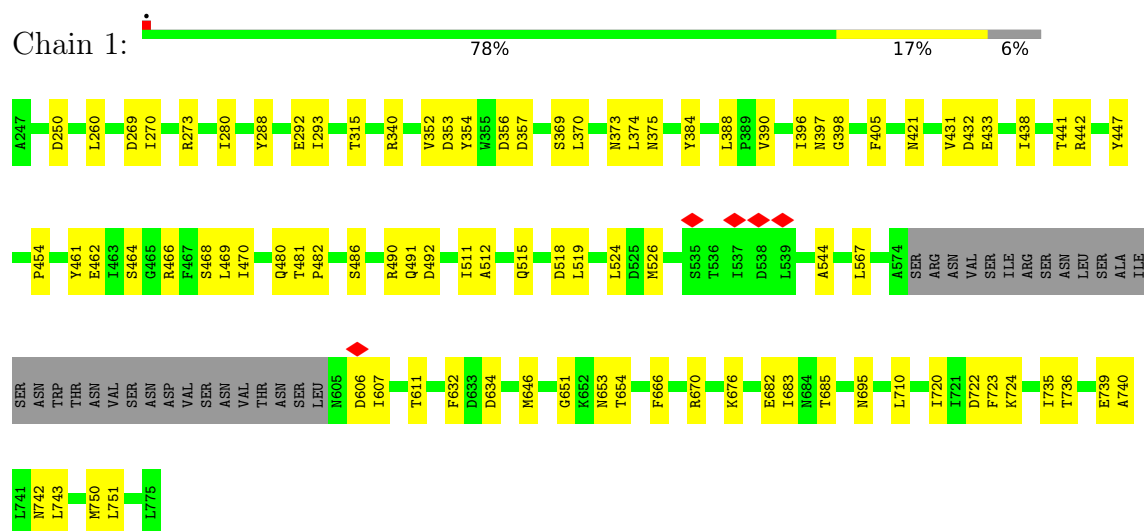
- Molecule 9 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		AltConf
9	a	4	Total	Ca	0
			4	4	
9	b	4	Total	Ca	0
			4	4	
9	c	4	Total	Ca	0
			4	4	
9	d	4	Total	Ca	0
			4	4	
9	e	4	Total	Ca	0
			4	4	
9	f	4	Total	Ca	0
			4	4	

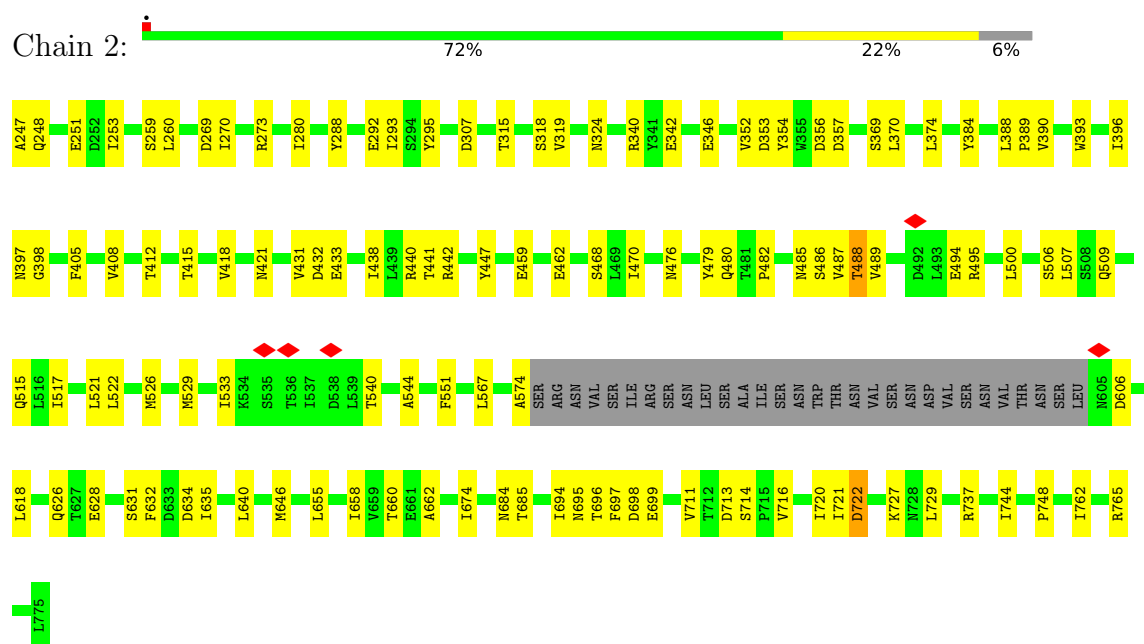
Continued on next page...

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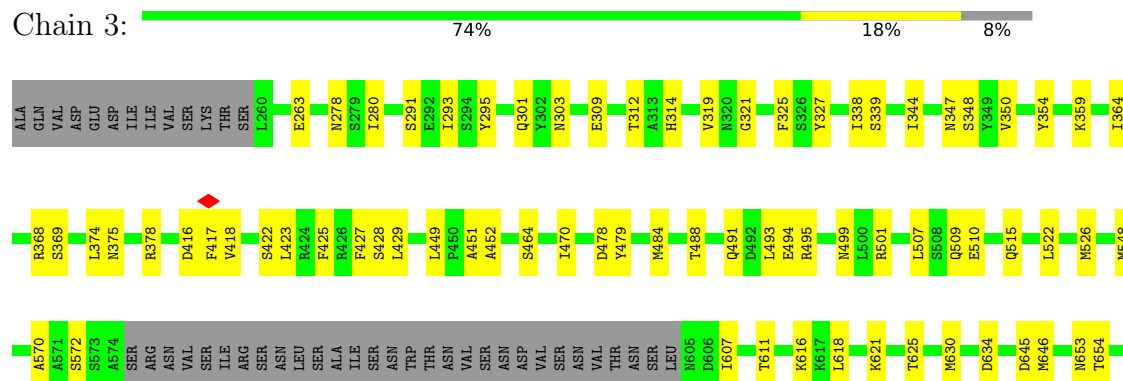
Mol	Chain	Residues	Atoms		AltConf
9	g	4	Total 4	Ca 4	0
9	h	4	Total 4	Ca 4	0
9	i	4	Total 4	Ca 4	0
9	j	4	Total 4	Ca 4	0
9	k	4	Total 4	Ca 4	0
9	l	4	Total 4	Ca 4	0
9	m	4	Total 4	Ca 4	0
9	n	4	Total 4	Ca 4	0
9	o	4	Total 4	Ca 4	0
9	p	4	Total 4	Ca 4	0
9	q	4	Total 4	Ca 4	0
9	r	4	Total 4	Ca 4	0



- Molecule 2: Outer capsid protein VP5*

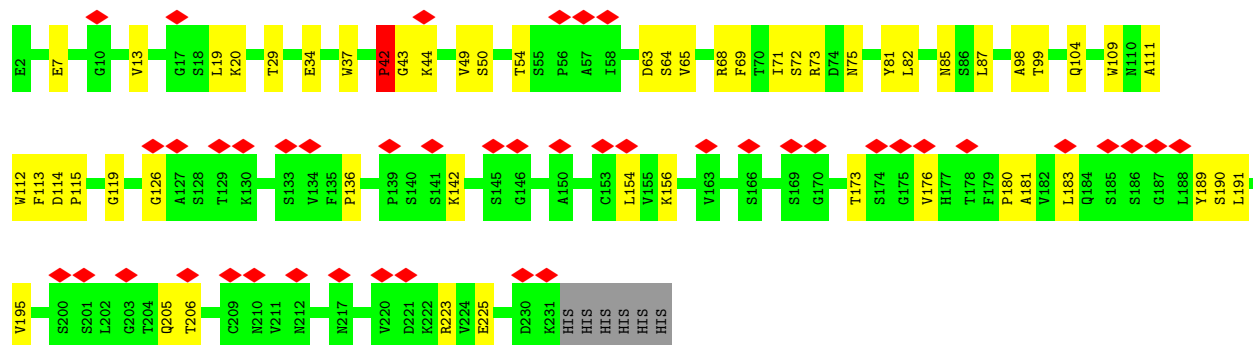
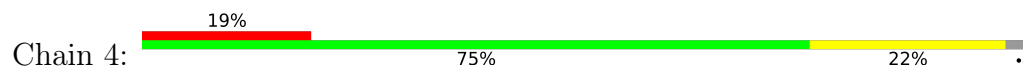


- Molecule 2: Outer capsid protein VP5*

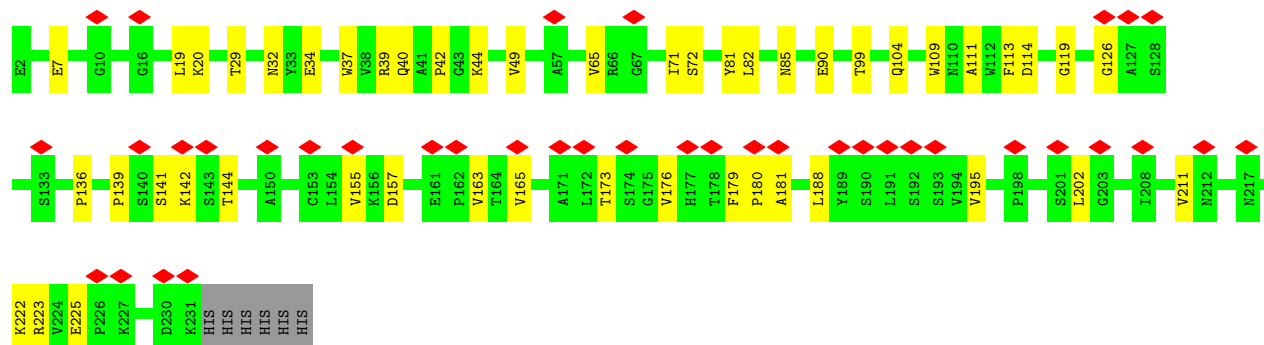
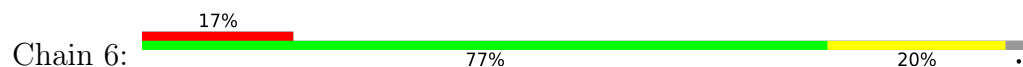




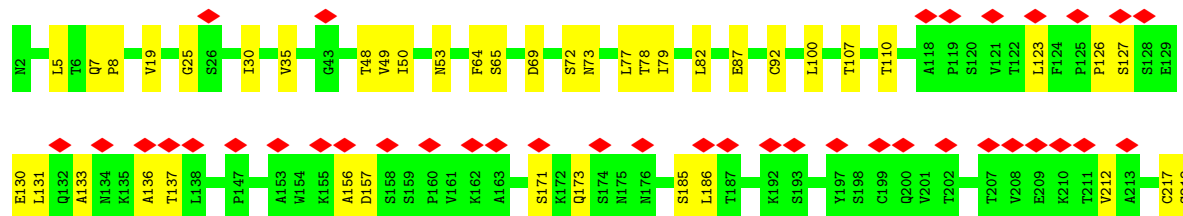
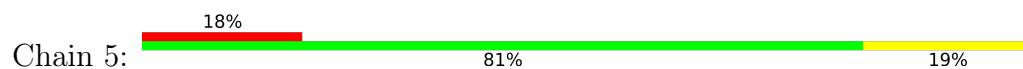
• Molecule 3: Fab 41 heavy chain



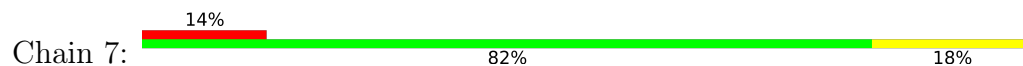
• Molecule 3: Fab 41 heavy chain

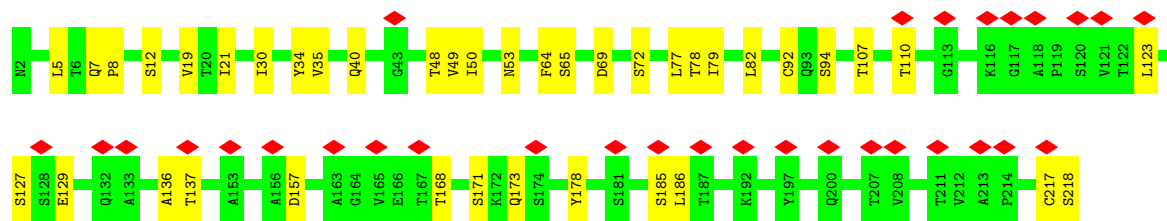


• Molecule 4: Fab 41 light chain



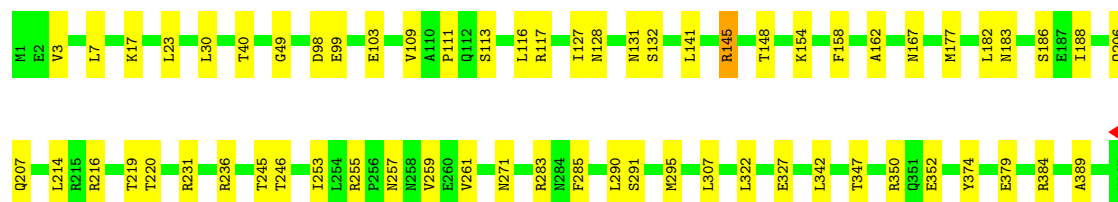
• Molecule 4: Fab 41 light chain





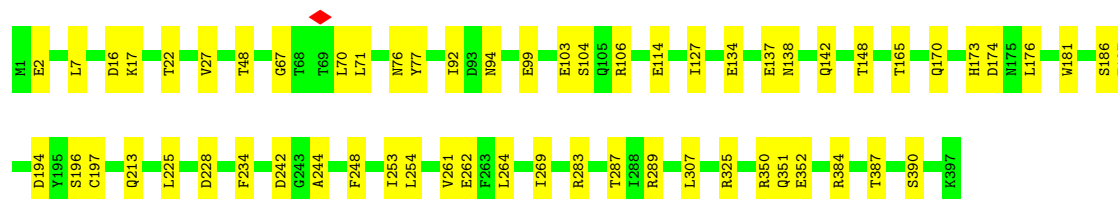
• Molecule 5: Intermediate capsid protein VP6

Chain A: 84% 16%



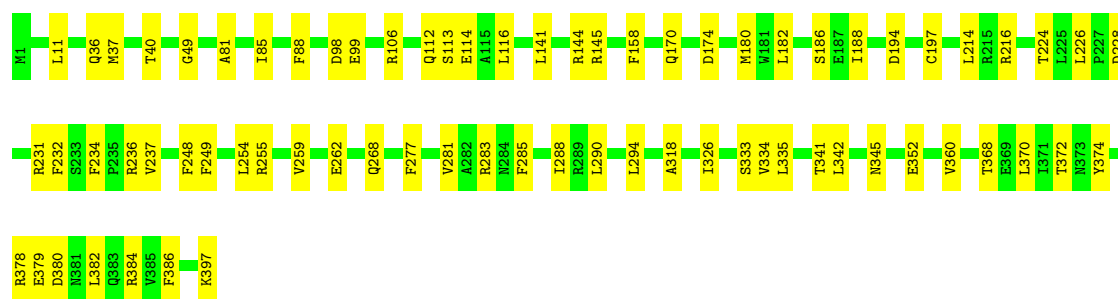
• Molecule 5: Intermediate capsid protein VP6

Chain B: 85% 15%



• Molecule 5: Intermediate capsid protein VP6

Chain C: 82% 18%



• Molecule 5: Intermediate capsid protein VP6

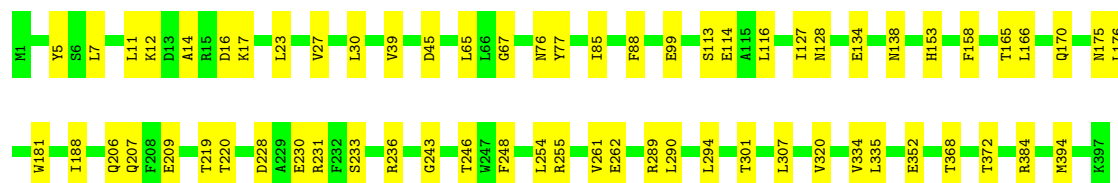
Chain D: 82% 18%





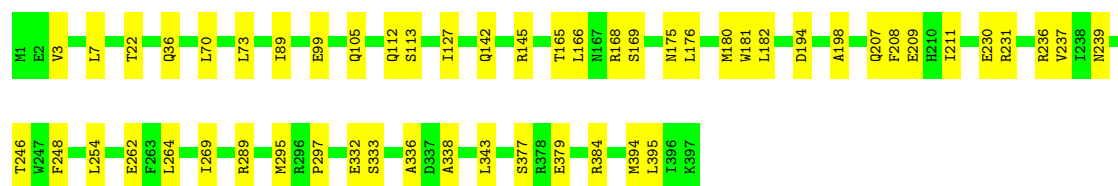
- Molecule 5: Intermediate capsid protein VP6

Chain E: 84% 16%



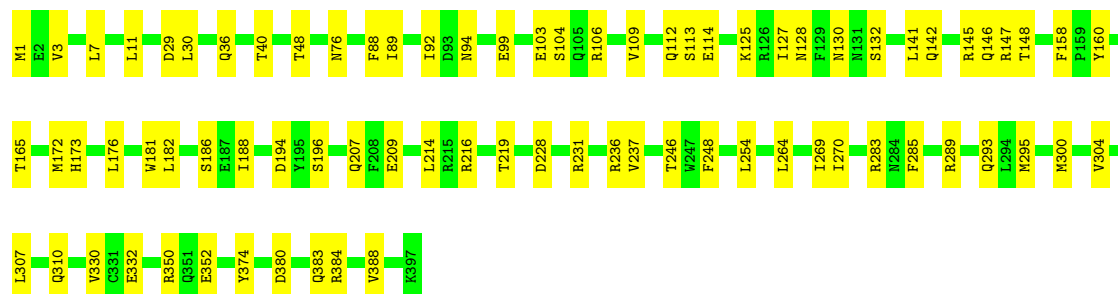
- Molecule 5: Intermediate capsid protein VP6

Chain F: 87% 13%



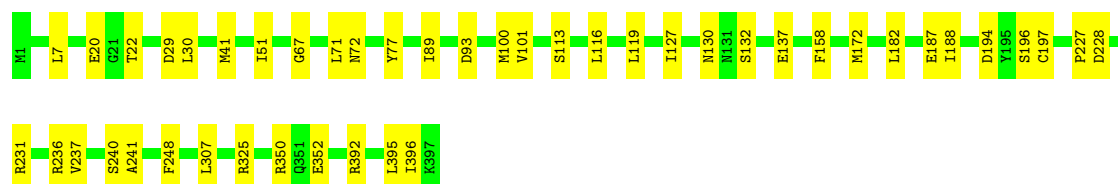
- Molecule 5: Intermediate capsid protein VP6

Chain G: 80% 20%

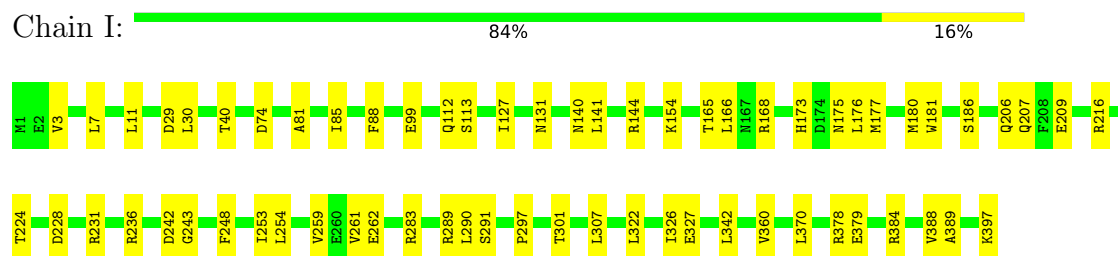


- Molecule 5: Intermediate capsid protein VP6

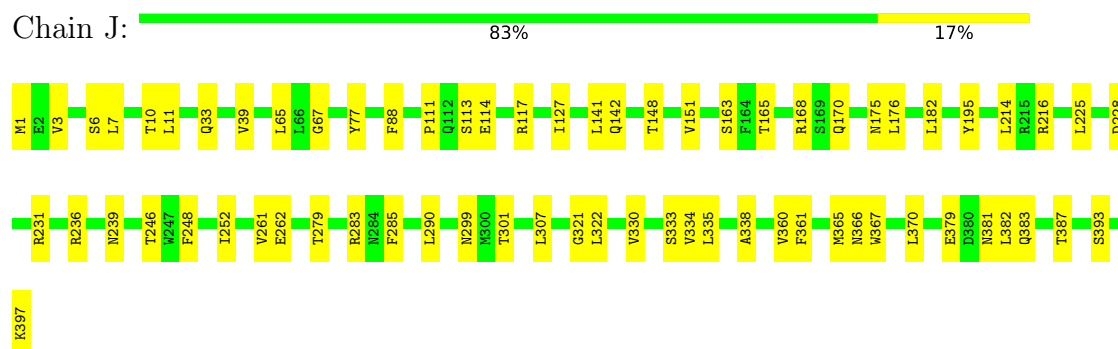
Chain H: 89% 11%



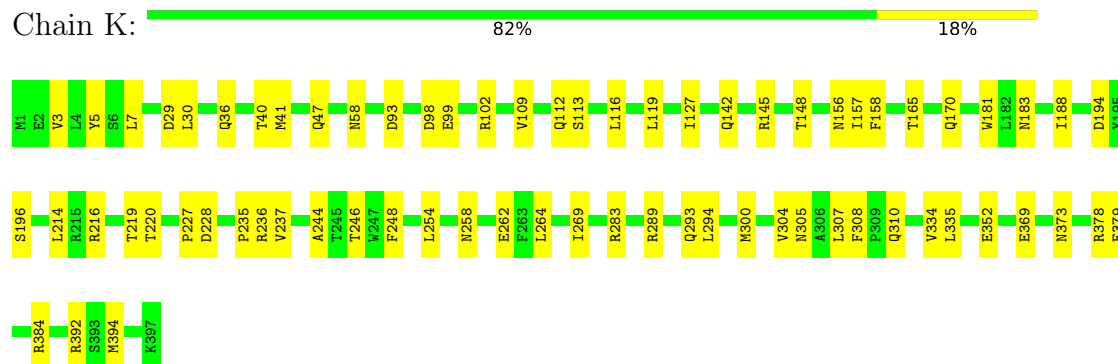
- Molecule 5: Intermediate capsid protein VP6



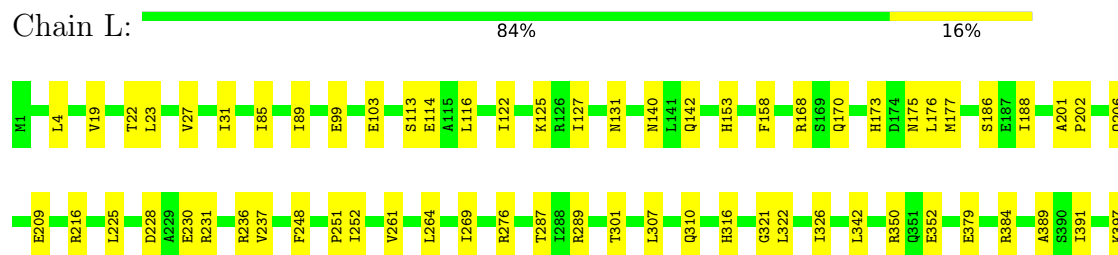
- Molecule 5: Intermediate capsid protein VP6



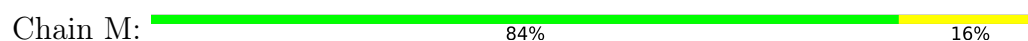
- Molecule 5: Intermediate capsid protein VP6

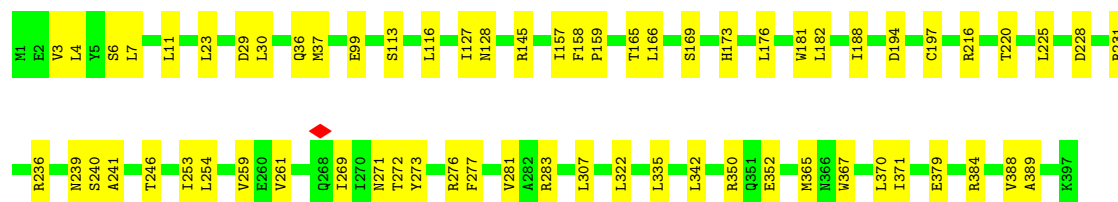


- Molecule 5: Intermediate capsid protein VP6



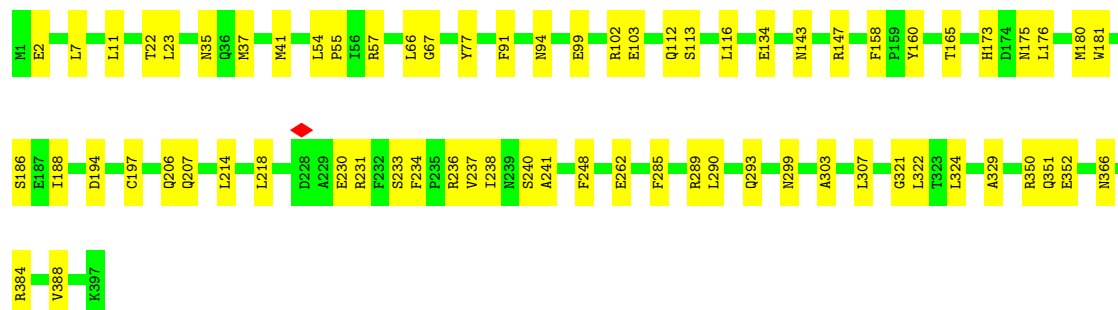
- Molecule 5: Intermediate capsid protein VP6





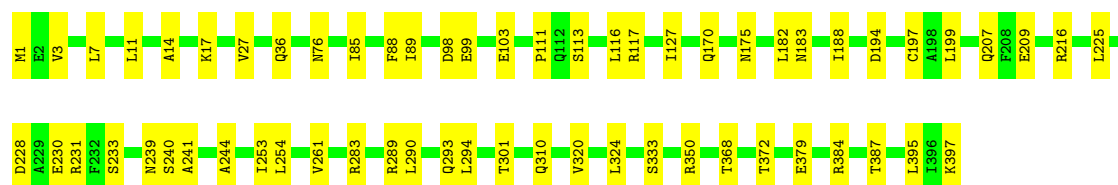
• Molecule 5: Intermediate capsid protein VP6

Chain N: 83% 17%



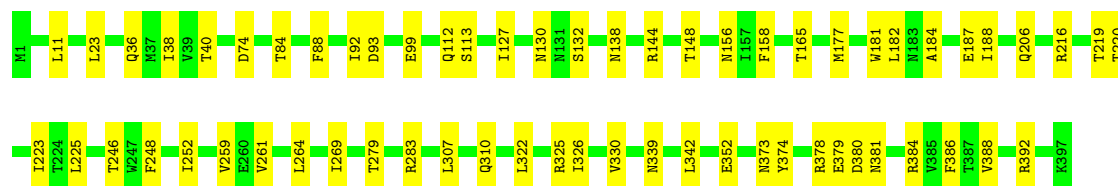
• Molecule 5: Intermediate capsid protein VP6

Chain O: 85% 15%



• Molecule 5: Intermediate capsid protein VP6

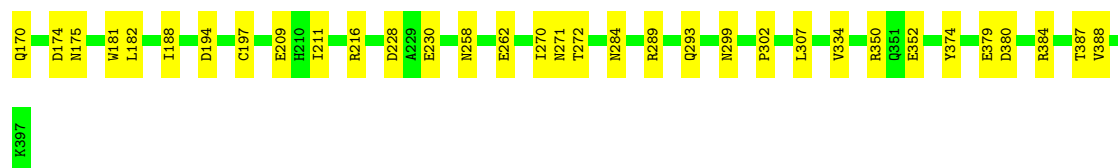
Chain P: 84% 16%



• Molecule 5: Intermediate capsid protein VP6

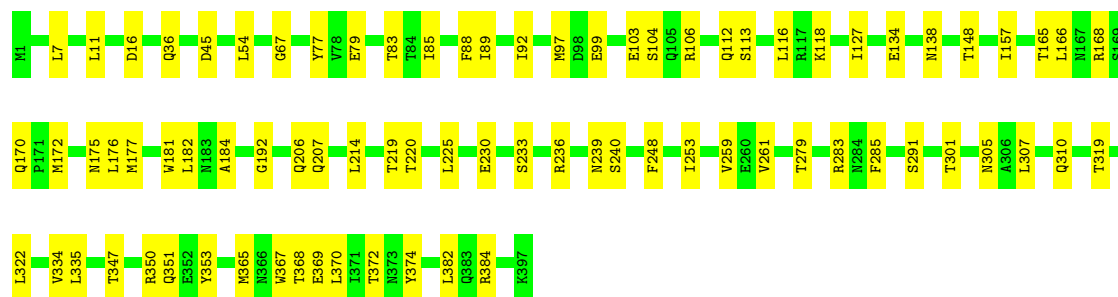
Chain Q: 83% 17%





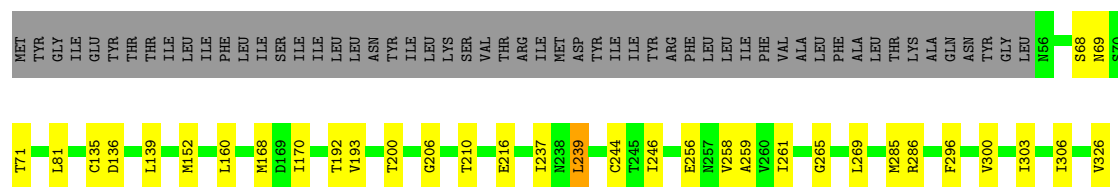
• Molecule 5: Intermediate capsid protein VP6

Chain R: 80% 20%



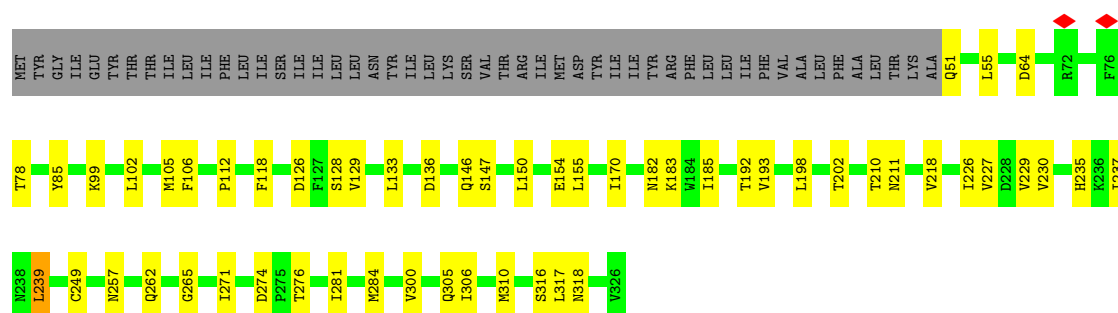
• Molecule 6: Outer capsid glycoprotein VP7

Chain a: 73% 10% 17%



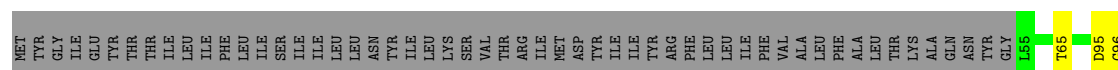
• Molecule 6: Outer capsid glycoprotein VP7

Chain b: 68% 17% 15%



• Molecule 6: Outer capsid glycoprotein VP7

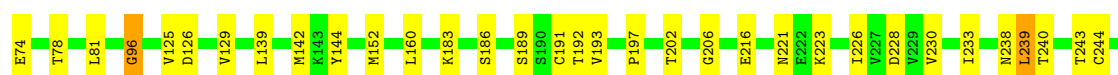
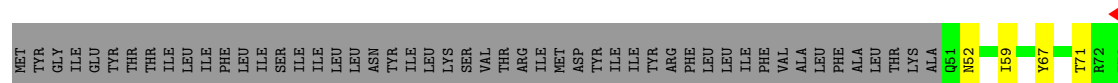
Chain c: 68% 12% 20%





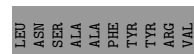
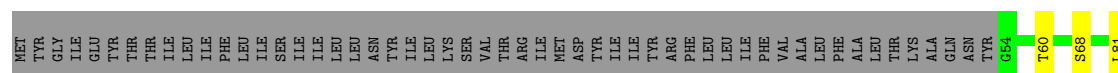
- Molecule 6: Outer capsid glycoprotein VP7

