



Full wwPDB EM Validation Report ⓘ

Jun 20, 2026 – 11:00 am BST

PDB ID : 9RHM / pdb_00009rhm
EMDB ID : EMD-53973
Title : Phospho-MCM double hexamer bound to Sld3-Sld7-Cdc45 on ARS1 DNA
Authors : Puehringer, T.; Butryn, A.; Couves, E.C.; Costa, A.
Deposited on : 2025-06-09
Resolution : 3.70 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 : 1.8.4, CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

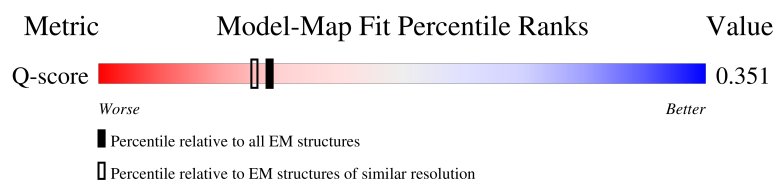
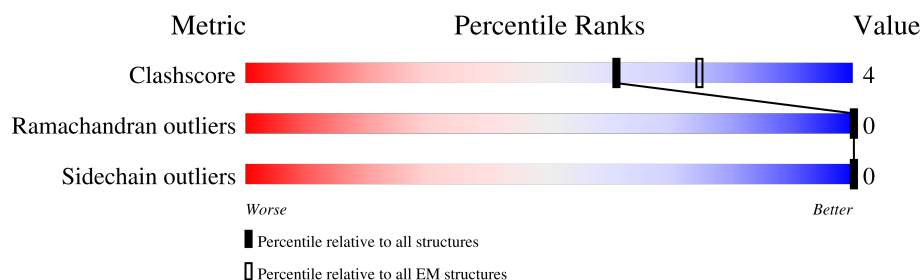
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.70 Å.

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



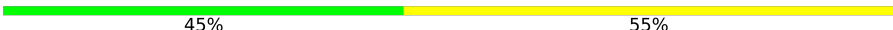
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	11569 (3.20 - 4.20)

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	A	257	97%
1	R	257	96%
2	H	704	98%
2	I	704	6% 10% 88%

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Mol	Chain	Length	Quality of chain
3	X	53	 45% 55%
4	Y	53	 51% 49%
5	E	657	 24% 51% 17% 33%
6	2	868	 8% 64% 8% 28%
6	a	868	 8% 64% 7% 28%
7	3	1006	 57% 6% 37%
7	b	1006	 58% 5% 37%
8	4	933	 5% 65% 7% 28%
8	c	933	 7% 64% 8% 28%
9	5	775	 5% 71% 13% 16%
9	d	775	 6% 75% 9% 16%
10	6	1017	 7% 54% 9% 37%
10	e	1017	 9% 54% 8% 37%
11	7	845	 76% 7% 17%
11	f	845	 6% 74% 9% 17%

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 68681 atoms, of which 0 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 Mitochondrial morphogenesis protein SLD7.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	A	7	Total	C	N	O	0	0
			50	29	8	13		
1	R	10	Total	C	N	O	0	0
			74	47	11	16		

- Molecule 2 is a protein called DNA replication regulator SLD3.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	15	Total	C	N	O	S	0	0
			124	76	24	23	1		
2	I	85	Total	C	N	O	S	0	0
			682	438	114	126	4		

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	669	GLU	-	expression tag	UNP P53135
H	670	ASN	-	expression tag	UNP P53135
H	671	LEU	-	expression tag	UNP P53135
H	672	TYR	-	expression tag	UNP P53135
H	673	PHE	-	expression tag	UNP P53135
H	674	GLN	-	expression tag	UNP P53135
H	675	GLY	-	expression tag	UNP P53135
H	676	TRP	-	expression tag	UNP P53135
H	677	SER	-	expression tag	UNP P53135
H	678	HIS	-	expression tag	UNP P53135
H	679	PRO	-	expression tag	UNP P53135
H	680	GLN	-	expression tag	UNP P53135
H	681	PHE	-	expression tag	UNP P53135
H	682	GLU	-	expression tag	UNP P53135
H	683	LYS	-	expression tag	UNP P53135
H	684	GLY	-	expression tag	UNP P53135
H	685	GLY	-	expression tag	UNP P53135

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Chain	Residue	Modelled	Actual	Comment	Reference
H	686	GLY	-	expression tag	UNP P53135
H	687	SER	-	expression tag	UNP P53135
H	688	GLY	-	expression tag	UNP P53135
H	689	GLY	-	expression tag	UNP P53135
H	690	GLY	-	expression tag	UNP P53135
H	691	SER	-	expression tag	UNP P53135
H	692	GLY	-	expression tag	UNP P53135
H	693	GLY	-	expression tag	UNP P53135
H	694	SER	-	expression tag	UNP P53135
H	695	SER	-	expression tag	UNP P53135
H	696	ALA	-	expression tag	UNP P53135
H	697	TRP	-	expression tag	UNP P53135
H	698	SER	-	expression tag	UNP P53135
H	699	HIS	-	expression tag	UNP P53135
H	700	PRO	-	expression tag	UNP P53135
H	701	GLN	-	expression tag	UNP P53135
H	702	PHE	-	expression tag	UNP P53135
H	703	GLU	-	expression tag	UNP P53135
H	704	LYS	-	expression tag	UNP P53135
I	669	GLU	-	expression tag	UNP P53135
I	670	ASN	-	expression tag	UNP P53135
I	671	LEU	-	expression tag	UNP P53135
I	672	TYR	-	expression tag	UNP P53135
I	673	PHE	-	expression tag	UNP P53135
I	674	GLN	-	expression tag	UNP P53135
I	675	GLY	-	expression tag	UNP P53135
I	676	TRP	-	expression tag	UNP P53135
I	677	SER	-	expression tag	UNP P53135
I	678	HIS	-	expression tag	UNP P53135
I	679	PRO	-	expression tag	UNP P53135
I	680	GLN	-	expression tag	UNP P53135
I	681	PHE	-	expression tag	UNP P53135
I	682	GLU	-	expression tag	UNP P53135
I	683	LYS	-	expression tag	UNP P53135
I	684	GLY	-	expression tag	UNP P53135
I	685	GLY	-	expression tag	UNP P53135
I	686	GLY	-	expression tag	UNP P53135
I	687	SER	-	expression tag	UNP P53135
I	688	GLY	-	expression tag	UNP P53135
I	689	GLY	-	expression tag	UNP P53135
I	690	GLY	-	expression tag	UNP P53135
I	691	SER	-	expression tag	UNP P53135

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Chain	Residue	Modelled	Actual	Comment	Reference
I	692	GLY	-	expression tag	UNP P53135
I	693	GLY	-	expression tag	UNP P53135
I	694	SER	-	expression tag	UNP P53135
I	695	SER	-	expression tag	UNP P53135
I	696	ALA	-	expression tag	UNP P53135
I	697	TRP	-	expression tag	UNP P53135
I	698	SER	-	expression tag	UNP P53135
I	699	HIS	-	expression tag	UNP P53135
I	700	PRO	-	expression tag	UNP P53135
I	701	GLN	-	expression tag	UNP P53135
I	702	PHE	-	expression tag	UNP P53135
I	703	GLU	-	expression tag	UNP P53135
I	704	LYS	-	expression tag	UNP P53135

- Molecule 3 is a DNA chain called DNA (53-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
3	X	53	Total	C	N	O	P	0	0
			1086	515	199	319	53		

- Molecule 4 is a DNA chain called DNA (53-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
4	Y	53	Total	C	N	O	P	0	0
			1087	515	202	317	53		

- Molecule 5 is a protein called Cell division control protein 45.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	442	Total	C	N	O	S	0	0
			3590	2309	610	658	13		

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	150W	ASP	-	linker	UNP Q08032
E	150X	TYR	-	linker	UNP Q08032
E	150Y	LYS	-	linker	UNP Q08032
E	150Z	ASP	-	linker	UNP Q08032
E	151A	ASP	-	linker	UNP Q08032
E	151B	ASP	-	linker	UNP Q08032
E	151C	GLY	-	linker	UNP Q08032

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Chain	Residue	Modelled	Actual	Comment	Reference
E	151D	ASP	-	linker	UNP Q08032
E	151E	TYR	-	linker	UNP Q08032
E	151F	LYS	-	linker	UNP Q08032
E	151G	ASP	-	linker	UNP Q08032
E	151H	ASP	-	linker	UNP Q08032
E	151I	ASP	-	linker	UNP Q08032

- Molecule 6 is a protein called DNA replication licensing factor MCM2.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	2	623	Total	C	N	O	S	0	0
			4932	3109	877	927	19		
6	a	623	Total	C	N	O	S	0	0
			4932	3109	877	927	19		

- Molecule 7 is a protein called DNA replication licensing factor MCM3.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	3	638	Total	C	N	O	S	0	0
			4994	3150	889	942	13		
7	b	638	Total	C	N	O	S	0	0
			4994	3150	889	942	13		

There are 70 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
3	-34	MET	-	initiating methionine	UNP P24279
3	-33	LYS	-	expression tag	UNP P24279
3	-32	ARG	-	expression tag	UNP P24279
3	-31	ARG	-	expression tag	UNP P24279
3	-30	TRP	-	expression tag	UNP P24279
3	-29	LYS	-	expression tag	UNP P24279
3	-28	LYS	-	expression tag	UNP P24279
3	-27	ASN	-	expression tag	UNP P24279
3	-26	PHE	-	expression tag	UNP P24279
3	-25	ILE	-	expression tag	UNP P24279
3	-24	ALA	-	expression tag	UNP P24279
3	-23	VAL	-	expression tag	UNP P24279
3	-22	SER	-	expression tag	UNP P24279
3	-21	ALA	-	expression tag	UNP P24279
3	-20	ALA	-	expression tag	UNP P24279
3	-19	ASN	-	expression tag	UNP P24279

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Chain	Residue	Modelled	Actual	Comment	Reference
3	-18	ARG	-	expression tag	UNP P24279
3	-17	PHE	-	expression tag	UNP P24279
3	-16	LYS	-	expression tag	UNP P24279
3	-15	LYS	-	expression tag	UNP P24279
3	-14	ILE	-	expression tag	UNP P24279
3	-13	SER	-	expression tag	UNP P24279
3	-12	SER	-	expression tag	UNP P24279
3	-11	SER	-	expression tag	UNP P24279
3	-10	GLY	-	expression tag	UNP P24279
3	-9	ALA	-	expression tag	UNP P24279
3	-8	LEU	-	expression tag	UNP P24279
3	-7	GLU	-	expression tag	UNP P24279
3	-6	ASN	-	expression tag	UNP P24279
3	-5	LEU	-	expression tag	UNP P24279
3	-4	TYR	-	expression tag	UNP P24279
3	-3	PHE	-	expression tag	UNP P24279
3	-2	GLN	-	expression tag	UNP P24279
3	-1	GLY	-	expression tag	UNP P24279
3	0	GLU	-	expression tag	UNP P24279
b	-34	MET	-	initiating methionine	UNP P24279
b	-33	LYS	-	expression tag	UNP P24279
b	-32	ARG	-	expression tag	UNP P24279
b	-31	ARG	-	expression tag	UNP P24279
b	-30	TRP	-	expression tag	UNP P24279
b	-29	LYS	-	expression tag	UNP P24279
b	-28	LYS	-	expression tag	UNP P24279
b	-27	ASN	-	expression tag	UNP P24279
b	-26	PHE	-	expression tag	UNP P24279
b	-25	ILE	-	expression tag	UNP P24279
b	-24	ALA	-	expression tag	UNP P24279
b	-23	VAL	-	expression tag	UNP P24279
b	-22	SER	-	expression tag	UNP P24279
b	-21	ALA	-	expression tag	UNP P24279
b	-20	ALA	-	expression tag	UNP P24279
b	-19	ASN	-	expression tag	UNP P24279
b	-18	ARG	-	expression tag	UNP P24279
b	-17	PHE	-	expression tag	UNP P24279
b	-16	LYS	-	expression tag	UNP P24279
b	-15	LYS	-	expression tag	UNP P24279
b	-14	ILE	-	expression tag	UNP P24279
b	-13	SER	-	expression tag	UNP P24279
b	-12	SER	-	expression tag	UNP P24279

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Chain	Residue	Modelled	Actual	Comment	Reference
b	-11	SER	-	expression tag	UNP P24279
b	-10	GLY	-	expression tag	UNP P24279
b	-9	ALA	-	expression tag	UNP P24279
b	-8	LEU	-	expression tag	UNP P24279
b	-7	GLU	-	expression tag	UNP P24279
b	-6	ASN	-	expression tag	UNP P24279
b	-5	LEU	-	expression tag	UNP P24279
b	-4	TYR	-	expression tag	UNP P24279
b	-3	PHE	-	expression tag	UNP P24279
b	-2	GLN	-	expression tag	UNP P24279
b	-1	GLY	-	expression tag	UNP P24279
b	0	GLU	-	expression tag	UNP P24279