Chain d: 71% 13% 15%



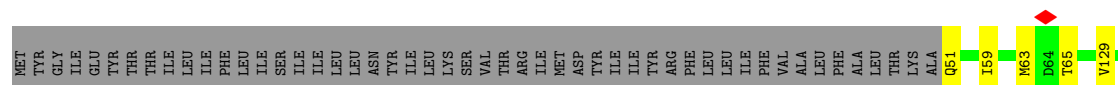
- Molecule 6: Outer capsid glycoprotein VP7

Chain e: 70% 10% 20%



- Molecule 6: Outer capsid glycoprotein VP7

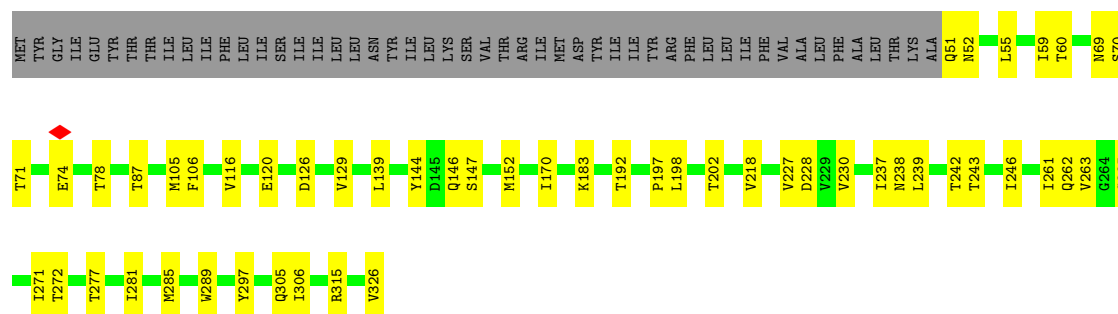
Chain f: 71% 14% 15%



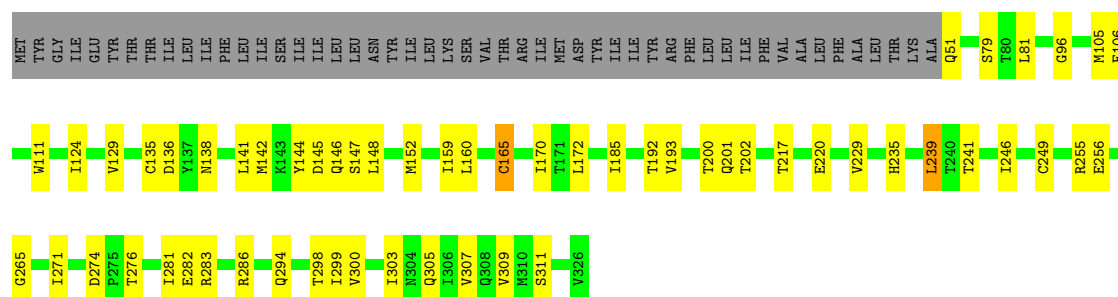
- Molecule 6: Outer capsid glycoprotein VP7

Chain g: 69% 15% 15%

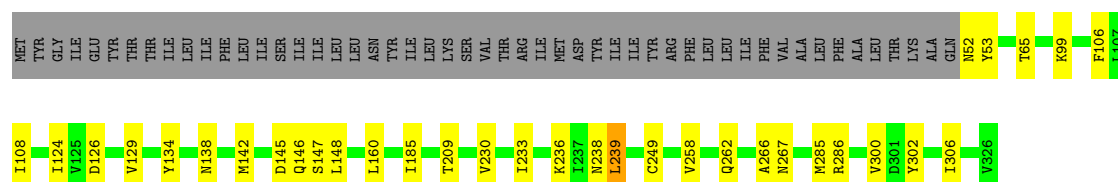
• Molecule 6: Outer capsid glycoprotein VP7

Chain k:  68% 16% 15%

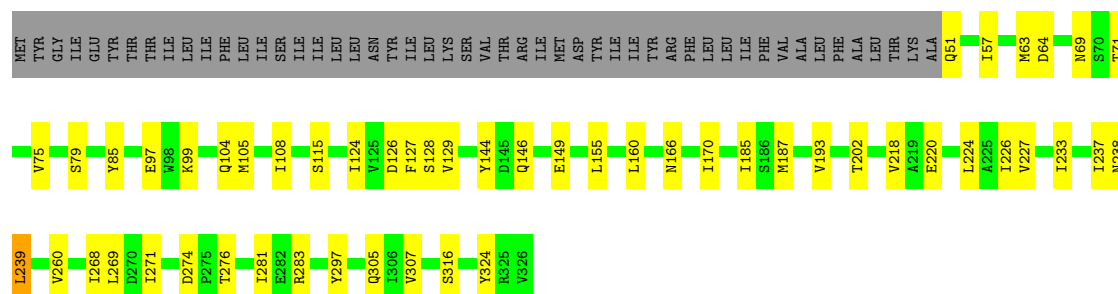
• Molecule 6: Outer capsid glycoprotein VP7

Chain l:  67% 17% 15%

• Molecule 6: Outer capsid glycoprotein VP7

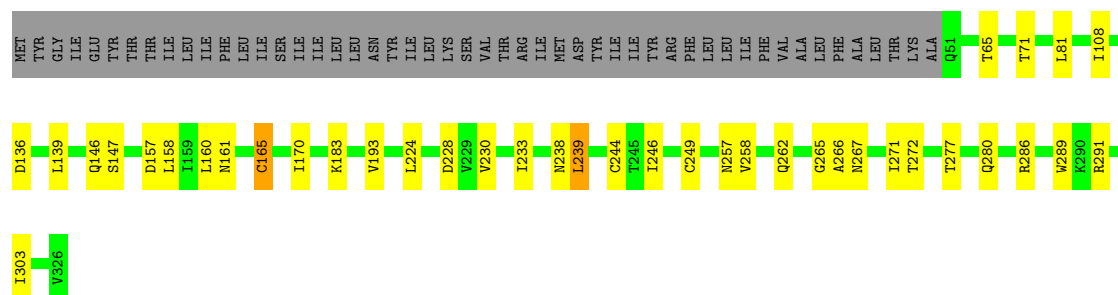
Chain m:  74% 10% 16%

• Molecule 6: Outer capsid glycoprotein VP7

Chain n:  68% 16% 15%

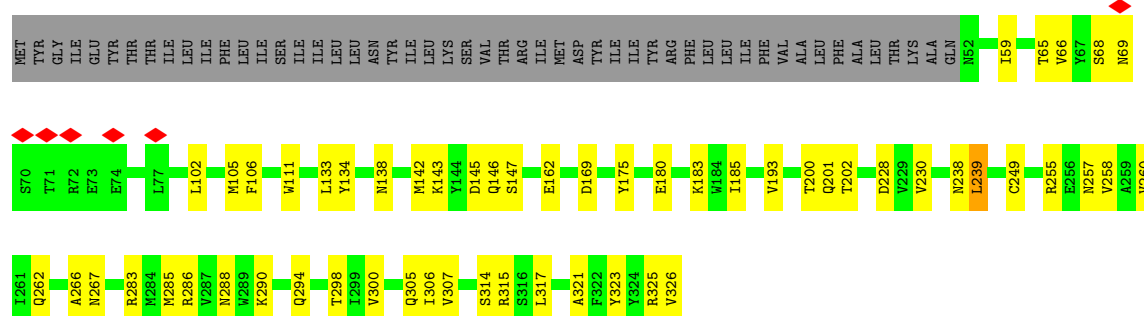
- Molecule 6: Outer capsid glycoprotein VP7

Chain o:  73% 11% 15%



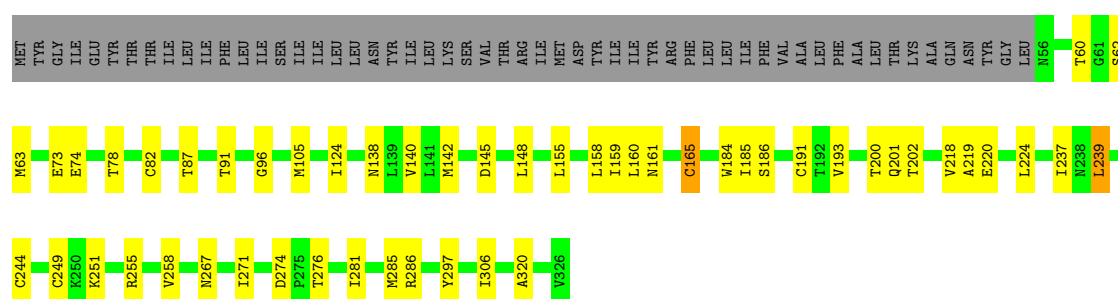
- Molecule 6: Outer capsid glycoprotein VP7

Chain p:  67% 17% 16%



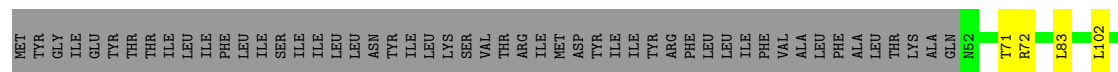
- Molecule 6: Outer capsid glycoprotein VP7

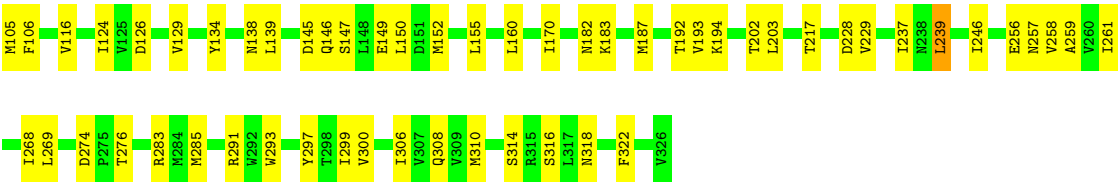
Chain q:  67% 15% 17%



- Molecule 6: Outer capsid glycoprotein VP7

Chain r:  66% 18% 16%





• Molecule 7: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	422920	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60.0	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3500	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.153	Depositor
Minimum map value	0.000	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.004	Depositor
Recommended contour level	0.01	Depositor
Map size ($\text{\AA}$)	393.91998, 393.91998, 393.91998	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.231, 1.231, 1.231	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, BMA, CA

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	T	0.20	0/1876	0.45	0/2565
1	U	0.20	0/1876	0.45	0/2565
1	V	0.24	0/206	0.41	0/279
2	1	0.21	0/4035	0.40	0/5459
2	2	0.21	0/4035	0.41	0/5459
2	3	0.23	0/3938	0.43	0/5327
3	4	0.26	0/1777	0.64	0/2425
3	6	0.26	0/1777	0.65	0/2425
4	5	0.24	0/1681	0.69	1/2295 (0.0%)
4	7	0.26	0/1681	0.70	0/2295
5	A	0.27	0/3238	0.40	0/4404
5	B	0.27	0/3238	0.41	0/4404
5	C	0.27	0/3238	0.44	0/4404
5	D	0.26	0/3238	0.40	0/4404
5	E	0.25	0/3238	0.39	0/4404
5	F	0.26	0/3238	0.41	0/4404
5	G	0.27	0/3238	0.42	0/4404
5	H	0.25	0/3238	0.39	0/4404
5	I	0.26	0/3238	0.43	0/4404
5	J	0.25	0/3238	0.39	0/4404
5	K	0.26	0/3238	0.40	0/4404
5	L	0.26	0/3238	0.40	0/4404
5	M	0.25	0/3238	0.40	0/4404
5	N	0.26	0/3238	0.40	0/4404
5	O	0.25	0/3238	0.40	0/4404
5	P	0.24	0/3238	0.41	0/4404
5	Q	0.25	0/3238	0.42	0/4404
5	R	0.25	0/3238	0.41	0/4404
6	a	0.23	0/2213	0.39	0/3013
6	b	0.27	0/2255	0.43	0/3070
6	c	0.24	0/2115	0.40	0/2882
6	d	0.24	0/2255	0.41	0/3070

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
6	e	0.22	0/2119	0.37	0/2887
6	f	0.26	0/2255	0.44	1/3070 (0.0%)
6	g	0.24	0/2255	0.41	0/3070
6	h	0.24	0/2213	0.44	1/3013 (0.0%)
6	i	0.26	0/2255	0.42	0/3070
6	j	0.23	0/2238	0.40	0/3047
6	k	0.23	0/2255	0.40	0/3070
6	l	0.25	0/2255	0.46	1/3070 (0.0%)
6	m	0.25	0/2246	0.40	0/3058
6	n	0.24	0/2255	0.40	0/3070
6	o	0.26	0/2255	0.46	1/3070 (0.0%)
6	p	0.26	0/2246	0.43	0/3058
6	q	0.25	0/2213	0.43	1/3013 (0.0%)
6	r	0.24	0/2246	0.40	0/3058
All	All	0.25	0/121310	0.43	6/165025 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	T	0	1
1	U	0	1
2	1	0	3
2	2	0	1
3	4	0	4
3	6	0	3
4	5	0	1
5	A	0	1
5	C	0	1
5	P	0	1
5	Q	0	1
6	a	0	1
6	b	0	1
6	c	0	1
6	d	0	2
6	e	0	1
6	f	0	1
6	g	0	2
6	h	0	1
6	i	0	2

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Mol	Chain	#Chirality outliers	#Planarity outliers
6	j	0	1
6	k	0	1
6	l	0	2
6	m	0	2
6	n	0	1
6	o	0	2
6	p	0	2
6	q	0	2
6	r	0	2
All	All	0	45

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	l	165	CYS	CA-CB-SG	8.26	133.40	114.40
6	o	165	CYS	CA-CB-SG	8.15	133.15	114.40
6	f	165	CYS	CA-CB-SG	8.10	133.02	114.40
6	q	165	CYS	CA-CB-SG	7.67	132.05	114.40
6	h	165	CYS	CA-CB-SG	7.53	131.72	114.40
4	5	156	ALA	N-CA-C	-6.18	98.44	108.52

There are no chirality outliers.

All (45) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	1	480	GLN	Peptide
2	1	481	THR	Peptide
2	1	482	PRO	Peptide
2	2	487	VAL	Peptide
3	4	113	PHE	Peptide
3	4	126	GLY	Peptide
3	4	42	PRO	Peptide
3	4	85	ASN	Peptide
4	5	100	LEU	Peptide
3	6	113	PHE	Peptide
3	6	126	GLY	Peptide
3	6	85	ASN	Peptide
5	A	145	ARG	Peptide
5	C	144	ARG	Peptide
5	P	144	ARG	Peptide

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Mol	Chain	Res	Type	Group
5	Q	144	ARG	Peptide
1	T	33	VAL	Peptide
1	U	36	ASN	Peptide
6	a	239	LEU	Peptide
6	b	239	LEU	Peptide
6	c	239	LEU	Peptide
6	d	239	LEU	Peptide
6	d	96	GLY	Peptide
6	e	239	LEU	Peptide
6	f	239	LEU	Peptide
6	g	239	LEU	Peptide
6	g	76	PHE	Peptide
6	h	239	LEU	Peptide
6	i	134	TYR	Peptide
6	i	239	LEU	Peptide
6	j	239	LEU	Peptide
6	k	239	LEU	Peptide
6	l	239	LEU	Peptide
6	l	96	GLY	Peptide
6	m	134	TYR	Peptide
6	m	239	LEU	Peptide
6	n	239	LEU	Peptide
6	o	239	LEU	Peptide
6	o	289	TRP	Peptide
6	p	134	TYR	Peptide
6	p	239	LEU	Peptide
6	q	239	LEU	Peptide
6	q	96	GLY	Peptide
6	r	134	TYR	Peptide
6	r	239	LEU	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	T	1827	1726	1726	40	0
1	U	1827	1726	1726	37	0
1	V	204	197	197	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	1	3963	3951	3951	65	0
2	2	3963	3951	3951	91	0
2	3	3866	3853	3854	65	0
3	4	1729	1667	1667	35	0
3	6	1729	1667	1667	30	0
4	5	1640	1562	1562	29	0
4	7	1640	1562	1562	29	0
5	A	3168	3127	3127	46	0
5	B	3168	3127	3127	44	0
5	C	3168	3127	3127	56	0
5	D	3168	3127	3127	52	0
5	E	3168	3127	3127	53	0
5	F	3168	3128	3127	42	0
5	G	3168	3127	3127	68	0
5	H	3168	3127	3127	38	0
5	I	3168	3127	3127	54	0
5	J	3168	3128	3127	54	0
5	K	3168	3127	3127	52	0
5	L	3168	3127	3127	48	0
5	M	3168	3128	3127	51	0
5	N	3168	3127	3127	53	0
5	O	3168	3127	3127	40	0
5	P	3168	3127	3127	41	0
5	Q	3168	3127	3127	50	0
5	R	3168	3127	3127	60	0
6	a	2168	2111	2111	30	0
6	b	2209	2148	2148	38	0
6	c	2073	2023	2023	26	0
6	d	2209	2148	2148	37	0
6	e	2077	2026	2026	26	0
6	f	2209	2148	2148	32	0
6	g	2209	2148	2148	40	0
6	h	2168	2111	2111	39	0
6	i	2209	2148	2148	38	0
6	j	2192	2134	2134	42	0
6	k	2209	2148	2148	40	0
6	l	2209	2148	2148	39	0
6	m	2200	2140	2140	27	0
6	n	2209	2148	2148	46	0
6	o	2209	2148	2148	32	0
6	p	2200	2140	2140	36	0
6	q	2168	2111	2111	42	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	r	2200	2140	2140	42	0
7	S	39	37	34	1	0
8	a	14	13	13	0	0
8	b	28	28	26	1	0
8	c	28	28	26	0	0
8	d	28	28	26	1	0
8	e	28	28	26	0	0
8	f	28	28	26	1	0
8	g	28	28	26	0	0
8	h	28	28	26	0	0
8	i	28	28	26	0	0
8	j	28	28	26	1	0
8	k	28	28	26	1	0
8	l	28	28	26	1	0
8	m	28	28	26	0	0
8	n	28	28	26	1	0
8	o	28	28	26	1	0
8	p	28	28	26	2	0
8	q	28	28	26	0	0
8	r	28	28	26	0	0
9	a	4	0	0	0	0
9	b	4	0	0	0	0
9	c	4	0	0	0	0
9	d	4	0	0	0	0
9	e	4	0	0	0	0
9	f	4	0	0	0	0
9	g	4	0	0	0	0
9	h	4	0	0	0	0
9	i	4	0	0	0	0
9	j	4	0	0	0	0
9	k	4	0	0	0	0
9	l	4	0	0	0	0
9	m	4	0	0	0	0
9	n	4	0	0	0	0
9	o	4	0	0	0	0
9	p	4	0	0	0	0
9	q	4	0	0	0	0
9	r	4	0	0	0	0
All	All	119340	116945	116906	1719	0