- Molecule 8 is a protein called DNA replication licensing factor MCM4.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	4	672	Total	C	N	O	S	0	0
			5356	3357	926	1043	30		
8	c	672	Total	C	N	O	S	0	0
			5356	3357	926	1043	30		

- Molecule 9 is a protein called Minichromosome maintenance protein 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	5	651	Total	C	N	O	S	0	0
			5097	3195	872	1005	25		
9	d	651	Total	C	N	O	S	0	0
			5097	3195	872	1005	25		

- Molecule 10 is a protein called DNA replication licensing factor MCM6.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	6	637	Total	C	N	O	S	0	0
			5041	3174	879	963	25		
10	e	637	Total	C	N	O	S	0	0
			5041	3174	879	963	25		

- Molecule 11 is a protein called DNA replication licensing factor MCM7.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	7	700	Total	C	N	O	S	0	0
			5513	3468	953	1060	32		

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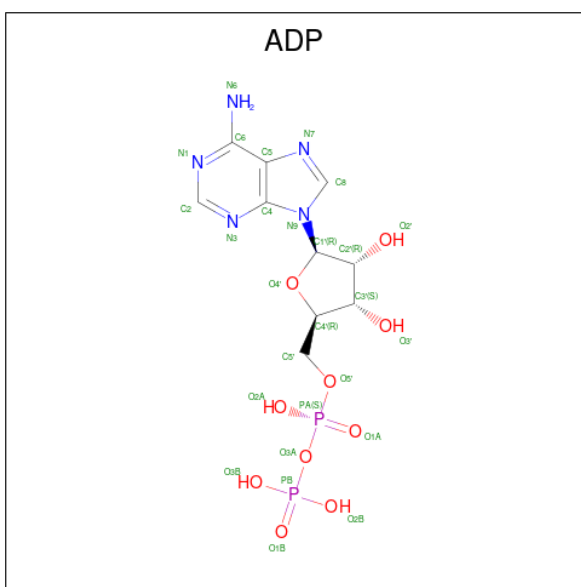
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Mol	Chain	Residues	Atoms					AltConf	Trace
11	f	700	Total	C	N	O	S	0	0
			5513	3468	953	1060	32		

- Molecule 12 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
12	2	1	Total	Zn	0
			1	1	
12	4	1	Total	Zn	0
			1	1	
12	5	1	Total	Zn	0
			1	1	
12	6	1	Total	Zn	0
			1	1	
12	7	1	Total	Zn	0
			1	1	
12	a	1	Total	Zn	0
			1	1	
12	c	1	Total	Zn	0
			1	1	
12	d	1	Total	Zn	0
			1	1	
12	e	1	Total	Zn	0
			1	1	
12	f	1	Total	Zn	0
			1	1	

- Molecule 13 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



Mol	Chain	Residues	Atoms					AltConf
13	3	1	Total 27	C 10	N 5	O 10	P 2	0
13	5	1	Total 27	C 10	N 5	O 10	P 2	0
13	b	1	Total 27	C 10	N 5	O 10	P 2	0
13	d	1	Total 27	C 10	N 5	O 10	P 2	0

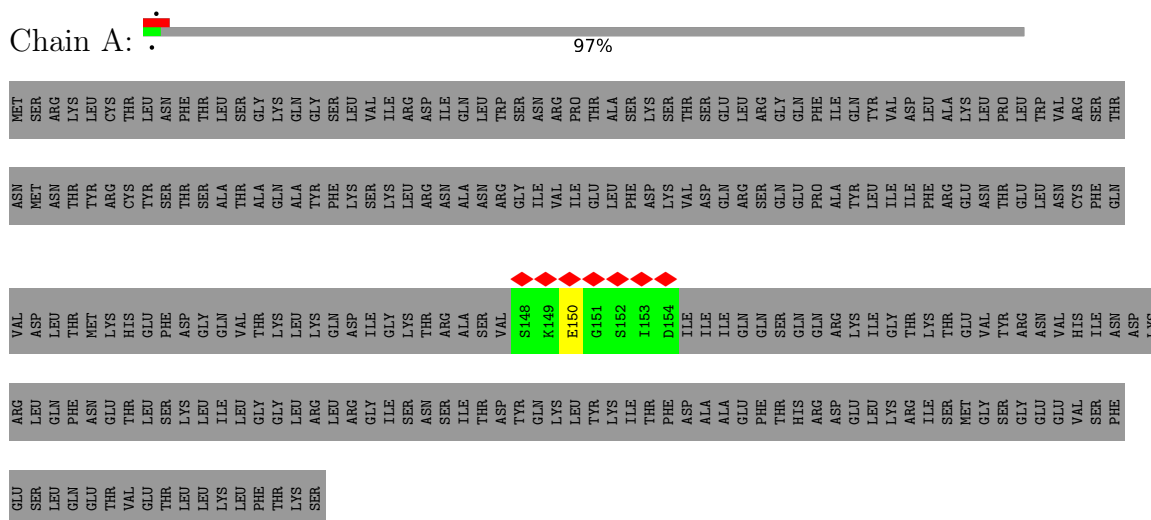
- Molecule 14 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	AltConf
14	3	1	Total Mg 1 1	0
14	5	1	Total Mg 1 1	0
14	a	1	Total Mg 1 1	0
14	b	1	Total Mg 1 1	0

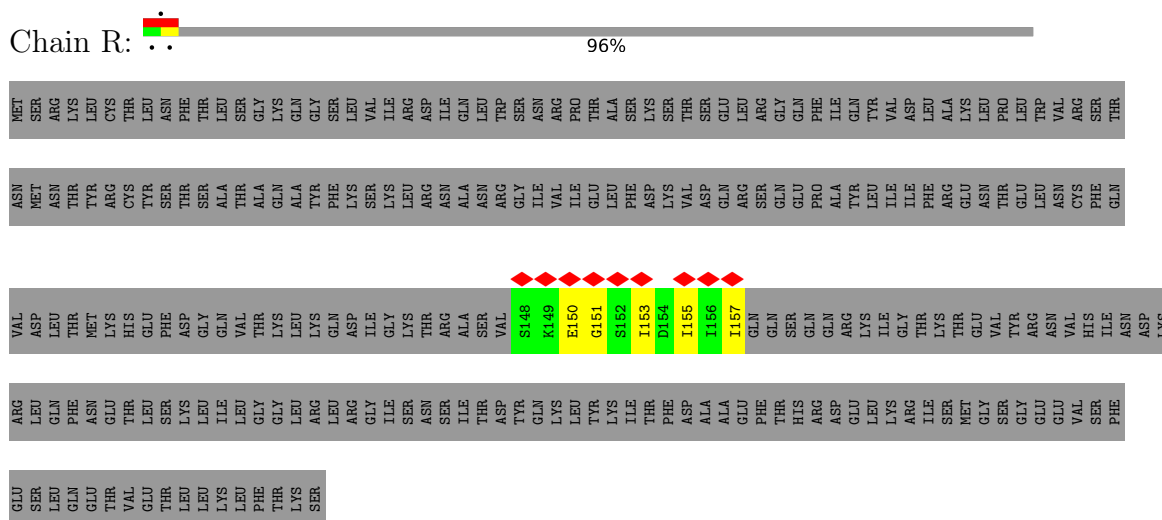
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Mitochondrial morphogenesis protein SLD7

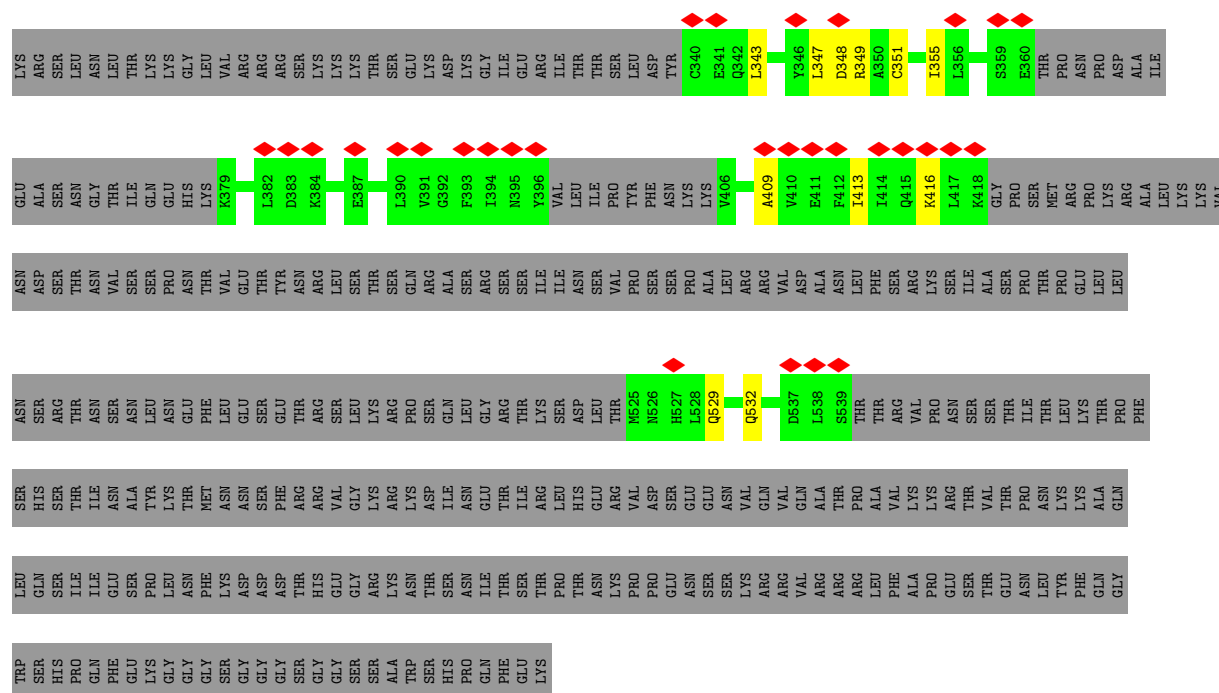


- Molecule 1: Mitochondrial morphogenesis protein SLD7



- Molecule 2: DNA replication regulator SLD3





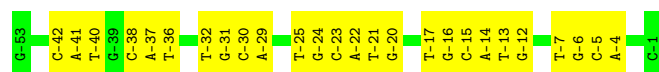
• Molecule 3: DNA (53-MER)

Chain X: 45% 55%



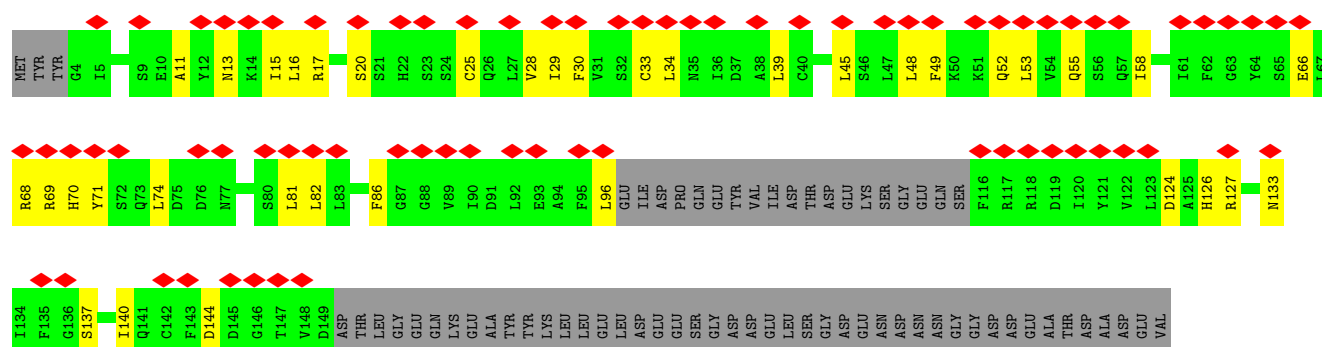
• Molecule 4: DNA (53-MER)

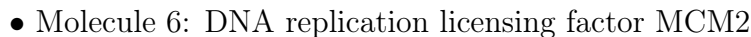
Chain Y: 51% 49%



• Molecule 5: Cell division control protein 45

Chain E: 24% 51% 17% 33%





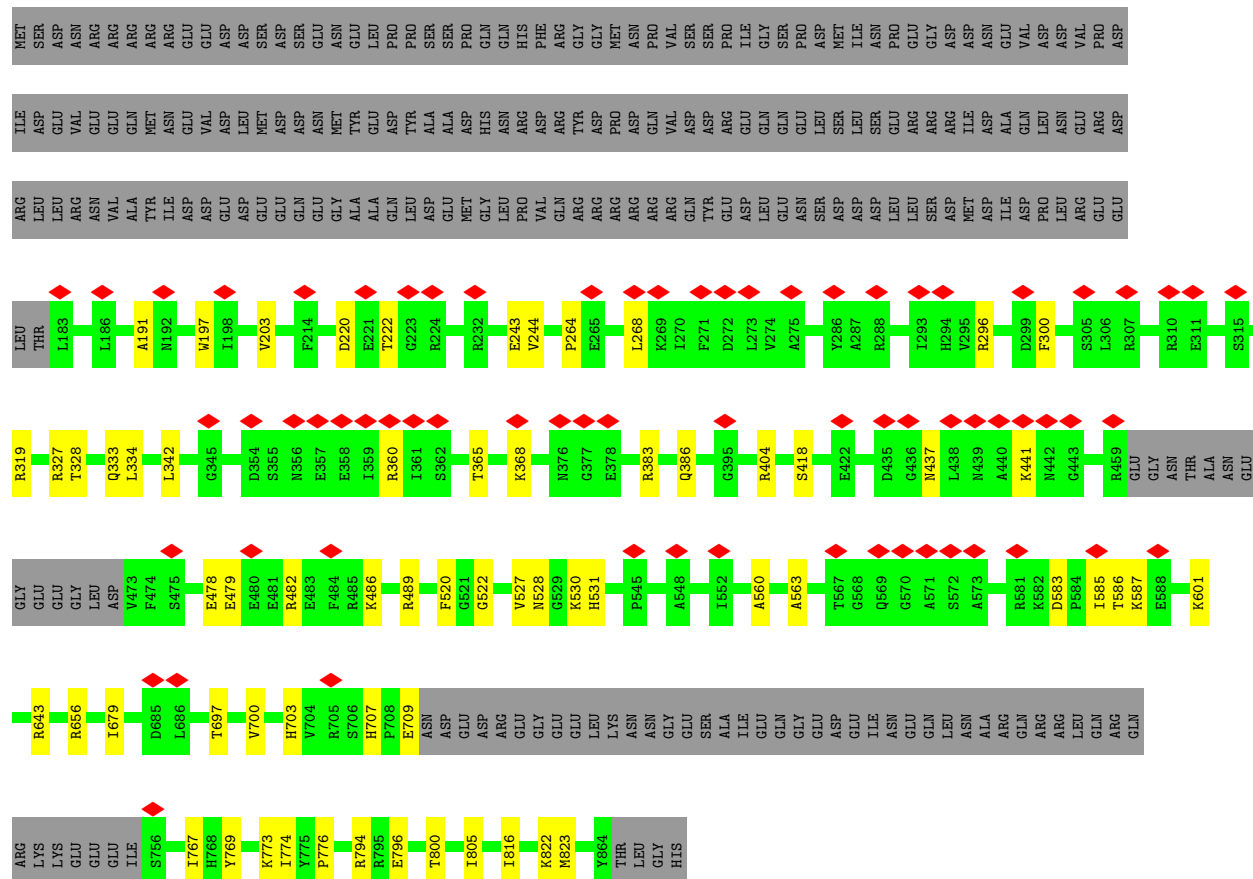
Frequency	Percentage
Daily	64%
Weekly	8%
Monthly	28%



LEU
GLY
HIS

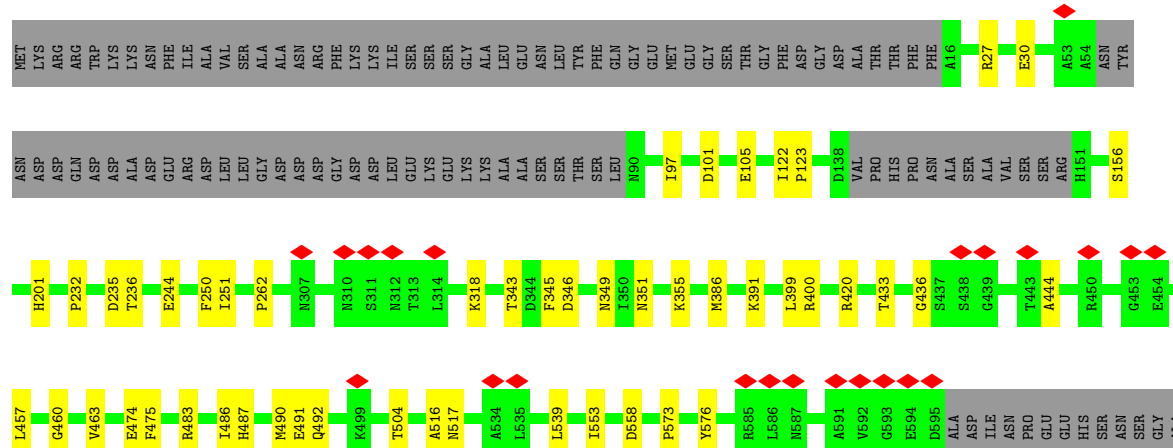
• Molecule 6: DNA replication licensing factor MCM2

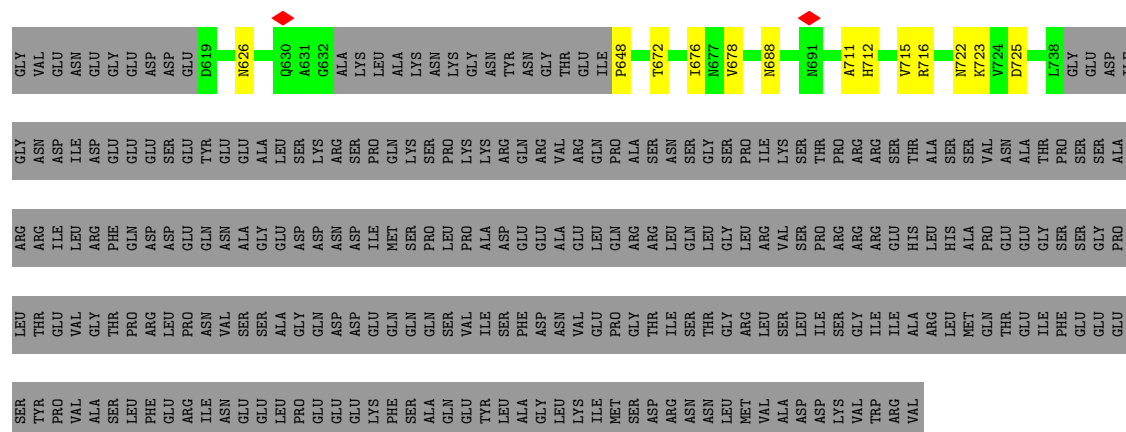
Chain a: 



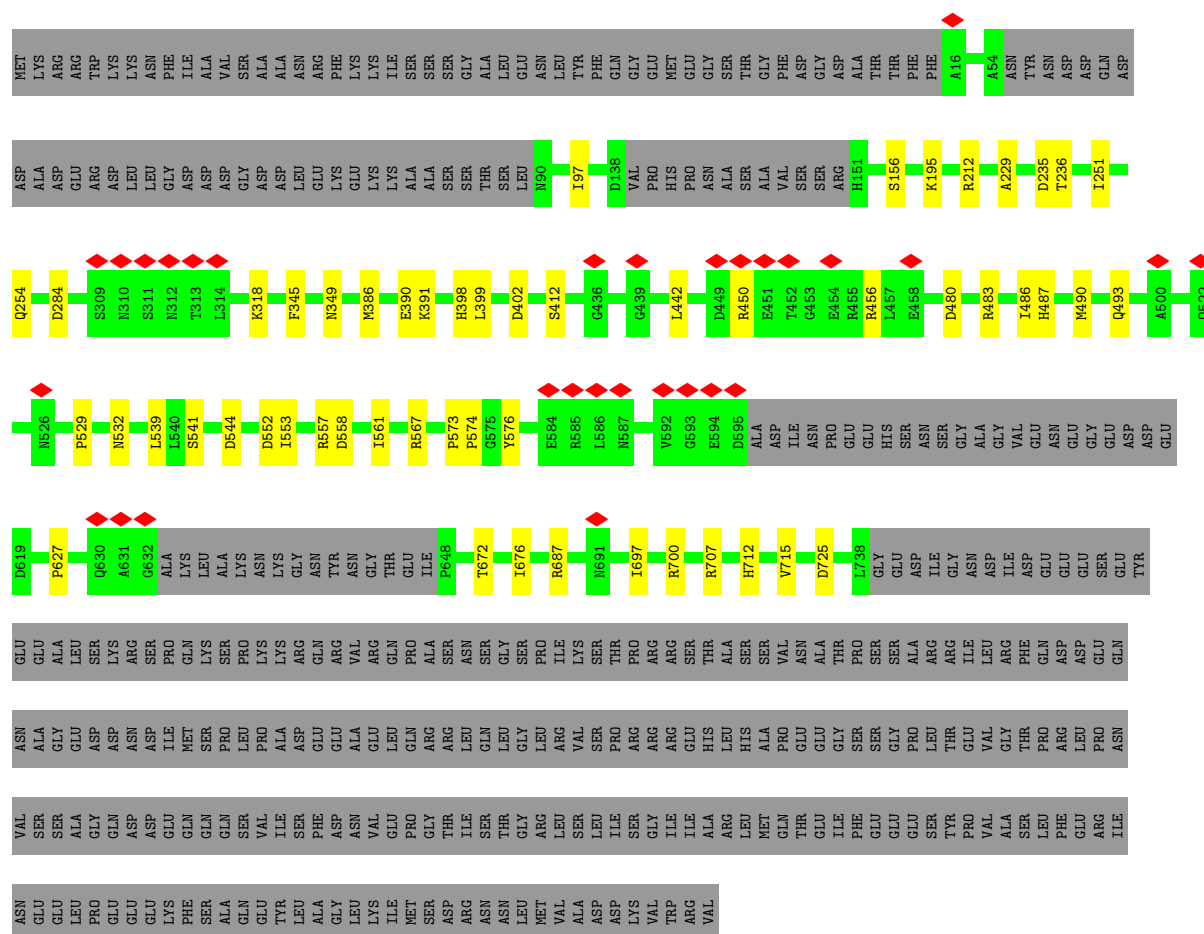
• Molecule 7: DNA replication licensing factor MCM3

Chain 3: 