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

All (1719) 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)
3:4:176:VAL:HG12	3:4:195:VAL:HG12	1.46	0.96
6:e:193:VAL:HG22	6:e:239:LEU:HD12	1.49	0.92
6:a:206:GLY:O	6:a:216:GLU:OE2	1.93	0.86
3:6:142:LYS:NZ	4:7:123:LEU:O	2.09	0.85
2:1:466:ARG:NH2	2:2:247:ALA:O	2.11	0.84
6:f:141:LEU:HD21	6:f:299:ILE:HD11	1.61	0.83
6:g:193:VAL:HG22	6:g:239:LEU:HD12	1.60	0.82
5:C:106:ARG:NH1	5:Q:379:GLU:OE1	2.13	0.82
3:4:136:PRO:O	4:5:127:SER:OG	1.96	0.81
3:6:176:VAL:HG12	3:6:195:VAL:HG12	1.62	0.81
5:G:283:ARG:NH2	5:H:352:GLU:OE2	2.12	0.81
6:p:183:LYS:NZ	6:p:228:ASP:OD1	2.13	0.81
6:f:165:CYS:HB3	6:f:249:CYS:HA	1.63	0.80
5:M:11:LEU:HD21	5:M:37:MET:HE1	1.64	0.80
5:M:216:ARG:NH2	5:M:379:GLU:OE2	2.15	0.80
6:l:256:GLU:OE1	6:l:311:SER:OG	1.98	0.80
5:C:11:LEU:HD12	5:C:85:ILE:HG23	1.65	0.79
5:M:276:ARG:NH2	5:N:366:ASN:OD1	2.16	0.79
5:D:283:ARG:NH1	5:E:352:GLU:OE2	2.16	0.78
5:I:99:GLU:OE2	5:I:384:ARG:NH1	2.16	0.78
5:G:300:MET:O	6:h:308:GLN:NE2	2.16	0.78
6:c:183:LYS:NZ	6:c:228:ASP:OD2	2.15	0.78
6:j:183:LYS:NZ	6:j:228:ASP:OD1	2.16	0.78
5:C:145:ARG:NH2	5:Q:111:PRO:O	2.16	0.78
5:M:236:ARG:NH2	5:O:228:ASP:OD2	2.17	0.78
5:K:216:ARG:NE	5:K:379:GLU:OE2	2.17	0.77
6:g:52:ASN:ND2	6:k:55:LEU:O	2.17	0.77
6:d:186:SER:OG	6:d:244:CYS:SG	2.42	0.77
5:P:130:ASN:OD1	5:P:132:SER:OG	2.03	0.77
6:c:274:ASP:OD2	6:c:276:THR:OG1	2.02	0.76
6:i:186:SER:OG	6:i:244:CYS:SG	2.43	0.76
2:3:495:ARG:NH2	2:3:499:ASN:OD1	2.18	0.76
5:N:262:GLU:OE1	5:N:289:ARG:NH2	2.17	0.76
5:A:132:SER:OG	5:B:16:ASP:OD2	2.01	0.76
5:A:103:GLU:OE2	5:A:350:ARG:NH2	2.18	0.76
5:M:6:SER:OG	5:M:127:ILE:O	2.04	0.76
2:2:567:LEU:HD12	2:3:522:LEU:HD11	1.68	0.76
5:G:228:ASP:OD2	5:H:236:ARG:NH2	2.19	0.76
5:L:216:ARG:NH2	5:L:379:GLU:OE1	2.18	0.76
6:o:183:LYS:NZ	6:o:228:ASP:OD1	2.18	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:g:186:SER:OG	6:g:244:CYS:SG	2.44	0.75
4:7:35:VAL:N	4:7:53:ASN:OD1	2.19	0.75
6:o:81:LEU:HD21	6:o:303:ILE:HD11	1.67	0.75
5:G:246:THR:O	5:I:301:THR:OG1	2.03	0.75
6:h:186:SER:OG	6:h:244:CYS:SG	2.43	0.75
5:Q:7:LEU:HD22	5:Q:127:ILE:HG21	1.67	0.75
5:G:130:ASN:OD1	5:G:132:SER:OG	2.03	0.75
6:h:216:GLU:OE1	6:i:293:TRP:NE1	2.19	0.75
3:4:156:LYS:NZ	3:4:190:SER:OG	2.20	0.75
6:g:291:ARG:NH1	6:i:220:GLU:OE1	2.19	0.75
6:k:183:LYS:NZ	6:k:228:ASP:OD2	2.20	0.75
4:5:35:VAL:N	4:5:53:ASN:OD1	2.20	0.74
5:H:231:ARG:O	5:H:236:ARG:NH1	2.20	0.74
4:5:137:THR:HG22	4:5:185:SER:HA	1.69	0.74
5:B:262:GLU:OE1	5:B:289:ARG:NH1	2.20	0.74
5:F:99:GLU:OE2	5:F:384:ARG:NH1	2.21	0.74
5:K:93:ASP:OD1	5:K:392:ARG:NH1	2.20	0.74
5:F:168:ARG:NH1	5:F:175:ASN:OD1	2.21	0.74
5:A:236:ARG:NH2	5:C:228:ASP:OD2	2.20	0.74
5:J:168:ARG:NH1	5:J:175:ASN:OD1	2.20	0.74
5:M:228:ASP:OD2	5:N:236:ARG:NH2	2.21	0.73
6:p:257:ASN:ND2	6:p:314:SER:O	2.21	0.73
2:1:280:ILE:HD13	2:1:293:ILE:HD13	1.69	0.73
5:D:168:ARG:NH1	5:D:175:ASN:OD1	2.22	0.73
6:f:265:GLY:O	6:f:286:ARG:NH1	2.22	0.73
5:A:255:ARG:NH1	6:c:65:THR:O	2.22	0.73
5:E:114:GLU:OE1	5:E:114:GLU:N	2.22	0.73
1:T:109:GLU:OE2	4:7:65:SER:OG	2.04	0.73
5:H:227:PRO:O	5:I:231:ARG:NH2	2.22	0.73
5:H:51:ILE:HD11	5:H:101:VAL:HG23	1.69	0.73
2:3:354:TYR:O	2:3:422:SER:OG	2.07	0.72
5:A:271:ASN:ND2	5:B:351:GLN:OE1	2.22	0.72
6:p:162:GLU:OE2	6:p:323:TYR:OH	2.06	0.72
5:L:142:GLN:O	5:M:145:ARG:NH1	2.22	0.72
1:U:179:GLU:O	1:U:183:ALA:N	2.22	0.72
3:6:223:ARG:NE	3:6:225:GLU:OE2	2.22	0.72
5:E:99:GLU:OE2	5:E:384:ARG:NH1	2.22	0.72
6:o:165:CYS:HB3	6:o:249:CYS:HA	1.71	0.72
3:4:142:LYS:NZ	4:5:123:LEU:O	2.21	0.72
5:B:76:ASN:ND2	5:D:76:ASN:OD1	2.22	0.72
5:D:229:ALA:O	5:D:233:SER:OG	2.08	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:K:373:ASN:O	5:K:378:ARG:NH1	2.23	0.72
1:T:158:THR:OG1	1:T:179:GLU:OE1	2.02	0.72
5:K:99:GLU:OE1	5:K:384:ARG:NH1	2.22	0.72
6:n:220:GLU:OE2	6:o:291:ARG:NH1	2.23	0.72
5:A:99:GLU:OE1	5:A:384:ARG:NH1	2.23	0.72
2:3:724:LYS:NZ	5:M:272:THR:OG1	2.22	0.72
5:J:163:SER:OG	6:j:61:GLY:O	2.06	0.72
2:1:454:PRO:O	2:1:461:TYR:OH	2.07	0.71
5:O:111:PRO:O	5:O:117:ARG:NH1	2.23	0.71
6:q:285:MET:HE1	6:q:306:ILE:HG23	1.72	0.71
5:P:373:ASN:O	5:P:378:ARG:NH1	2.23	0.71
2:3:278:ASN:ND2	2:3:452:ALA:O	2.23	0.71
5:R:172:MET:HE3	6:r:256:GLU:HG2	1.72	0.71
6:k:51:GLN:N	6:k:51:GLN:OE1	2.23	0.71
5:Q:271:ASN:ND2	5:R:351:GLN:OE1	2.24	0.71
6:j:109:LYS:NZ	6:j:301:ASP:O	2.23	0.71
2:2:315:THR:OG1	2:2:353:ASP:O	2.06	0.70
5:N:103:GLU:OE1	5:N:350:ARG:NH2	2.23	0.70
5:R:370:LEU:HD21	5:R:382:LEU:HD13	1.73	0.70
6:d:206:GLY:O	6:d:216:GLU:OE2	2.08	0.70
5:K:283:ARG:NH2	5:L:352:GLU:OE2	2.23	0.70
2:1:492:ASP:OD2	6:i:234:ASN:ND2	2.24	0.70
2:3:321:GLY:O	2:3:347:ASN:ND2	2.24	0.70
5:C:341:THR:O	5:C:345:ASN:ND2	2.25	0.70
5:P:138:ASN:OD1	5:P:148:THR:OG1	2.10	0.70
6:p:255:ARG:O	6:p:283:ARG:NH1	2.25	0.70
6:q:220:GLU:OE1	6:r:291:ARG:NH1	2.24	0.70
1:U:109:GLU:OE2	4:5:65:SER:OG	2.06	0.70
5:B:137:GLU:OE2	5:C:397:LYS:NZ	2.24	0.70
5:C:141:LEU:O	5:C:145:ARG:HA	1.92	0.70
5:H:130:ASN:OD1	5:H:132:SER:OG	2.08	0.70
5:O:76:ASN:ND2	5:Q:76:ASN:OD1	2.24	0.70
5:A:219:THR:HG22	5:A:220:THR:HG23	1.74	0.70
5:R:192:GLY:O	5:R:319:THR:OG1	2.10	0.70
3:4:34:GLU:O	3:4:99:THR:OG1	2.09	0.69
5:A:131:ASN:N	5:B:16:ASP:OD1	2.25	0.69
5:G:352:GLU:OE2	5:I:283:ARG:NE	2.24	0.69
2:1:736:THR:OG1	2:1:739:GLU:OE1	2.11	0.69
5:B:228:ASP:OD2	5:C:236:ARG:NH2	2.24	0.69
5:P:283:ARG:NH1	5:Q:352:GLU:OE1	2.25	0.69
5:R:67:GLY:O	5:R:77:TYR:OH	2.10	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:1:486:SER:OG	6:i:172:LEU:O	2.07	0.69
5:E:76:ASN:OD1	5:G:76:ASN:ND2	2.25	0.69
6:h:144:TYR:OH	6:h:146:GLN:OE1	2.09	0.69
2:3:291:SER:O	2:3:339:SER:OG	2.11	0.69
5:A:216:ARG:NH2	5:A:374:TYR:O	2.24	0.69
5:G:104:SER:OG	5:G:106:ARG:O	2.10	0.69
5:J:114:GLU:OE1	5:J:114:GLU:N	2.26	0.69
2:2:251:GLU:OE2	2:2:253:ILE:HG23	1.93	0.69
5:O:225:LEU:HD11	5:O:261:VAL:HG21	1.74	0.68
5:J:67:GLY:O	5:J:77:TYR:OH	2.08	0.68
5:O:209:GLU:OE2	5:O:289:ARG:NE	2.26	0.68
6:p:193:VAL:HG22	6:p:239:LEU:HD13	1.76	0.68
5:C:194:ASP:OD2	5:C:197:CYS:N	2.27	0.68
5:G:147:ARG:NH1	5:G:332:GLU:OE1	2.27	0.68
5:I:7:LEU:HD22	5:I:127:ILE:HG21	1.75	0.68
6:o:265:GLY:O	6:o:286:ARG:NH1	2.27	0.68
2:2:494:GLU:OE1	2:2:494:GLU:N	2.24	0.68
6:p:200:THR:OG1	6:p:201:GLN:OE1	2.08	0.68
5:H:187:GLU:OE2	5:H:325:ARG:NH1	2.27	0.68
4:5:77:LEU:HD13	4:5:79:ILE:HD11	1.75	0.68
5:G:7:LEU:HD22	5:G:127:ILE:HG21	1.76	0.68
5:N:57:ARG:NH1	5:N:94:ASN:OD1	2.27	0.68
2:1:375:ASN:OD1	2:1:464:SER:OG	2.11	0.68
6:m:267:ASN:OD1	6:m:286:ARG:NE	2.25	0.68
4:5:87:GLU:OE2	4:5:110:THR:OG1	2.09	0.67
5:D:147:ARG:NH1	5:D:148:THR:O	2.26	0.67
5:E:39:VAL:HG23	5:E:65:LEU:HD21	1.75	0.67
5:L:170:GLN:OE1	5:L:170:GLN:N	2.28	0.67
6:j:281:ILE:HG22	6:j:284:MET:HE3	1.75	0.67
2:1:724:LYS:NZ	5:I:262:GLU:OE1	2.26	0.67
2:3:364:ILE:HD11	2:3:423:LEU:HD21	1.75	0.67
6:d:189:SER:OG	6:d:243:THR:OG1	2.11	0.67
5:K:369:GLU:OE2	5:K:373:ASN:ND2	2.27	0.67
6:n:75:VAL:HG11	6:n:307:VAL:HG11	1.76	0.67
2:2:494:GLU:OE2	6:a:200:THR:OG1	2.13	0.67
2:3:618:LEU:HD22	2:3:716:VAL:HG21	1.76	0.67
5:Q:98:ASP:OD2	5:Q:113:SER:OG	2.04	0.67
5:E:12:LYS:NZ	5:E:394:MET:O	2.28	0.67
6:g:157:ASP:OD2	6:g:161:ASN:ND2	2.28	0.67
5:G:11:LEU:HD11	5:G:88:PHE:HB3	1.76	0.66
5:L:99:GLU:OE1	5:L:384:ARG:NH1	2.29	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:f:238:ASN:OD1	8:f:504:NAG:N2	2.28	0.66
6:g:149:GLU:HG2	6:g:268:ILE:HD12	1.77	0.66
2:2:292:GLU:OE2	2:2:340:ARG:NH1	2.28	0.66
5:G:11:LEU:HD12	5:G:92:ILE:HD12	1.76	0.66
6:o:158:LEU:HD22	6:o:224:LEU:HD11	1.75	0.66
4:5:123:LEU:HD23	4:5:212:VAL:HG13	1.76	0.66
5:F:105:GLN:O	5:F:377:SER:OG	2.14	0.66
5:K:170:GLN:N	5:K:170:GLN:OE1	2.29	0.66
5:Q:7:LEU:CD1	5:Q:37:MET:HE3	2.26	0.66
2:3:301:GLN:OE1	2:3:314:HIS:ND1	2.28	0.66
5:H:100:MET:O	5:H:350:ARG:NH1	2.28	0.66
5:J:7:LEU:HD22	5:J:127:ILE:HG21	1.78	0.66
5:A:216:ARG:NH1	5:A:379:GLU:OE2	2.29	0.66
5:D:240:SER:OG	5:D:245:THR:O	2.11	0.66
2:1:315:THR:OG1	2:1:353:ASP:O	2.07	0.65
5:I:11:LEU:HD12	5:I:85:ILE:HG23	1.78	0.65
5:J:301:THR:OG1	5:K:246:THR:O	2.13	0.65
5:R:157:ILE:HD13	5:R:335:LEU:HD21	1.78	0.65
2:1:743:LEU:HG	2:1:750:MET:HE1	1.78	0.65
5:A:182:LEU:HD13	5:A:231:ARG:NE	2.12	0.65
5:J:111:PRO:O	5:J:117:ARG:NH1	2.28	0.65
5:N:160:TYR:OH	6:n:64:ASP:OD2	2.13	0.65
1:U:97:ASN:ND2	1:U:100:ASP:OD1	2.29	0.65
1:U:144:ARG:NH1	1:U:146:SER:O	2.29	0.65
5:M:194:ASP:OD2	5:M:197:CYS:N	2.29	0.65
5:E:134:GLU:O	5:E:138:ASN:ND2	2.30	0.65
2:3:493:LEU:HD13	6:n:108:ILE:HD11	1.77	0.65
6:l:144:TYR:HA	6:l:152:MET:HE1	1.77	0.65
6:b:51:GLN:OE1	6:f:59:ILE:HG21	1.96	0.65
6:g:164:LEU:HD11	6:g:325:ARG:HG3	1.78	0.65
5:A:7:LEU:HD22	5:A:127:ILE:HG21	1.78	0.65
5:L:177:MET:SD	5:L:206:GLN:NE2	2.70	0.65
5:I:40:THR:HG21	5:I:127:ILE:HD11	1.79	0.64
6:m:126:ASP:O	6:m:129:VAL:HG12	1.96	0.64
6:n:218:VAL:HG11	6:n:227:VAL:HG23	1.78	0.64
5:M:350:ARG:HE	5:M:365:MET:HE3	1.61	0.64
5:N:194:ASP:OD2	5:N:197:CYS:N	2.31	0.64
5:D:228:ASP:OD2	5:E:236:ARG:NH2	2.30	0.64
6:g:285:MET:HE1	6:g:306:ILE:HG23	1.79	0.64
2:2:521:LEU:HD11	2:2:748:PRO:CB	2.28	0.64
3:4:154:LEU:HD12	3:4:191:LEU:O	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:G:109:VAL:HG21	5:K:145:ARG:HD3	1.77	0.64
5:J:165:THR:HG22	6:j:60:THR:HG22	1.79	0.64
5:M:182:LEU:HD13	5:M:231:ARG:NE	2.13	0.64
6:e:206:GLY:O	6:e:216:GLU:OE2	2.16	0.64
3:6:222:LYS:NZ	4:7:129:GLU:OE1	2.26	0.64
5:C:216:ARG:NH2	5:C:379:GLU:OE1	2.30	0.64
3:6:155:VAL:HG11	3:6:163:VAL:HG21	1.78	0.64
5:A:352:GLU:OE2	5:C:283:ARG:NH2	2.31	0.64
5:M:220:THR:N	5:M:283:ARG:O	2.30	0.64
5:R:166:LEU:HD23	5:R:176:LEU:HD22	1.79	0.64
6:b:305:GLN:OE1	6:b:305:GLN:N	2.31	0.64
6:l:274:ASP:OD2	6:l:276:THR:OG1	2.14	0.64
6:n:271:ILE:HD11	6:n:281:ILE:HD12	1.79	0.64
5:O:98:ASP:OD2	5:O:113:SER:OG	2.11	0.64
6:f:183:LYS:NZ	6:f:228:ASP:OD1	2.21	0.64
6:j:166:ASN:ND2	6:j:326:VAL:O	2.31	0.64
2:2:488:THR:OG1	2:2:489:VAL:N	2.26	0.64
5:I:209:GLU:OE1	5:I:289:ARG:NH1	2.31	0.64
6:i:133:LEU:HD12	6:i:133:LEU:O	1.98	0.64
6:i:274:ASP:OD2	6:i:276:THR:OG1	2.14	0.64
6:k:305:GLN:OE1	6:k:305:GLN:N	2.31	0.64
6:p:133:LEU:HD12	6:p:133:LEU:O	1.98	0.64
2:2:389:PRO:O	3:6:32:ASN:ND2	2.31	0.64
3:4:173:THR:O	3:4:176:VAL:HG22	1.98	0.64
5:R:170:GLN:OE1	5:R:170:GLN:N	2.30	0.64
3:6:7:GLU:OE2	3:6:119:GLY:N	2.32	0.63
2:2:476:ASN:ND2	2:2:479:TYR:O	2.30	0.63
5:Q:99:GLU:OE2	5:Q:384:ARG:NH1	2.31	0.63
6:a:160:LEU:HD23	6:a:258:VAL:HG23	1.80	0.63
6:c:209:THR:O	6:c:236:LYS:NZ	2.30	0.63
3:4:7:GLU:OE2	3:4:119:GLY:N	2.31	0.63
5:C:180:MET:HE1	5:C:249:PHE:CZ	2.33	0.63
5:G:216:ARG:NH2	5:G:374:TYR:O	2.31	0.63
5:H:67:GLY:O	5:H:77:TYR:OH	2.12	0.63
6:g:256:GLU:OE1	6:g:314:SER:OG	2.14	0.63
1:T:73:THR:OG1	1:T:199:THR:OG1	2.08	0.63
5:N:11:LEU:HD21	5:N:37:MET:HE1	1.80	0.63
6:h:111:TRP:HZ2	6:h:307:VAL:HG21	1.64	0.63
5:O:254:LEU:HD12	5:O:294:LEU:HD11	1.79	0.63
2:1:292:GLU:OE2	2:1:340:ARG:NH1	2.32	0.63
4:5:7:GLN:NE2	4:5:92:CYS:SG	2.71	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:o:193:VAL:HG12	6:o:239:LEU:HD22	1.81	0.63
1:U:21:GLU:O	1:U:25:ILE:HG22	1.99	0.63
5:F:333:SER:OG	5:F:379:GLU:OE2	2.17	0.63
4:7:137:THR:HG22	4:7:185:SER:HA	1.81	0.63
6:r:274:ASP:OD2	6:r:276:THR:OG1	2.12	0.63
2:3:691:ALA:O	2:3:700:VAL:N	2.32	0.62
5:A:49:GLY:N	5:A:98:ASP:OD2	2.31	0.62
6:k:242:THR:HG23	6:k:243:THR:HG23	1.81	0.62
6:n:274:ASP:OD2	6:n:276:THR:N	2.32	0.62
6:p:305:GLN:OE1	6:p:305:GLN:N	2.30	0.62
6:r:170:ILE:HD11	6:r:246:ILE:HG21	1.80	0.62
6:b:133:LEU:HD12	6:b:133:LEU:O	1.99	0.62
6:r:138:ASN:HB2	6:r:258:VAL:HG12	1.81	0.62
5:M:99:GLU:OE2	5:M:384:ARG:NH1	2.31	0.62
6:e:274:ASP:OD1	6:e:276:THR:OG1	2.17	0.62
2:1:370:LEU:HD12	2:1:468:SER:O	1.99	0.62
2:3:263:GLU:OE2	2:3:368:ARG:NH1	2.32	0.62
5:N:102:ARG:O	5:N:384:ARG:NH2	2.31	0.62
5:G:158:PHE:CZ	5:G:188:ILE:HD11	2.35	0.62
5:O:207:GLN:OE1	5:O:293:GLN:NE2	2.32	0.62
2:3:319:VAL:HG22	2:3:350:VAL:HG12	1.82	0.62
5:E:113:SER:O	5:E:116:LEU:N	2.32	0.62
2:3:309:GLU:OE1	2:3:309:GLU:N	2.32	0.62
2:3:493:LEU:HD23	2:3:494:GLU:N	2.15	0.62
5:O:368:THR:O	5:O:372:THR:HG23	2.00	0.62
3:4:68:ARG:HB2	3:4:87:LEU:HD21	1.81	0.62
5:I:141:LEU:O	5:I:144:ARG:O	2.18	0.62
5:L:225:LEU:CD1	5:L:261:VAL:HG21	2.30	0.62
6:g:105:MET:HE1	6:g:297:TYR:CE1	2.35	0.62
6:j:220:GLU:OE1	6:j:220:GLU:N	2.31	0.62
6:b:126:ASP:O	6:b:129:VAL:HG12	1.99	0.61
5:L:168:ARG:NH1	5:L:175:ASN:OD1	2.33	0.61
5:N:7:LEU:HD22	5:N:37:MET:HE3	1.82	0.61
6:h:124:ILE:HD11	6:h:155:LEU:HD23	1.82	0.61
6:m:233:ILE:CD1	6:n:108:ILE:HD13	2.29	0.61
6:i:193:VAL:HG22	6:i:239:LEU:HD12	1.81	0.61
6:r:105:MET:HE1	6:r:297:TYR:CE2	2.35	0.61
3:4:223:ARG:NE	3:4:225:GLU:OE2	2.33	0.61
5:C:180:MET:HE3	5:C:232:PHE:CD2	2.35	0.61
5:D:182:LEU:HD21	5:D:184:ALA:HB2	1.83	0.61
5:J:170:GLN:OE1	5:J:170:GLN:N	2.34	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:618:LEU:HD21	2:2:711:VAL:CG2	2.31	0.61
2:1:442:ARG:NH2	1:U:205:TYR:OH	2.34	0.61
2:2:713:ASP:OD1	2:2:714:SER:N	2.33	0.61
6:n:149:GLU:N	6:n:149:GLU:OE2	2.34	0.61
3:6:40:GLN:NE2	4:7:40:GLN:OE1	2.33	0.61
4:7:8:PRO:O	4:7:107:THR:HG23	2.01	0.61
5:M:246:THR:O	5:O:301:THR:OG1	2.12	0.61
6:d:305:GLN:OE1	6:d:305:GLN:N	2.33	0.61
1:T:218:TYR:O	1:T:222:GLY:N	2.34	0.61
5:L:4:LEU:HD23	5:L:391:ILE:HG21	1.81	0.61
5:O:103:GLU:OE2	5:O:350:ARG:NH1	2.33	0.61
1:U:168:LYS:HD2	1:U:196:ILE:HD11	1.82	0.60
5:A:141:LEU:O	5:A:145:ARG:HA	2.01	0.60
6:j:105:MET:HE1	6:j:297:TYR:CE1	2.36	0.60
6:n:51:GLN:N	6:n:51:GLN:OE1	2.33	0.60
1:U:35:ILE:HD13	2:2:482:PRO:HB3	1.82	0.60
5:E:254:LEU:HD12	5:E:294:LEU:CD1	2.32	0.60
6:r:316:SER:OG	6:r:318:ASN:O	2.17	0.60
2:2:388:LEU:O	3:6:29:THR:OG1	2.18	0.60
5:A:23:LEU:HD11	5:C:36:GLN:HG2	1.82	0.60
5:E:368:THR:O	5:E:372:THR:HG23	2.00	0.60
5:G:214:LEU:HD11	5:G:285:PHE:CE2	2.36	0.60
5:L:209:GLU:OE1	5:L:289:ARG:NE	2.34	0.60
5:Q:228:ASP:OD2	5:R:236:ARG:NH2	2.34	0.60
6:f:305:GLN:N	6:f:305:GLN:OE1	2.34	0.60
2:3:416:ASP:OD1	2:3:417:PHE:N	2.34	0.60
2:3:618:LEU:HD21	2:3:711:VAL:CG2	2.32	0.60
3:6:173:THR:O	3:6:176:VAL:HG22	2.02	0.60
6:i:242:THR:HG23	6:i:243:THR:HG23	1.82	0.60
6:j:170:ILE:HD12	6:j:237:ILE:HB	1.84	0.60
6:o:239:LEU:HD21	6:o:244:CYS:SG	2.42	0.60
1:T:228:ASN:OD1	2:2:440:ARG:NE	2.35	0.60
2:1:388:LEU:O	3:4:29:THR:OG1	2.18	0.60
2:2:515:GLN:OE1	2:2:515:GLN:N	2.34	0.60
6:j:230:VAL:CG2	6:k:105:MET:HE2	2.31	0.60
6:l:201:GLN:OE1	6:l:201:GLN:N	2.35	0.60
5:J:393:SER:O	5:J:397:LYS:NZ	2.34	0.60
5:R:370:LEU:HD21	5:R:382:LEU:CD1	2.32	0.60
6:i:230:VAL:CG1	6:i:233:ILE:HD12	2.32	0.60
2:3:374:LEU:HD11	2:3:425:PHE:CD1	2.36	0.60
5:B:225:LEU:HD11	5:B:261:VAL:HG21	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:O:182:LEU:HD13	5:O:231:ARG:NE	2.17	0.60
6:h:157:ASP:OD1	6:h:161:ASN:ND2	2.35	0.60
5:P:246:THR:O	5:R:301:THR:OG1	2.07	0.59
6:f:161:ASN:OD1	6:f:251:LYS:NZ	2.24	0.59
2:2:393:TRP:O	2:2:440:ARG:NH1	2.35	0.59
5:G:236:ARG:NH2	5:I:228:ASP:OD2	2.34	0.59
5:L:310:GLN:OE1	5:L:310:GLN:N	2.35	0.59
5:R:45:ASP:O	5:R:118:LYS:NZ	2.35	0.59
6:a:285:MET:HE1	6:a:306:ILE:HG23	1.83	0.59
6:e:305:GLN:OE1	6:e:305:GLN:N	2.35	0.59
5:L:131:ASN:ND2	5:L:140:ASN:OD1	2.35	0.59
5:O:230:GLU:N	5:O:230:GLU:OE2	2.35	0.59
6:k:87:THR:OG1	6:k:120:GLU:OE2	2.08	0.59
1:T:149:ASN:OD1	1:T:150:GLU:N	2.35	0.59
5:K:98:ASP:OD1	5:K:102:ARG:NH1	2.34	0.59
6:m:160:LEU:HD23	6:m:258:VAL:HG23	1.84	0.59
1:U:149:ASN:OD1	1:U:150:GLU:N	2.34	0.59
5:C:170:GLN:NE2	5:C:174:ASP:OD2	2.35	0.59
2:2:495:ARG:HD3	2:2:500:LEU:HD13	1.83	0.59
5:E:219:THR:HG22	5:E:220:THR:HG23	1.85	0.59
1:U:2:ALA:N	2:2:634:ASP:OD1	2.35	0.59
2:2:640:LEU:HD13	2:2:662:ALA:HB2	1.83	0.59
2:3:369:SER:OG	2:3:470:ILE:HD12	2.01	0.59
5:A:111:PRO:O	5:A:117:ARG:NH1	2.36	0.59
5:A:145:ARG:NH1	5:F:142:GLN:O	2.34	0.59
6:l:185:ILE:HD12	6:l:249:CYS:SG	2.42	0.59
2:2:260:LEU:HD13	2:2:480:GLN:HE22	1.67	0.59
2:2:273:ARG:NH2	2:2:462:GLU:OE2	2.35	0.59
5:I:131:ASN:ND2	5:I:140:ASN:OD1	2.36	0.59
6:f:285:MET:HE1	6:f:306:ILE:HG23	1.83	0.59
6:l:159:ILE:HG13	6:l:160:LEU:HD12	1.83	0.59
6:n:144:TYR:OH	6:n:146:GLN:OE1	2.21	0.59
2:3:699:GLU:OE2	5:M:269:ILE:N	2.36	0.59
5:E:14:ALA:HA	5:E:30:LEU:HD21	1.83	0.59
5:G:248:PHE:CD1	5:I:307:LEU:HD11	2.38	0.59
5:P:99:GLU:OE1	5:P:384:ARG:NE	2.32	0.59
6:b:154:GLU:OE2	6:b:226:ILE:N	2.31	0.59
5:D:227:PRO:O	5:E:231:ARG:NH1	2.35	0.59
6:a:193:VAL:HG12	6:a:239:LEU:HD22	1.84	0.59
6:p:290:LYS:HB2	6:r:150:LEU:HD23	1.85	0.59
2:2:618:LEU:HD22	2:2:716:VAL:HG11	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:628:GLU:OE1	2:2:628:GLU:N	2.36	0.58
5:J:236:ARG:NH2	5:L:228:ASP:OD2	2.35	0.58
5:K:142:GLN:HB2	5:K:148:THR:HG21	1.83	0.58
6:l:271:ILE:HD11	6:l:281:ILE:HG13	1.85	0.58
2:1:722:ASP:OD1	2:1:723:PHE:N	2.35	0.58
5:E:134:GLU:OE1	5:E:134:GLU:N	2.36	0.58
5:J:216:ARG:NH1	5:J:379:GLU:OE2	2.36	0.58
6:c:135:CYS:O	6:c:138:ASN:ND2	2.36	0.58
5:B:225:LEU:CD1	5:B:261:VAL:HG21	2.33	0.58
5:F:182:LEU:HD13	5:F:231:ARG:NE	2.18	0.58
2:3:488:THR:O	2:3:491:GLN:N	2.36	0.58
5:B:187:GLU:OE2	5:B:325:ARG:NE	2.36	0.58
5:Q:11:LEU:HD21	5:Q:37:MET:HE1	1.85	0.58
5:F:207:GLN:OE1	5:F:208:PHE:N	2.37	0.58
5:J:1:MET:HE1	5:J:387:THR:HG23	1.86	0.58
3:6:34:GLU:O	3:6:99:THR:OG1	2.09	0.58
5:P:225:LEU:HD11	5:P:261:VAL:HG11	1.84	0.58
6:d:186:SER:OG	6:d:191:CYS:SG	2.61	0.58
6:k:285:MET:HE3	6:k:306:ILE:CD1	2.34	0.58
2:1:646:MET:HE2	2:3:572:SER:HB2	1.86	0.58
1:U:97:ASN:HB3	1:U:196:ILE:HG22	1.86	0.58
2:2:631:SER:O	2:2:634:ASP:N	2.34	0.58
5:H:89:ILE:HD13	5:H:395:LEU:HD13	1.85	0.58
6:j:193:VAL:HG22	6:j:239:LEU:HD13	1.85	0.58
5:J:39:VAL:HG23	5:J:65:LEU:HD11	1.86	0.58
5:M:352:GLU:OE1	5:O:283:ARG:NH2	2.36	0.58
6:b:55:LEU:HD11	6:f:319:SER:HA	1.85	0.58
6:q:74:GLU:O	6:q:78:THR:HG22	2.04	0.58
6:q:159:ILE:O	6:q:258:VAL:HG21	2.04	0.58
2:3:364:ILE:HD11	2:3:423:LEU:CD2	2.34	0.58
4:7:69:ASP:OD1	4:7:72:SER:OG	2.17	0.58
5:D:182:LEU:HD13	5:D:231:ARG:NE	2.19	0.58
6:b:306:ILE:HG22	6:b:310:MET:HE2	1.86	0.58
5:L:158:PHE:CZ	5:L:188:ILE:HD11	2.39	0.57
5:M:4:LEU:HD22	5:M:116:LEU:HD13	1.86	0.57
6:d:239:LEU:HD21	6:d:244:CYS:SG	2.44	0.57
1:T:135:ASN:O	1:T:135:ASN:ND2	2.37	0.57
5:K:307:LEU:HD11	5:L:248:PHE:CD1	2.38	0.57
6:d:74:GLU:O	6:d:78:THR:HG22	2.04	0.57
2:1:611:THR:HG23	2:1:683:ILE:HD12	1.86	0.57
5:E:254:LEU:HD12	5:E:294:LEU:HD11	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:d:183:LYS:NZ	6:d:228:ASP:OD2	2.23	0.57
6:h:106:PHE:CE2	6:h:300:VAL:HG22	2.40	0.57
6:l:165:CYS:HB3	6:l:249:CYS:HA	1.85	0.57
6:o:160:LEU:HD23	6:o:258:VAL:HG23	1.86	0.57
1:T:87:ASN:ND2	3:6:104:GLN:OE1	2.36	0.57
2:1:356:ASP:OD1	2:1:357:ASP:N	2.37	0.57
5:A:261:VAL:HG23	5:A:290:LEU:HD23	1.86	0.57
6:l:294:GLN:O	6:l:298:THR:HG23	2.04	0.57
2:2:432:ASP:OD1	2:2:433:GLU:N	2.37	0.57
2:2:699:GLU:O	5:C:268:GLN:NE2	2.38	0.57
5:M:283:ARG:NH1	5:N:352:GLU:OE1	2.37	0.57
5:R:112:GLN:O	5:R:113:SER:OG	2.22	0.57
6:e:81:LEU:HD11	6:e:303:ILE:HD11	1.84	0.57
6:q:161:ASN:OD1	6:q:251:LYS:NZ	2.21	0.57
5:R:347:THR:HG22	5:R:365:MET:HB2	1.86	0.57
6:l:51:GLN:OE1	6:l:51:GLN:N	2.37	0.57
1:T:9:LEU:HD23	2:1:526:MET:HE1	1.87	0.57
6:b:78:THR:HG21	6:b:112:PRO:HG2	1.87	0.57
6:g:168:MET:HB2	6:g:246:ILE:HG23	1.87	0.57
6:q:184:TRP:CZ3	6:q:237:ILE:HD11	2.39	0.57
5:B:170:GLN:NE2	5:B:174:ASP:OD2	2.37	0.57
5:D:395:LEU:HB2	5:D:396:ILE:HD12	1.86	0.57
6:l:135:CYS:O	6:l:138:ASN:ND2	2.38	0.57
2:1:515:GLN:NE2	2:3:570:ALA:O	2.37	0.57
3:6:49:VAL:HG13	3:6:65:VAL:HG11	1.86	0.57
5:G:264:LEU:HD23	5:G:269:ILE:HA	1.87	0.57
5:H:113:SER:O	5:H:116:LEU:N	2.38	0.57
5:I:259:VAL:CG2	5:I:322:LEU:HD11	2.34	0.57
5:L:173:HIS:HA	5:L:176:LEU:HD11	1.86	0.57
5:Q:113:SER:O	5:Q:116:LEU:N	2.37	0.57
6:b:218:VAL:HG21	6:b:227:VAL:HG21	1.85	0.57
6:h:165:CYS:HB3	6:h:249:CYS:HA	1.87	0.57
6:k:263:VAL:HG12	6:k:289:TRP:CD1	2.40	0.57
2:1:432:ASP:OD1	2:1:433:GLU:N	2.37	0.56
3:4:37:TRP:CD1	3:4:82:LEU:HD11	2.40	0.56
6:a:81:LEU:HD21	6:a:303:ILE:HD11	1.87	0.56
6:g:128:SER:OG	6:g:155:LEU:HD12	2.05	0.56
6:k:74:GLU:O	6:k:78:THR:OG1	2.10	0.56
6:m:285:MET:HE2	6:m:306:ILE:HG12	1.87	0.56
1:V:2:ALA:N	2:3:522:LEU:O	2.38	0.56
5:L:4:LEU:CD2	5:L:391:ILE:HG21	2.34	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:N:321:GLY:O	5:N:322:LEU:HD12	2.05	0.56
6:a:239:LEU:HD21	6:a:244:CYS:SG	2.45	0.56
1:U:93:TYR:CD2	1:U:198:ILE:HD11	2.40	0.56
5:I:168:ARG:NH1	5:I:175:ASN:OD1	2.38	0.56
1:U:218:TYR:O	1:U:222:GLY:N	2.38	0.56
3:6:165:VAL:HG22	3:6:211:VAL:HG22	1.85	0.56
5:C:106:ARG:NE	5:Q:380:ASP:OD1	2.34	0.56
5:C:180:MET:HE1	5:C:249:PHE:HZ	1.70	0.56
5:D:219:THR:HG22	5:D:220:THR:HG23	1.86	0.56
5:J:151:VAL:HG22	5:J:330:VAL:HG22	1.87	0.56
5:M:113:SER:O	5:M:116:LEU:N	2.38	0.56
6:j:81:LEU:CD1	6:j:303:ILE:HD11	2.36	0.56
6:k:106:PHE:CD1	6:k:116:VAL:HG21	2.41	0.56
3:6:157:ASP:CG	3:6:188:LEU:HD13	2.30	0.56
5:F:180:MET:HE2	6:f:63:MET:HE1	1.87	0.56
5:L:225:LEU:HD11	5:L:261:VAL:HG21	1.86	0.56
5:L:252:ILE:HD11	5:L:316:HIS:ND1	2.20	0.56
6:p:185:ILE:HD12	6:p:249:CYS:SG	2.46	0.56
6:q:185:ILE:HD12	6:q:249:CYS:SG	2.46	0.56
2:1:611:THR:CG2	2:1:683:ILE:HD12	2.35	0.56
5:G:109:VAL:HG23	5:G:380:ASP:HB3	1.87	0.56
6:g:161:ASN:O	6:g:255:ARG:NH1	2.39	0.56
5:A:109:VAL:HG21	5:F:145:ARG:HD3	1.87	0.56
5:A:257:ASN:ND2	5:A:295:MET:HE3	2.20	0.56
5:J:225:LEU:CD1	5:J:261:VAL:HG21	2.36	0.56
5:M:239:ASN:OD1	6:m:65:THR:OG1	2.23	0.56
6:h:102:LEU:HD23	6:h:118:PHE:CE1	2.41	0.56
6:j:139:LEU:HD22	6:j:261:ILE:HD11	1.88	0.56
6:r:106:PHE:CG	6:r:116:VAL:HG21	2.40	0.56
5:A:283:ARG:NH1	5:B:352:GLU:OE1	2.39	0.56
5:P:187:GLU:OE2	5:P:325:ARG:NE	2.39	0.56
5:Q:170:GLN:NE2	5:Q:175:ASN:O	2.37	0.56
5:F:166:LEU:HD22	5:F:176:LEU:HD22	1.87	0.56
5:I:141:LEU:O	5:I:144:ARG:C	2.48	0.56
5:O:113:SER:O	5:O:116:LEU:N	2.39	0.56
5:R:11:LEU:HD11	5:R:88:PHE:HB2	1.87	0.56
6:a:193:VAL:CG1	6:a:239:LEU:HD22	2.36	0.56
6:i:169:ASP:OD2	6:i:172:LEU:HD12	2.06	0.56
1:T:3:SER:OG	1:T:7:ARG:NH2	2.39	0.56
6:n:305:GLN:OE1	6:n:305:GLN:N	2.38	0.56
6:p:262:GLN:NE2	6:p:266:ALA:O	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:4:63:ASP:OD2	3:4:64:SER:N	2.39	0.55
4:5:19:VAL:HG12	4:5:82:LEU:HD21	1.88	0.55
5:C:81:ALA:O	5:C:85:ILE:HD12	2.06	0.55
5:L:103:GLU:OE1	5:L:350:ARG:NH2	2.39	0.55
5:R:103:GLU:OE2	5:R:350:ARG:NH2	2.40	0.55
6:i:161:ASN:OD1	6:i:251:LYS:NZ	2.22	0.55
2:2:295:TYR:N	2:2:342:GLU:OE1	2.36	0.55
5:I:81:ALA:O	5:I:85:ILE:HD12	2.05	0.55
5:N:134:GLU:OE2	5:O:397:LYS:NZ	2.40	0.55
6:q:105:MET:HE1	6:q:297:TYR:CE1	2.42	0.55
5:F:295:MET:HE3	6:d:67:TYR:HD2	1.71	0.55
5:G:99:GLU:OE2	5:G:113:SER:OG	2.25	0.55
6:j:146:GLN:O	6:j:147:SER:OG	2.21	0.55
1:T:168:LYS:HD2	1:T:196:ILE:HD11	1.88	0.55
4:5:5:LEU:HD11	4:5:30:ILE:HD11	1.87	0.55
5:M:273:TYR:OH	5:M:281:VAL:O	2.18	0.55
5:Q:216:ARG:NE	5:Q:379:GLU:OE2	2.33	0.55
6:j:144:TYR:HA	6:j:152:MET:HE1	1.87	0.55
2:2:695:ASN:OD1	2:2:696:THR:N	2.39	0.55
5:C:158:PHE:CE2	5:C:188:ILE:HD11	2.42	0.55
5:L:251:PRO:O	5:L:252:ILE:HD13	2.06	0.55
3:6:109:TRP:CZ3	3:6:111:ALA:HB2	2.42	0.55
5:D:214:LEU:HD23	5:D:218:LEU:HD13	1.89	0.55
6:b:262:GLN:NE2	6:b:265:GLY:O	2.40	0.55
6:b:316:SER:OG	6:b:318:ASN:O	2.24	0.55
2:2:521:LEU:HD11	2:2:748:PRO:HB3	1.89	0.55
5:B:254:LEU:HD23	5:C:237:VAL:CG2	2.36	0.55
5:C:182:LEU:HD22	5:C:231:ARG:NE	2.21	0.55
6:a:168:MET:HE2	6:a:170:ILE:HD11	1.88	0.55
6:m:138:ASN:HB2	6:m:258:VAL:HG12	1.88	0.55
2:1:666:PHE:O	2:1:670:ARG:NH2	2.39	0.55
2:3:515:GLN:N	2:3:515:GLN:OE1	2.39	0.55
3:4:109:TRP:CZ3	3:4:111:ALA:HB2	2.42	0.55
5:C:262:GLU:OE1	5:C:262:GLU:N	2.40	0.55
5:G:11:LEU:HD13	5:G:89:ILE:HG13	1.87	0.55
6:g:256:GLU:OE1	6:g:256:GLU:N	2.39	0.55
6:o:262:GLN:NE2	6:o:266:ALA:O	2.38	0.55
2:2:488:THR:O	2:2:489:VAL:HG23	2.07	0.55
2:3:295:TYR:OH	2:3:449:LEU:O	2.25	0.55
5:G:295:MET:HE3	6:h:67:TYR:HD2	1.72	0.55
5:N:180:MET:HE2	6:n:63:MET:CE	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:R:207:GLN:NE2	5:R:291:SER:OG	2.40	0.55
5:J:239:ASN:OD1	6:j:65:THR:OG1	2.23	0.54
5:J:283:ARG:NE	5:K:352:GLU:OE1	2.40	0.54
5:O:99:GLU:OE2	5:O:384:ARG:NH1	2.39	0.54
6:d:144:TYR:HA	6:d:152:MET:HE1	1.88	0.54
6:d:191:CYS:O	6:d:221:ASN:N	2.39	0.54
6:h:111:TRP:CZ2	6:h:307:VAL:HG21	2.41	0.54
6:q:193:VAL:HG22	6:q:219:ALA:HB3	1.89	0.54
1:T:84:ILE:HD13	1:T:200:ILE:HD12	1.90	0.54
5:J:370:LEU:HD21	5:J:382:LEU:HD13	1.89	0.54
5:M:7:LEU:HD22	5:M:127:ILE:HG21	1.88	0.54
6:d:193:VAL:HG22	6:d:239:LEU:HD22	1.88	0.54
2:2:722:ASP:OD1	2:2:765:ARG:NH1	2.41	0.54
2:2:369:SER:HB3	2:2:470:ILE:HD12	1.88	0.54
5:J:228:ASP:OD2	5:K:236:ARG:NH1	2.38	0.54
6:d:285:MET:HE1	6:d:306:ILE:HG23	1.89	0.54
6:l:81:LEU:HD11	6:l:303:ILE:HD11	1.87	0.54
6:p:202:THR:O	6:p:202:THR:HG22	2.08	0.54
1:U:3:SER:N	2:2:634:ASP:OD1	2.38	0.54
2:3:478:ASP:OD1	2:3:479:TYR:N	2.40	0.54
5:C:334:VAL:O	5:C:335:LEU:HD12	2.07	0.54
5:C:380:ASP:OD2	5:Q:106:ARG:NE	2.40	0.54
5:H:307:LEU:HD11	5:I:248:PHE:CD1	2.43	0.54
6:m:233:ILE:HD11	6:n:108:ILE:HD13	1.89	0.54
6:r:202:THR:O	6:r:203:LEU:HD12	2.08	0.54
5:G:310:GLN:OE1	5:G:310:GLN:N	2.41	0.54
5:Q:307:LEU:HD11	5:R:248:PHE:CD2	2.42	0.54
5:R:134:GLU:O	5:R:138:ASN:ND2	2.41	0.54
5:F:262:GLU:OE1	5:F:289:ARG:NH1	2.40	0.54
6:i:170:ILE:HD11	6:i:246:ILE:HG21	1.90	0.54
6:i:202:THR:HG22	6:i:202:THR:O	2.07	0.54
6:r:126:ASP:O	6:r:129:VAL:HG12	2.08	0.54
5:D:182:LEU:HD13	5:D:231:ARG:HE	1.71	0.54
5:R:97:MET:HE3	5:R:353:TYR:CE1	2.42	0.54
6:i:166:ASN:ND2	6:i:326:VAL:O	2.39	0.54
6:o:157:ASP:OD2	6:o:161:ASN:ND2	2.41	0.54
1:T:22:ILE:HD11	1:U:22:ILE:CD1	2.38	0.54
2:2:529:MET:HE1	2:2:640:LEU:HD23	1.89	0.54
5:I:216:ARG:NH2	5:I:379:GLU:OE1	2.39	0.54
6:b:185:ILE:HD12	6:b:249:CYS:SG	2.48	0.54