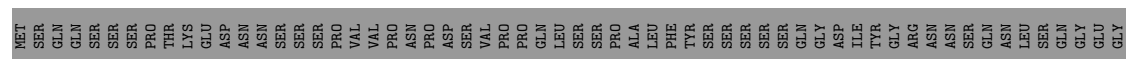


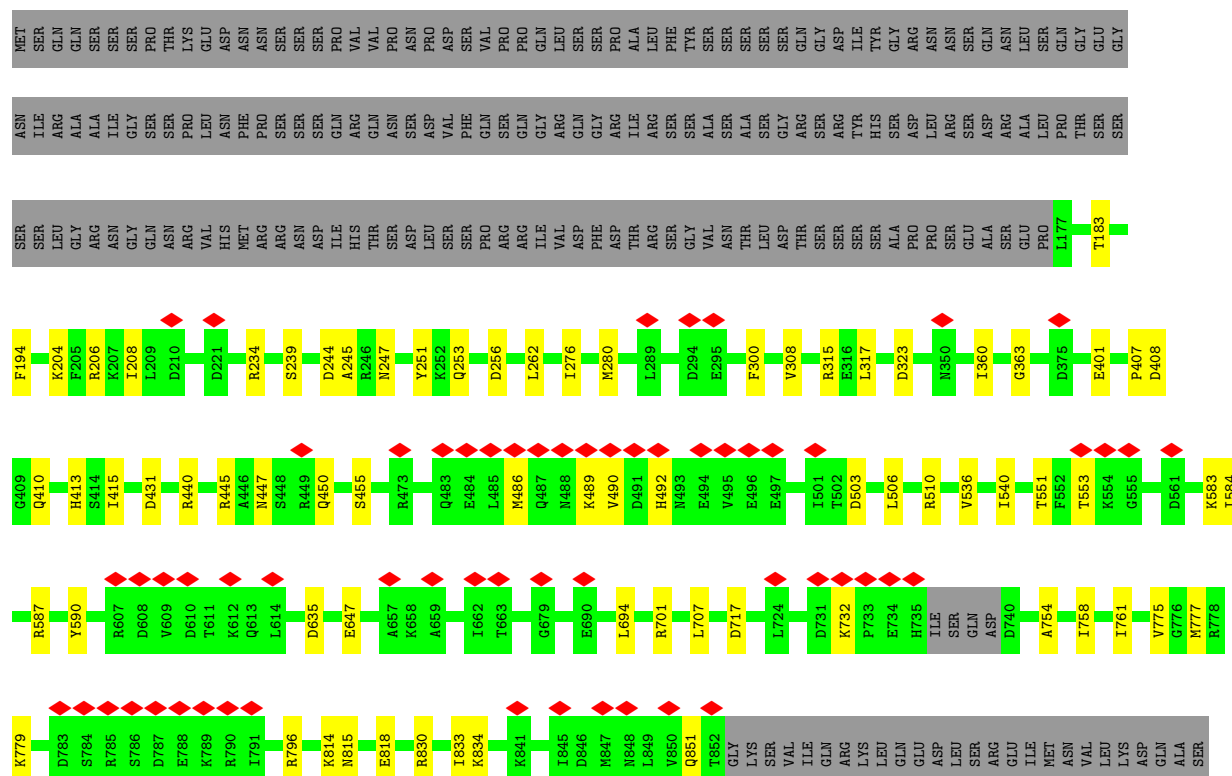


- Molecule 7: DNA replication licensing factor MCM3

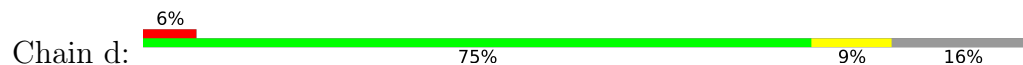


- Molecule 8: DNA replication licensing factor MCM4





Chain 5: 5% 71% 13% 16%



HIS
GLU
THR
ILE
GLN
LEU
ARG
HIS
GLY
GLN
ASN
ILE
TYR
ARG
SER
GLY
VAL

• Molecule 10: DNA replication licensing factor MCM6



MET
SER
SER
PRO
PRO
PHE
PRO
ALA
ASP
THR
SER
PRO
SER
SER
ASN
ARG
PRO
SER
SER
ASN
SER
SER
PRO
PRO
PRO
MET
SER
SER
SER
ILE
GLY
ALA
GLY
PHE
GLY
SER
SER
SER
SER
GLY
LEU
ASP
ALA
SER
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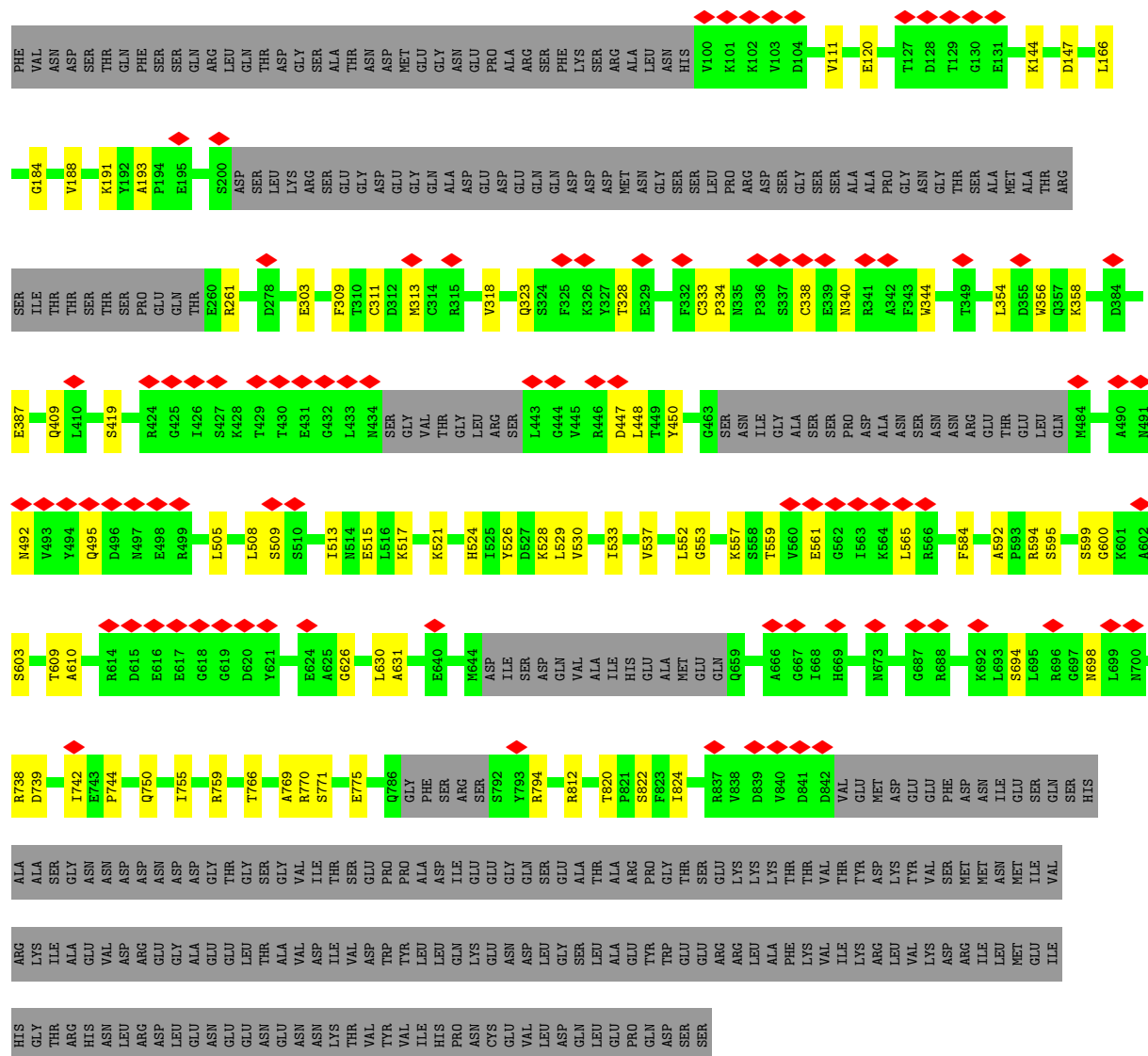
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• Molecule 10: DNA replication licensing factor MCM6

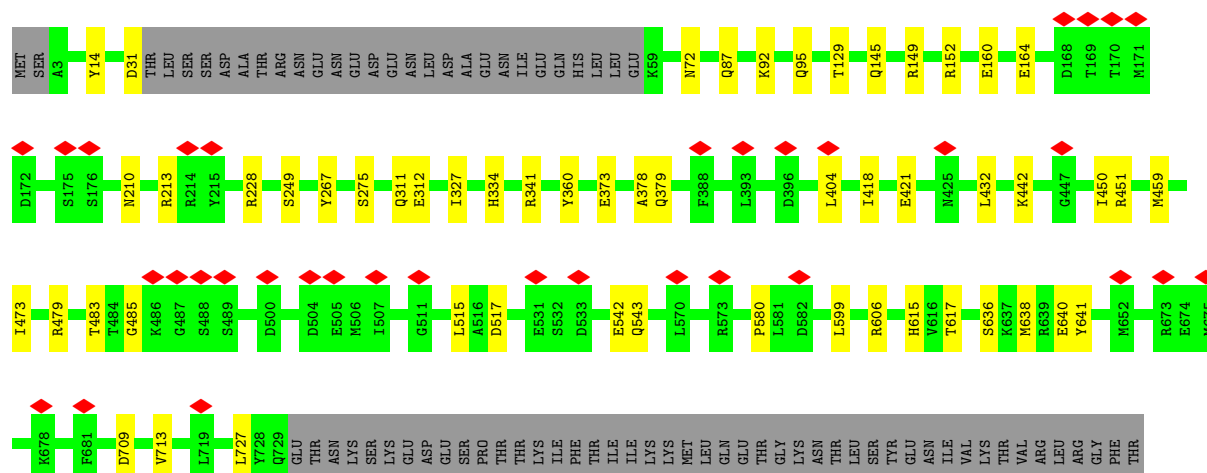


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• Molecule 11: DNA replication licensing factor MCM7

Chain 7: 76% 7% 17%



MET	GLN
LEU	GLN
GLN	SER
LEU	ASP
ASP	ILE
SER	ASN
ASN	CYS
CYS	ILE
ILE	GLN
GLN	GLU
GLU	THR
THR	TYR
TYR	SER
SER	TYR
TYR	LEU
LEU	ASN
ASN	VAL
VAL	TRP
TRP	HIS
HIS	LEU
LEU	ILE
ILE	ASN
ASN	GLY
GLY	GLU
GLU	ASN
ASN	THR
THR	LEU
LEU	LYS
LYS	PHE
PHE	VAL
VAL	ASP
ASP	ASP
ASP	GLY
GLY	THR
THR	LEU
LEU	MET
MET	GLU
GLU	ASP
ASP	THR
THR	ASP
ASP	GLN
GLN	GLU
GLU	ASP
ASP	SER
SER	LEU
LEU	VAL
VAL	SER
SER	THR
THR	PRO
PRO	LYS
LYS	LEU
LEU	ALA
ALA	PRO
PRO	GLN
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THR	ALA
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SER	ALA
ALA	ASN
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SER	ALA

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● Molecule 11: DNA replication licensing factor MCM7



THR	LEU	LYS	PHE	VAL	ASP	ASP	GLY	THR	MET	ASP	THR	THR	ASP	GLN	GLU	ASP	SER	LEU	VAL	SER	THR	THR	LYS	PRO	THR	LEU	ALA	ALA	SER	ARG	LEU	ALA	ASN	VAL	GLY	SER	PHE	THR	ALA	ASP	SER	ASP	LEU	GLN	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR	THR
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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	72693	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	49.8	Depositor
Minimum defocus (nm)	1100	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	130000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.782	Depositor
Minimum map value	-0.303	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.024	Depositor
Recommended contour level	0.15	Depositor
Map size ($\text{\AA}$)	483.84003, 483.84003, 483.84003	wwPDB
Map dimensions	448, 448, 448	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.08, 1.08, 1.08	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, MG, ADP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.06	0/49	0.15	0/63
1	R	0.07	0/73	0.18	0/96
2	H	0.06	0/125	0.22	0/165
2	I	0.08	0/685	0.21	0/914
3	X	0.17	0/1217	0.35	0/1876
4	Y	0.17	0/1219	0.34	0/1879
5	E	0.10	0/3651	0.27	0/4929
6	2	0.10	0/5019	0.29	0/6781
6	a	0.10	0/5019	0.29	0/6781
7	3	0.11	0/5081	0.30	0/6889
7	b	0.11	0/5081	0.29	0/6889
8	4	0.11	0/5434	0.32	0/7345
8	c	0.10	0/5434	0.29	0/7345
9	5	0.11	0/5170	0.30	0/6992
9	d	0.11	0/5170	0.30	0/6992
10	6	0.10	0/5119	0.28	0/6906
10	e	0.10	0/5119	0.28	0/6906
11	7	0.11	0/5599	0.27	0/7569
11	f	0.10	0/5599	0.28	0/7569
All	All	0.11	0/69863	0.29	0/94886

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	50	0	46	1	0
1	R	74	0	79	3	0
2	H	124	0	122	3	0
2	I	682	0	717	8	0
3	X	1086	0	596	17	0
4	Y	1087	0	595	17	0
5	E	3590	0	3619	71	0
6	2	4932	0	4985	43	0
6	a	4932	0	4985	38	0
7	3	4994	0	5058	39	0
7	b	4994	0	5058	33	0
8	4	5356	0	5410	44	0
8	c	5356	0	5410	46	0
9	5	5097	0	5142	65	0
9	d	5097	0	5142	43	0
10	6	5041	0	5065	52	0
10	e	5041	0	5065	51	0
11	7	5513	0	5575	34	0
11	f	5513	0	5575	45	0
12	2	1	0	0	0	0
12	4	1	0	0	0	0
12	5	1	0	0	0	0
12	6	1	0	0	0	0
12	7	1	0	0	0	0
12	a	1	0	0	0	0
12	c	1	0	0	0	0
12	d	1	0	0	0	0
12	e	1	0	0	0	0
12	f	1	0	0	0	0
13	3	27	0	12	1	0
13	5	27	0	12	3	0
13	b	27	0	12	1	0
13	d	27	0	12	3	0
14	3	1	0	0	0	0
14	5	1	0	0	0	0
14	a	1	0	0	0	0
14	b	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	68681	0	68292	598	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 4.