6:g:255:ARG:NH2	6:g:321:ALA:O	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:l:255:ARG:O	6:l:283:ARG:NH1	2.38	0.54
6:q:160:LEU:HD23	6:q:258:VAL:HG23	1.89	0.54
2:2:259:SER:C	2:2:260:LEU:HD12	2.33	0.54
2:3:375:ASN:ND2	2:3:464:SER:OG	2.41	0.54
5:F:295:MET:HE3	6:d:67:TYR:CD2	2.42	0.54
5:N:147:ARG:NH2	5:O:397:LYS:OXT	2.41	0.54
6:j:146:GLN:OE1	6:j:146:GLN:N	2.41	0.54
6:r:194:LYS:HG2	6:r:217:THR:HG22	1.89	0.54
2:3:645:ASP:OD1	2:3:646:MET:N	2.41	0.53
5:P:158:PHE:CE1	5:P:188:ILE:HD11	2.43	0.53
5:Q:1:MET:HE1	5:Q:387:THR:HG21	1.90	0.53
6:d:238:ASN:OD1	8:d:402:NAG:N2	2.41	0.53
6:k:139:LEU:HD22	6:k:261:ILE:HD12	1.90	0.53
6:p:294:GLN:O	6:p:298:THR:HG23	2.08	0.53
5:D:342:LEU:HD12	5:D:389:ALA:HB3	1.90	0.53
2:2:356:ASP:OD1	2:2:357:ASP:N	2.41	0.53
4:7:7:GLN:NE2	4:7:92:CYS:SG	2.81	0.53
5:A:167:ASN:O	6:f:51:GLN:N	2.41	0.53
5:N:214:LEU:HD13	5:N:218:LEU:HD13	1.90	0.53
6:l:265:GLY:O	6:l:286:ARG:NH1	2.41	0.53
2:2:438:ILE:HG23	2:2:441:THR:HB	1.90	0.53
6:b:146:GLN:O	6:b:147:SER:OG	2.27	0.53
5:P:23:LEU:HD11	5:R:36:GLN:OE1	2.08	0.53
6:d:139:LEU:HD21	6:d:306:ILE:HG21	1.91	0.53
6:i:230:VAL:HG13	6:i:233:ILE:HD12	1.90	0.53
6:l:229:VAL:HG21	6:l:235:HIS:CD2	2.43	0.53
1:U:25:ILE:HG21	1:V:22:ILE:HG23	1.89	0.53
2:3:327:TYR:O	2:3:338:ILE:N	2.41	0.53
5:C:99:GLU:OE2	5:C:384:ARG:NE	2.39	0.53
5:G:1:MET:HE1	5:G:388:VAL:CG2	2.38	0.53
5:O:239:ASN:ND2	6:o:65:THR:O	2.42	0.53
6:a:296:PHE:O	6:a:300:VAL:HG23	2.08	0.53
1:U:138:LYS:HD3	1:U:158:THR:HG22	1.90	0.53
5:J:338:ALA:O	5:L:153:HIS:ND1	2.41	0.53
5:Q:194:ASP:OD2	5:Q:197:CYS:N	2.42	0.53
6:b:106:PHE:CZ	6:b:300:VAL:HG22	2.43	0.53
6:g:81:LEU:HD12	6:g:137:TYR:O	2.08	0.53
6:k:197:PRO:O	6:k:198:LEU:HD23	2.08	0.53
6:l:111:TRP:HZ2	6:l:307:VAL:HG21	1.74	0.53
6:p:180:GLU:OE1	6:p:180:GLU:N	2.39	0.53
6:q:138:ASN:HB2	6:q:258:VAL:HG12	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:3:378:ARG:NH1	6:a:210:THR:OG1	2.42	0.53
3:4:72:SER:OG	3:4:81:TYR:O	2.26	0.53
5:K:219:THR:HG22	5:K:220:THR:HG23	1.90	0.53
5:R:310:GLN:N	5:R:310:GLN:OE1	2.42	0.53
6:e:201:GLN:O	6:e:202:THR:OG1	2.26	0.53
6:k:238:ASN:OD1	8:k:402:NAG:N2	2.42	0.53
6:k:285:MET:HE3	6:k:306:ILE:HD12	1.91	0.53
6:m:239:LEU:HD23	6:m:239:LEU:O	2.09	0.53
5:Q:67:GLY:O	5:Q:77:TYR:OH	2.24	0.53
6:c:102:LEU:HD23	6:c:105:MET:HE2	1.91	0.53
6:g:109:LYS:NZ	6:g:301:ASP:O	2.41	0.53
4:7:64:PHE:CD1	4:7:79:ILE:HD12	2.44	0.53
4:7:77:LEU:HD13	4:7:79:ILE:HD11	1.90	0.53
5:K:300:MET:HE1	5:K:308:PHE:HD2	1.74	0.53
5:M:271:ASN:ND2	5:N:351:GLN:OE1	2.40	0.53
5:M:307:LEU:HD11	5:N:248:PHE:CD1	2.44	0.53
5:Q:209:GLU:OE2	5:Q:211:ILE:HD11	2.09	0.53
6:g:305:GLN:OE1	6:g:305:GLN:N	2.42	0.53
6:l:165:CYS:CB	6:l:249:CYS:HA	2.38	0.53
6:p:267:ASN:OD1	6:p:286:ARG:NE	2.41	0.53
6:r:160:LEU:HD23	6:r:258:VAL:HG23	1.90	0.53
2:1:646:MET:HE2	2:3:572:SER:CB	2.38	0.52
1:U:168:LYS:HG3	1:U:193:LEU:HD21	1.91	0.52
3:4:225:GLU:OE1	3:4:225:GLU:N	2.43	0.52
5:B:70:LEU:C	5:B:71:LEU:HD12	2.34	0.52
5:N:299:ASN:ND2	5:O:244:ALA:O	2.41	0.52
6:a:81:LEU:HD11	6:a:139:LEU:HD13	1.90	0.52
6:b:85:TYR:OH	6:b:99:LYS:NZ	2.33	0.52
6:m:124:ILE:HG21	6:m:148:LEU:HD23	1.91	0.52
6:n:97:GLU:N	6:n:97:GLU:OE2	2.42	0.52
5:E:209:GLU:OE2	5:E:289:ARG:NH1	2.41	0.52
5:G:207:GLN:OE1	5:G:293:GLN:NE2	2.41	0.52
5:I:40:THR:CG2	5:I:127:ILE:HD11	2.39	0.52
5:K:29:ASP:OD2	5:K:30:LEU:N	2.43	0.52
8:b:401:NAG:O7	8:b:401:NAG:O3	2.19	0.52
2:1:676:LYS:HE3	2:1:710:LEU:HD22	1.91	0.52
5:R:225:LEU:HD22	5:R:261:VAL:HG21	1.92	0.52
6:b:193:VAL:HG22	6:b:239:LEU:HD13	1.89	0.52
6:j:142:MET:HE3	6:j:152:MET:CG	2.39	0.52
6:k:105:MET:HE1	6:k:297:TYR:CE1	2.44	0.52
5:J:225:LEU:HD11	5:J:261:VAL:HG21	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:e:106:PHE:CZ	6:e:300:VAL:HG22	2.45	0.52
6:g:127:PHE:HD1	6:g:155:LEU:HD11	1.75	0.52
6:j:106:PHE:CE1	6:j:300:VAL:HG22	2.45	0.52
6:p:169:ASP:N	6:p:175:TYR:OH	2.43	0.52
6:q:165:CYS:CB	6:q:249:CYS:HA	2.40	0.52
1:T:2:ALA:CB	2:1:524:LEU:HD13	2.39	0.52
5:H:93:ASP:OD2	5:H:392:ARG:NE	2.42	0.52
5:O:1:MET:HE1	5:O:387:THR:OG1	2.10	0.52
5:R:104:SER:OG	5:R:106:ARG:O	2.28	0.52
6:n:220:GLU:OE1	6:n:220:GLU:N	2.43	0.52
3:6:179:PHE:CE1	4:7:168:THR:HG21	2.45	0.52
4:7:34:TYR:O	4:7:94:SER:OG	2.20	0.52
5:O:7:LEU:HD22	5:O:127:ILE:HG21	1.92	0.52
6:c:218:VAL:HG11	6:c:227:VAL:HG23	1.90	0.52
6:e:186:SER:HB2	6:e:239:LEU:HD11	1.92	0.52
5:G:209:GLU:OE2	5:G:289:ARG:NH1	2.41	0.52
6:b:274:ASP:OD2	6:b:276:THR:N	2.43	0.52
6:p:69:ASN:OD1	8:p:401:NAG:N2	2.42	0.52
1:T:64:ILE:HG21	2:2:442:ARG:HB3	1.92	0.52
5:G:142:GLN:OE1	5:G:383:GLN:NE2	2.42	0.52
5:G:172:MET:HE3	6:g:256:GLU:OE2	2.10	0.52
5:J:141:LEU:HB3	5:J:148:THR:HG22	1.91	0.52
5:L:122:ILE:HD12	5:L:125:LYS:HD3	1.91	0.52
6:a:152:MET:HE3	6:a:269:LEU:HD11	1.90	0.52
6:h:165:CYS:CB	6:h:249:CYS:HA	2.39	0.52
6:q:239:LEU:O	6:q:239:LEU:HD23	2.09	0.52
6:r:139:LEU:HD21	6:r:306:ILE:HG21	1.91	0.52
3:4:49:VAL:HG13	3:4:65:VAL:HG11	1.92	0.52
5:A:128:ASN:O	5:B:22:THR:HG22	2.10	0.52
5:A:154:LYS:N	5:A:327:GLU:O	2.40	0.52
5:H:240:SER:OG	5:H:241:ALA:N	2.42	0.52
5:M:23:LEU:HD11	5:O:36:GLN:HG2	1.92	0.52
6:b:150:LEU:HD23	6:c:290:LYS:CG	2.39	0.52
1:U:35:ILE:HD12	2:3:325:PHE:HE2	1.74	0.52
4:5:64:PHE:CD1	4:5:79:ILE:HD12	2.45	0.52
5:I:154:LYS:N	5:I:327:GLU:O	2.42	0.52
5:J:262:GLU:N	5:J:262:GLU:OE1	2.43	0.52
6:a:326:VAL:HG12	6:q:320:ALA:HA	1.92	0.52
6:j:124:ILE:HD11	6:j:155:LEU:HD23	1.92	0.52
2:3:767:GLU:O	2:3:771:LEU:HD23	2.10	0.51
5:J:361:PHE:CD2	5:J:365:MET:HE3	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:K:227:PRO:O	5:L:231:ARG:NH2	2.42	0.51
6:d:81:LEU:HD21	6:d:303:ILE:HD11	1.92	0.51
1:U:100:ASP:OD1	1:U:100:ASP:N	2.42	0.51
3:6:71:ILE:HG22	3:6:82:LEU:CD2	2.40	0.51
3:6:71:ILE:HG22	3:6:82:LEU:HD22	1.92	0.51
5:E:262:GLU:OE1	5:E:289:ARG:NH2	2.43	0.51
6:e:239:LEU:HD21	6:e:244:CYS:SG	2.51	0.51
6:h:262:GLN:OE1	6:h:286:ARG:NH1	2.43	0.51
6:i:295:VAL:O	6:i:298:THR:OG1	2.26	0.51
6:n:105:MET:HE1	6:n:297:TYR:CE1	2.45	0.51
6:r:160:LEU:O	6:r:283:ARG:NH2	2.43	0.51
2:2:431:VAL:HG21	2:2:447:TYR:HB3	1.92	0.51
2:2:517:ILE:O	2:2:521:LEU:HD23	2.11	0.51
5:I:360:VAL:O	5:I:378:ARG:NH2	2.43	0.51
5:M:259:VAL:CG2	5:M:322:LEU:HD11	2.40	0.51
6:d:142:MET:HB3	6:d:152:MET:HE2	1.93	0.51
6:h:305:GLN:O	6:h:309:VAL:HG23	2.10	0.51
6:i:281:ILE:HG22	6:i:283:ARG:H	1.75	0.51
5:J:360:VAL:HG21	5:J:381:ASN:HD22	1.75	0.51
6:n:274:ASP:OD2	6:n:276:THR:OG1	2.21	0.51
1:T:2:ALA:N	2:1:634:ASP:OD1	2.44	0.51
2:1:490:ARG:NE	2:1:491:GLN:O	2.43	0.51
5:G:103:GLU:OE1	5:G:350:ARG:NH1	2.41	0.51
6:a:71:THR:O	6:a:71:THR:HG22	2.11	0.51
6:k:146:GLN:O	6:k:147:SER:OG	2.25	0.51
5:G:1:MET:HE1	5:G:388:VAL:HG22	1.93	0.51
5:M:128:ASN:O	5:N:22:THR:HG22	2.11	0.51
5:R:54:LEU:HD13	5:R:353:TYR:CE1	2.46	0.51
6:e:185:ILE:HD12	6:e:249:CYS:SG	2.50	0.51
6:h:276:THR:HG21	6:i:305:GLN:HB3	1.93	0.51
6:j:238:ASN:OD1	8:j:402:NAG:N2	2.44	0.51
1:T:133:ASP:OD1	1:T:134:SER:N	2.44	0.51
3:4:71:ILE:HG22	3:4:82:LEU:CD2	2.41	0.51
5:A:245:THR:HG22	5:A:246:THR:H	1.76	0.51
5:N:180:MET:HE2	6:n:63:MET:HE2	1.92	0.51
5:R:369:GLU:OE1	5:R:369:GLU:N	2.44	0.51
6:o:238:ASN:OD1	8:o:504:NAG:N2	2.43	0.51
6:q:200:THR:OG1	6:q:201:GLN:OE1	2.29	0.51
1:T:93:TYR:CD1	1:T:198:ILE:HD11	2.46	0.51
1:U:7:ARG:NH2	2:2:626:GLN:OE1	2.44	0.51
5:M:3:VAL:O	5:M:7:LEU:HD23	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:N:230:GLU:O	5:N:233:SER:OG	2.25	0.51
6:e:230:VAL:HG11	6:e:233:ILE:HD12	1.91	0.51
6:f:295:VAL:O	6:f:298:THR:OG1	2.22	0.51
6:i:126:ASP:O	6:i:129:VAL:HG12	2.10	0.51
4:7:19:VAL:HG12	4:7:82:LEU:HD21	1.92	0.51
5:F:332:GLU:N	5:F:332:GLU:OE2	2.44	0.51
5:G:145:ARG:HH12	5:K:109:VAL:HG11	1.75	0.51
5:L:321:GLY:O	5:L:322:LEU:HD12	2.11	0.51
5:P:223:ILE:HG12	5:P:326:ILE:HD13	1.93	0.51
6:n:124:ILE:HD11	6:n:155:LEU:HD23	1.92	0.51
6:n:126:ASP:O	6:n:129:VAL:HG12	2.11	0.51
4:5:8:PRO:O	4:5:107:THR:HG23	2.11	0.51
5:B:103:GLU:OE2	5:B:350:ARG:NH2	2.44	0.51
5:M:166:LEU:HD23	5:M:169:SER:HB2	1.93	0.51
5:N:165:THR:HG23	5:N:181:TRP:CH2	2.46	0.51
5:R:172:MET:HE2	5:R:172:MET:HA	1.92	0.51
6:f:169:ASP:N	6:f:175:TYR:OH	2.41	0.51
6:h:274:ASP:OD2	6:h:276:THR:N	2.42	0.51
6:l:141:LEU:HD11	6:l:299:ILE:HD11	1.91	0.51
6:l:193:VAL:HG22	6:l:239:LEU:HD12	1.92	0.51
6:m:302:TYR:OH	6:o:183:LYS:NZ	2.44	0.51
6:r:83:LEU:CD2	6:r:139:LEU:HD12	2.41	0.51
5:H:51:ILE:HD11	5:H:101:VAL:CG2	2.40	0.50
6:a:261:ILE:CG1	6:a:285:MET:HE3	2.40	0.50
6:i:229:VAL:HG11	6:i:235:HIS:NE2	2.26	0.50
6:p:255:ARG:NH2	6:p:321:ALA:O	2.44	0.50
1:T:193:LEU:O	1:T:196:ILE:HG23	2.12	0.50
2:3:491:GLN:OE1	6:n:104:GLN:NE2	2.44	0.50
5:A:214:LEU:HD11	5:A:285:PHE:CE2	2.47	0.50
5:F:165:THR:HG23	5:F:181:TRP:CH2	2.47	0.50
5:K:228:ASP:OD2	5:L:236:ARG:NH2	2.41	0.50
5:O:14:ALA:HB1	5:O:85:ILE:HD11	1.93	0.50
5:R:230:GLU:O	5:R:233:SER:OG	2.29	0.50
6:r:187:MET:HE1	6:r:322:PHE:HE2	1.77	0.50
5:A:307:LEU:HD11	5:B:248:PHE:CE1	2.45	0.50
6:e:130:ASP:OD2	6:e:187:MET:HE3	2.11	0.50
6:j:85:TYR:OH	6:j:99:LYS:NZ	2.44	0.50
2:3:703:ASP:OD1	2:3:703:ASP:N	2.44	0.50
3:4:98:ALA:HB1	3:4:115:PRO:O	2.12	0.50
3:6:141:SER:HA	3:6:144:THR:HG22	1.92	0.50
5:E:334:VAL:O	5:E:335:LEU:HD12	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:K:158:PHE:CE2	5:K:214:LEU:HD21	2.46	0.50
5:N:11:LEU:CD2	5:N:37:MET:HE1	2.41	0.50
6:g:208:GLN:N	6:g:208:GLN:OE1	2.45	0.50
6:o:165:CYS:CB	6:o:249:CYS:HA	2.40	0.50
1:V:13:SER:O	1:V:16:VAL:HG22	2.11	0.50
2:3:417:PHE:CE2	2:3:418:VAL:HG22	2.47	0.50
2:3:616:LYS:NZ	2:3:685:THR:O	2.45	0.50
5:B:283:ARG:NE	5:C:352:GLU:OE2	2.44	0.50
5:G:165:THR:HG23	5:G:181:TRP:CH2	2.46	0.50
5:P:88:PHE:O	5:P:92:ILE:HD12	2.11	0.50
5:R:239:ASN:ND2	5:R:240:SER:O	2.44	0.50
5:R:368:THR:O	5:R:372:THR:HG23	2.11	0.50
6:o:136:ASP:O	6:o:257:ASN:ND2	2.43	0.50
2:2:415:THR:O	2:2:418:VAL:HG12	2.12	0.50
6:q:274:ASP:OD2	6:q:276:THR:N	2.45	0.50
5:D:99:GLU:CG	5:D:388:VAL:HG21	2.42	0.50
5:D:259:VAL:CG2	5:D:322:LEU:HD11	2.42	0.50
5:E:11:LEU:HD12	5:E:85:ILE:HG23	1.94	0.50
5:O:3:VAL:O	5:O:7:LEU:HD23	2.11	0.50
5:P:216:ARG:NH2	5:P:374:TYR:O	2.44	0.50
6:b:136:ASP:O	6:b:257:ASN:ND2	2.40	0.50
6:p:65:THR:HB	6:p:66:VAL:HG13	1.93	0.50
2:1:288:TYR:CE1	2:1:390:VAL:HG12	2.47	0.50
3:4:42:PRO:O	3:4:44:LYS:N	2.45	0.50
3:6:136:PRO:O	4:7:127:SER:OG	2.28	0.50
5:F:22:THR:OG1	5:F:73:LEU:HD22	2.12	0.50
5:I:242:ASP:OD2	5:I:243:GLY:N	2.44	0.50
6:m:230:VAL:CG2	6:n:105:MET:HE2	2.41	0.50
6:q:124:ILE:HD11	6:q:142:MET:HE2	1.93	0.50
2:3:359:LYS:NZ	2:3:484:MET:SD	2.85	0.50
5:J:307:LEU:HD11	5:K:248:PHE:CD2	2.47	0.50
5:L:113:SER:O	5:L:116:LEU:N	2.45	0.50
5:P:11:LEU:HD11	5:P:88:PHE:HB2	1.92	0.50
6:a:326:VAL:HG12	6:q:320:ALA:HB2	1.94	0.50
2:1:739:GLU:OE1	2:1:739:GLU:N	2.45	0.49
5:Q:3:VAL:O	5:Q:7:LEU:HD23	2.12	0.49
5:R:367:TRP:CE3	5:R:370:LEU:HD23	2.46	0.49
6:b:182:ASN:OD1	6:b:183:LYS:N	2.45	0.49
6:g:138:ASN:N	6:g:257:ASN:O	2.43	0.49
6:p:102:LEU:HD23	6:p:105:MET:HE3	1.93	0.49
6:q:87:THR:O	6:q:91:THR:HG23	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:4:183:LEU:HD13	3:4:189:TYR:CE1	2.47	0.49
4:5:5:LEU:O	4:5:7:GLN:NE2	2.45	0.49
5:G:128:ASN:O	5:H:22:THR:HG22	2.12	0.49
5:H:182:LEU:HD22	5:H:231:ARG:NE	2.27	0.49
5:M:342:LEU:HD12	5:M:389:ALA:HB3	1.94	0.49
4:5:30:ILE:HG23	4:5:35:VAL:HG21	1.94	0.49
5:E:206:GLN:OE1	5:E:207:GLN:N	2.44	0.49
5:F:3:VAL:O	5:F:7:LEU:HD23	2.12	0.49
5:K:194:ASP:OD1	5:K:196:SER:N	2.43	0.49
5:P:23:LEU:HD11	5:R:36:GLN:CD	2.37	0.49
6:k:202:THR:HG22	6:k:202:THR:O	2.12	0.49
6:n:170:ILE:HG23	6:n:237:ILE:HB	1.93	0.49
2:1:273:ARG:NH2	2:1:462:GLU:OE2	2.46	0.49
2:2:529:MET:HE1	2:2:640:LEU:CD2	2.42	0.49
2:2:697:PHE:CE1	2:2:744:ILE:HD11	2.47	0.49
5:I:180:MET:HE2	6:i:63:MET:HE1	1.95	0.49
5:L:27:VAL:HG12	5:L:31:ILE:HG12	1.93	0.49
5:M:158:PHE:CZ	5:M:188:ILE:HD11	2.48	0.49
6:f:315:ARG:HB3	6:f:326:VAL:HG12	1.95	0.49
6:q:145:ASP:OD1	6:q:148:LEU:HD12	2.13	0.49
6:r:138:ASN:CB	6:r:258:VAL:HG12	2.41	0.49
1:U:158:THR:OG1	1:U:179:GLU:OE1	2.15	0.49
2:2:674:ILE:HD11	2:2:716:VAL:HG12	1.95	0.49
3:4:71:ILE:HG22	3:4:82:LEU:HD22	1.94	0.49
4:7:217:CYS:SG	4:7:218:SER:N	2.86	0.49
5:D:173:HIS:HA	5:D:176:LEU:HD11	1.94	0.49
5:G:48:THR:HG21	5:G:94:ASN:HB3	1.93	0.49
5:P:219:THR:HG22	5:P:220:THR:HG23	1.94	0.49
6:b:271:ILE:HG23	6:b:284:MET:HE2	1.94	0.49
6:h:135:CYS:SG	6:h:136:ASP:N	2.84	0.49
6:q:165:CYS:HB2	6:q:249:CYS:HA	1.94	0.49
2:1:526:MET:HE3	2:1:544:ALA:HB1	1.94	0.49
5:E:230:GLU:O	5:E:233:SER:OG	2.28	0.49
5:F:209:GLU:OE2	5:F:211:ILE:HD11	2.12	0.49
5:G:182:LEU:HD22	5:G:231:ARG:NE	2.28	0.49
5:H:7:LEU:CD2	5:H:127:ILE:HG21	2.42	0.49
6:f:274:ASP:OD1	6:f:276:THR:N	2.41	0.49
6:h:230:VAL:HB	6:h:233:ILE:HD12	1.94	0.49
6:m:262:GLN:NE2	6:m:266:ALA:O	2.44	0.49
5:B:48:THR:HG21	5:B:94:ASN:HB3	1.95	0.49
5:D:301:THR:OG1	5:E:246:THR:O	2.21	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:P:182:LEU:HD21	5:P:184:ALA:HB2	1.95	0.49
6:j:139:LEU:HD22	6:j:261:ILE:CD1	2.41	0.49
1:V:21:GLU:OE1	1:V:24:GLN:NE2	2.45	0.49
5:F:165:THR:HG23	5:F:181:TRP:HH2	1.76	0.49
5:J:333:SER:OG	5:J:379:GLU:OE1	2.28	0.49
5:N:207:GLN:OE1	5:N:293:GLN:NE2	2.46	0.49
5:Q:27:VAL:HG23	5:Q:30:LEU:HD23	1.95	0.49
5:C:158:PHE:CZ	5:C:188:ILE:HD11	2.48	0.49
5:E:5:TYR:CE1	5:E:394:MET:HE1	2.48	0.49
5:E:301:THR:OG1	5:F:246:THR:O	2.25	0.49
5:J:321:GLY:O	5:J:322:LEU:HD12	2.13	0.49
5:M:159:PRO:HG3	5:M:371:ILE:HG23	1.95	0.49
5:P:112:GLN:O	5:P:113:SER:OG	2.31	0.49
6:o:170:ILE:HD11	6:o:246:ILE:HG21	1.93	0.49
2:2:618:LEU:HD21	2:2:711:VAL:HG22	1.93	0.49
5:A:207:GLN:NE2	5:A:291:SER:OG	2.46	0.49
5:B:170:GLN:OE1	5:B:170:GLN:N	2.46	0.49
5:J:3:VAL:O	5:J:7:LEU:HD23	2.12	0.49
5:L:114:GLU:OE1	5:L:114:GLU:N	2.44	0.49
6:h:126:ASP:O	6:h:129:VAL:HG12	2.13	0.49
6:k:230:VAL:CG2	6:l:105:MET:HE2	2.43	0.49
5:P:342:LEU:HD12	5:P:386:PHE:O	2.13	0.48
6:e:151:ASP:OD2	6:f:290:LYS:NZ	2.46	0.48
6:j:230:VAL:HG21	6:k:105:MET:HE2	1.95	0.48
6:j:268:ILE:HG13	6:j:269:LEU:HD12	1.93	0.48
6:r:149:GLU:HB2	6:r:268:ILE:HD12	1.95	0.48
6:r:261:ILE:HD13	6:r:299:ILE:HD12	1.95	0.48
3:6:179:PHE:HE1	4:7:168:THR:HG21	1.78	0.48
5:K:300:MET:HE3	5:K:305:ASN:HA	1.94	0.48
5:P:38:ILE:HD11	5:P:84:THR:HG21	1.95	0.48
6:d:160:LEU:HD23	6:d:258:VAL:HG13	1.94	0.48
2:3:428:SER:C	2:3:429:LEU:HD12	2.38	0.48
5:D:261:VAL:HG23	5:D:290:LEU:HD23	1.95	0.48
5:E:294:LEU:HD12	5:E:320:VAL:HG13	1.95	0.48
5:K:36:GLN:HG2	5:L:23:LEU:HD11	1.94	0.48
6:e:274:ASP:OD1	6:e:276:THR:N	2.44	0.48
6:h:80:THR:CG2	6:h:326:VAL:HG22	2.43	0.48
6:l:241:THR:OG1	8:l:504:NAG:H83	2.13	0.48
6:q:186:SER:OG	6:q:191:CYS:SG	2.71	0.48
5:F:89:ILE:HD12	5:F:395:LEU:HD12	1.94	0.48
5:H:29:ASP:OD1	5:H:30:LEU:N	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:M:173:HIS:HA	5:M:176:LEU:HD21	1.95	0.48
1:U:93:TYR:HD2	1:U:198:ILE:HD11	1.78	0.48
2:2:526:MET:HE3	2:2:544:ALA:HB1	1.95	0.48
5:E:255:ARG:NH2	6:f:65:THR:O	2.45	0.48
5:I:166:LEU:O	6:d:52:ASN:ND2	2.47	0.48
5:J:182:LEU:HD13	5:J:231:ARG:CD	2.44	0.48
5:P:165:THR:HG23	5:P:181:TRP:CH2	2.48	0.48
5:P:352:GLU:OE1	5:R:283:ARG:NH1	2.46	0.48
5:Q:230:GLU:N	5:Q:230:GLU:OE2	2.46	0.48
6:i:261:ILE:HD13	6:i:299:ILE:CD1	2.44	0.48
6:m:145:ASP:OD2	6:m:145:ASP:N	2.45	0.48
6:q:62:SER:OG	6:q:63:MET:N	2.47	0.48
5:N:173:HIS:HA	5:N:176:LEU:HD11	1.96	0.48
6:h:233:ILE:HG12	6:i:108:ILE:HD13	1.94	0.48
6:k:126:ASP:O	6:k:129:VAL:HG12	2.14	0.48
6:m:209:THR:O	6:m:236:LYS:NZ	2.29	0.48
5:I:29:ASP:OD1	5:I:30:LEU:N	2.46	0.48
5:J:142:GLN:HG3	5:J:148:THR:HG21	1.96	0.48
5:R:165:THR:HG23	5:R:181:TRP:CH2	2.49	0.48
6:g:105:MET:HE1	6:g:297:TYR:CD1	2.48	0.48
2:2:280:ILE:HD13	2:2:293:ILE:HD13	1.96	0.48
4:7:49:VAL:O	4:7:50:ILE:HD13	2.13	0.48
5:A:158:PHE:CZ	5:A:188:ILE:HD11	2.49	0.48
5:C:49:GLY:N	5:C:98:ASP:OD2	2.46	0.48
5:C:259:VAL:HG11	5:C:277:PHE:CE1	2.49	0.48
5:F:7:LEU:HD22	5:F:127:ILE:HG21	1.96	0.48
5:M:225:LEU:HD22	5:M:261:VAL:HG21	1.96	0.48
6:m:185:ILE:HD12	6:m:249:CYS:SG	2.54	0.48
4:7:12:SER:OG	4:7:110:THR:O	2.30	0.48
5:K:112:GLN:O	5:K:113:SER:OG	2.30	0.48
5:M:367:TRP:O	5:M:371:ILE:HG22	2.13	0.48
5:N:180:MET:HE1	5:N:238:ILE:HD13	1.94	0.48
5:N:188:ILE:HD12	5:N:324:LEU:HD23	1.96	0.48
5:R:214:LEU:HD11	5:R:285:PHE:CE1	2.49	0.48
6:g:144:TYR:HD1	6:g:152:MET:HE1	1.78	0.48
6:l:220:GLU:OE1	6:l:220:GLU:N	2.47	0.48
6:q:202:THR:HG22	6:q:202:THR:O	2.14	0.48
6:q:267:ASN:OD1	6:q:286:ARG:NE	2.43	0.48
6:q:271:ILE:CD1	6:q:281:ILE:HD11	2.44	0.48
1:T:39:PRO:HB3	2:1:260:LEU:HD13	1.95	0.48
2:2:405:PHE:HZ	2:2:408:VAL:HG23	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:4:114:ASP:OD1	4:5:48:THR:HG22	2.14	0.48
5:K:3:VAL:O	5:K:7:LEU:HD23	2.13	0.48
5:L:201:ALA:HB1	5:L:202:PRO:HD2	1.96	0.48
5:N:67:GLY:O	5:N:77:TYR:OH	2.28	0.48
5:O:188:ILE:HD12	5:O:324:LEU:HD22	1.95	0.48
5:P:380:ASP:OD1	5:P:381:ASN:N	2.47	0.48
6:b:230:VAL:HG21	6:c:105:MET:SD	2.53	0.48
6:d:230:VAL:HG11	6:d:233:ILE:HD12	1.96	0.48
6:r:124:ILE:HD11	6:r:155:LEU:HD23	1.96	0.48
1:T:73:THR:HG1	1:T:199:THR:HG1	1.34	0.47
2:2:470:ILE:HG22	2:2:470:ILE:O	2.13	0.47
5:E:39:VAL:CG2	5:E:65:LEU:HD21	2.42	0.47
5:I:99:GLU:HG3	5:I:388:VAL:HG21	1.96	0.47
6:a:135:CYS:SG	6:a:136:ASP:N	2.83	0.47
6:a:259:ALA:HB1	6:a:285:MET:HE2	1.95	0.47
6:r:146:GLN:O	6:r:147:SER:OG	2.27	0.47
4:5:25:GLY:O	4:5:73:ASN:ND2	2.47	0.47
5:I:259:VAL:HG21	5:I:322:LEU:HD11	1.95	0.47
5:K:157:ILE:HD13	5:K:335:LEU:HD21	1.95	0.47
5:L:264:LEU:HD22	5:L:269:ILE:HD13	1.96	0.47
5:Q:334:VAL:O	5:Q:374:TYR:OH	2.31	0.47
6:a:326:VAL:HG12	6:q:320:ALA:CA	2.44	0.47
6:g:164:LEU:HD11	6:g:325:ARG:CG	2.44	0.47
2:1:742:ASN:HB2	2:1:750:MET:HE2	1.96	0.47
2:2:260:LEU:HD13	2:2:480:GLN:NE2	2.30	0.47
5:G:112:GLN:O	5:G:113:SER:OG	2.31	0.47
5:K:254:LEU:HD12	5:K:294:LEU:CD1	2.44	0.47
5:P:310:GLN:OE1	5:P:310:GLN:N	2.47	0.47
5:Q:262:GLU:OE1	5:Q:272:THR:OG1	2.19	0.47
5:R:79:GLU:O	5:R:83:THR:HG23	2.14	0.47
6:a:139:LEU:HD21	6:a:306:ILE:HG21	1.95	0.47
6:p:138:ASN:HB2	6:p:258:VAL:HG12	1.96	0.47
6:r:187:MET:HE1	6:r:322:PHE:CE2	2.49	0.47
5:D:103:GLU:OE2	5:D:350:ARG:NH2	2.47	0.47
5:D:165:THR:HG23	5:D:181:TRP:CH2	2.50	0.47
5:G:300:MET:CG	5:G:304:VAL:HG13	2.45	0.47
6:b:198:LEU:HD12	6:b:202:THR:C	2.38	0.47
2:2:370:LEU:HD12	2:2:468:SER:O	2.15	0.47
2:2:697:PHE:CZ	2:2:744:ILE:HD11	2.49	0.47
3:4:65:VAL:HG12	3:4:69:PHE:CE1	2.49	0.47
5:B:307:LEU:HD11	5:C:248:PHE:CD1	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:G:173:HIS:CA	5:G:176:LEU:HD11	2.44	0.47
5:N:186:SER:O	5:N:186:SER:OG	2.32	0.47
6:n:160:LEU:HD11	6:n:260:VAL:HG23	1.96	0.47
6:o:71:THR:O	6:o:71:THR:HG22	2.15	0.47
6:o:267:ASN:ND2	6:o:280:GLN:OE1	2.47	0.47
6:r:259:ALA:HB3	6:r:310:MET:HE2	1.97	0.47
1:T:13:SER:O	1:T:16:VAL:HG22	2.14	0.47
2:3:621:LYS:O	2:3:625:THR:HG22	2.15	0.47
5:A:183:ASN:OD1	5:A:188:ILE:HG23	2.14	0.47
5:D:259:VAL:HG21	5:D:322:LEU:HD11	1.96	0.47
5:F:394:MET:C	5:F:395:LEU:HD22	2.38	0.47
5:G:186:SER:O	5:G:186:SER:OG	2.28	0.47
5:O:11:LEU:HD11	5:O:88:PHE:HB3	1.96	0.47
5:O:233:SER:HA	5:O:253:ILE:HD11	1.97	0.47
6:r:193:VAL:HG22	6:r:239:LEU:HD13	1.95	0.47
1:U:22:ILE:CD1	1:V:22:ILE:HD11	2.44	0.47
2:2:318:SER:OG	2:2:319:VAL:N	2.47	0.47
5:A:177:MET:HE2	5:A:206:GLN:NE2	2.28	0.47
5:D:4:LEU:HD22	5:D:116:LEU:HD12	1.96	0.47
5:D:252:ILE:HD13	5:D:307:LEU:HD13	1.96	0.47
5:G:300:MET:HG2	5:G:304:VAL:HG13	1.96	0.47
5:N:35:ASN:OD1	5:N:66:LEU:N	2.46	0.47
5:P:99:GLU:HG3	5:P:388:VAL:HG21	1.97	0.47
5:Q:174:ASP:OD2	5:Q:175:ASN:N	2.47	0.47
5:Q:302:PRO:CD	5:R:172:MET:HE1	2.45	0.47
6:i:230:VAL:HG11	6:i:233:ILE:HD12	1.96	0.47
6:o:165:CYS:HB3	6:o:249:CYS:CA	2.38	0.47
6:r:72:ARG:NH1	6:r:308:GLN:OE1	2.48	0.47
1:T:156:THR:OG1	1:T:157:LEU:N	2.48	0.47
5:A:257:ASN:HD22	5:A:295:MET:HE3	1.80	0.47
5:G:7:LEU:HB3	5:G:92:ILE:HD11	1.97	0.47
5:H:137:GLU:OE1	5:I:397:LYS:NZ	2.43	0.47
5:I:165:THR:HG22	6:i:60:THR:HA	1.97	0.47
5:K:7:LEU:HD22	5:K:127:ILE:HG21	1.97	0.47
5:M:37:MET:HA	5:M:127:ILE:HD12	1.97	0.47
6:a:256:GLU:OE1	6:a:256:GLU:N	2.47	0.47
6:f:145:ASP:OD1	6:f:145:ASP:N	2.47	0.47
5:E:7:LEU:O	5:E:11:LEU:HD23	2.14	0.47
5:E:11:LEU:HD11	5:E:88:PHE:CB	2.45	0.47
5:J:176:LEU:HD12	5:J:195:TYR:CZ	2.49	0.47
5:K:47:GLN:NE2	5:K:58:ASN:OD1	2.42	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:L:230:GLU:N	5:L:230:GLU:OE2	2.48	0.47
5:N:158:PHE:CZ	5:N:188:ILE:HD11	2.48	0.47
6:c:105:MET:HE3	6:c:300:VAL:HG11	1.96	0.47
6:c:142:MET:HE1	6:c:155:LEU:HD23	1.97	0.47
6:h:145:ASP:N	6:h:145:ASP:OD1	2.46	0.47
2:1:676:LYS:CE	2:1:710:LEU:HD22	2.45	0.47
2:2:397:ASN:OD1	2:2:398:GLY:N	2.48	0.47
5:A:186:SER:O	5:A:186:SER:OG	2.33	0.47
5:D:99:GLU:OE2	5:D:384:ARG:NE	2.44	0.47
5:I:224:THR:OG1	5:I:327:GLU:OE2	2.22	0.47
5:R:177:MET:SD	5:R:206:GLN:NE2	2.88	0.47
6:i:136:ASP:O	6:i:257:ASN:ND2	2.47	0.47
1:T:2:ALA:HB2	2:1:524:LEU:HD13	1.97	0.46
2:1:606:ASP:OD1	2:1:607:ILE:N	2.46	0.46
5:B:7:LEU:HB3	5:B:92:ILE:HD11	1.96	0.46
5:E:334:VAL:C	5:E:335:LEU:HD12	2.40	0.46
5:R:219:THR:HG22	5:R:220:THR:HG23	1.98	0.46
2:2:346:GLU:OE1	2:2:346:GLU:N	2.47	0.46
2:2:640:LEU:HD13	2:2:662:ALA:CB	2.43	0.46
2:3:278:ASN:ND2	2:3:451:ALA:O	2.49	0.46
3:4:176:VAL:HG12	3:4:195:VAL:CG1	2.30	0.46
5:K:113:SER:O	5:K:116:LEU:N	2.49	0.46
5:R:157:ILE:CD1	5:R:335:LEU:HD21	2.44	0.46
6:b:128:SER:HA	6:b:155:LEU:HD13	1.97	0.46
6:e:147:SER:N	6:e:149:GLU:OE2	2.48	0.46
6:l:274:ASP:OD2	6:l:276:THR:N	2.45	0.46
2:3:348:SER:O	2:3:429:LEU:HD13	2.15	0.46
3:6:37:TRP:CD1	3:6:82:LEU:HD11	2.50	0.46
6:n:85:TYR:OH	6:n:99:LYS:NZ	2.47	0.46
1:U:162:ARG:O	1:U:163:LEU:HD22	2.15	0.46
4:7:21:ILE:HD13	4:7:77:LEU:HD11	1.97	0.46
5:A:3:VAL:O	5:A:7:LEU:HD23	2.14	0.46
5:D:85:ILE:HG22	5:D:89:ILE:HD12	1.98	0.46
5:D:166:LEU:HD12	6:d:59:ILE:HD11	1.96	0.46
5:I:186:SER:OG	5:I:326:ILE:O	2.31	0.46
5:M:36:GLN:HG2	5:N:23:LEU:HD21	1.97	0.46
6:c:185:ILE:HD12	6:c:249:CYS:SG	2.55	0.46
6:h:71:THR:HG22	6:h:71:THR:O	2.14	0.46
6:j:81:LEU:HD11	6:j:303:ILE:HD11	1.95	0.46
6:k:262:GLN:NE2	6:k:265:GLY:O	2.49	0.46
6:p:230:VAL:HG21	6:q:105:MET:HE2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:T:72:THR:HG22	1:T:73:THR:N	2.30	0.46
2:1:370:LEU:HD13	2:1:469:LEU:CD1	2.46	0.46
5:B:104:SER:OG	5:B:106:ARG:O	2.33	0.46
5:E:165:THR:HG23	5:E:181:TRP:CH2	2.50	0.46
5:F:239:ASN:ND2	6:f:65:THR:O	2.49	0.46
5:I:177:MET:HE3	5:I:206:GLN:CD	2.41	0.46
5:N:2:GLU:OE1	5:N:143:ASN:ND2	2.48	0.46
5:O:170:GLN:NE2	5:O:175:ASN:O	2.45	0.46
6:d:71:THR:O	6:d:71:THR:HG22	2.15	0.46
6:g:126:ASP:O	6:g:129:VAL:HG12	2.16	0.46
6:i:138:ASN:HB2	6:i:258:VAL:HG23	1.98	0.46
6:j:185:ILE:HD12	6:j:249:CYS:SG	2.55	0.46
6:k:170:ILE:HG23	6:k:237:ILE:HB	1.97	0.46
6:k:271:ILE:HD11	6:k:281:ILE:HD12	1.96	0.46
6:l:124:ILE:HD12	6:l:142:MET:HG3	1.97	0.46
6:q:73:GLU:N	6:q:73:GLU:OE1	2.49	0.46
2:3:721:ILE:HD11	2:3:751:LEU:HD11	1.97	0.46
5:B:253:ILE:HD13	5:C:234:PHE:CD2	2.51	0.46
5:D:237:VAL:CG2	5:F:254:LEU:HD13	2.45	0.46
5:I:206:GLN:NE2	5:I:207:GLN:O	2.48	0.46
5:O:89:ILE:HG12	5:O:395:LEU:HD22	1.98	0.46
6:e:239:LEU:HD23	6:e:239:LEU:O	2.16	0.46
6:j:170:ILE:HD11	6:j:246:ILE:HG21	1.97	0.46
6:n:202:THR:HG22	6:n:202:THR:O	2.15	0.46
1:U:179:GLU:O	1:U:182:ARG:C	2.59	0.46
5:D:307:LEU:HD11	5:E:248:PHE:CD2	2.50	0.46
5:G:29:ASP:OD1	5:G:30:LEU:N	2.48	0.46
5:I:7:LEU:CD2	5:I:127:ILE:HG21	2.44	0.46
5:J:334:VAL:C	5:J:335:LEU:HD12	2.41	0.46
5:K:41:MET:HE1	5:K:119:LEU:HD21	1.98	0.46
5:L:85:ILE:HG22	5:L:89:ILE:HD12	1.97	0.46
2:1:269:ASP:OD2	2:2:248:GLN:NE2	2.45	0.46
4:5:123:LEU:CD2	4:5:212:VAL:HG13	2.45	0.46
5:H:41:MET:HE1	5:H:119:LEU:HD21	1.98	0.46
5:I:11:LEU:HD11	5:I:88:PHE:CB	2.45	0.46
5:J:1:MET:CE	5:J:387:THR:HG23	2.46	0.46
5:M:7:LEU:CD1	5:M:37:MET:HE3	2.46	0.46
5:N:54:LEU:HD12	5:N:55:PRO:HD2	1.98	0.46
5:R:253:ILE:HD13	5:R:319:THR:HG23	1.98	0.46
6:d:239:LEU:HD12	6:d:240:THR:HG22	1.96	0.46
6:e:149:GLU:OE2	6:e:149:GLU:N	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:m:124:ILE:CG2	6:m:148:LEU:HD23	2.46	0.46
6:r:106:PHE:CE2	6:r:300:VAL:HG22	2.50	0.46
1:T:48:VAL:HG13	2:1:421:ASN:OD1	2.16	0.46
2:1:293:ILE:HG23	2:1:293:ILE:O	2.15	0.46
4:5:126:PRO:CB	4:5:131:LEU:HD21	2.46	0.46
5:C:113:SER:O	5:C:116:LEU:N	2.47	0.46
5:D:239:ASN:O	5:D:247:TRP:NE1	2.49	0.46
5:E:165:THR:HG22	6:e:60:THR:HA	1.98	0.46
5:G:3:VAL:O	5:G:7:LEU:HD23	2.15	0.46
5:I:168:ARG:NH2	5:I:177:MET:SD	2.89	0.46
5:I:261:VAL:HG22	5:I:290:LEU:HD23	1.98	0.46
5:J:261:VAL:HG22	5:J:290:LEU:HD23	1.97	0.46
5:P:93:ASP:OD1	5:P:392:ARG:NE	2.49	0.46
5:R:7:LEU:HD12	5:R:127:ILE:HG21	1.97	0.46
6:d:276:THR:HG21	6:e:305:GLN:HB3	1.98	0.46
6:p:230:VAL:CG2	6:q:105:MET:HE2	2.46	0.46
2:1:670:ARG:O	2:1:685:THR:HG23	2.16	0.46
1:U:72:THR:HG22	1:U:73:THR:N	2.31	0.46
2:2:635:ILE:HD11	2:2:720:ILE:CD1	2.45	0.46
4:5:69:ASP:OD1	4:5:72:SER:OG	2.19	0.46
5:G:106:ARG:NH2	5:L:397:LYS:OXT	2.44	0.46
5:H:71:LEU:O	5:H:71:LEU:HD23	2.15	0.46
5:N:112:GLN:O	5:N:113:SER:OG	2.34	0.46
5:R:7:LEU:HB3	5:R:92:ILE:HD11	1.97	0.46
6:k:170:ILE:HD11	6:k:246:ILE:HG21	1.98	0.46
6:n:233:ILE:HD11	6:o:108:ILE:HG21	1.98	0.46
5:A:162:ALA:HB2	5:A:182:LEU:HD12	1.97	0.45
5:B:165:THR:HG23	5:B:181:TRP:HH2	1.80	0.45
5:E:166:LEU:HD22	5:E:176:LEU:HD22	1.97	0.45
5:F:169:SER:HA	5:F:176:LEU:HD23	1.97	0.45
5:K:165:THR:HG23	5:K:181:TRP:CH2	2.51	0.45
5:M:231:ARG:O	5:M:236:ARG:NH1	2.49	0.45
6:r:285:MET:HE1	6:r:306:ILE:HG23	1.98	0.45
2:2:551:PHE:HA	2:2:660:THR:HG22	1.98	0.45
4:5:49:VAL:O	4:5:50:ILE:HD13	2.16	0.45
5:K:237:VAL:O	5:K:237:VAL:HG13	2.15	0.45
5:M:29:ASP:OD1	5:M:30:LEU:N	2.49	0.45
5:N:214:LEU:HD11	5:N:285:PHE:CE2	2.51	0.45
5:Q:14:ALA:HA	5:Q:30:LEU:HD21	1.98	0.45
5:R:113:SER:O	5:R:116:LEU:N	2.49	0.45
6:d:192:THR:HG23	6:d:192:THR:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:r:202:THR:O	6:r:202:THR:HG22	2.16	0.45
2:2:655:LEU:HA	2:2:658:ILE:HG22	1.99	0.45
5:A:113:SER:O	5:A:116:LEU:N	2.49	0.45
5:C:342:LEU:HD12	5:C:386:PHE:O	2.16	0.45
5:N:99:GLU:OE1	5:N:384:ARG:NE	2.40	0.45
6:c:102:LEU:HD23	6:c:105:MET:CE	2.46	0.45
6:c:111:TRP:HZ2	6:c:307:VAL:HG21	1.80	0.45
6:l:106:PHE:CE2	6:l:300:VAL:HG22	2.52	0.45
6:m:106:PHE:CZ	6:m:300:VAL:HG22	2.52	0.45
6:p:111:TRP:CZ2	6:p:307:VAL:HG11	2.51	0.45
6:p:288:ASN:ND2	6:r:149:GLU:OE1	2.45	0.45
1:T:84:ILE:O	1:T:163:LEU:HD13	2.17	0.45
1:T:93:TYR:HB2	1:T:200:ILE:HD13	1.98	0.45
1:U:129:ASN:OD1	1:U:130:VAL:N	2.49	0.45
3:4:37:TRP:CG	3:4:82:LEU:HD11	2.51	0.45
3:6:225:GLU:N	3:6:225:GLU:OE1	2.50	0.45
5:C:180:MET:HE3	5:C:232:PHE:HD2	1.80	0.45
5:G:173:HIS:C	5:G:176:LEU:HD11	2.42	0.45
5:K:165:THR:HG22	6:k:60:THR:HA	1.98	0.45
5:K:334:VAL:C	5:K:335:LEU:HD12	2.42	0.45
5:M:165:THR:HG23	5:M:181:TRP:CH2	2.52	0.45
6:c:261:ILE:HD13	6:c:299:ILE:CD1	2.47	0.45
6:j:124:ILE:HD13	6:j:142:MET:HG3	1.97	0.45
6:l:192:THR:HG23	6:l:192:THR:O	2.16	0.45
6:r:182:ASN:O	6:r:183:LYS:NZ	2.38	0.45
2:2:533:ILE:HG23	2:2:533:ILE:O	2.17	0.45
2:3:765:ARG:O	2:3:769:LEU:HD23	2.17	0.45
5:F:194:ASP:OD2	5:F:198:ALA:N	2.47	0.45
5:H:89:ILE:CG2	5:H:396:ILE:HD11	2.46	0.45
5:L:27:VAL:HG12	5:L:27:VAL:O	2.15	0.45
5:Q:99:GLU:HG2	5:Q:388:VAL:HG21	1.99	0.45
6:d:126:ASP:O	6:d:129:VAL:HG12	2.16	0.45
6:h:193:VAL:HG22	6:h:239:LEU:HD12	1.98	0.45
5:D:113:SER:O	5:D:113:SER:OG	2.32	0.45
5:I:112:GLN:O	5:I:113:SER:OG	2.34	0.45
5:L:264:LEU:CD2	5:L:269:ILE:HD13	2.46	0.45
5:R:138:ASN:OD1	5:R:148:THR:OG1	2.35	0.45
6:f:178:SER:OG	6:f:179:GLY:N	2.49	0.45
6:j:81:LEU:HD13	6:j:303:ILE:HD11	1.97	0.45
6:k:315:ARG:HB2	6:k:326:VAL:HG12	1.97	0.45
2:1:431:VAL:HG21	2:1:447:TYR:HB3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:1:518:ASP:OD1	2:1:519:LEU:N	2.49	0.45
2:2:288:TYR:CZ	2:2:390:VAL:HG12	2.52	0.45
2:2:721:ILE:HD11	2:2:762:ILE:CD1	2.47	0.45
5:A:347:THR:HB	5:C:281:VAL:HG21	1.97	0.45
5:C:231:ARG:O	5:C:236:ARG:NH1	2.50	0.45
5:E:228:ASP:OD1	5:F:236:ARG:NH2	2.45	0.45
5:G:142:GLN:HB2	5:G:148:THR:HG21	1.99	0.45
5:N:240:SER:OG	5:N:241:ALA:N	2.50	0.45
5:P:156:ASN:HD22	5:R:279:THR:HG21	1.81	0.45
6:l:305:GLN:O	6:l:309:VAL:HG23	2.16	0.45
1:U:211:GLN:N	1:U:211:GLN:OE1	2.50	0.45
4:7:30:ILE:HG23	4:7:35:VAL:HG21	1.97	0.45
5:E:127:ILE:HG22	5:E:127:ILE:O	2.17	0.45
5:O:199:LEU:HD11	5:O:310:GLN:OE1	2.16	0.45
6:a:326:VAL:HG12	6:q:320:ALA:CB	2.47	0.45
6:c:142:MET:CE	6:c:155:LEU:HD23	2.47	0.45
6:q:142:MET:HE1	6:q:155:LEU:HD22	1.98	0.45
1:T:100:ASP:OD2	1:T:100:ASP:N	2.50	0.45
1:U:156:THR:OG1	1:U:157:LEU:N	2.49	0.45
5:A:182:LEU:HD13	5:A:231:ARG:HE	1.82	0.45
5:C:158:PHE:CE1	5:C:326:ILE:HD12	2.52	0.45
5:C:368:THR:O	5:C:372:THR:OG1	2.31	0.45
5:E:128:ASN:O	5:F:22:THR:HG22	2.16	0.45
5:F:230:GLU:OE1	5:F:230:GLU:N	2.49	0.45
5:I:74:ASP:OD2	5:I:74:ASP:N	2.50	0.45
5:J:11:LEU:HD11	5:J:88:PHE:HB2	1.99	0.45
5:P:36:GLN:O	5:P:40:THR:HG22	2.17	0.45
6:g:124:ILE:HG23	6:g:155:LEU:HD13	1.97	0.45
1:T:205:TYR:OH	2:2:442:ARG:NH2	2.50	0.45
2:2:270:ILE:HD11	2:2:354:TYR:CE2	2.51	0.45
2:2:574:ALA:N	2:3:510:GLU:OE2	2.49	0.45
2:3:280:ILE:HG23	2:3:293:ILE:CD1	2.47	0.45
2:3:428:SER:O	2:3:429:LEU:HD12	2.17	0.45
4:7:110:THR:HG21	4:7:178:TYR:OH	2.16	0.45
5:C:186:SER:O	5:C:186:SER:OG	2.29	0.45
5:F:70:LEU:HD23	5:F:73:LEU:HA	1.98	0.45
5:G:141:LEU:O	5:G:146:GLN:N	2.49	0.45
5:G:270:ILE:HG23	5:G:283:ARG:HH11	1.82	0.45
5:I:207:GLN:NE2	5:I:291:SER:OG	2.49	0.45
5:J:299:ASN:ND2	5:K:244:ALA:O	2.47	0.45
6:b:150:LEU:HD23	6:c:290:LYS:HG3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:k:144:TYR:HB2	6:k:152:MET:HE1	1.98	0.45
6:p:285:MET:HE2	6:p:306:ILE:CD1	2.47	0.45
2:1:369:SER:HB3	2:1:470:ILE:HD12	1.98	0.44
2:1:438:ILE:HG23	2:1:441:THR:HB	1.99	0.44
4:7:136:ALA:HB3	4:7:186:LEU:O	2.18	0.44
5:B:7:LEU:HD12	5:B:127:ILE:HG21	1.99	0.44
5:C:112:GLN:O	5:C:113:SER:OG	2.32	0.44
5:D:165:THR:HG23	5:D:181:TRP:HH2	1.82	0.44
5:J:261:VAL:HG22	5:J:290:LEU:CD2	2.46	0.44
5:P:219:THR:N	5:P:330:VAL:O	2.50	0.44
5:P:384:ARG:O	5:P:388:VAL:HG23	2.16	0.44
5:Q:127:ILE:O	5:Q:127:ILE:HG22	2.17	0.44
6:a:261:ILE:HG13	6:a:285:MET:HE3	1.98	0.44
6:h:164:LEU:HB2	6:h:252:LEU:HD11	1.99	0.44
6:j:142:MET:HE3	6:j:152:MET:HG3	1.98	0.44
6:r:145:ASP:N	6:r:145:ASP:OD1	2.49	0.44
2:1:441:THR:HG22	2:1:442:ARG:N	2.32	0.44
5:C:288:ILE:HG22	5:C:290:LEU:CD1	2.47	0.44
5:K:262:GLU:OE1	5:K:289:ARG:NH2	2.50	0.44
6:g:75:VAL:HG13	6:g:75:VAL:O	2.16	0.44
6:h:85:TYR:OH	6:h:99:LYS:NZ	2.50	0.44
6:i:71:THR:O	6:i:71:THR:HG22	2.17	0.44
6:m:108:ILE:HD13	6:o:233:ILE:HD11	1.99	0.44
5:A:17:LYS:HD2	5:A:30:LEU:HD22	2.00	0.44
5:Q:75:ALA:O	5:Q:78:VAL:N	2.47	0.44
5:Q:165:THR:HG22	6:q:60:THR:HA	1.99	0.44
6:a:265:GLY:O	6:a:286:ARG:NH1	2.50	0.44
6:k:218:VAL:HG21	6:k:227:VAL:HG21	1.99	0.44
6:l:160:LEU:O	6:l:283:ARG:NH1	2.49	0.44
6:m:108:ILE:HD11	6:o:230:VAL:CG1	2.47	0.44
2:2:618:LEU:CD2	2:2:716:VAL:HG11	2.45	0.44
5:B:67:GLY:O	5:B:77:TYR:OH	2.28	0.44
5:D:222:THR:HG22	5:D:328:SER:OG	2.18	0.44
5:E:254:LEU:HD23	5:F:237:VAL:HG22	2.00	0.44
5:G:40:THR:HG21	5:G:127:ILE:HG13	1.98	0.44
5:H:89:ILE:CD1	5:H:395:LEU:HD22	2.47	0.44
5:J:252:ILE:HD13	5:K:235:PRO:HB2	1.98	0.44
6:d:160:LEU:HD22	6:d:283:ARG:HH11	1.82	0.44
6:j:187:MET:HE1	6:j:322:PHE:CE2	2.52	0.44
6:k:192:THR:O	6:k:192:THR:HG23	2.18	0.44
6:l:142:MET:HG2	6:l:152:MET:HE2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:l:146:GLN:O	6:l:147:SER:OG	2.28	0.44
6:l:172:LEU:HD11	6:m:99:LYS:HE3	2.00	0.44
6:m:230:VAL:HG21	6:n:105:MET:HE2	1.99	0.44
6:n:268:ILE:HG13	6:n:269:LEU:HD12	2.00	0.44
6:r:228:ASP:OD2	6:r:229:VAL:N	2.51	0.44
5:C:334:VAL:O	5:C:374:TYR:OH	2.32	0.44
5:E:11:LEU:HD11	5:E:88:PHE:HB2	2.00	0.44
5:G:237:VAL:CG2	5:I:254:LEU:HD13	2.47	0.44
5:J:366:ASN:OD1	5:L:276:ARG:NH2	2.50	0.44
6:j:139:LEU:HD21	6:j:306:ILE:HG21	1.99	0.44
2:1:735:ILE:HD12	2:1:740:ALA:HA	2.00	0.44
2:3:611:THR:O	2:3:683:ILE:HD11	2.18	0.44
3:6:139:PRO:HG2	3:6:202:LEU:HD21	2.00	0.44
5:B:2:GLU:OE1	5:B:2:GLU:N	2.47	0.44
5:B:134:GLU:O	5:B:138:ASN:ND2	2.50	0.44
5:H:194:ASP:OD2	5:H:197:CYS:N	2.50	0.44
5:J:383:GLN:O	5:J:387:THR:HG22	2.17	0.44
5:N:231:ARG:O	5:N:236:ARG:NH1	2.51	0.44
6:p:68:SER:OG	6:p:69:ASN:N	2.50	0.44
5:G:236:ARG:HG2	5:I:253:ILE:HD11	2.00	0.44
5:I:3:VAL:O	5:I:7:LEU:HD23	2.17	0.44
5:I:342:LEU:HD12	5:I:389:ALA:HB3	2.00	0.44
5:P:264:LEU:CD2	5:P:269:ILE:HD13	2.48	0.44
5:R:99:GLU:OE1	5:R:384:ARG:NE	2.50	0.44
6:o:146:GLN:O	6:o:147:SER:OG	2.29	0.44
6:p:143:LYS:NZ	6:p:145:ASP:OD1	2.48	0.44
2:1:397:ASN:OD1	2:1:398:GLY:N	2.48	0.44
2:1:611:THR:CG2	2:1:683:ILE:HG23	2.48	0.44
2:1:611:THR:HG22	2:1:683:ILE:HG23	1.98	0.44
2:2:695:ASN:O	2:2:737:ARG:NH1	2.51	0.44
3:4:19:LEU:HD12	3:4:20:LYS:N	2.33	0.44
4:5:171:SER:O	4:5:173:GLN:NE2	2.51	0.44
5:B:173:HIS:HA	5:B:176:LEU:HD11	1.99	0.44
5:D:246:THR:HG21	5:F:297:PRO:HB3	2.00	0.44
5:G:384:ARG:O	5:G:388:VAL:HG23	2.18	0.44
5:M:157:ILE:HD13	5:M:335:LEU:HD11	2.00	0.44
5:N:165:THR:HG23	5:N:181:TRP:CZ3	2.53	0.44
6:c:146:GLN:O	6:c:147:SER:OG	2.30	0.44
6:h:267:ASN:OD1	6:h:286:ARG:NH2	2.51	0.44
2:2:694:ILE:HG13	2:2:744:ILE:HD13	1.99	0.44
4:5:78:THR:O	4:5:79:ILE:HD13	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:17:LYS:HB3	5:E:27:VAL:HG12	2.00	0.44
5:N:180:MET:HE1	5:N:238:ILE:CD1	2.48	0.44
5:O:127:ILE:O	5:O:127:ILE:HG22	2.17	0.44
6:c:142:MET:SD	6:c:260:VAL:HG13	2.58	0.44
1:T:194:ASN:OD1	3:4:54:THR:HG21	2.18	0.43
2:2:729:LEU:CD1	2:2:762:ILE:HD11	2.48	0.43
3:4:205:GLN:NE2	3:4:206:THR:O	2.46	0.43
5:H:89:ILE:CD1	5:H:395:LEU:HD13	2.48	0.43
5:K:183:ASN:OD1	5:K:188:ILE:HD12	2.18	0.43
5:N:158:PHE:CE1	5:N:188:ILE:HD11	2.53	0.43
6:f:165:CYS:HB3	6:f:249:CYS:CA	2.30	0.43
6:g:129:VAL:HG13	6:g:129:VAL:O	2.18	0.43
6:j:135:CYS:O	6:j:315:ARG:NE	2.51	0.43
6:n:233:ILE:CD1	6:o:108:ILE:HG21	2.48	0.43
2:2:698:ASP:OD1	2:2:698:ASP:N	2.52	0.43
5:E:243:GLY:O	6:e:68:SER:OG	2.33	0.43
5:H:22:THR:O	5:H:72:ASN:ND2	2.51	0.43
5:H:127:ILE:HG22	5:H:127:ILE:O	2.17	0.43
5:J:113:SER:O	5:J:113:SER:OG	2.34	0.43
5:R:182:LEU:HD11	5:R:184:ALA:HB2	2.00	0.43
6:h:74:GLU:O	6:h:78:THR:HG22	2.18	0.43
6:h:129:VAL:HG13	6:h:129:VAL:O	2.18	0.43
6:o:81:LEU:HD11	6:o:139:LEU:HD13	2.00	0.43
1:T:132:ASN:OD1	1:T:133:ASP:N	2.52	0.43
1:U:72:THR:HG22	1:U:73:THR:H	1.83	0.43
2:2:459:GLU:OE1	2:2:459:GLU:N	2.49	0.43
3:6:114:ASP:OD1	4:7:48:THR:HG22	2.18	0.43
5:A:141:LEU:HB3	5:A:148:THR:HG22	2.00	0.43
5:E:170:GLN:NE2	5:E:175:ASN:O	2.47	0.43
5:I:165:THR:HG23	5:I:181:TRP:CH2	2.53	0.43
5:N:290:LEU:HD21	5:N:324:LEU:HD22	2.00	0.43
6:c:137:TYR:CE2	6:c:307:VAL:HG13	2.54	0.43
6:i:233:ILE:HG22	6:i:234:ASN:O	2.18	0.43
6:k:272:THR:HG21	6:k:277:THR:CG2	2.49	0.43
6:n:69:ASN:ND2	8:n:401:NAG:O7	2.51	0.43
2:1:270:ILE:HD11	2:1:354:TYR:CE2	2.53	0.43
2:2:506:SER:O	2:2:509:GLN:NE2	2.51	0.43
2:3:630:MET:HG2	2:3:720:ILE:HD13	2.00	0.43
5:G:307:LEU:HD11	5:H:248:PHE:CE1	2.54	0.43
5:R:165:THR:C	5:R:166:LEU:HD12	2.44	0.43
6:d:125:VAL:HG12	6:d:223:LYS:CG	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:f:144:TYR:OH	6:f:146:GLN:OE1	2.22	0.43
6:f:263:VAL:HG12	6:f:289:TRP:HB3	2.00	0.43
6:g:193:VAL:HG22	6:g:239:LEU:CD1	2.40	0.43
6:g:230:VAL:HG21	6:h:105:MET:HG2	1.99	0.43
6:h:139:LEU:HD12	6:h:261:ILE:HD12	2.00	0.43
6:l:217:THR:HG23	6:l:217:THR:O	2.19	0.43
6:m:108:ILE:HD11	6:o:230:VAL:HG11	1.99	0.43
2:1:651:GLY:N	2:1:654:THR:OG1	2.46	0.43
4:7:5:LEU:O	4:7:7:GLN:NE2	2.51	0.43
5:F:336:ALA:HB2	5:F:343:LEU:HD22	2.00	0.43
5:H:228:ASP:OD2	5:I:236:ARG:NH2	2.45	0.43
5:N:99:GLU:HG2	5:N:388:VAL:HG21	2.00	0.43
5:P:279:THR:HG21	5:Q:156:ASN:HD22	1.83	0.43
6:m:146:GLN:O	6:m:147:SER:OG	2.25	0.43
2:3:699:GLU:OE2	5:M:269:ILE:HG22	2.19	0.43
4:7:5:LEU:HD11	4:7:30:ILE:HD11	1.99	0.43
5:J:10:THR:HG21	5:J:33:GLN:OE1	2.18	0.43
5:O:17:LYS:HB3	5:O:27:VAL:HG12	2.01	0.43
5:P:216:ARG:NH2	5:P:379:GLU:OE1	2.50	0.43
5:Q:19:VAL:O	5:Q:22:THR:HG22	2.18	0.43
6:a:170:ILE:HD12	6:a:237:ILE:HB	2.00	0.43
6:b:105:MET:HB3	6:b:300:VAL:HG11	2.01	0.43
6:d:183:LYS:NZ	6:e:302:TYR:OH	2.50	0.43
6:m:108:ILE:HD13	6:o:233:ILE:CD1	2.49	0.43
6:n:185:ILE:HD13	6:n:226:ILE:HG23	2.01	0.43
7:S:1:NAG:O6	7:S:2:NAG:N2	2.52	0.43
3:4:54:THR:O	3:4:54:THR:HG22	2.18	0.43
5:D:23:LEU:HD11	5:F:36:GLN:HB2	2.00	0.43
5:D:259:VAL:HG23	5:D:292:PHE:HB3	2.00	0.43
5:F:264:LEU:HD23	5:F:269:ILE:HA	2.00	0.43
5:N:303:ALA:O	5:N:307:LEU:HD23	2.19	0.43
6:b:271:ILE:HD11	6:b:281:ILE:HD11	2.01	0.43
6:n:187:MET:HG2	6:n:224:LEU:HD13	2.01	0.43
6:q:271:ILE:HD11	6:q:281:ILE:HD11	2.01	0.43
2:1:567:LEU:HD12	2:2:522:LEU:HD22	2.01	0.43
2:2:374:LEU:HD23	2:2:405:PHE:CE1	2.54	0.43
5:L:19:VAL:O	5:L:22:THR:HG22	2.19	0.43
6:q:124:ILE:CD1	6:q:142:MET:HE2	2.49	0.43
2:2:485:ASN:O	2:2:486:SER:OG	2.30	0.43
5:C:36:GLN:O	5:C:40:THR:HG22	2.19	0.43
5:D:131:ASN:N	5:E:16:ASP:OD1	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:N:206:GLN:OE1	5:N:207:GLN:N	2.52	0.43
5:O:216:ARG:NH2	5:O:379:GLU:OE1	2.48	0.43
5:R:7:LEU:CD1	5:R:127:ILE:HG21	2.49	0.43
6:g:108:ILE:HG13	6:i:233:ILE:HD11	2.00	0.43
6:j:126:ASP:O	6:j:129:VAL:HG12	2.19	0.43
6:l:79:SER:OG	6:l:136:ASP:OD1	2.14	0.43
6:l:202:THR:O	6:l:202:THR:HG22	2.19	0.43
6:n:57:ILE:HD13	6:p:59:ILE:HG22	2.01	0.43
6:n:281:ILE:HG22	6:n:283:ARG:H	1.84	0.43
6:q:255:ARG:O	6:q:258:VAL:HG22	2.19	0.43
1:T:211:GLN:OE1	1:T:211:GLN:N	2.52	0.43
4:5:131:LEU:HD12	4:5:218:SER:OXT	2.19	0.43
5:F:112:GLN:O	5:F:113:SER:OG	2.37	0.43
6:d:197:PRO:CG	6:e:105:MET:HE1	2.49	0.43
6:g:59:ILE:HD12	6:k:52:ASN:ND2	2.34	0.43
2:1:374:LEU:HD23	2:1:405:PHE:CE1	2.54	0.42
2:2:697:PHE:O	2:2:727:LYS:NZ	2.43	0.42
2:3:607:ILE:HG23	2:3:704:VAL:HG11	2.02	0.42
5:C:214:LEU:HD11	5:C:285:PHE:CE2	2.54	0.42
5:E:67:GLY:O	5:E:77:TYR:OH	2.33	0.42
5:Q:163:SER:OG	5:Q:181:TRP:NE1	2.52	0.42
5:R:127:ILE:O	5:R:127:ILE:HG22	2.19	0.42
6:a:68:SER:OG	6:a:69:ASN:N	2.52	0.42
6:c:126:ASP:O	6:c:129:VAL:HG12	2.19	0.42
6:i:193:VAL:HG22	6:i:239:LEU:CD1	2.49	0.42
6:q:140:VAL:CG1	6:q:142:MET:HE3	2.49	0.42
6:r:170:ILE:HD12	6:r:237:ILE:HB	2.00	0.42
1:T:76:PRO:HD3	1:T:198:ILE:HG22	2.00	0.42
1:U:87:ASN:N	1:U:202:SER:OG	2.51	0.42
5:D:6:SER:OG	5:D:127:ILE:O	2.32	0.42
5:E:45:ASP:OD2	5:E:45:ASP:N	2.50	0.42
5:E:307:LEU:HD21	5:F:248:PHE:CE1	2.54	0.42
5:I:370:LEU:O	5:I:370:LEU:HD23	2.19	0.42
5:J:360:VAL:HG23	5:J:361:PHE:N	2.34	0.42
5:L:4:LEU:HD12	5:L:116:LEU:HD13	2.00	0.42
5:L:186:SER:OG	5:L:326:ILE:N	2.52	0.42
6:i:129:VAL:O	6:i:129:VAL:HG13	2.19	0.42
6:i:228:ASP:OD2	6:i:229:VAL:N	2.52	0.42
6:k:129:VAL:HG13	6:k:129:VAL:O	2.19	0.42
1:V:3:SER:N	2:3:634:ASP:OD1	2.45	0.42
5:B:7:LEU:CD1	5:B:127:ILE:HG21	2.48	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:C:37:MET:HE3	5:C:88:PHE:CE1	2.54	0.42
5:C:360:VAL:O	5:C:378:ARG:NH2	2.53	0.42
5:D:99:GLU:HG2	5:D:388:VAL:HG21	2.01	0.42
5:G:36:GLN:O	5:G:40:THR:HG22	2.19	0.42
5:J:127:ILE:O	5:J:127:ILE:HG22	2.17	0.42
5:P:248:PHE:CE2	5:R:307:LEU:HD21	2.55	0.42
5:Q:158:PHE:CE1	5:Q:188:ILE:HD11	2.54	0.42
6:b:317:LEU:HD12	6:f:324:TYR:HB3	2.01	0.42
6:p:106:PHE:CZ	6:p:300:VAL:HG22	2.55	0.42
5:D:38:ILE:HD11	5:D:84:THR:HG21	2.02	0.42
5:N:41:MET:HE1	5:N:91:PHE:CE2	2.55	0.42
5:Q:103:GLU:OE1	5:Q:350:ARG:NH1	2.47	0.42
6:r:152:MET:SD	6:r:269:LEU:HD11	2.60	0.42
1:T:138:LYS:HD3	1:T:158:THR:HG22	2.02	0.42
2:3:618:LEU:HD21	2:3:711:VAL:HG22	2.01	0.42
5:K:258:ASN:ND2	5:K:293:GLN:OE1	2.48	0.42
5:L:264:LEU:HD12	5:L:287:THR:OG1	2.20	0.42
1:T:111:HIS:CG	4:7:78:THR:HG21	2.55	0.42
2:1:384:TYR:HB3	2:1:396:ILE:HG22	2.01	0.42
2:1:653:ASN:OD1	2:1:654:THR:N	2.52	0.42
2:2:269:ASP:OD1	2:2:269:ASP:N	2.52	0.42
5:C:170:GLN:N	5:C:170:GLN:OE1	2.53	0.42
5:K:5:TYR:CE2	5:K:394:MET:HE1	2.55	0.42
5:M:99:GLU:CG	5:M:388:VAL:HG21	2.50	0.42
5:P:252:ILE:HD13	5:P:307:LEU:HD23	2.00	0.42
5:R:165:THR:HG23	5:R:181:TRP:HH2	1.85	0.42
6:b:317:LEU:HD21	6:f:162:GLU:HB3	2.01	0.42
6:i:123:ASN:OD1	6:i:123:ASN:N	2.53	0.42
6:o:272:THR:HG21	6:o:277:THR:OG1	2.20	0.42
2:1:288:TYR:CZ	2:1:390:VAL:HG12	2.54	0.42
2:2:412:THR:HG22	2:2:421:ASN:ND2	2.34	0.42
3:6:72:SER:OG	3:6:81:TYR:O	2.36	0.42
5:B:99:GLU:OE2	5:B:384:ARG:NH1	2.53	0.42
5:D:215:ARG:O	5:D:216:ARG:NH1	2.48	0.42
5:F:127:ILE:O	5:F:127:ILE:HG22	2.19	0.42
5:G:160:TYR:OH	6:g:64:ASP:OD2	2.20	0.42
5:H:158:PHE:CZ	5:H:188:ILE:HD11	2.55	0.42
5:L:127:ILE:O	5:L:127:ILE:HG22	2.20	0.42
5:P:177:MET:SD	5:P:206:GLN:NE2	2.93	0.42
6:d:226:ILE:HG23	6:d:226:ILE:O	2.20	0.42
6:f:129:VAL:O	6:f:129:VAL:HG13	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:m:52:ASN:OD1	6:m:53:TYR:N	2.53	0.42
6:o:81:LEU:HD21	6:o:303:ILE:CD1	2.45	0.42
6:q:82:CYS:HB2	6:q:138:ASN:HA	2.00	0.42
4:5:217:CYS:SG	4:5:218:SER:N	2.93	0.42
5:A:307:LEU:HD11	5:B:248:PHE:CD1	2.55	0.42
5:B:127:ILE:O	5:B:127:ILE:HG22	2.20	0.42
5:G:236:ARG:CZ	5:I:253:ILE:HD11	2.50	0.42
5:I:40:THR:CB	5:I:127:ILE:HD11	2.50	0.42
5:J:6:SER:O	5:J:10:THR:HG23	2.20	0.42
5:M:259:VAL:HG23	5:M:322:LEU:HD11	2.01	0.42
5:R:85:ILE:HG22	5:R:89:ILE:HD12	2.01	0.42
6:b:202:THR:O	6:b:202:THR:HG22	2.19	0.42
6:l:145:ASP:OD1	6:l:148:LEU:HD22	2.20	0.42
6:r:192:THR:HG23	6:r:192:THR:O	2.20	0.42
2:1:720:ILE:HD11	2:1:751:LEU:HD13	2.02	0.42
2:2:324:ASN:OD1	2:2:324:ASN:N	2.52	0.42
2:2:606:ASP:OD2	2:2:606:ASP:N	2.52	0.42
2:2:684:ASN:OD1	2:2:685:THR:N	2.53	0.42
5:G:112:GLN:OE1	5:G:112:GLN:N	2.53	0.42
5:I:384:ARG:O	5:I:388:VAL:HG23	2.20	0.42
5:M:254:LEU:HD23	5:N:237:VAL:HB	2.02	0.42
6:b:129:VAL:HG13	6:b:129:VAL:O	2.19	0.42
6:c:145:ASP:OD1	6:c:148:LEU:HD22	2.20	0.42
6:e:106:PHE:CE1	6:e:300:VAL:HG22	2.54	0.42
6:f:137:TYR:HD1	6:f:310:MET:HE3	1.85	0.42
6:n:127:PHE:HD1	6:n:155:LEU:HD21	1.84	0.42
6:n:129:VAL:O	6:n:129:VAL:HG13	2.20	0.42
6:r:129:VAL:HG13	6:r:129:VAL:O	2.19	0.42
1:T:168:LYS:CD	1:T:196:ILE:HD11	2.49	0.42
2:1:682:GLU:O	2:1:683:ILE:HD13	2.20	0.42
2:1:695:ASN:OD1	2:1:695:ASN:N	2.52	0.42
1:U:87:ASN:ND2	3:4:104:GLN:OE1	2.48	0.42
1:U:133:ASP:OD1	1:U:134:SER:N	2.53	0.42
2:3:507:LEU:HD22	2:3:646:MET:HE3	2.01	0.42
4:7:136:ALA:HB3	4:7:186:LEU:C	2.45	0.42
5:C:255:ARG:NH2	6:b:64:ASP:O	2.53	0.42
5:D:251:PRO:HB3	5:D:319:THR:HG21	2.02	0.42
5:J:367:TRP:CE3	5:J:370:LEU:HD23	2.55	0.42
5:M:99:GLU:HG2	5:M:388:VAL:HG21	2.01	0.42
5:O:194:ASP:OD2	5:O:197:CYS:N	2.52	0.42
5:R:233:SER:HB3	5:R:253:ILE:HD11	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:k:139:LEU:HD21	6:k:306:ILE:HG21	2.02	0.42
6:l:170:ILE:HD11	6:l:246:ILE:HG21	2.02	0.42
6:r:102:LEU:HD23	6:r:105:MET:HE3	2.02	0.42
1:T:22:ILE:HD11	1:U:22:ILE:HD13	2.01	0.41
1:U:138:LYS:CD	1:U:158:THR:HG22	2.50	0.41
2:3:526:MET:HE2	2:3:548:MET:HE3	2.02	0.41
3:4:112:TRP:CZ2	4:5:48:THR:HG21	2.55	0.41
5:D:113:SER:O	5:D:116:LEU:N	2.53	0.41
5:G:125:LYS:NZ	5:H:20:GLU:OE2	2.29	0.41
5:K:254:LEU:HD12	5:K:294:LEU:HD11	2.00	0.41
5:M:240:SER:OG	5:M:241:ALA:N	2.52	0.41
5:R:168:ARG:NE	5:R:175:ASN:OD1	2.53	0.41
5:R:334:VAL:O	5:R:374:TYR:OH	2.38	0.41
6:b:192:THR:HG23	6:b:192:THR:O	2.20	0.41
6:f:233:ILE:HG22	6:f:234:ASN:O	2.20	0.41
6:g:150:LEU:HD23	6:h:290:LYS:HB2	2.03	0.41
2:1:250:ASP:N	2:2:307:ASP:OD2	2.50	0.41
4:5:87:GLU:OE1	4:5:110:THR:HG23	2.20	0.41
5:D:307:LEU:HD11	5:E:248:PHE:CG	2.55	0.41
5:O:294:LEU:HD12	5:O:320:VAL:HG13	2.02	0.41
5:P:40:THR:HG21	5:P:127:ILE:HG13	2.02	0.41
5:Q:134:GLU:OE1	5:Q:134:GLU:N	2.51	0.41
5:Q:258:ASN:ND2	5:Q:293:GLN:OE1	2.49	0.41
6:b:229:VAL:HG21	6:b:235:HIS:CD2	2.55	0.41
6:d:230:VAL:CG1	6:d:233:ILE:HD12	2.51	0.41
6:j:142:MET:HG2	6:j:152:MET:HE2	2.02	0.41
6:n:193:VAL:HG22	6:n:239:LEU:HD13	2.01	0.41
2:3:509:GLN:OE1	6:n:71:THR:HG22	2.20	0.41
2:3:684:ASN:OD1	2:3:685:THR:N	2.53	0.41
4:5:130:GLU:O	4:5:133:ALA:HB3	2.19	0.41
5:A:259:VAL:CG2	5:A:322:LEU:HD11	2.51	0.41
5:B:17:LYS:HB3	5:B:27:VAL:HG12	2.03	0.41
5:B:142:GLN:HB3	5:B:148:THR:HG21	2.03	0.41
5:G:165:THR:HG23	5:G:181:TRP:HH2	1.85	0.41
5:G:254:LEU:HD13	5:H:237:VAL:CG2	2.50	0.41
5:H:7:LEU:HD21	5:H:127:ILE:HG21	2.01	0.41
6:d:129:VAL:O	6:d:129:VAL:HG13	2.20	0.41
6:i:243:THR:OG1	6:i:244:CYS:N	2.53	0.41
6:n:128:SER:OG	6:n:155:LEU:HD22	2.20	0.41
6:n:316:SER:OG	6:p:325:ARG:NH2	2.53	0.41
2:1:352:VAL:HG23	2:1:352:VAL:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:264:LEU:HD23	5:B:269:ILE:HA	2.03	0.41
5:C:333:SER:OG	5:C:379:GLU:OE2	2.33	0.41
5:I:173:HIS:HA	5:I:176:LEU:HD21	2.01	0.41
5:J:248:PHE:CZ	5:L:307:LEU:HD11	2.56	0.41
5:K:40:THR:OG1	5:K:127:ILE:HD11	2.21	0.41
5:M:276:ARG:NH1	5:M:277:PHE:O	2.53	0.41
5:N:113:SER:O	5:N:116:LEU:N	2.53	0.41
5:N:175:ASN:C	5:N:176:LEU:HD12	2.45	0.41
5:O:261:VAL:HG13	5:O:290:LEU:HD23	2.01	0.41
6:c:233:ILE:HG22	6:c:234:ASN:O	2.20	0.41
6:d:240:THR:HG23	6:d:240:THR:O	2.20	0.41
6:h:285:MET:SD	6:h:306:ILE:HD12	2.59	0.41
6:p:146:GLN:OE1	6:p:147:SER:N	2.53	0.41
2:2:293:ILE:HG23	2:2:293:ILE:O	2.20	0.41
5:G:194:ASP:OD1	5:G:196:SER:N	2.52	0.41
5:K:307:LEU:HD11	5:L:248:PHE:CE1	2.55	0.41
5:O:240:SER:OG	5:O:241:ALA:N	2.51	0.41
5:Q:165:THR:HG23	5:Q:181:TRP:HH2	1.86	0.41
6:a:170:ILE:HD11	6:a:246:ILE:HG21	2.02	0.41
6:h:162:GLU:OE1	6:h:313:ARG:NH2	2.53	0.41
6:h:239:LEU:HD23	6:h:239:LEU:O	2.19	0.41
6:i:124:ILE:HD12	6:i:124:ILE:H	1.86	0.41
6:j:160:LEU:O	6:j:255:ARG:N	2.49	0.41
2:2:384:TYR:HB3	2:2:396:ILE:HG22	2.02	0.41
4:5:136:ALA:HB3	4:5:186:LEU:HB2	2.02	0.41
3:6:19:LEU:HD12	3:6:20:LYS:N	2.35	0.41
5:B:242:ASP:OD2	5:B:244:ALA:N	2.52	0.41
5:C:158:PHE:HE1	5:C:326:ILE:HD12	1.85	0.41
5:C:224:THR:HG22	5:C:226:LEU:HD13	2.03	0.41
5:G:11:LEU:CD1	5:G:92:ILE:HD12	2.46	0.41
6:q:220:GLU:OE1	6:q:220:GLU:N	2.53	0.41
1:T:129:ASN:OD1	1:T:130:VAL:N	2.53	0.41
2:3:303:ASN:OD1	2:3:312:THR:HG22	2.21	0.41
2:3:501:ARG:NH1	2:3:661:GLU:OE2	2.52	0.41
3:4:13:VAL:HG22	3:4:19:LEU:CD2	2.50	0.41
3:4:50:SER:HB2	3:4:71:ILE:HG21	2.03	0.41
5:B:186:SER:O	5:B:186:SER:OG	2.38	0.41
5:B:194:ASP:OD1	5:B:197:CYS:N	2.54	0.41
5:C:254:LEU:HD12	5:C:318:ALA:HB1	2.03	0.41
5:H:172:MET:HE2	6:h:256:GLU:HG2	2.02	0.41
6:e:189:SER:O	6:e:223:LYS:NZ	2.48	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:g:185:ILE:HD13	6:g:249:CYS:SG	2.61	0.41
6:k:139:LEU:HD22	6:k:261:ILE:CD1	2.51	0.41
6:k:261:ILE:HG13	6:k:285:MET:HE2	2.01	0.41
3:6:39:ARG:NH2	3:6:90:GLU:O	2.54	0.41
5:J:11:LEU:HD11	5:J:88:PHE:CB	2.51	0.41
5:K:310:GLN:OE1	5:K:310:GLN:N	2.53	0.41
5:O:183:ASN:OD1	5:O:188:ILE:HG23	2.21	0.41
5:Q:29:ASP:OD1	5:Q:30:LEU:N	2.54	0.41
5:Q:131:ASN:N	5:R:16:ASP:OD2	2.47	0.41
6:b:51:GLN:OE1	6:f:59:ILE:HD13	2.21	0.41
6:j:256:GLU:N	6:j:256:GLU:OE2	2.54	0.41
6:p:142:MET:SD	6:p:260:VAL:HG13	2.61	0.41
1:T:159:SER:HA	1:T:180:THR:HG22	2.02	0.41
2:3:374:LEU:HD12	2:3:427:PHE:CE2	2.56	0.41
3:6:42:PRO:O	3:6:44:LYS:N	2.54	0.41
5:C:370:LEU:HD21	5:C:382:LEU:HD22	2.02	0.41
5:G:219:THR:N	5:G:330:VAL:O	2.48	0.41
5:G:246:THR:HG21	5:I:297:PRO:HB3	2.02	0.41
5:J:321:GLY:C	5:J:322:LEU:HD12	2.46	0.41
5:K:264:LEU:HD23	5:K:269:ILE:HA	2.02	0.41
5:K:304:VAL:HG12	5:L:237:VAL:CG1	2.49	0.41
5:N:218:LEU:HD23	5:N:329:ALA:CB	2.51	0.41
5:Q:99:GLU:CG	5:Q:388:VAL:HG21	2.50	0.41
5:Q:162:ALA:HB2	5:Q:182:LEU:HD12	2.03	0.41
5:Q:211:ILE:HD12	5:Q:289:ARG:HG2	2.01	0.41
5:R:259:VAL:HG23	5:R:322:LEU:HD11	2.03	0.41
6:a:170:ILE:HG23	6:a:237:ILE:HB	2.03	0.41
6:b:210:THR:OG1	6:b:211:ASN:N	2.54	0.41
6:e:141:LEU:HD11	6:e:299:ILE:HD11	2.03	0.41
6:j:142:MET:HE1	6:j:269:LEU:HD21	2.03	0.41
6:p:325:ARG:O	6:p:326:VAL:HG12	2.20	0.41
8:p:401:NAG:O7	8:p:401:NAG:O3	2.36	0.41
6:q:158:LEU:HD13	6:q:224:LEU:HD11	2.02	0.41
6:q:218:VAL:HG12	6:r:293:TRP:HB3	2.02	0.41
6:r:268:ILE:HG13	6:r:269:LEU:HD12	2.03	0.41
2:2:507:LEU:HD21	2:2:646:MET:SD	2.61	0.41
5:A:40:THR:CG2	5:A:127:ILE:HD11	2.51	0.41
5:B:194:ASP:OD1	5:B:196:SER:N	2.54	0.41
5:C:37:MET:HE3	5:C:88:PHE:CZ	2.56	0.41
5:C:254:LEU:HD13	5:C:294:LEU:HD11	2.02	0.41
5:E:153:HIS:ND1	5:F:338:ALA:O	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:158:PHE:CE2	5:E:188:ILE:HD11	2.56	0.41
5:H:194:ASP:OD2	5:H:196:SER:N	2.52	0.41
5:P:248:PHE:CZ	5:R:307:LEU:HD21	2.55	0.41
5:Q:270:ILE:HG21	5:Q:284:ASN:O	2.21	0.41
6:a:192:THR:HG23	6:a:192:THR:O	2.20	0.41
6:c:95:ASP:OD1	6:c:96:GLY:N	2.53	0.41
6:d:59:ILE:HD13	6:i:51:GLN:HB3	2.02	0.41
6:g:81:LEU:HD11	6:g:139:LEU:HD13	2.03	0.41
6:r:257:ASN:ND2	6:r:314:SER:O	2.54	0.41
2:1:373:ASN:C	2:1:374:LEU:HD22	2.46	0.40
3:4:73:ARG:NH2	3:4:75:ASN:OD1	2.53	0.40
5:D:253:ILE:HD13	5:D:319:THR:OG1	2.21	0.40
5:D:254:LEU:HD12	5:D:318:ALA:HB1	2.02	0.40
5:E:17:LYS:HD2	5:E:30:LEU:HD22	2.04	0.40
5:E:261:VAL:HG13	5:E:290:LEU:HD23	2.02	0.40
5:G:304:VAL:HG23	5:H:237:VAL:HG11	2.03	0.40
5:J:279:THR:HG21	5:K:156:ASN:HD22	1.85	0.40
5:K:40:THR:CB	5:K:127:ILE:HD11	2.51	0.40
5:P:74:ASP:N	5:P:74:ASP:OD1	2.53	0.40
5:P:127:ILE:O	5:P:127:ILE:HG22	2.21	0.40
6:d:202:THR:O	6:d:202:THR:HG22	2.21	0.40
6:f:256:GLU:OE1	6:f:256:GLU:N	2.53	0.40
6:n:233:ILE:HD12	6:o:108:ILE:HD13	2.03	0.40
2:1:511:ILE:HG22	2:1:512:ALA:O	2.21	0.40
2:2:540:THR:HG22	2:2:540:THR:O	2.21	0.40
2:3:344:ILE:HD12	2:3:429:LEU:HD21	2.03	0.40
5:A:253:ILE:HG13	5:B:234:PHE:CE2	2.56	0.40
5:D:4:LEU:HD22	5:D:116:LEU:CD1	2.51	0.40
5:D:36:GLN:NE2	5:E:23:LEU:HD21	2.36	0.40
5:D:276:ARG:NH1	5:D:280:ILE:HG22	2.36	0.40
5:M:253:ILE:HD12	5:N:234:PHE:CE1	2.56	0.40
6:b:170:ILE:HD11	6:b:237:ILE:C	2.45	0.40
6:c:100:ASP:O	6:c:103:SER:OG	2.26	0.40
6:g:314:SER:OG	6:g:314:SER:O	2.38	0.40
6:l:200:THR:OG1	6:l:201:GLN:OE1	2.22	0.40
6:n:166:ASN:OD1	6:p:315:ARG:NH2	2.47	0.40
6:n:324:TYR:HB3	6:p:317:LEU:HD12	2.02	0.40
1:U:141:GLU:O	1:U:155:ARG:NH2	2.54	0.40
4:7:171:SER:O	4:7:173:GLN:NE2	2.54	0.40
5:A:342:LEU:HD12	5:A:389:ALA:HB3	2.04	0.40
5:B:387:THR:O	5:B:390:SER:OG	2.24	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:F:89:ILE:CD1	5:F:395:LEU:HD12	2.52	0.40
5:F:113:SER:OG	5:F:113:SER:O	2.37	0.40
5:M:370:LEU:HD23	5:M:370:LEU:O	2.22	0.40
5:Q:134:GLU:O	5:Q:138:ASN:ND2	2.45	0.40
6:g:53:TYR:C	6:k:59:ILE:HD11	2.47	0.40
6:h:186:SER:OG	6:h:191:CYS:SG	2.78	0.40
6:k:71:THR:HG22	6:k:71:THR:O	2.21	0.40
6:m:124:ILE:HD13	6:m:142:MET:HG3	2.04	0.40
2:2:352:VAL:HG23	2:2:352:VAL:O	2.21	0.40
5:B:213:GLN:HA	5:B:287:THR:HG22	2.03	0.40
5:D:98:ASP:OD2	5:D:102:ARG:NH2	2.55	0.40
5:D:360:VAL:HG23	5:D:361:PHE:N	2.36	0.40
5:I:127:ILE:O	5:I:127:ILE:HG22	2.21	0.40
5:J:214:LEU:HD11	5:J:285:PHE:CE2	2.57	0.40
5:J:246:THR:O	5:L:301:THR:OG1	2.33	0.40
5:O:216:ARG:NH1	5:O:333:SER:OG	2.54	0.40
5:Q:209:GLU:OE2	5:Q:289:ARG:NH2	2.55	0.40
5:Q:299:ASN:OD1	6:r:71:THR:HG22	2.22	0.40
6:b:102:LEU:HD13	6:b:118:PHE:HE2	1.87	0.40
6:c:218:VAL:HG12	6:c:218:VAL:O	2.21	0.40
6:g:285:MET:HE1	6:g:306:ILE:CG2	2.51	0.40
6:h:263:VAL:HG12	6:h:289:TRP:HB3	2.04	0.40
6:l:129:VAL:O	6:l:129:VAL:HG13	2.21	0.40
6:l:282:GLU:HB3	6:l:309:VAL:HG13	2.03	0.40
6:n:79:SER:O	6:n:115:SER:OG	2.38	0.40
6:o:271:ILE:H	6:o:271:ILE:HD12	1.87	0.40
6:q:186:SER:OG	6:q:244:CYS:SG	2.80	0.40
2:3:653:ASN:O	2:3:654:THR:HG22	2.22	0.40
5:G:173:HIS:HA	5:G:176:LEU:HD11	2.03	0.40
5:G:254:LEU:HD13	5:H:237:VAL:HG23	2.02	0.40
5:K:158:PHE:HE2	5:K:214:LEU:HD21	1.85	0.40
5:L:342:LEU:HD12	5:L:389:ALA:HB3	2.03	0.40
5:P:259:VAL:CG2	5:P:322:LEU:HD11	2.52	0.40
6:e:169:ASP:OD2	6:e:172:LEU:HD13	2.21	0.40
6:f:152:MET:HE2	6:f:269:LEU:HD21	2.03	0.40
6:j:139:LEU:HD21	6:j:306:ILE:CG2	2.52	0.40
6:j:230:VAL:HG23	6:k:105:MET:HE2	2.02	0.40
6:j:261:ILE:HD13	6:j:299:ILE:CD1	2.52	0.40
6:k:69:ASN:OD1	6:k:70:SER:N	2.47	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 PDB entries followed by that with respect to all EM entries.