All (598) 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)
5:E:16:LEU:O	5:E:20:SER:HB2	1.76	0.85
7:3:444:ALA:HB1	7:3:463:VAL:HG12	1.62	0.79
2:I:349:ARG:HH21	5:E:646:LEU:HD23	1.52	0.74
5:E:526:ARG:NH1	5:E:556:CYS:SG	2.62	0.72
4:Y:-37:DA:H2"	4:Y:-36:DT:H5"	1.72	0.71
11:7:442:LYS:HB3	11:7:450:ILE:HB	1.74	0.70
8:c:445:ARG:NH1	10:e:447:ASP:OD2	2.28	0.67
9:d:77:LYS:NZ	9:d:81:ASP:OD2	2.24	0.66
5:E:29:ILE:HB	5:E:58:ILE:HG22	1.77	0.66
10:6:553:GLY:O	10:6:812:ARG:NH1	2.29	0.66
6:a:794:ARG:NH1	6:a:805:ILE:O	2.28	0.65
8:c:431:ASP:OD1	8:c:587:ARG:NH1	2.30	0.64
10:e:592:ALA:HB3	10:e:595:SER:HB3	1.78	0.64
8:c:851:GLN:N	8:c:851:GLN:OE1	2.30	0.64
10:6:270:LEU:HD12	10:6:289:SER:HB2	1.80	0.63
8:4:431:ASP:OD1	8:4:587:ARG:NH1	2.32	0.63
5:E:124:ASP:OD1	5:E:126:HIS:ND1	2.31	0.62
11:f:453:ASP:OD2	11:f:544:GLN:NE2	2.32	0.62
7:3:716:ARG:NH2	7:3:722:ASN:OD1	2.32	0.62
7:b:573:PRO:HG2	7:b:576:TYR:HB2	1.80	0.62
10:6:642:ASP:OD2	10:6:643:LYS:NZ	2.29	0.62
6:2:344:CYS:HB3	6:2:367:CYS:SG	2.39	0.62
6:2:292:GLU:OE2	6:2:335:LYS:NZ	2.32	0.62
11:7:483:THR:HG22	11:7:485:GLY:H	1.64	0.62
1:R:150:GLU:HG3	10:6:517:LYS:HD2	1.82	0.61
10:6:742:ILE:HG22	10:6:744:PRO:HD3	1.82	0.61
6:a:478:GLU:OE1	6:a:482:ARG:NH2	2.32	0.61
10:6:310:THR:HG22	10:6:317:ILE:HG12	1.82	0.61
7:b:687:ARG:NH1	7:b:697:ILE:O	2.34	0.61
8:c:814:LYS:NZ	8:c:818:GLU:OE1	2.27	0.61
11:f:625:GLN:OE1	11:f:627:ASP:N	2.33	0.60
6:a:328:THR:O	6:a:386:GLN:NE2	2.35	0.60
8:c:204:LYS:O	8:c:208:ILE:HG12	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:349:SER:O	5:E:352:ASN:ND2	2.35	0.60
9:5:77:LYS:NZ	9:5:81:ASP:OD2	2.26	0.59
8:c:551:THR:O	10:e:738:ARG:NH2	2.32	0.59
7:b:254:GLN:OE1	7:b:284:ASP:N	2.34	0.59
8:c:262:LEU:HD11	8:c:308:VAL:HG23	1.84	0.59
10:e:594:ARG:HB3	10:e:631:ALA:HA	1.84	0.59
11:7:479:ARG:NH2	11:7:515:LEU:O	2.36	0.59
5:E:270:LEU:HD22	5:E:302:LEU:HD12	1.85	0.59
7:b:558:ASP:OD2	9:d:630:ARG:NE	2.34	0.59
1:A:150:GLU:HG3	10:e:517:LYS:HD2	1.85	0.58
9:d:498:GLU:OE1	9:d:651:ARG:NH1	2.36	0.58
7:b:712:HIS:ND1	7:b:725:ASP:OD1	2.36	0.58
10:e:742:ILE:HG22	10:e:744:PRO:HD3	1.85	0.58
3:X:38:DG:OP2	8:4:612:LYS:NZ	2.36	0.58
8:c:777:MET:HG2	8:c:830:ARG:HH21	1.67	0.58
5:E:643:LYS:HA	5:E:646:LEU:HD12	1.86	0.58
7:b:390:GLU:OE1	7:b:398:HIS:NE2	2.32	0.58
8:4:294:ASP:O	8:4:298:THR:HG23	2.04	0.58
11:f:290:SER:HA	11:f:295:LYS:HE3	1.85	0.58
10:e:609:THR:OG1	10:e:610:ALA:N	2.35	0.58
11:f:160:GLU:O	11:f:164:GLU:HG2	2.04	0.58
5:E:388:LEU:HG	5:E:392:PHE:CE2	2.39	0.58
10:6:532:SER:HB2	10:6:745:PRO:HG2	1.85	0.58
7:3:475:PHE:HB3	7:3:516:ALA:HB2	1.85	0.58
6:a:528:ASN:HD22	6:a:530:LYS:HE2	1.69	0.58
10:e:557:LYS:HB2	10:e:565:LEU:HD12	1.84	0.57
2:I:343:LEU:O	2:I:347:LEU:HG	2.04	0.57
11:7:432:LEU:HB3	11:7:473:ILE:HD11	1.85	0.57
9:d:444:SER:HB2	9:d:447:ALA:HB3	1.87	0.57
7:3:244:GLU:OE2	11:7:14:TYR:OH	2.23	0.57
10:6:609:THR:OG1	10:6:610:ALA:N	2.37	0.57
6:a:585:ILE:HG13	6:a:586:THR:HG23	1.85	0.57
2:I:348:ASP:OD1	2:I:416:LYS:NZ	2.37	0.57
8:c:239:SER:OG	8:c:300:PHE:O	2.21	0.57
5:E:277:THR:HG22	5:E:288:TYR:HE1	1.70	0.56
6:2:520:PHE:HE1	6:2:767:ILE:HG12	1.69	0.56
9:5:149:ARG:NH1	9:5:260:GLU:OE1	2.38	0.56
9:5:588:GLU:HG3	9:5:593:GLU:HB2	1.87	0.56
7:3:487:HIS:CD2	7:3:539:LEU:HD13	2.40	0.56
8:c:244:ASP:OD1	8:c:245:ALA:N	2.38	0.56
9:d:50:LEU:HD13	9:d:61:LEU:HD12	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:e:553:GLY:O	10:e:812:ARG:NH1	2.37	0.56
7:b:97:ILE:HG12	7:b:156:SER:HB2	1.87	0.56
9:d:580:ALA:HA	9:d:584:GLN:HB2	1.87	0.56
11:f:138:VAL:HG21	11:f:303:ARG:HH21	1.71	0.56
3:X:17:DA:H2"	3:X:18:DT:H5"	1.86	0.56
7:b:552:ASP:O	7:b:557:ARG:NH2	2.36	0.56
3:X:16:DC:H2"	3:X:17:DA:C8	2.41	0.56
11:f:122:ASP:OD2	11:f:198:ARG:NH1	2.37	0.56
6:a:243:GLU:HA	6:a:296:ARG:HB2	1.88	0.56
10:e:492:ASN:HB2	10:e:495:GLN:HG3	1.88	0.56
5:E:16:LEU:O	5:E:20:SER:CB	2.50	0.55
8:4:212:ARG:HD2	8:4:215:PHE:HE2	1.71	0.55
7:3:97:ILE:HG12	7:3:156:SER:HB2	1.89	0.55
7:3:672:THR:O	7:3:676:ILE:HG12	2.06	0.55
11:f:444:VAL:HB	11:f:448:MET:HB2	1.89	0.55
8:4:292:ASP:OD1	8:4:292:ASP:N	2.39	0.55
8:4:437:GLY:HA2	8:4:464:VAL:HG22	1.89	0.55
5:E:467:THR:HG22	5:E:471:LYS:HE3	1.87	0.55
8:4:602:THR:HG21	8:4:656:ILE:HD11	1.88	0.55
6:a:601:LYS:N	6:a:643:ARG:O	2.34	0.55
5:E:526:ARG:HH21	5:E:567:MET:HG3	1.71	0.55
9:5:72:ASN:OD1	9:5:73:GLU:N	2.39	0.55
9:5:580:ALA:HA	9:5:584:GLN:HB2	1.88	0.55
10:e:409:GLN:NE2	10:e:450:TYR:O	2.30	0.55
8:c:775:VAL:O	8:c:779:LYS:HG2	2.06	0.55
9:d:168:SER:HB3	9:d:257:LYS:HB2	1.89	0.55
6:a:191:ALA:HB3	6:a:197:TRP:HB2	1.89	0.54
7:b:672:THR:O	7:b:676:ILE:HG12	2.07	0.54
5:E:580:LEU:HA	5:E:630:ILE:O	2.07	0.54
10:6:125:GLN:NE2	10:6:129:THR:O	2.37	0.54
6:2:260:LEU:HA	6:2:267:MET:HG3	1.90	0.54
3:X:43:DC:H2"	3:X:44:DG:C8	2.43	0.54
8:4:350:ASN:OD1	8:4:381:SER:OG	2.22	0.54
8:c:536:VAL:O	8:c:540:ILE:HG12	2.08	0.54
10:6:816:VAL:HG12	10:6:818:GLU:H	1.72	0.54
11:7:459:MET:HE3	11:7:599:LEU:HD13	1.90	0.54
4:Y:-25:DT:H2"	4:Y:-24:DG:C8	2.43	0.53
5:E:250:SER:HB2	5:E:280:LEU:HD11	1.90	0.53
9:5:50:LEU:HD13	9:5:61:LEU:HD12	1.90	0.53
11:f:262:CYS:HB3	11:f:265:CYS:SG	2.48	0.53
2:H:529:GLN:HA	2:H:532:GLN:NE2	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:e:505:LEU:HD23	10:e:508:LEU:HD12	1.90	0.53
4:Y:-13:DT:H2''	4:Y:-12:DG:C8	2.44	0.53
11:7:312:GLU:OE1	11:7:334:HIS:NE2	2.39	0.53
8:c:707:LEU:O	8:c:834:LYS:NZ	2.38	0.53
6:a:333:GLN:OE1	6:a:383:ARG:NH1	2.41	0.53
10:e:303:GLU:HB2	10:e:356:TRP:HB2	1.90	0.53
11:f:459:MET:HE3	11:f:599:LEU:HD13	1.91	0.53
4:Y:-30:DC:H2''	4:Y:-29:DA:C8	2.44	0.53
4:Y:-15:DC:H2''	4:Y:-14:DA:C8	2.43	0.53
5:E:29:ILE:HG12	5:E:82:LEU:HD13	1.91	0.53
9:5:559:ASP:O	9:5:564:ARG:NH2	2.39	0.53
8:4:209:LEU:HG	8:4:210:ASP:H	1.73	0.52
11:7:580:PRO:HB3	11:7:727:LEU:HD11	1.90	0.52
4:Y:-37:DA:H2'	4:Y:-36:DT:H71	1.91	0.52
9:5:168:SER:HB3	9:5:257:LYS:HB2	1.91	0.52
10:6:107:THR:OG1	10:6:176:ARG:NH2	2.42	0.52
7:b:483:ARG:HA	7:b:486:ILE:HG22	1.91	0.52
9:5:625:ASN:O	9:5:629:ILE:HG13	2.09	0.52
8:c:317:LEU:O	11:f:341:ARG:NH2	2.42	0.52
10:e:515:GLU:OE2	10:e:750:GLN:NE2	2.39	0.52
10:6:524:HIS:HB3	10:6:528:LYS:HE3	1.91	0.52
5:E:71:TYR:HA	5:E:74:LEU:HG	1.92	0.52
6:2:693:GLU:OE2	10:6:778:LYS:NZ	2.31	0.52
7:3:688:ASN:OD1	11:7:606:ARG:NH2	2.43	0.52
6:2:774:ILE:HG22	6:2:776:PRO:HD3	1.92	0.52
11:f:432:LEU:HB3	11:f:473:ILE:HD11	1.90	0.52
7:3:400:ARG:NH1	7:3:490:MET:O	2.43	0.52
7:3:626:ASN:ND2	7:3:648:PRO:O	2.42	0.52
8:4:310:SER:OG	8:4:327:ASN:OD1	2.22	0.52
11:f:479:ARG:NH2	11:f:515:LEU:O	2.42	0.52
5:E:328:LEU:HD22	5:E:496:ILE:HD11	1.92	0.52
9:5:65:MET:SD	9:5:161:ARG:NH2	2.83	0.52
10:6:313:MET:HB2	10:6:338:CYS:SG	2.50	0.52
8:c:635:ASP:HA	8:c:694:LEU:HD21	1.92	0.52
8:4:213:GLU:OE2	8:4:213:GLU:N	2.24	0.51
7:b:486:ILE:HG12	7:b:490:MET:HE2	1.92	0.51
10:e:552:LEU:HD21	10:e:755:ILE:HG23	1.90	0.51
5:E:270:LEU:O	5:E:274:ILE:HG13	2.10	0.51
9:5:149:ARG:HH21	9:5:269:GLU:CD	2.19	0.51
8:4:455:SER:OG	11:7:275:SER:O	2.26	0.51
10:6:147:ASP:OD2	10:6:261:ARG:NH2	2.42	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:2:198:ILE:HG21	6:2:266:GLU:HB3	1.93	0.51
5:E:334:LEU:HB2	5:E:380:MET:HE1	1.92	0.51
7:3:712:HIS:ND1	7:3:725:ASP:OD1	2.41	0.51
8:4:536:VAL:O	8:4:540:ILE:HG12	2.09	0.51