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	T	225/230 (98%)	202 (90%)	23 (10%)	0	100	100
1	U	225/230 (98%)	200 (89%)	25 (11%)	0	100	100
1	V	23/230 (10%)	23 (100%)	0	0	100	100
2	1	495/529 (94%)	444 (90%)	50 (10%)	1 (0%)	43	74
2	2	495/529 (94%)	452 (91%)	40 (8%)	3 (1%)	21	54
2	3	482/529 (91%)	438 (91%)	44 (9%)	0	100	100
3	4	228/236 (97%)	204 (90%)	20 (9%)	4 (2%)	6	34
3	6	228/236 (97%)	204 (90%)	22 (10%)	2 (1%)	14	47
4	5	215/217 (99%)	196 (91%)	18 (8%)	1 (0%)	24	57
4	7	215/217 (99%)	193 (90%)	21 (10%)	1 (0%)	24	57
5	A	395/397 (100%)	370 (94%)	25 (6%)	0	100	100
5	B	395/397 (100%)	365 (92%)	29 (7%)	1 (0%)	36	67
5	C	395/397 (100%)	373 (94%)	21 (5%)	1 (0%)	36	67
5	D	395/397 (100%)	374 (95%)	21 (5%)	0	100	100
5	E	395/397 (100%)	364 (92%)	31 (8%)	0	100	100
5	F	395/397 (100%)	366 (93%)	29 (7%)	0	100	100
5	G	395/397 (100%)	372 (94%)	22 (6%)	1 (0%)	36	67
5	H	395/397 (100%)	374 (95%)	21 (5%)	0	100	100
5	I	395/397 (100%)	372 (94%)	23 (6%)	0	100	100
5	J	395/397 (100%)	372 (94%)	23 (6%)	0	100	100
5	K	395/397 (100%)	369 (93%)	26 (7%)	0	100	100
5	L	395/397 (100%)	376 (95%)	19 (5%)	0	100	100
5	M	395/397 (100%)	367 (93%)	28 (7%)	0	100	100
5	N	395/397 (100%)	369 (93%)	26 (7%)	0	100	100
5	O	395/397 (100%)	372 (94%)	23 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
5	P	395/397 (100%)	366 (93%)	29 (7%)	0	100	100
5	Q	395/397 (100%)	372 (94%)	23 (6%)	0	100	100
5	R	395/397 (100%)	364 (92%)	31 (8%)	0	100	100
6	a	269/326 (82%)	254 (94%)	15 (6%)	0	100	100
6	b	274/326 (84%)	248 (90%)	26 (10%)	0	100	100
6	c	258/326 (79%)	239 (93%)	19 (7%)	0	100	100
6	d	274/326 (84%)	254 (93%)	19 (7%)	1 (0%)	30	62
6	e	259/326 (79%)	243 (94%)	16 (6%)	0	100	100
6	f	274/326 (84%)	253 (92%)	21 (8%)	0	100	100
6	g	274/326 (84%)	257 (94%)	17 (6%)	0	100	100
6	h	269/326 (82%)	255 (95%)	14 (5%)	0	100	100
6	i	274/326 (84%)	250 (91%)	24 (9%)	0	100	100
6	j	272/326 (83%)	253 (93%)	19 (7%)	0	100	100
6	k	274/326 (84%)	256 (93%)	18 (7%)	0	100	100
6	l	274/326 (84%)	255 (93%)	19 (7%)	0	100	100
6	m	273/326 (84%)	256 (94%)	17 (6%)	0	100	100
6	n	274/326 (84%)	250 (91%)	24 (9%)	0	100	100
6	o	274/326 (84%)	252 (92%)	22 (8%)	0	100	100
6	p	273/326 (84%)	249 (91%)	24 (9%)	0	100	100
6	q	269/326 (82%)	248 (92%)	21 (8%)	0	100	100
6	r	273/326 (84%)	241 (88%)	32 (12%)	0	100	100
All	All	14822/16197 (92%)	13726 (93%)	1080 (7%)	16 (0%)	49	79

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	4	181	ALA
3	6	181	ALA
4	7	157	ASP
2	1	632	PHE
3	4	180	PRO
3	6	180	PRO
2	2	488	THR
2	2	632	PHE
4	5	157	ASP

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Mol	Chain	Res	Type
5	C	114	GLU
2	2	722	ASP
5	B	114	GLU
5	G	114	GLU
3	4	43	GLY
3	4	42	PRO
6	d	96	GLY

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	T	208/211 (99%)	208 (100%)	0	100	100
1	U	208/211 (99%)	208 (100%)	0	100	100
1	V	23/211 (11%)	23 (100%)	0	100	100
2	1	448/477 (94%)	448 (100%)	0	100	100
2	2	448/477 (94%)	448 (100%)	0	100	100
2	3	436/477 (91%)	436 (100%)	0	100	100
3	4	193/199 (97%)	193 (100%)	0	100	100
3	6	193/199 (97%)	193 (100%)	0	100	100
4	5	187/187 (100%)	187 (100%)	0	100	100
4	7	187/187 (100%)	187 (100%)	0	100	100
5	A	348/348 (100%)	348 (100%)	0	100	100
5	B	348/348 (100%)	348 (100%)	0	100	100
5	C	348/348 (100%)	348 (100%)	0	100	100
5	D	348/348 (100%)	348 (100%)	0	100	100
5	E	348/348 (100%)	348 (100%)	0	100	100
5	F	348/348 (100%)	348 (100%)	0	100	100
5	G	348/348 (100%)	348 (100%)	0	100	100
5	H	348/348 (100%)	348 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
5	I	348/348 (100%)	348 (100%)	0	100	100
5	J	348/348 (100%)	348 (100%)	0	100	100
5	K	348/348 (100%)	348 (100%)	0	100	100
5	L	348/348 (100%)	348 (100%)	0	100	100
5	M	348/348 (100%)	348 (100%)	0	100	100
5	N	348/348 (100%)	348 (100%)	0	100	100
5	O	348/348 (100%)	348 (100%)	0	100	100
5	P	348/348 (100%)	347 (100%)	1 (0%)	86	83
5	Q	348/348 (100%)	348 (100%)	0	100	100
5	R	348/348 (100%)	347 (100%)	1 (0%)	86	83
6	a	250/300 (83%)	250 (100%)	0	100	100
6	b	254/300 (85%)	254 (100%)	0	100	100
6	c	241/300 (80%)	240 (100%)	1 (0%)	84	81
6	d	254/300 (85%)	254 (100%)	0	100	100
6	e	241/300 (80%)	241 (100%)	0	100	100
6	f	254/300 (85%)	254 (100%)	0	100	100
6	g	254/300 (85%)	253 (100%)	1 (0%)	84	81
6	h	250/300 (83%)	250 (100%)	0	100	100
6	i	254/300 (85%)	254 (100%)	0	100	100
6	j	252/300 (84%)	251 (100%)	1 (0%)	84	81
6	k	254/300 (85%)	254 (100%)	0	100	100
6	l	254/300 (85%)	254 (100%)	0	100	100
6	m	253/300 (84%)	252 (100%)	1 (0%)	84	81
6	n	254/300 (85%)	253 (100%)	1 (0%)	84	81
6	o	254/300 (85%)	254 (100%)	0	100	100
6	p	253/300 (84%)	252 (100%)	1 (0%)	84	81
6	q	250/300 (83%)	250 (100%)	0	100	100
6	r	253/300 (84%)	253 (100%)	0	100	100
All	All	13324/14500 (92%)	13316 (100%)	8 (0%)	87	87

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
5	P	339	ASN
5	R	305	ASN
6	c	238	ASN
6	g	238	ASN
6	j	69	ASN
6	m	238	ASN
6	n	238	ASN
6	p	238	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (99) such sidechains are listed below:

Mol	Chain	Res	Type
2	1	385	HIS
2	1	730	ASN
1	U	135	ASN
2	2	480	GLN
2	2	742	ASN
2	2	757	GLN
2	3	445	ASN
2	3	753	ASN
2	3	758	ASN
3	4	213	HIS
3	6	110	ASN
3	6	210	ASN
3	6	212	ASN
4	7	55	GLN
4	7	93	GLN
5	A	173	HIS
5	B	204	ASN
5	C	173	HIS
5	C	339	ASN
5	C	345	ASN
5	D	72	ASN
5	D	112	GLN
5	D	131	ASN
5	D	140	ASN
5	D	153	HIS
5	D	206	GLN
5	D	266	ASN
5	D	274	GLN
5	D	345	ASN
5	E	173	HIS
5	E	183	ASN

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Mol	Chain	Res	Type
5	E	316	HIS
5	F	140	ASN
5	F	183	ASN
5	F	204	ASN
5	F	206	GLN
5	H	26	ASN
5	H	72	ASN
5	H	173	HIS
5	H	317	HIS
5	I	42	ASN
5	I	268	GLN
5	I	305	ASN
5	J	140	ASN
5	J	183	ASN
5	J	189	GLN
5	J	200	ASN
5	J	316	HIS
5	J	345	ASN
5	J	383	GLN
5	K	239	ASN
5	K	250	ASN
5	K	373	ASN
5	L	26	ASN
5	M	167	ASN
5	M	183	ASN
5	M	210	HIS
5	O	36	GLN
5	O	105	GLN
5	O	204	ASN
5	O	206	GLN
5	O	207	GLN
5	O	293	GLN
5	P	183	ASN
5	P	239	ASN
5	P	345	ASN
5	Q	146	GLN
5	Q	173	HIS
5	Q	183	ASN
5	Q	210	HIS
5	R	105	GLN
5	R	131	ASN
5	R	140	ASN

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Mol	Chain	Res	Type
5	R	146	GLN
5	R	305	ASN
5	R	383	GLN
6	a	161	ASN
6	a	166	ASN
6	b	235	HIS
6	b	262	GLN
6	b	294	GLN
6	c	267	ASN
6	d	288	ASN
6	e	294	GLN
6	f	56	ASN
6	g	308	GLN
6	i	56	ASN
6	i	234	ASN
6	k	51	GLN
6	k	235	HIS
6	k	262	GLN
6	k	308	GLN
6	l	52	ASN
6	l	176	GLN
6	l	267	ASN
6	l	305	GLN
6	n	294	GLN
6	p	234	ASN
6	p	257	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

3 monosaccharides 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	$\# Z > 2$
7	NAG	S	1	7,6	14,14,15	0.61	0	17,19,21	1.12	1 (5%)
7	NAG	S	2	7	14,14,15	0.24	0	17,19,21	0.39	0
7	BMA	S	3	7	11,11,12	0.64	0	15,15,17	0.68	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
7	NAG	S	1	7,6	-	2/6/23/26	0/1/1/1
7	NAG	S	2	7	-	2/6/23/26	0/1/1/1
7	BMA	S	3	7	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	S	1	NAG	C1-O5-C5	3.66	117.09	112.19

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	S	1	NAG	C4-C5-C6-O6
7	S	2	NAG	O5-C5-C6-O6
7	S	1	NAG	O5-C5-C6-O6
7	S	2	NAG	C4-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 1 short contact:

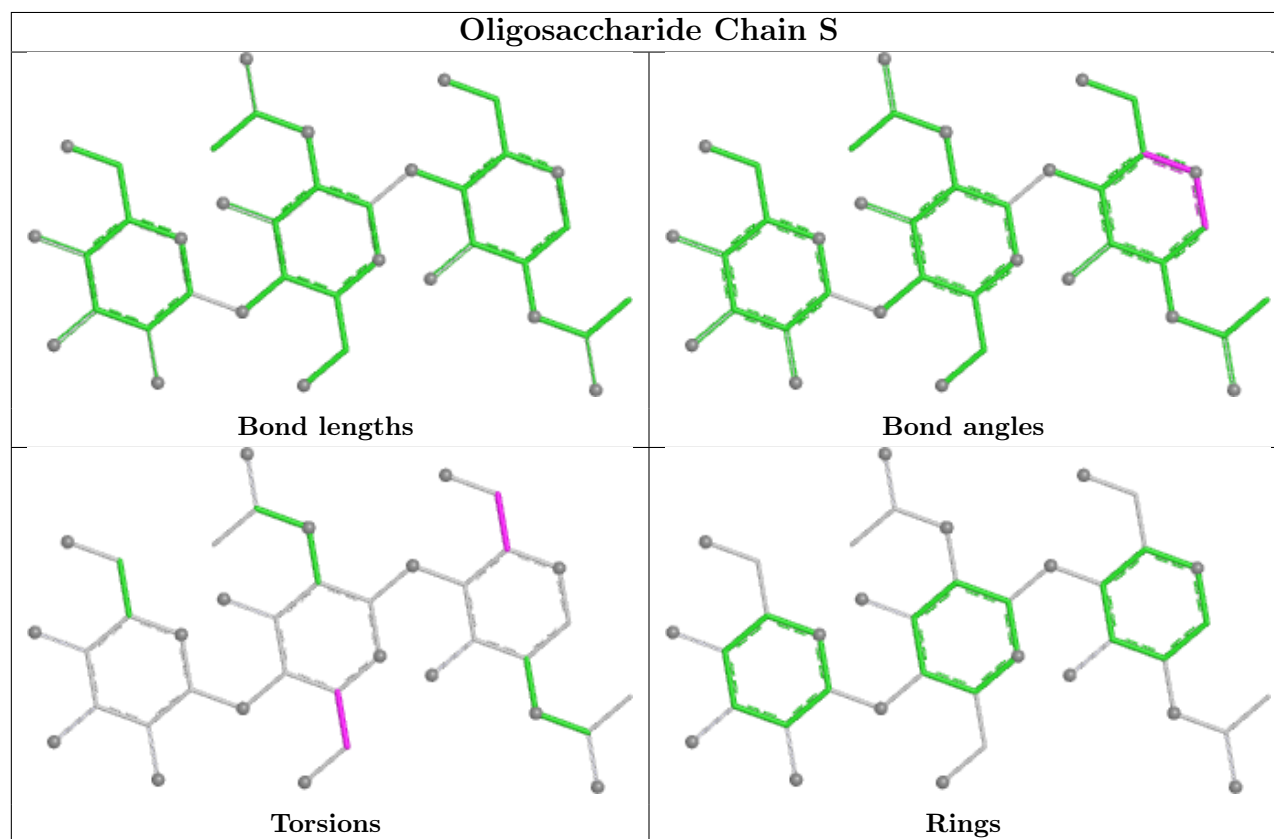
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	S	1	NAG	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	S	2	NAG	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



5.6 Ligand geometry [i](#)

Of 107 ligands modelled in this entry, 72 are monoatomic - leaving 35 for Mogul analysis.

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	$\# Z > 2$
8	NAG	e	402	6	14,14,15	0.34	0	17,19,21	0.39	0
8	NAG	h	402	6	14,14,15	0.22	0	17,19,21	0.41	0
8	NAG	o	503	6	14,14,15	0.42	0	17,19,21	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	NAG	q	402	6	14,14,15	0.33	0	17,19,21	0.41	0
8	NAG	j	402	6	14,14,15	0.44	0	17,19,21	0.43	0
8	NAG	h	401	6	14,14,15	0.24	0	17,19,21	0.45	0
8	NAG	b	402	6	14,14,15	0.27	0	17,19,21	0.42	0
8	NAG	c	503	6	14,14,15	0.18	0	17,19,21	0.49	0
8	NAG	l	503	6	14,14,15	0.30	0	17,19,21	0.47	0
8	NAG	l	504	6	14,14,15	0.35	0	17,19,21	0.38	0
8	NAG	b	401	6	14,14,15	0.38	0	17,19,21	0.42	0
8	NAG	d	402	6	14,14,15	0.52	0	17,19,21	0.50	0
8	NAG	m	401	6	14,14,15	0.21	0	17,19,21	0.41	0
8	NAG	f	503	6	14,14,15	0.27	0	17,19,21	0.46	0
8	NAG	q	401	6	14,14,15	0.19	0	17,19,21	0.44	0
8	NAG	p	402	6	14,14,15	0.34	0	17,19,21	0.43	0
8	NAG	g	402	6	14,14,15	0.30	0	17,19,21	0.39	0
8	NAG	g	401	6	14,14,15	0.31	0	17,19,21	1.58	2 (11%)
8	NAG	p	401	6	14,14,15	0.63	1 (7%)	17,19,21	0.51	0
8	NAG	c	504	6	14,14,15	0.29	0	17,19,21	0.37	0
8	NAG	o	504	6	14,14,15	0.42	0	17,19,21	0.60	0
8	NAG	k	402	6	14,14,15	0.43	0	17,19,21	0.57	0
8	NAG	n	402	6	14,14,15	0.20	0	17,19,21	0.39	0
8	NAG	r	503	6	14,14,15	0.28	0	17,19,21	0.40	0
8	NAG	j	401	6	14,14,15	0.28	0	17,19,21	0.51	0
8	NAG	i	504	6	14,14,15	0.24	0	17,19,21	0.38	0
8	NAG	f	504	6	14,14,15	0.55	0	17,19,21	0.52	0
8	NAG	d	401	6	14,14,15	0.24	0	17,19,21	0.41	0
8	NAG	i	503	6	14,14,15	0.20	0	17,19,21	0.34	0
8	NAG	a	401	6	14,14,15	0.26	0	17,19,21	0.55	0
8	NAG	k	401	6	14,14,15	0.30	0	17,19,21	0.67	1 (5%)
8	NAG	n	401	6	14,14,15	0.38	0	17,19,21	0.53	0
8	NAG	m	402	6	14,14,15	0.25	0	17,19,21	0.42	0
8	NAG	r	504	6	14,14,15	0.38	0	17,19,21	0.45	0
8	NAG	e	401	6	14,14,15	0.22	0	17,19,21	0.50	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
8	NAG	e	402	6	-	2/6/23/26	0/1/1/1
8	NAG	h	402	6	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	NAG	o	503	6	-	2/6/23/26	0/1/1/1
8	NAG	q	402	6	-	0/6/23/26	0/1/1/1
8	NAG	j	402	6	-	2/6/23/26	0/1/1/1
8	NAG	h	401	6	-	2/6/23/26	0/1/1/1
8	NAG	b	402	6	-	0/6/23/26	0/1/1/1
8	NAG	c	503	6	-	2/6/23/26	0/1/1/1
8	NAG	l	503	6	-	2/6/23/26	0/1/1/1
8	NAG	l	504	6	-	0/6/23/26	0/1/1/1
8	NAG	b	401	6	-	3/6/23/26	0/1/1/1
8	NAG	d	402	6	-	0/6/23/26	0/1/1/1
8	NAG	m	401	6	-	2/6/23/26	0/1/1/1
8	NAG	f	503	6	-	2/6/23/26	0/1/1/1
8	NAG	q	401	6	-	2/6/23/26	0/1/1/1
8	NAG	p	402	6	-	2/6/23/26	0/1/1/1
8	NAG	g	402	6	-	2/6/23/26	0/1/1/1
8	NAG	g	401	6	-	5/6/23/26	0/1/1/1
8	NAG	p	401	6	-	3/6/23/26	0/1/1/1
8	NAG	c	504	6	-	1/6/23/26	0/1/1/1
8	NAG	o	504	6	-	2/6/23/26	0/1/1/1
8	NAG	k	402	6	-	0/6/23/26	0/1/1/1
8	NAG	n	402	6	-	0/6/23/26	0/1/1/1
8	NAG	r	503	6	-	1/6/23/26	0/1/1/1
8	NAG	j	401	6	-	2/6/23/26	0/1/1/1
8	NAG	i	504	6	-	1/6/23/26	0/1/1/1
8	NAG	f	504	6	-	2/6/23/26	0/1/1/1
8	NAG	d	401	6	-	2/6/23/26	0/1/1/1
8	NAG	i	503	6	-	1/6/23/26	0/1/1/1
8	NAG	a	401	6	-	0/6/23/26	0/1/1/1
8	NAG	k	401	6	-	2/6/23/26	0/1/1/1
8	NAG	n	401	6	-	2/6/23/26	0/1/1/1
8	NAG	m	402	6	-	1/6/23/26	0/1/1/1
8	NAG	r	504	6	-	2/6/23/26	0/1/1/1
8	NAG	e	401	6	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	p	401	NAG	C1-C2	2.22	1.55	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	g	401	NAG	C2-N2-C7	4.94	129.52	122.90
8	g	401	NAG	C1-C2-N2	3.28	115.61	110.43
8	k	401	NAG	C1-O5-C5	2.25	115.20	112.19

There are no chirality outliers.

All (56) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	g	401	NAG	C1-C2-N2-C7
8	j	401	NAG	C4-C5-C6-O6
8	e	401	NAG	O5-C5-C6-O6
8	m	401	NAG	O5-C5-C6-O6
8	h	402	NAG	O5-C5-C6-O6
8	n	401	NAG	O5-C5-C6-O6
8	p	402	NAG	O5-C5-C6-O6
8	o	503	NAG	C4-C5-C6-O6
8	j	401	NAG	O5-C5-C6-O6
8	h	401	NAG	O5-C5-C6-O6
8	o	504	NAG	O5-C5-C6-O6
8	h	402	NAG	C4-C5-C6-O6
8	n	401	NAG	C4-C5-C6-O6
8	p	402	NAG	C4-C5-C6-O6
8	d	401	NAG	O5-C5-C6-O6
8	q	401	NAG	O5-C5-C6-O6
8	f	504	NAG	C4-C5-C6-O6
8	g	402	NAG	C4-C5-C6-O6
8	e	401	NAG	C4-C5-C6-O6
8	f	504	NAG	O5-C5-C6-O6
8	g	402	NAG	O5-C5-C6-O6
8	o	503	NAG	O5-C5-C6-O6
8	m	401	NAG	C4-C5-C6-O6
8	h	401	NAG	C4-C5-C6-O6
8	r	504	NAG	O5-C5-C6-O6
8	l	503	NAG	O5-C5-C6-O6
8	e	402	NAG	O5-C5-C6-O6
8	j	402	NAG	O5-C5-C6-O6
8	g	401	NAG	O5-C5-C6-O6
8	g	401	NAG	C8-C7-N2-C2
8	g	401	NAG	O7-C7-N2-C2
8	d	401	NAG	C4-C5-C6-O6
8	f	503	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
8	j	402	NAG	C4-C5-C6-O6
8	p	401	NAG	O5-C5-C6-O6
8	r	504	NAG	C4-C5-C6-O6
8	l	503	NAG	C4-C5-C6-O6
8	o	504	NAG	C4-C5-C6-O6
8	q	401	NAG	C4-C5-C6-O6
8	e	402	NAG	C4-C5-C6-O6
8	i	503	NAG	O5-C5-C6-O6
8	c	503	NAG	C4-C5-C6-O6
8	c	503	NAG	O5-C5-C6-O6
8	i	504	NAG	O5-C5-C6-O6
8	m	402	NAG	O5-C5-C6-O6
8	g	401	NAG	C4-C5-C6-O6
8	r	503	NAG	O5-C5-C6-O6
8	f	503	NAG	C4-C5-C6-O6
8	c	504	NAG	O5-C5-C6-O6
8	b	401	NAG	C1-C2-N2-C7
8	p	401	NAG	C4-C5-C6-O6
8	k	401	NAG	C4-C5-C6-O6
8	b	401	NAG	C3-C2-N2-C7
8	p	401	NAG	C3-C2-N2-C7
8	b	401	NAG	C4-C5-C6-O6
8	k	401	NAG	O5-C5-C6-O6

There are no ring outliers.

9 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	j	402	NAG	1	0
8	l	504	NAG	1	0
8	b	401	NAG	1	0
8	d	402	NAG	1	0
8	p	401	NAG	2	0
8	o	504	NAG	1	0
8	k	402	NAG	1	0
8	f	504	NAG	1	0
8	n	401	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

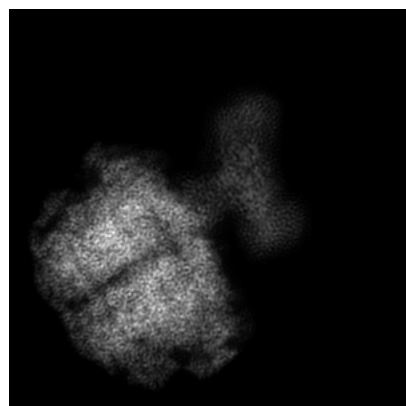
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-26608. These allow visual inspection of the internal detail of the map and identification of artifacts.

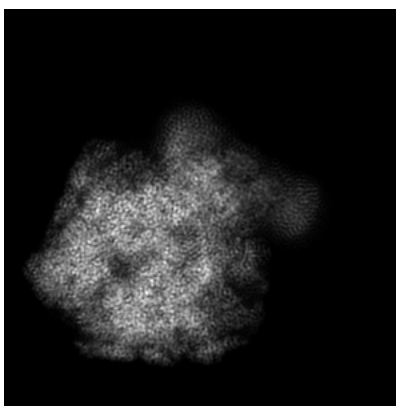
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

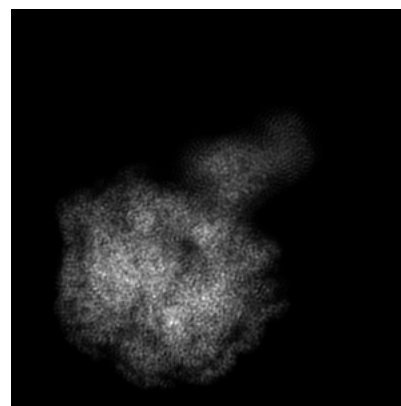
6.1.1 Primary map



X

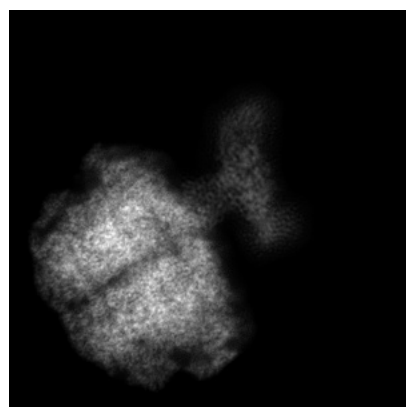


Y

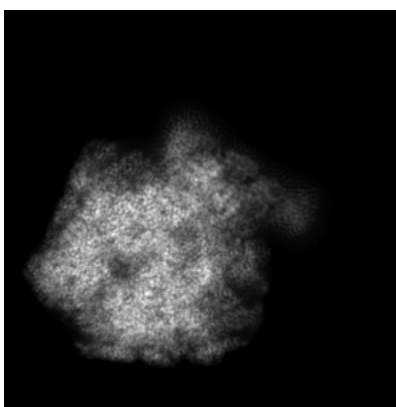


Z

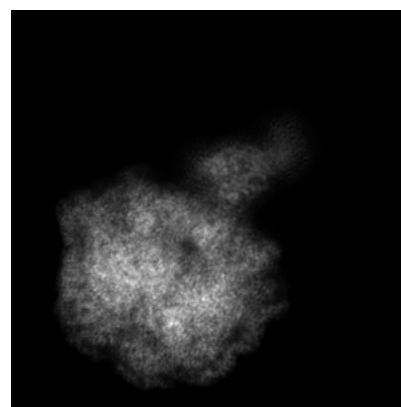
6.1.2 Raw map



X



Y

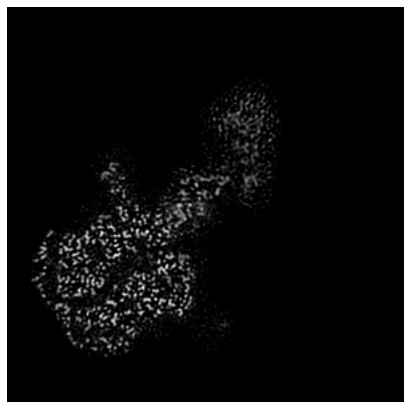


Z

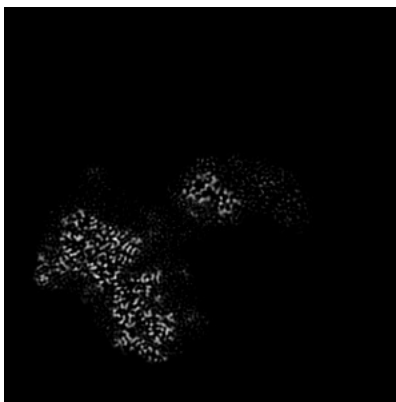
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 160

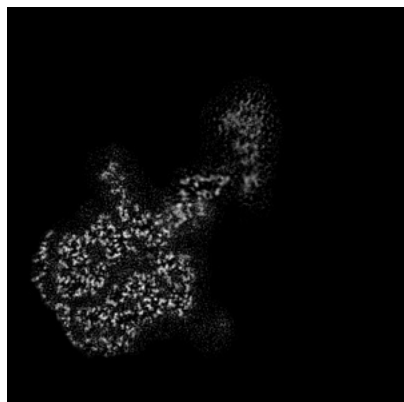


Y Index: 160

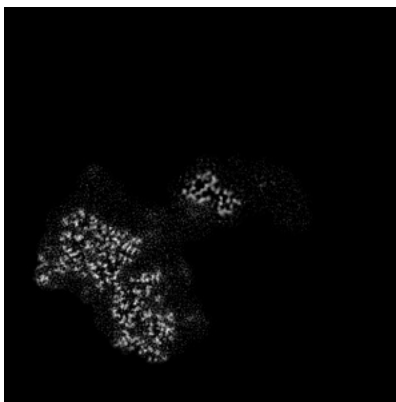


Z Index: 160

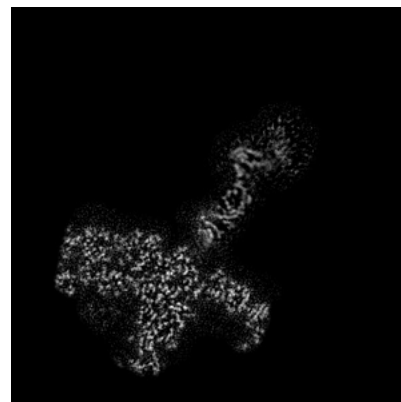
6.2.2 Raw map



X Index: 160



Y Index: 160



Z Index: 160

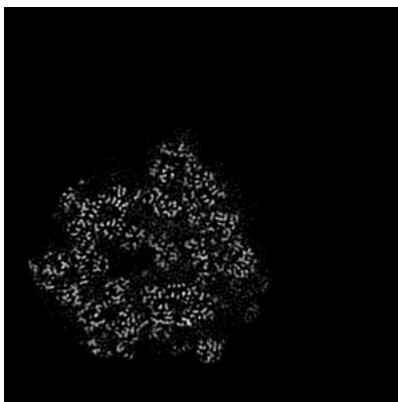
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

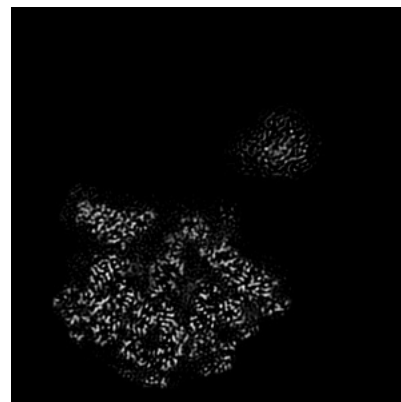
6.3.1 Primary map



X Index: 111

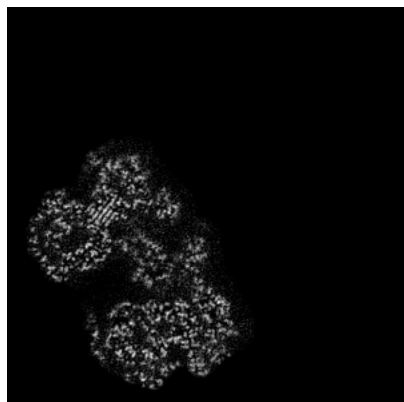


Y Index: 96

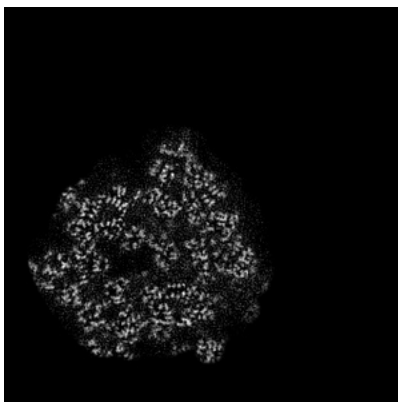


Z Index: 134

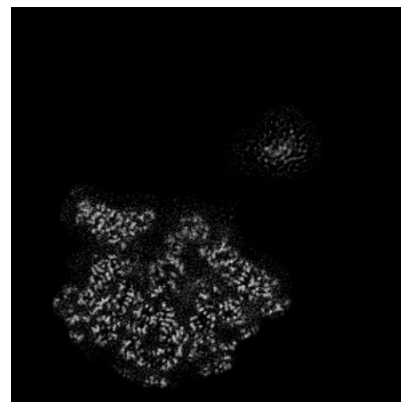
6.3.2 Raw map



X Index: 111



Y Index: 96

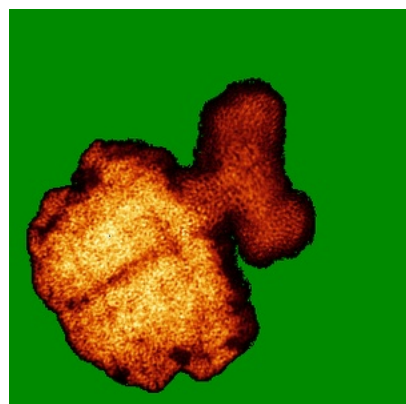


Z Index: 134

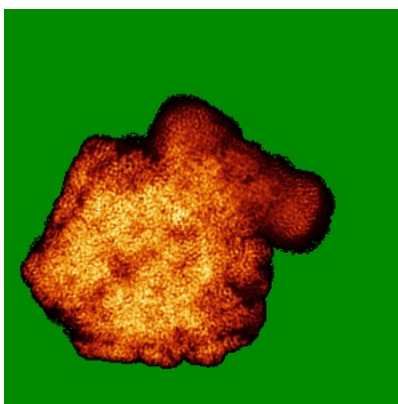
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