11:7:160:GLU:O	11:7:164:GLU:HG2	2.10	0.51
10:e:309:PHE:N	10:e:318:VAL:O	2.41	0.51
7:b:235:ASP:OD1	7:b:236:THR:N	2.42	0.51
10:e:559:THR:HG22	10:e:561:GLU:H	1.76	0.51
8:4:360:ILE:HD11	8:4:363:GLY:HA2	1.92	0.51
7:b:544:ASP:OD2	7:b:707:ARG:NH1	2.43	0.51
5:E:388:LEU:HG	5:E:392:PHE:HE2	1.74	0.51
7:3:235:ASP:OD1	7:3:236:THR:N	2.43	0.51
10:6:581:LYS:HA	10:6:584:PHE:HD2	1.76	0.51
9:d:421:ALA:HB2	13:d:801:ADP:C8	2.46	0.51
7:3:491:GLU:HG3	7:3:492:GLN:HG3	1.94	0.50
8:4:552:PHE:HB2	8:4:556:GLY:HA3	1.94	0.50
7:b:195:LYS:HE2	7:b:251:ILE:HD11	1.93	0.50
10:e:771:SER:O	10:e:775:GLU:HG2	2.12	0.50
11:f:256:GLU:OE1	11:f:306:LYS:NZ	2.44	0.50
4:Y:-23:DC:H2"	4:Y:-22:DA:C8	2.47	0.50
10:6:303:GLU:HB2	10:6:356:TRP:HB2	1.93	0.50
7:b:391:LYS:HB2	7:b:399:LEU:HD12	1.93	0.50
6:2:490:ASP:OD1	6:2:491:ARG:N	2.44	0.50
6:a:404:ARG:NH2	10:e:387:GLU:OE2	2.44	0.50
11:f:442:LYS:HB3	11:f:450:ILE:HB	1.93	0.50
6:2:519:LEU:O	6:2:771:ARG:NH1	2.44	0.50
8:4:481:ILE:HG22	8:4:484:GLU:HB2	1.94	0.50
7:b:386:MET:HG3	7:b:715:VAL:HG22	1.93	0.50
7:b:553:ILE:HG21	9:d:634:LEU:HB2	1.93	0.50
7:3:345:PHE:O	7:3:349:ASN:ND2	2.44	0.50
6:a:479:GLU:OE1	6:a:482:ARG:NH1	2.43	0.50
6:2:682:VAL:HG12	6:2:684:ARG:H	1.75	0.50
7:b:345:PHE:O	7:b:349:ASN:ND2	2.44	0.50
7:b:557:ARG:O	7:b:561:ILE:HG12	2.12	0.50
6:2:394:PRO:O	10:6:673:ASN:ND2	2.35	0.50
9:d:365:LYS:O	9:d:369:ILE:HG12	2.12	0.50
7:3:101:ASP:O	7:3:105:GLU:HG2	2.12	0.49
9:d:419:GLY:N	13:d:801:ADP:O3B	2.30	0.49
9:d:455:ARG:NH1	9:d:460:ARG:O	2.43	0.49
3:X:6:DC:H2"	3:X:7:DA:C8	2.47	0.49
7:3:27:ARG:NH1	7:3:30:GLU:OE1	2.38	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:d:559:ASP:OD2	9:d:564:ARG:NH1	2.44	0.49
3:X:45:DC:H2"	3:X:46:DA:C8	2.46	0.49
6:a:560:ALA:HB3	6:a:563:ALA:HB2	1.95	0.49
9:d:625:ASN:O	9:d:629:ILE:HG13	2.11	0.49
5:E:386:ARG:HH22	9:5:77:LYS:HD3	1.77	0.49
6:a:327:ARG:NH1	6:a:418:SER:O	2.39	0.49
5:E:34:LEU:HD22	5:E:432:LEU:HD13	1.93	0.49
5:E:48:LEU:O	5:E:52:GLN:HG2	2.13	0.49
7:b:487:HIS:CD2	7:b:539:LEU:HD13	2.48	0.49
9:d:480:ASP:OD1	9:d:481:GLU:N	2.45	0.49
10:6:771:SER:O	10:6:775:GLU:HG2	2.13	0.49
9:5:476:VAL:HG22	9:5:518:SER:HB2	1.94	0.49
9:d:66:GLU:HA	9:d:69:ILE:HG22	1.94	0.49
5:E:365:ARG:O	5:E:365:ARG:NH1	2.45	0.49
10:6:699:LEU:HD12	10:6:706:MET:HE3	1.95	0.49
9:d:624:SER:O	9:d:628:THR:HG23	2.13	0.49
7:3:444:ALA:HA	7:3:460:GLY:HA3	1.93	0.49
4:Y:-17:DT:H2"	4:Y:-16:DG:C8	2.48	0.48
11:f:378:ALA:O	11:f:379:GLN:NE2	2.38	0.48
1:R:153:ILE:O	1:R:157:ILE:HG13	2.12	0.48
6:2:264:PRO:O	6:2:268:LEU:HG	2.13	0.48
6:2:520:PHE:CE1	6:2:767:ILE:HG12	2.48	0.48
6:2:660:THR:HA	6:2:851:VAL:HG23	1.95	0.48
6:a:486:LYS:HD3	6:a:489:ARG:HH21	1.78	0.48
5:E:553:ILE:HA	5:E:556:CYS:SG	2.54	0.48
8:4:223:GLU:OE1	8:4:228:LYS:NZ	2.40	0.48
4:Y:-21:DT:H2"	4:Y:-20:DG:H8	1.78	0.48
9:5:379:PHE:HB3	13:5:801:ADP:HN62	1.79	0.48
8:c:447:ASN:HD21	8:c:450:GLN:HB2	1.78	0.48
9:d:182:MET:HG3	9:d:189:THR:HG22	1.95	0.48
10:e:318:VAL:HG22	10:e:334:PRO:HG3	1.96	0.48
9:5:497:MET:HE2	9:5:519:VAL:HG11	1.95	0.48
3:X:37:DT:H2"	3:X:38:DG:C8	2.48	0.48
3:X:49:DC:H2"	3:X:50:DA:C8	2.49	0.48
7:b:412:SER:HB2	9:d:651:ARG:HH21	1.78	0.48
10:e:311:CYS:SG	10:e:340:ASN:ND2	2.87	0.48
10:e:419:SER:HB2	10:e:448:LEU:HD11	1.94	0.48
10:6:559:THR:HG22	10:6:561:GLU:H	1.78	0.48
11:7:149:ARG:HE	9:d:13:VAL:HA	1.78	0.48
9:d:456:ASP:HB3	9:d:461:GLU:H	1.79	0.48
10:6:303:GLU:N	10:6:354:LEU:O	2.37	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:6:694:SER:O	10:6:698:ASN:ND2	2.38	0.48
3:X:31:DC:H2"	3:X:32:DA:C8	2.49	0.48
3:X:47:DT:H2"	3:X:48:DG:C8	2.49	0.48
7:3:386:MET:HG3	7:3:715:VAL:HG22	1.94	0.48
9:5:674:GLU:HA	9:5:677:VAL:HB	1.95	0.48
8:c:486:MET:O	8:c:490:VAL:HG23	2.14	0.48
11:f:653:SER:HB3	11:f:707:MET:HE2	1.96	0.48
6:2:243:GLU:HA	6:2:296:ARG:HB2	1.95	0.48
6:2:435:ASP:CG	6:2:437:ASN:HD22	2.22	0.48
9:5:298:TYR:HA	9:5:328:ILE:HD13	1.96	0.48
11:7:267:TYR:CZ	9:d:12:PRO:HB3	2.49	0.48
11:f:158:THR:HG23	11:f:181:LEU:HD21	1.95	0.48
5:E:382:HIS:CD2	9:5:73:GLU:HB3	2.49	0.47
7:3:420:ARG:NH1	9:5:501:THR:OG1	2.47	0.47
6:a:220:ASP:OD1	6:a:222:THR:OG1	2.28	0.47
6:2:527:VAL:HG13	9:5:581:ASN:HB2	1.96	0.47
1:R:151:GLY:O	1:R:155:ILE:HG13	2.15	0.47
6:2:397:VAL:HG21	6:2:403:PRO:HG3	1.96	0.47
6:2:703:HIS:CE1	10:6:565:LEU:HD13	2.48	0.47
10:e:769:ALA:HB1	10:e:824:ILE:HD11	1.95	0.47
6:2:688:ASP:OD1	6:2:689:GLU:N	2.47	0.47
9:5:624:SER:O	9:5:628:THR:HG23	2.14	0.47
11:7:421:GLU:C	11:7:615:HIS:HE2	2.21	0.47
11:f:92:LYS:HD3	11:f:95:GLN:NE2	2.30	0.47
13:3:1001:ADP:O3B	9:5:549:ARG:NH2	2.47	0.47
8:c:401:GLU:HB2	8:c:415:ILE:HD12	1.97	0.47
5:E:566:PRO:HA	5:E:586:PRO:HD3	1.96	0.47
9:5:182:MET:HG3	9:5:189:THR:HG22	1.95	0.47
11:7:228:ARG:NH2	11:7:327:ILE:O	2.46	0.47
8:c:647:GLU:OE1	10:e:599:SER:OG	2.22	0.47
7:3:483:ARG:HA	7:3:486:ILE:HG22	1.96	0.47
8:4:447:ASN:HD21	8:4:450:GLN:HB2	1.79	0.47
9:5:66:GLU:HA	9:5:69:ILE:HG22	1.96	0.47
10:6:598:THR:HG23	10:6:638:ILE:HG23	1.96	0.47
6:a:774:ILE:HG22	6:a:776:PRO:HD3	1.96	0.47
11:f:29:LYS:HE2	11:f:61:PRO:HG3	1.96	0.47
11:f:361:THR:N	11:f:373:GLU:OE2	2.47	0.47
11:f:479:ARG:NE	11:f:517:ASP:O	2.41	0.47
10:6:795:ILE:HG23	10:6:799:GLN:HG3	1.97	0.47
10:e:147:ASP:OD2	10:e:261:ARG:NH2	2.48	0.47
9:5:438:TYR:OH	9:5:480:ASP:OD2	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:e:333:CYS:HB2	10:e:340:ASN:HB3	1.97	0.47
4:Y:-7:DT:H2"	4:Y:-6:DG:C8	2.50	0.47
6:2:821:ALA:HB2	6:2:829:VAL:HG22	1.97	0.47
9:5:618:ALA:HB1	9:5:677:VAL:HG21	1.97	0.47
9:d:555:ILE:HD12	9:d:694:GLN:HE22	1.79	0.47
11:f:483:THR:HG22	11:f:485:GLY:H	1.80	0.47
8:c:830:ARG:HA	8:c:833:ILE:HG22	1.97	0.46
9:d:485:MET:SD	9:d:490:ARG:NH1	2.88	0.46
9:d:536:PRO:HA	9:d:539:ASN:HB2	1.97	0.46
3:X:22:DT:H2"	3:X:23:DG:C8	2.51	0.46
8:4:512:VAL:O	8:4:515:ARG:HG2	2.15	0.46
9:5:378:ILE:HD13	9:5:425:LEU:HD21	1.97	0.46
9:5:421:ALA:HB2	13:5:801:ADP:C8	2.51	0.46
7:b:553:ILE:HB	9:d:634:LEU:HD13	1.98	0.46
9:d:298:TYR:HA	9:d:328:ILE:HD13	1.96	0.46
6:a:365:THR:O	6:a:368:LYS:NZ	2.38	0.46
7:b:480:ASP:HA	7:b:483:ARG:HB2	1.96	0.46
2:I:268:SER:O	2:I:272:LYS:HG3	2.15	0.46
9:5:53:ASN:ND2	9:5:60:SER:O	2.35	0.46
9:5:455:ARG:HA	9:5:461:GLU:O	2.16	0.46
9:d:32:LYS:HD3	9:d:100:ARG:NH2	2.30	0.46
11:f:92:LYS:HD3	11:f:95:GLN:HE21	1.80	0.46
5:E:266:ASN:OD1	5:E:267:LEU:N	2.48	0.46
10:e:333:CYS:HA	10:e:344:TRP:HH2	1.81	0.46
10:e:759:ARG:O	10:e:812:ARG:NH2	2.48	0.46
3:X:26:DT:H2"	3:X:27:DG:C8	2.51	0.46
5:E:318:LEU:HD13	5:E:411:ARG:HB2	1.97	0.46
8:4:486:MET:O	8:4:490:VAL:HG23	2.16	0.46
9:5:559:ASP:N	9:5:559:ASP:OD1	2.48	0.46
10:6:557:LYS:HB2	10:6:565:LEU:HD12	1.97	0.46
10:6:592:ALA:HB3	10:6:595:SER:OG	2.16	0.46
6:a:656:ARG:HG3	10:e:794:ARG:HH21	1.80	0.46
4:Y:-38:DC:H2"	4:Y:-37:DA:C8	2.51	0.46
5:E:525:TYR:HB3	5:E:605:PHE:CZ	2.51	0.46
8:c:553:THR:OG1	10:e:739:ASP:OD2	2.27	0.46
8:c:761:ILE:HG12	8:c:815:ASN:O	2.16	0.46
11:f:435:LEU:HD12	11:f:454:ILE:HB	1.98	0.46
7:3:573:PRO:HG2	7:3:576:TYR:HB2	1.98	0.46
8:4:194:PHE:HE2	8:4:276:ILE:HG13	1.80	0.46
10:6:194:PRO:O	10:6:197:LEU:HB2	2.16	0.46
5:E:133:ASN:HB3	5:E:140:ILE:HG21	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:563:GLN:HG2	5:E:564:LEU:H	1.79	0.45
7:3:351:ASN:O	7:3:355:LYS:HG2	2.15	0.45
7:3:558:ASP:OD2	9:5:630:ARG:NE	2.39	0.45
7:3:676:ILE:HD12	11:7:617:THR:OG1	2.16	0.45
9:5:551:ASP:CG	9:5:655:ALA:HB1	2.41	0.45
11:f:636:SER:O	11:f:640:GLU:HG3	2.16	0.45
6:2:528:ASN:HB3	6:2:530:LYS:HG2	1.98	0.45
9:d:13:VAL:HG12	9:d:14:LEU:HD13	1.97	0.45
5:E:66:GLU:HB3	5:E:70:HIS:CD2	2.51	0.45
5:E:127:ARG:NH2	5:E:144:ASP:OD1	2.49	0.45
11:7:378:ALA:O	11:7:379:GLN:NE2	2.38	0.45
3:X:29:DT:H2"	3:X:30:DG:C8	2.51	0.45
7:3:474:GLU:OE1	7:3:517:ASN:ND2	2.42	0.45
11:f:312:GLU:OE1	11:f:334:HIS:NE2	2.49	0.45
3:X:51:DT:H2"	3:X:52:DG:C8	2.51	0.45
4:Y:-5:DC:H2"	4:Y:-4:DA:C8	2.51	0.45
8:4:318:ASN:OD1	11:7:341:ARG:NH1	2.32	0.45
6:a:527:VAL:HG23	6:a:531:HIS:HB3	1.99	0.45
8:c:256:ASP:HB3	11:f:134:TYR:CD2	2.52	0.45
9:d:259:GLN:HG2	9:d:260:GLU:N	2.31	0.45
5:E:541:ASN:OD1	5:E:544:THR:N	2.48	0.45
6:a:197:TRP:O	6:a:203:VAL:HG11	2.17	0.45
9:d:411:ASN:OD1	9:d:519:VAL:N	2.47	0.45
5:E:386:ARG:NH1	9:5:74:ASP:OD1	2.49	0.45
10:e:524:HIS:HB3	10:e:528:LYS:HE3	1.99	0.45
5:E:377:TRP:HA	5:E:380:MET:HB2	1.99	0.45
7:3:343:THR:OG1	7:3:346:ASP:OD2	2.27	0.45
9:5:485:MET:SD	9:5:490:ARG:NH1	2.90	0.45
3:X:16:DC:OP1	6:a:587:LYS:HD2	2.17	0.45
5:E:11:ALA:O	5:E:15:ILE:HG13	2.17	0.45
7:3:262:PRO:HG3	9:5:462:PHE:HE1	1.82	0.45
9:5:444:SER:HB2	9:5:447:ALA:HB3	1.99	0.45
10:6:509:SER:O	10:6:513:ILE:HG13	2.17	0.45
10:6:696:ARG:HA	10:6:706:MET:HE1	1.98	0.45
5:E:25:CYS:HB2	5:E:55:GLN:HG2	1.99	0.45
9:5:525:PRO:HG2	9:5:530:TYR:HB2	1.99	0.45
6:a:583:ASP:O	6:a:587:LYS:N	2.43	0.45
10:e:529:LEU:O	10:e:533:ILE:HG12	2.17	0.45
11:f:526:PHE:CD1	11:f:529:MET:HG3	2.52	0.45
5:E:277:THR:HG22	5:E:288:TYR:CE1	2.51	0.44
7:b:402:ASP:OD2	7:b:493:GLN:NE2	2.37	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:Y:-21:DT:H4'	4:Y:-20:DG:OP1	2.16	0.44
11:7:360:TYR:HA	11:7:373:GLU:HG2	1.98	0.44
9:d:82:GLU:O	9:d:86:ILE:HG12	2.17	0.44
9:d:674:GLU:HA	9:d:677:VAL:HB	1.97	0.44
5:E:330:ARG:HA	5:E:377:TRP:CE3	2.52	0.44
9:5:341:SER:N	9:5:344:ASN:OD1	2.43	0.44
9:d:379:PHE:N	13:d:801:ADP:N1	2.65	0.44
10:6:766:THR:O	10:6:770:ARG:HG3	2.17	0.44
11:7:249:SER:O	11:7:311:GLN:NE2	2.50	0.44
8:c:503:ASP:OD1	8:c:503:ASP:N	2.51	0.44
10:e:526:TYR:O	10:e:530:VAL:HG23	2.18	0.44
4:Y:-32:DT:H2''	4:Y:-31:DG:C8	2.52	0.44
2:I:409:ALA:O	2:I:413:ILE:HG13	2.17	0.44
2:I:529:GLN:HA	2:I:532:GLN:NE2	2.31	0.44
6:2:404:ARG:NH2	10:6:387:GLU:OE2	2.51	0.44
6:a:520:PHE:CE1	6:a:767:ILE:HG12	2.52	0.44
6:a:703:HIS:CE1	10:e:565:LEU:HD13	2.53	0.44
10:e:120:GLU:OE2	10:e:191:LYS:NZ	2.46	0.44
6:2:437:ASN:OD1	6:2:440:ALA:N	2.49	0.44
6:a:300:PHE:HB3	6:a:319:ARG:HD2	1.99	0.44
8:c:251:TYR:HD2	8:c:253:GLN:HB2	1.83	0.44
7:3:553:ILE:HB	9:5:634:LEU:HD13	2.00	0.44
9:5:12:PRO:HB3	11:f:267:TYR:CZ	2.53	0.44
5:E:485:ASP:OD2	5:E:487:ARG:HB2	2.18	0.44
8:4:481:ILE:O	8:4:485:LEU:HG	2.18	0.44
11:7:31:ASP:CG	11:7:87:GLN:HE22	2.26	0.44
6:a:679:ILE:HD11	6:a:816:ILE:HD11	1.98	0.44
7:b:212:ARG:HD3	7:b:229:ALA:HB1	2.00	0.44
11:f:193:PRO:HD2	11:f:196:LEU:HD12	2.00	0.44
11:f:388:PHE:HE1	11:f:639:ARG:HG2	1.83	0.44
3:X:4:DT:H2''	3:X:5:DG:C8	2.53	0.44
6:2:690:GLU:CD	6:2:690:GLU:H	2.25	0.44
9:5:685:GLN:O	9:5:689:MET:HB2	2.18	0.44
8:c:407:PRO:HG2	8:c:410:GLN:HB2	1.99	0.44
8:c:701:ARG:HH21	8:c:796:ARG:HH21	1.65	0.44
5:E:330:ARG:HG3	5:E:331:HIS:NE2	2.33	0.43
6:2:227:TYR:OH	6:2:248:HIS:ND1	2.41	0.43
8:4:767:LYS:HG3	10:6:732:VAL:HG11	2.00	0.43
9:5:224:GLU:O	9:5:227:ILE:HG12	2.18	0.43
6:2:520:PHE:CD2	6:2:823:MET:HG2	2.53	0.43
8:4:262:LEU:HD12	8:4:262:LEU:HA	1.87	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:5:642:GLU:O	9:5:697:ILE:HG12	2.18	0.43
7:b:529:PRO:HA	7:b:532:ASN:HB2	2.00	0.43
8:c:194:PHE:HE2	8:c:276:ILE:HG13	1.83	0.43
4:Y:-17:DT:H2"	4:Y:-16:DG:N7	2.34	0.43
5:E:71:TYR:CE2	5:E:96:LEU:HB3	2.53	0.43
9:5:230:GLU:OE2	7:b:318:LYS:NZ	2.37	0.43
11:7:210:ASN:OD1	11:7:213:ARG:NH2	2.51	0.43
6:2:314:LEU:HD13	6:2:432:ASN:HD21	1.83	0.43
7:3:318:LYS:NZ	9:d:230:GLU:OE2	2.31	0.43
7:3:678:VAL:HG21	7:3:723:LYS:HG3	1.99	0.43
9:5:535:SER:O	9:5:539:ASN:ND2	2.52	0.43
10:6:737:LYS:HB2	10:6:741:ALA:HB2	1.99	0.43
6:a:244:VAL:N	6:a:296:ARG:O	2.51	0.43
6:2:520:PHE:CE2	6:2:823:MET:HG2	2.53	0.43
7:b:442:LEU:HD23	7:b:442:LEU:HA	1.83	0.43
10:e:323:GLN:O	10:e:328:THR:OG1	2.37	0.43
10:e:766:THR:O	10:e:770:ARG:HG3	2.18	0.43
5:E:33:CYS:SG	5:E:34:LEU:HG	2.59	0.43
6:2:474:PHE:CZ	6:2:561:HIS:HA	2.54	0.43
8:4:212:ARG:HD3	8:4:212:ARG:HA	1.83	0.43
10:6:330:PRO:HD2	10:6:344:TRP:CD1	2.54	0.43
11:f:439:GLY:O	11:f:701:LYS:NZ	2.39	0.43
5:E:13:ASN:O	5:E:17:ARG:HG3	2.18	0.43
5:E:425:VAL:O	5:E:429:THR:HG23	2.19	0.43
10:6:529:LEU:O	10:6:533:ILE:HG12	2.19	0.43
9:d:152:ASP:OD1	9:d:155:HIS:ND1	2.43	0.43
9:d:568:ILE:O	9:d:572:VAL:HG23	2.18	0.43
5:E:30:PHE:HB3	5:E:86:PHE:HZ	1.83	0.43
8:4:503:ASP:N	8:4:503:ASP:OD1	2.50	0.43
10:6:526:TYR:O	10:6:530:VAL:HG23	2.18	0.43
10:6:641:PHE:O	10:6:644:MET:HG2	2.19	0.43
10:e:303:GLU:N	10:e:354:LEU:O	2.39	0.43
10:e:313:MET:HB2	10:e:338:CYS:SG	2.58	0.43
10:e:537:VAL:HG11	10:e:584:PHE:CZ	2.53	0.43
5:E:66:GLU:H	5:E:69:ARG:HH21	1.67	0.43
5:E:638:SER:O	5:E:642:GLU:HG3	2.19	0.43
8:4:784:SER:HB2	8:4:787:ASP:HB2	2.00	0.43
8:c:315:ARG:HD2	8:c:413:HIS:CD2	2.54	0.43
4:Y:-42:DC:H2"	4:Y:-41:DA:C8	2.54	0.42
5:E:68:ARG:HA	5:E:71:TYR:HB3	2.01	0.42
6:2:632:SER:HA	6:2:636:ILE:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:5:456:ASP:HB2	9:5:461:GLU:H	1.84	0.42
7:b:541:SER:OG	7:b:700:ARG:NH1	2.50	0.42
8:c:183:THR:O	11:f:145:GLN:NE2	2.48	0.42
8:c:489:LYS:O	8:c:492:HIS:ND1	2.49	0.42
10:e:820:THR:OG1	10:e:822:SER:OG	2.31	0.42
5:E:39:LEU:HD11	5:E:481:TRP:HE3	1.84	0.42
5:E:370:LEU:HG	5:E:374:GLN:NE2	2.34	0.42
6:2:597:VAL:HG11	6:2:640:LEU:HD13	2.00	0.42
6:a:264:PRO:O	6:a:268:LEU:HG	2.18	0.42
6:a:520:PHE:CD2	6:a:823:MET:HG2	2.54	0.42
8:4:262:LEU:HD11	8:4:308:VAL:HG23	2.00	0.42
6:a:769:TYR:CE1	6:a:773:LYS:HD2	2.54	0.42
9:d:610:CYS:HB3	9:d:665:LYS:HG3	2.01	0.42
10:e:492:ASN:ND2	10:e:495:GLN:OE1	2.44	0.42
10:e:509:SER:O	10:e:513:ILE:HG13	2.19	0.42
2:I:416:LYS:HE3	5:E:522:LEU:HD23	2.01	0.42
6:2:583:ASP:OD2	6:2:586:THR:HG22	2.19	0.42
10:6:832:ARG:HA	10:6:835:ILE:HD12	2.01	0.42
8:c:590:TYR:HE1	11:f:543:GLN:HE22	1.66	0.42
8:c:732:LYS:HE3	11:f:445:GLY:HA2	2.00	0.42
11:f:378:ALA:C	11:f:379:GLN:HE21	2.26	0.42
9:5:373:SER:O	9:5:376:PRO:HD3	2.20	0.42
10:6:521:LYS:HE2	10:6:521:LYS:HA	2.00	0.42
7:b:450:ARG:CZ	7:b:456:ARG:HB3	2.49	0.42
9:d:536:PRO:HD2	9:d:695:ASP:OD1	2.19	0.42
5:E:277:THR:HG21	5:E:295:LEU:HD11	2.02	0.42
6:2:560:ALA:HB3	6:2:563:ALA:HB2	2.01	0.42
9:5:82:GLU:O	9:5:86:ILE:HG12	2.19	0.42
9:5:610:CYS:HB3	9:5:665:LYS:HG3	2.02	0.42
11:7:451:ARG:NH1	11:7:542:GLU:O	2.53	0.42
6:a:707:HIS:HE1	6:a:709:GLU:HB3	1.85	0.42
10:e:111:VAL:HG13	10:e:166:LEU:HD11	2.01	0.42
11:f:404:LEU:HD23	11:f:641:TYR:CD2	2.55	0.42
11:f:414:LEU:O	11:f:418:ILE:HG13	2.20	0.42
11:f:459:MET:SD	11:f:584:ILE:HD13	2.60	0.42
5:E:519:ILE:HG12	5:E:528:CYS:SG	2.59	0.42
10:6:584:PHE:O	10:6:588:VAL:HG23	2.19	0.42
5:E:334:LEU:HD11	5:E:338:PHE:HE1	1.85	0.42
6:2:344:CYS:SG	6:2:346:SER:HB3	2.60	0.42
8:4:238:THR:HG22	8:4:240:ASN:H	1.84	0.42