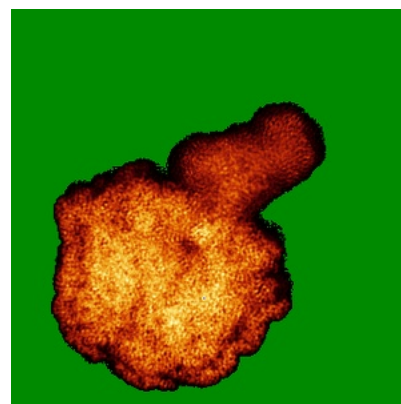
6.4.1 Primary map



X

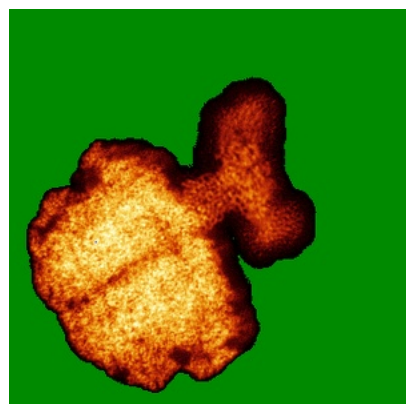


Y

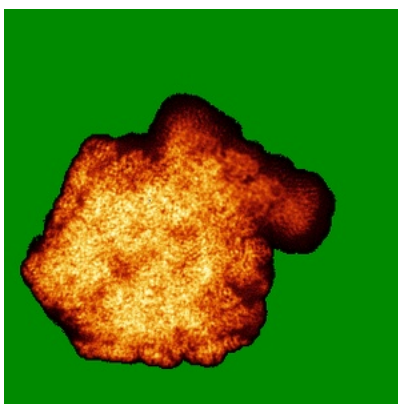


Z

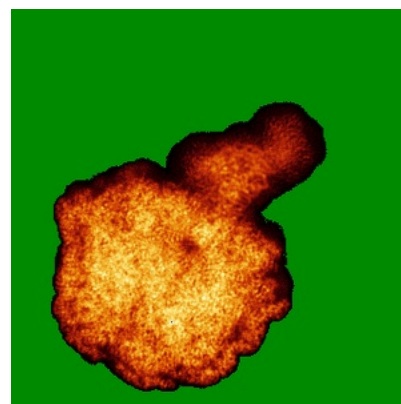
6.4.2 Raw map



X



Y



Z

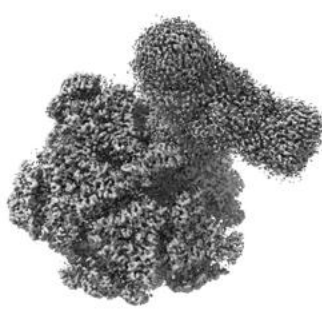
The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



X



Y



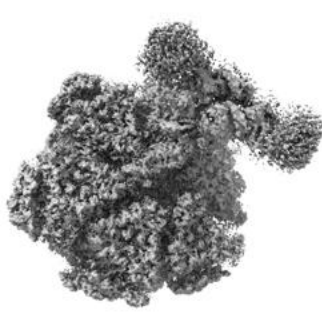
Z

The images above show the 3D surface view of the map at the recommended contour level 0.01. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

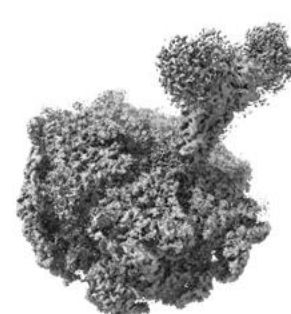
6.5.2 Raw map



X



Y



Z

These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

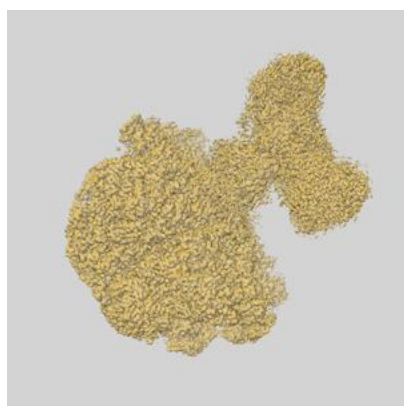
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

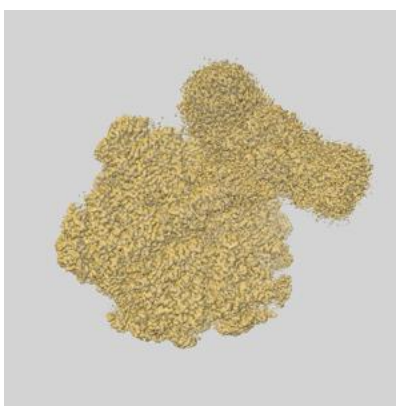
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

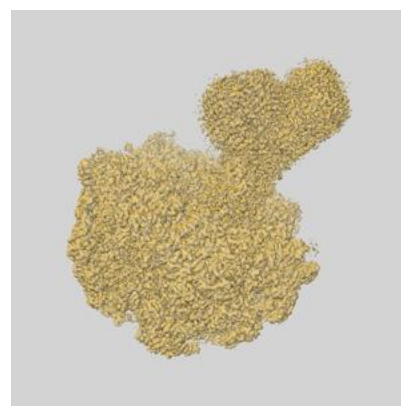
6.6.1 emd_26608_msk_1.map [i](#)



X



Y

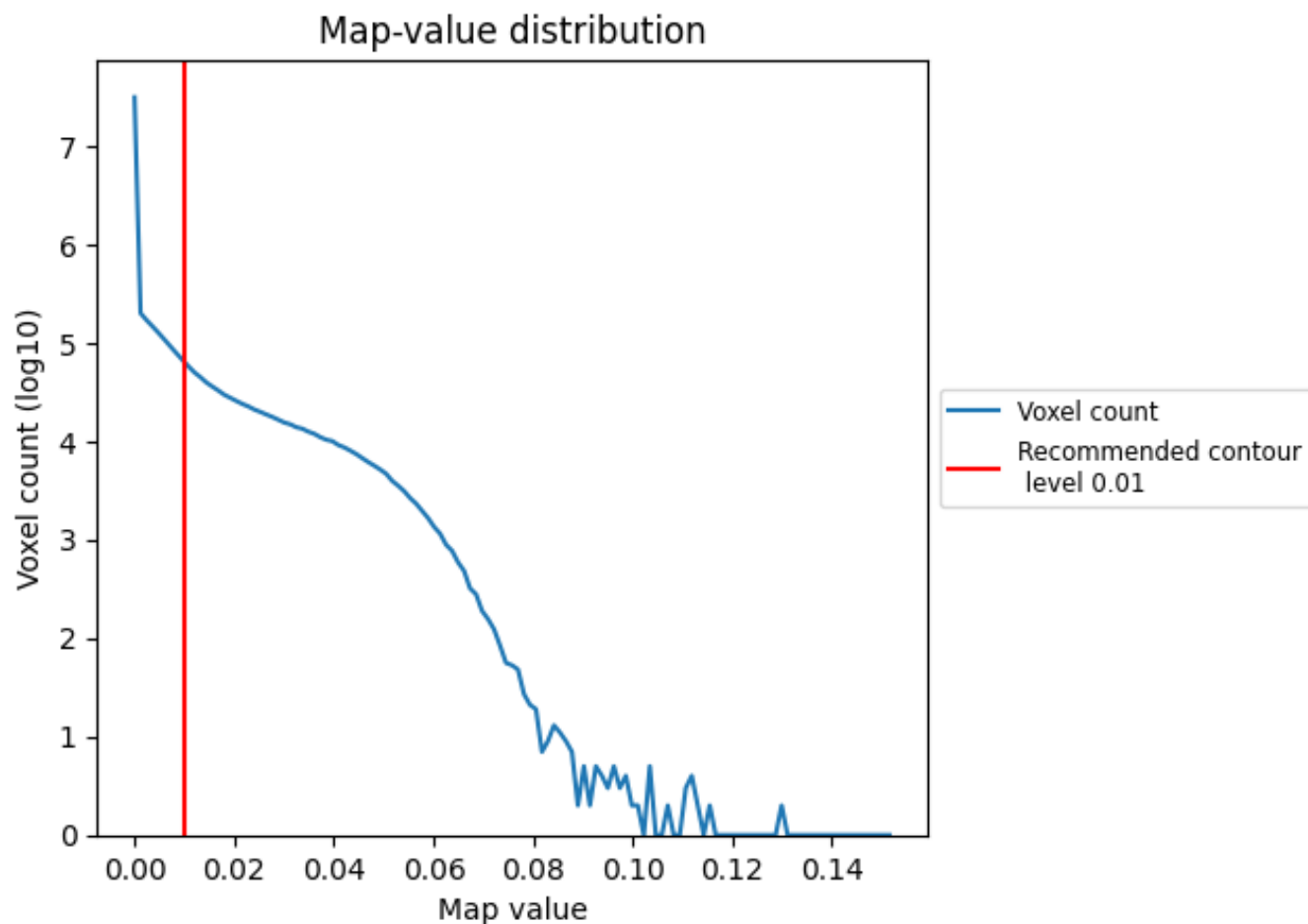


Z

7 Map analysis [i](#)

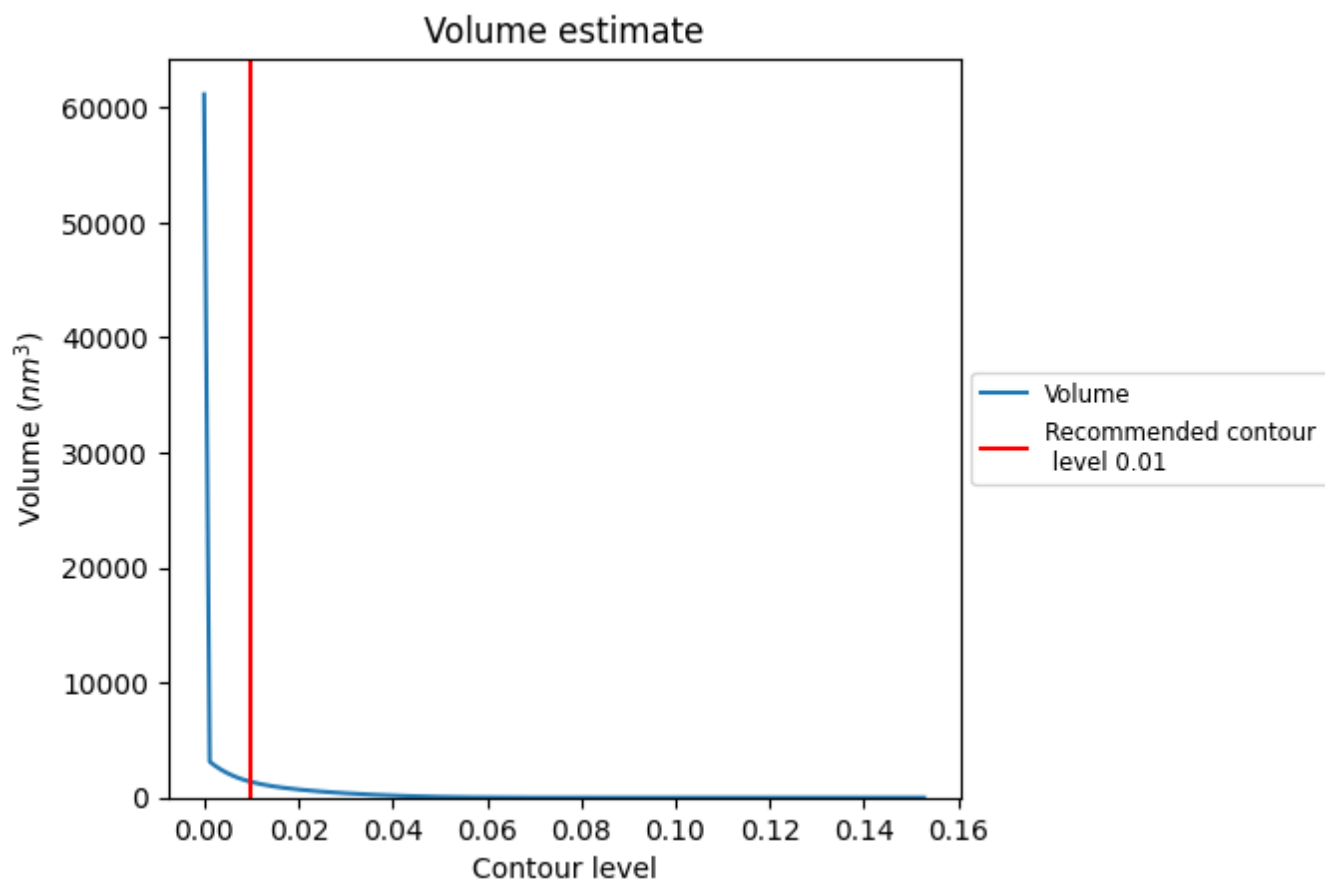
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

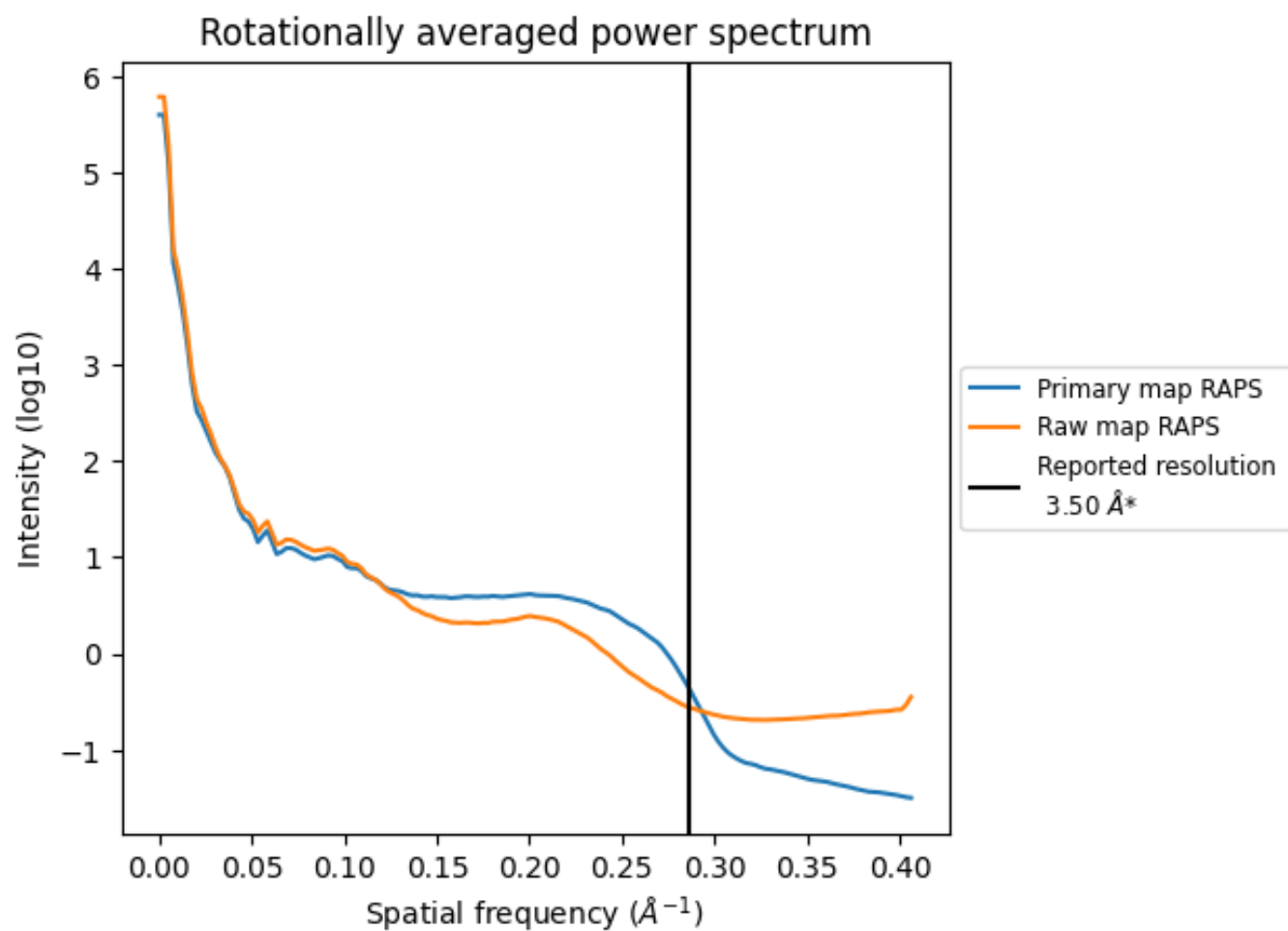
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1360 nm³; this corresponds to an approximate mass of 1229 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

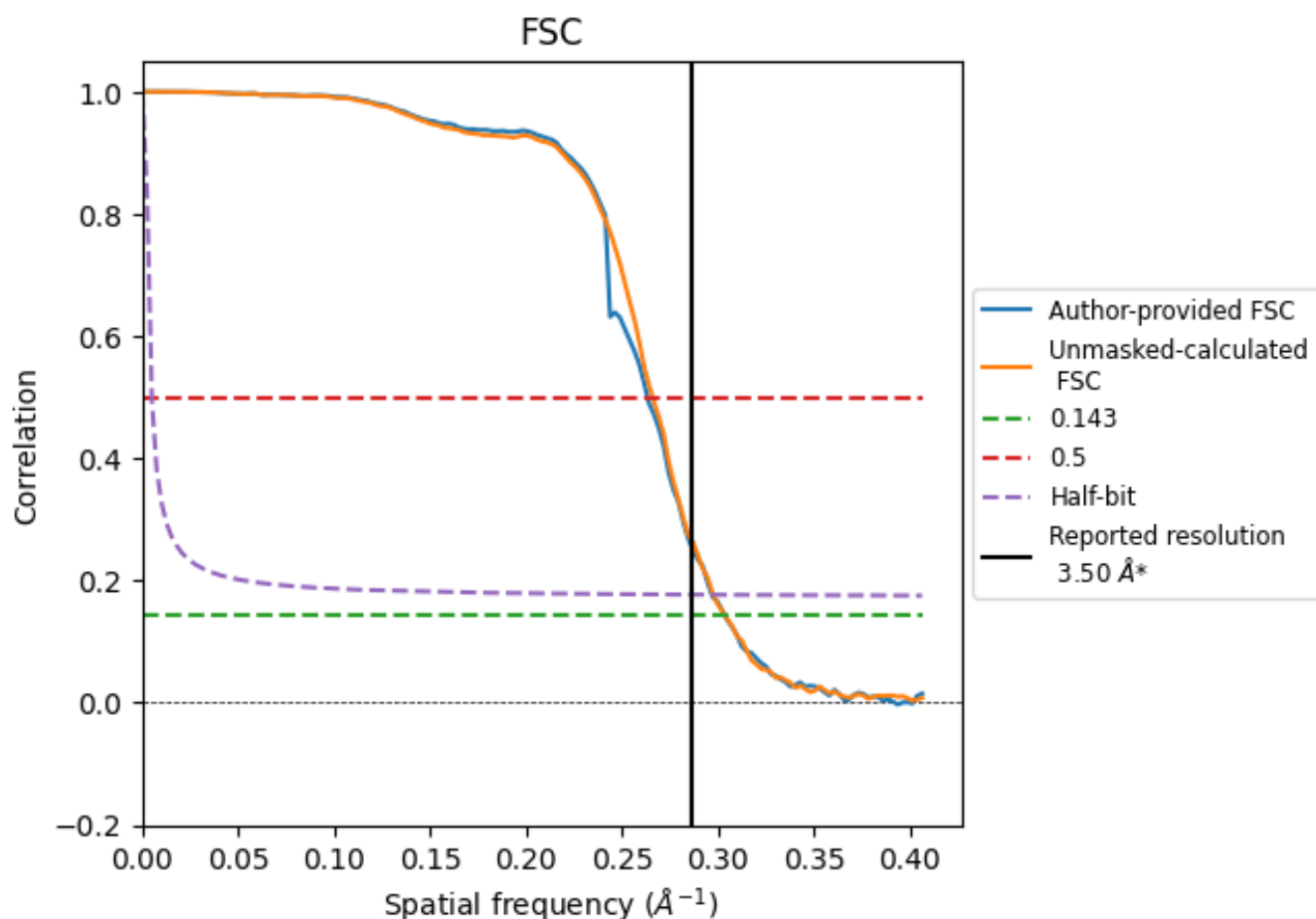


*Reported resolution corresponds to spatial frequency of 0.286 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.286 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

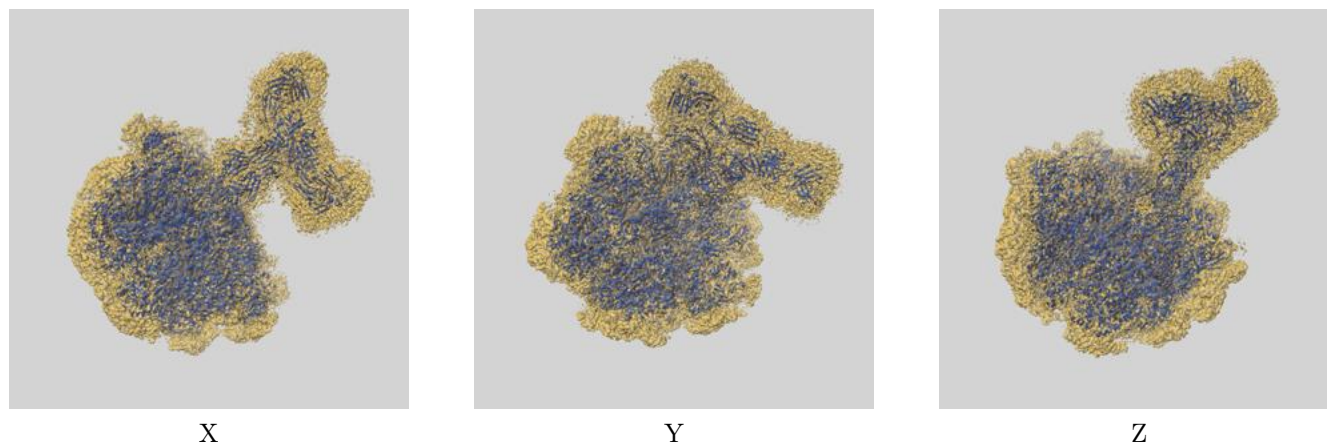
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.50	-	-
Author-provided FSC curve	3.30	3.80	3.37
Unmasked-calculated*	3.29	3.76	3.36

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit [i](#)

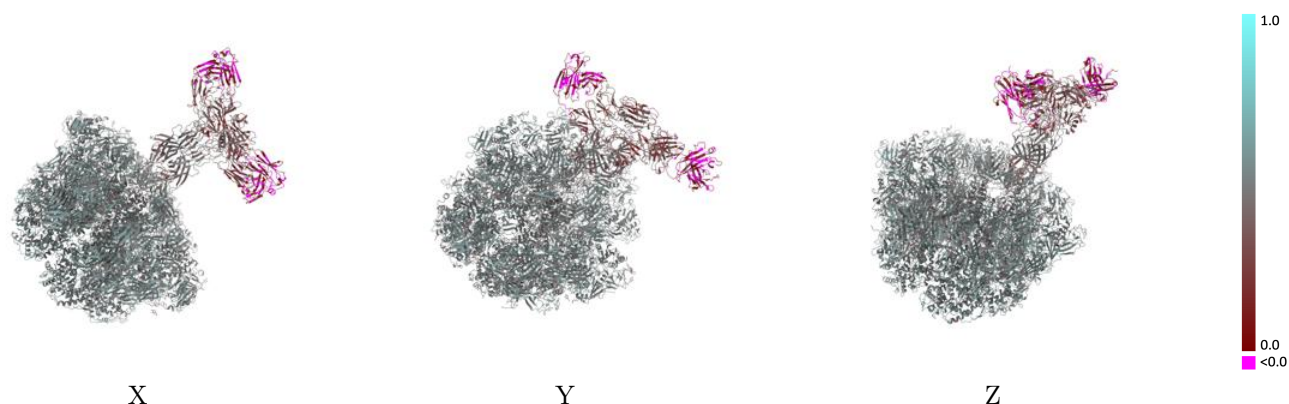
This section contains information regarding the fit between EMDB map EMD-26608 and PDB model 7UMS. Per-residue inclusion information can be found in section [3](#) on page [14](#).

9.1 Map-model overlay [i](#)



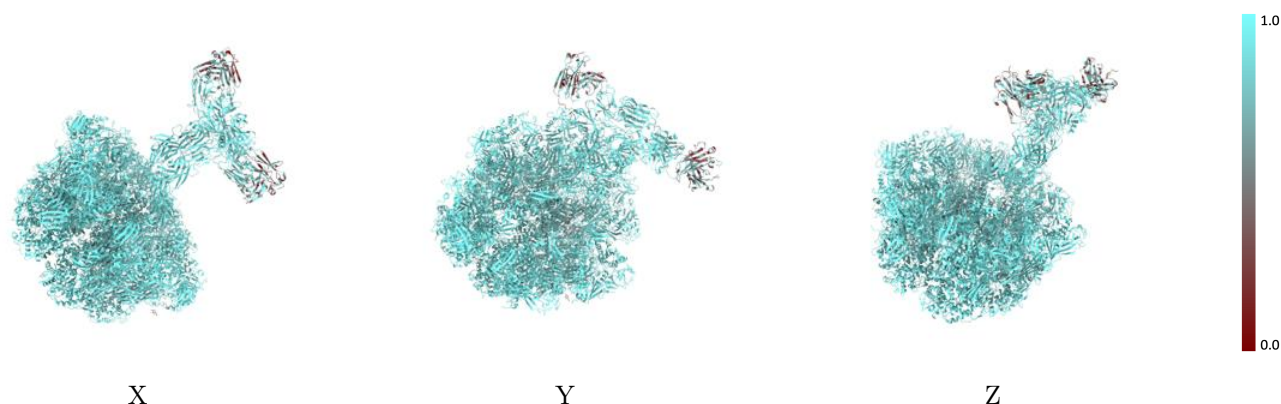
The images above show the 3D surface view of the map at the recommended contour level 0.01 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



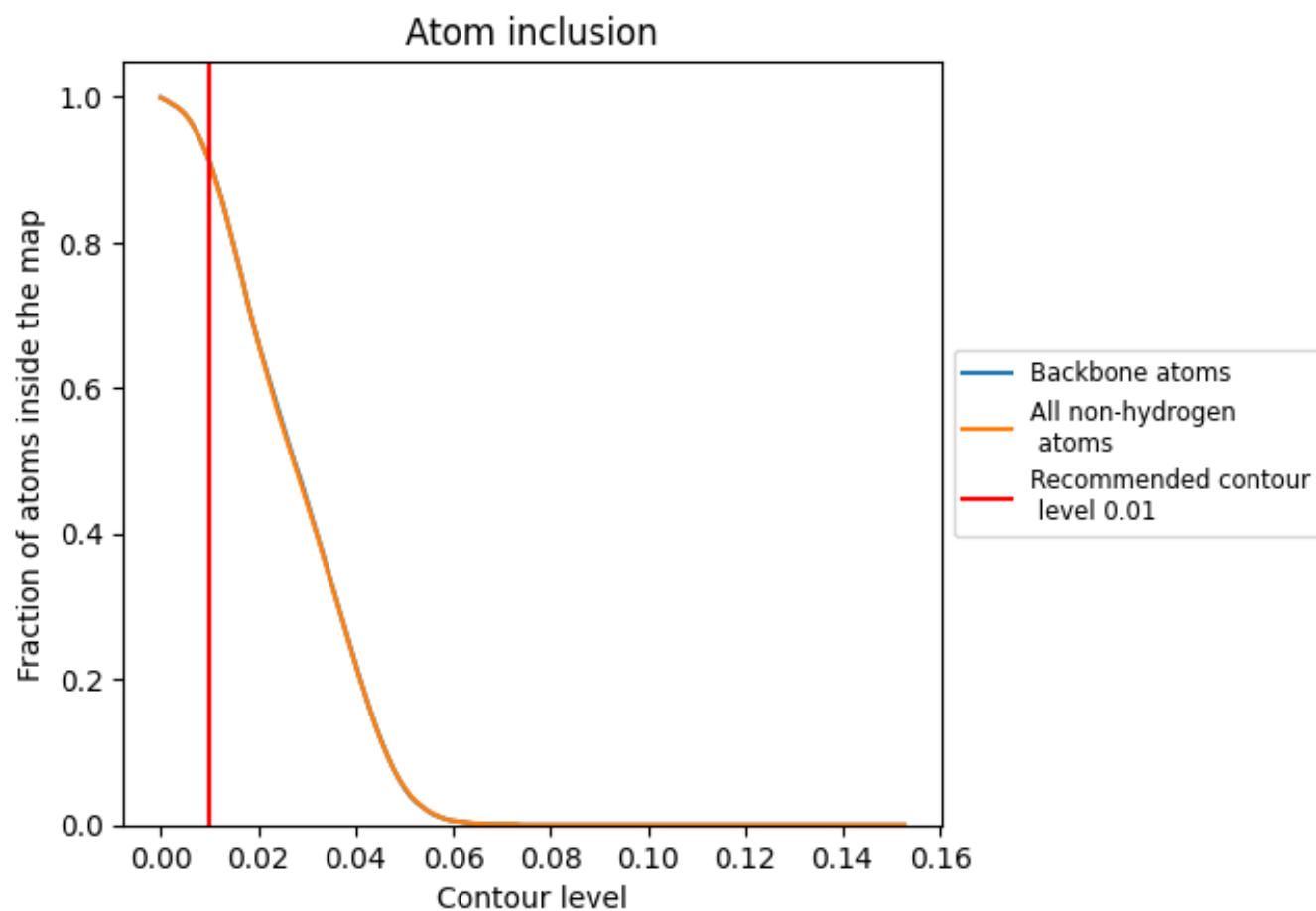
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.01).

9.4 Atom inclusion ⓘ



At the recommended contour level, 91% of all backbone atoms, 91% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (0.01) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9130	 0.4930
1	 0.8620	 0.4540
2	 0.8640	 0.4520
3	 0.8920	 0.4830
4	 0.6790	 0.1990
5	 0.6730	 0.1990
6	 0.6920	 0.2080
7	 0.6910	 0.2010
A	 0.9350	 0.5300
B	 0.9470	 0.5270
C	 0.9410	 0.5250
D	 0.9360	 0.5330
E	 0.9430	 0.5310
F	 0.9450	 0.5360
G	 0.9350	 0.5260
H	 0.9390	 0.5270
I	 0.9390	 0.5300
J	 0.9350	 0.5280
K	 0.9390	 0.5280
L	 0.9320	 0.5230
M	 0.9340	 0.5230
N	 0.9350	 0.5250
O	 0.9370	 0.5240
P	 0.9430	 0.5260
Q	 0.9360	 0.5240
R	 0.9400	 0.5240
S	 0.7690	 0.4220
T	 0.8630	 0.3750
U	 0.8700	 0.3810
V	 0.8890	 0.4790
a	 0.9390	 0.5200
b	 0.9330	 0.5100
c	 0.9380	 0.5160
d	 0.9400	 0.5170
e	 0.9390	 0.5180



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Chain	Atom inclusion	Q-score
f	 0.9370	 0.5120
g	 0.9280	 0.5090
h	 0.9350	 0.5120
i	 0.9380	 0.5160
j	 0.9350	 0.5130
k	 0.9330	 0.5070
l	 0.9320	 0.5170
m	 0.9360	 0.5190
n	 0.9470	 0.5160
o	 0.9330	 0.5120
p	 0.9290	 0.5100
q	 0.9400	 0.5090
r	 0.9330	 0.5050