11:7:72:ASN:OD1	11:7:129:THR:OG1	2.24	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:a:334:LEU:O	6:a:383:ARG:NH1	2.52	0.42
10:e:521:LYS:HE2	10:e:521:LYS:HA	2.02	0.42
5:E:49:PHE:O	5:E:53:LEU:N	2.53	0.42
5:E:511:VAL:O	5:E:515:GLU:HG2	2.20	0.42
5:E:564:LEU:HD22	5:E:586:PRO:HB2	2.02	0.42
8:4:204:LYS:O	8:4:208:ILE:HG12	2.19	0.42
9:5:456:ASP:N	9:5:461:GLU:O	2.52	0.42
10:6:356:TRP:CH2	10:6:358:LYS:HB2	2.54	0.42
10:e:694:SER:O	10:e:698:ASN:ND2	2.45	0.42
5:E:480:SER:O	5:E:484:LEU:HG	2.20	0.42
6:2:220:ASP:OD1	6:2:222:THR:OG1	2.29	0.42
6:2:265:GLU:OE1	6:2:431:LYS:NZ	2.48	0.42
9:5:419:GLY:H	13:5:801:ADP:PB	2.42	0.42
9:5:441:GLY:N	9:5:480:ASP:O	2.53	0.42
11:7:418:ILE:HG12	11:7:638:MET:HE1	2.02	0.42
6:a:360:ARG:HA	6:a:360:ARG:NE	2.35	0.42
7:b:412:SER:H	13:b:1001:ADP:PB	2.43	0.42
8:c:206:ARG:NH1	8:c:247:ASN:OD1	2.51	0.42
8:c:583:LYS:NZ	11:f:448:MET:SD	2.93	0.42
9:d:369:ILE:HD12	9:d:592:SER:HB3	2.02	0.42
3:X:24:DC:H2''	3:X:25:DA:C8	2.54	0.41
5:E:133:ASN:HA	5:E:137:SER:HB3	2.01	0.41
6:2:392:GLU:HB2	6:2:403:PRO:HB3	2.02	0.41
8:4:280:MET:O	8:4:284:ILE:HG12	2.20	0.41
8:4:593:GLY:N	8:4:632:ASP:O	2.53	0.41
9:d:618:ALA:HB1	9:d:677:VAL:HG21	2.02	0.41
5:E:330:ARG:NH1	5:E:418:SER:HB2	2.35	0.41
6:2:422:GLU:CD	6:2:458:ARG:HE	2.28	0.41
8:4:584:ILE:HD13	8:4:584:ILE:HA	1.92	0.41
8:4:830:ARG:HA	8:4:833:ILE:HG22	2.02	0.41
11:7:479:ARG:NE	11:7:517:ASP:O	2.48	0.41
11:7:636:SER:O	11:7:640:GLU:HG3	2.19	0.41
10:e:356:TRP:CH2	10:e:358:LYS:HB2	2.55	0.41
7:3:457:LEU:HD11	7:3:504:THR:HG21	2.02	0.41
7:3:553:ILE:HG21	9:5:634:LEU:HB2	2.02	0.41
9:5:455:ARG:HH12	9:5:462:PHE:HB3	1.84	0.41
9:5:551:ASP:OD1	9:5:655:ALA:HB1	2.20	0.41
9:5:568:ILE:O	9:5:572:VAL:HG23	2.20	0.41
6:2:695:LEU:HD11	10:6:797:VAL:HG11	2.01	0.41
8:4:432:ARG:O	8:4:473:ARG:NH2	2.53	0.41
11:7:92:LYS:HD3	11:7:95:GLN:NE2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:a:522:GLY:O	6:a:822:LYS:NZ	2.39	0.41
8:c:506:LEU:O	8:c:510:ARG:HG2	2.20	0.41
8:4:183:THR:O	11:7:145:GLN:NE2	2.54	0.41
10:6:556:HIS:CD2	10:6:567:GLY:H	2.38	0.41
6:a:437:ASN:ND2	6:a:441:LYS:HD2	2.35	0.41
8:c:234:ARG:HB2	8:c:280:MET:HE3	2.02	0.41
8:c:262:LEU:HD12	8:c:262:LEU:HA	1.84	0.41
9:d:483:ASP:HA	9:d:490:ARG:HH22	1.86	0.41
6:2:707:HIS:HE1	6:2:709:GLU:HB3	1.85	0.41
8:4:635:ASP:HA	8:4:694:LEU:HD21	2.02	0.41
9:5:415:LEU:O	9:5:555:ILE:HA	2.21	0.41
10:6:199:THR:HG23	10:6:260:GLU:HG3	2.02	0.41
10:6:749:GLU:CD	10:6:749:GLU:H	2.29	0.41
11:7:378:ALA:C	11:7:379:GLN:HE21	2.25	0.41
7:b:567:ARG:NE	7:b:627:PRO:O	2.54	0.41
10:e:626:GLY:O	10:e:630:LEU:HD13	2.21	0.41
11:f:317:GLU:OE1	11:f:329:ARG:N	2.44	0.41
6:2:195:SER:O	6:2:199:THR:HG23	2.20	0.41
7:3:262:PRO:HG3	9:5:462:PHE:CE1	2.55	0.41
8:4:698:LEU:HA	8:4:701:ARG:HG3	2.01	0.41
11:f:621:MET:HE3	11:f:622:HIS:CE1	2.55	0.41
11:f:674:GLU:OE1	11:f:728:TYR:OH	2.26	0.41
7:3:201:HIS:NE2	7:3:232:PRO:HG2	2.35	0.41
8:4:239:SER:O	8:4:239:SER:OG	2.38	0.41
8:4:579:GLN:NE2	11:7:543:GLN:OE1	2.54	0.41
9:5:397:LYS:C	9:5:398:LYS:HD3	2.46	0.41
10:6:433:LEU:HG	6:a:342:LEU:HB3	2.02	0.41
10:6:769:ALA:HB1	10:6:824:ILE:HD11	2.03	0.41
2:H:537:ASP:O	11:7:152:ARG:NE	2.49	0.41
5:E:268:SER:O	5:E:272:LEU:HG	2.21	0.41
5:E:494:ARG:O	5:E:498:LEU:HG	2.21	0.41
7:3:250:PHE:O	7:3:251:ILE:HD13	2.20	0.41
7:3:391:LYS:HB2	7:3:399:LEU:HD12	2.03	0.41
8:4:204:LYS:HE2	8:4:204:LYS:HB3	1.86	0.41
9:5:346:VAL:HG22	9:5:348:MET:H	1.85	0.41
9:5:374:ILE:HG23	9:5:428:PHE:HE2	1.86	0.41
9:5:621:LYS:NZ	9:5:678:ASP:OD1	2.54	0.41
11:7:404:LEU:HD23	11:7:641:TYR:CD2	2.55	0.41
11:7:709:ASP:O	11:7:713:VAL:HG23	2.21	0.41
8:c:455:SER:OG	11:f:275:SER:O	2.35	0.41
10:e:144:LYS:HD3	10:e:193:ALA:HB1	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:583:GLY:HA3	5:E:609:PHE:HZ	1.84	0.41
6:2:287:ALA:HA	6:2:290:HIS:O	2.21	0.41
7:3:711:ALA:O	7:3:715:VAL:HG23	2.20	0.41
10:6:610:ALA:HB3	10:6:663:ILE:HD11	2.03	0.41
7:b:573:PRO:HA	7:b:574:PRO:HD3	2.00	0.41
8:c:360:ILE:HD11	8:c:363:GLY:HA2	2.03	0.41
2:H:529:GLN:HA	2:H:532:GLN:HE21	1.86	0.40
5:E:45:LEU:HD11	5:E:49:PHE:CZ	2.56	0.40
10:6:154:ASP:OD1	10:6:156:GLN:N	2.50	0.40
10:e:184:GLY:O	10:e:188:VAL:HG23	2.21	0.40
2:I:351:CYS:O	2:I:355:ILE:HG13	2.22	0.40
6:a:796:GLU:O	6:a:800:THR:OG1	2.29	0.40
5:E:363:PHE:CE1	5:E:368:ILE:HB	2.57	0.40
7:3:122:ILE:HB	7:3:123:PRO:HD3	2.02	0.40
8:4:194:PHE:HB3	8:4:275:THR:HG21	2.03	0.40
9:5:31:PHE:O	9:5:35:ILE:HG12	2.22	0.40
10:6:571:ILE:HG12	10:6:711:LEU:HB2	2.03	0.40
6:a:697:THR:HA	6:a:700:VAL:HG12	2.02	0.40
8:c:323:ASP:OD1	8:c:440:ARG:NH1	2.51	0.40
8:c:584:ILE:HD13	8:c:584:ILE:HA	1.88	0.40
8:c:754:ALA:HA	8:c:758:ILE:HD13	2.03	0.40
10:e:600:GLY:O	10:e:603:SER:OG	2.36	0.40
5:E:28:VAL:O	5:E:81:LEU:HA	2.21	0.40
5:E:387:GLU:O	5:E:391:ILE:HG13	2.21	0.40
8:c:408:ASP:OD1	8:c:408:ASP:N	2.53	0.40
8:c:717:ASP:CG	11:f:668:ARG:HE	2.30	0.40
8:c:851:GLN:N	8:c:851:GLN:CD	2.79	0.40
9:d:663:LEU:O	9:d:666:LEU:HG	2.20	0.40
4:Y:-41:DA:H2'	4:Y:-40:DT:H71	2.04	0.40
7:3:433:THR:HG23	7:3:436:GLY:H	1.86	0.40
8:4:758:ILE:HG12	8:4:813:LEU:HA	2.04	0.40
9:5:536:PRO:HD2	9:5:695:ASP:OD1	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	A	5/257 (2%)	5 (100%)	0	0	100	100
1	R	8/257 (3%)	8 (100%)	0	0	100	100
2	H	13/704 (2%)	13 (100%)	0	0	100	100
2	I	75/704 (11%)	75 (100%)	0	0	100	100
5	E	420/657 (64%)	413 (98%)	7 (2%)	0	100	100
6	2	617/868 (71%)	611 (99%)	6 (1%)	0	100	100
6	a	617/868 (71%)	610 (99%)	7 (1%)	0	100	100
7	3	628/1006 (62%)	615 (98%)	13 (2%)	0	100	100
7	b	628/1006 (62%)	615 (98%)	13 (2%)	0	100	100
8	4	668/933 (72%)	648 (97%)	20 (3%)	0	100	100
8	c	668/933 (72%)	655 (98%)	13 (2%)	0	100	100
9	5	643/775 (83%)	632 (98%)	11 (2%)	0	100	100
9	d	643/775 (83%)	632 (98%)	11 (2%)	0	100	100
10	6	625/1017 (62%)	612 (98%)	13 (2%)	0	100	100
10	e	625/1017 (62%)	609 (97%)	16 (3%)	0	100	100
11	7	696/845 (82%)	679 (98%)	17 (2%)	0	100	100
11	f	696/845 (82%)	682 (98%)	14 (2%)	0	100	100
All	All	8275/13467 (61%)	8114 (98%)	161 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	A	6/234 (3%)	6 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	R	9/234 (4%)	9 (100%)	0	100	100
2	H	15/654 (2%)	15 (100%)	0	100	100
2	I	80/654 (12%)	80 (100%)	0	100	100
5	E	400/592 (68%)	400 (100%)	0	100	100
6	2	547/770 (71%)	547 (100%)	0	100	100
6	a	547/770 (71%)	547 (100%)	0	100	100
7	3	550/864 (64%)	550 (100%)	0	100	100
7	b	550/864 (64%)	550 (100%)	0	100	100
8	4	611/848 (72%)	611 (100%)	0	100	100
8	c	611/848 (72%)	611 (100%)	0	100	100
9	5	583/688 (85%)	583 (100%)	0	100	100
9	d	583/688 (85%)	583 (100%)	0	100	100
10	6	557/886 (63%)	557 (100%)	0	100	100
10	e	557/886 (63%)	557 (100%)	0	100	100
11	7	619/753 (82%)	619 (100%)	0	100	100
11	f	619/753 (82%)	619 (100%)	0	100	100
All	All	7444/11986 (62%)	7444 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (53) such sidechains are listed below:

Mol	Chain	Res	Type
5	E	18	ASN
5	E	77	ASN
5	E	321	GLN
5	E	500	GLN
6	2	282	HIS
6	2	294	HIS
6	2	432	ASN
6	2	760	GLN
6	2	849	GLN
7	3	177	ASN
7	3	259	GLN
7	3	374	HIS
7	3	487	HIS

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Mol	Chain	Res	Type
8	4	386	HIS
8	4	563	ASN
8	4	723	HIS
8	4	726	ASN
9	5	223	ASN
9	5	411	ASN
9	5	571	HIS
9	5	579	ASN
10	6	156	GLN
10	6	633	ASN
11	7	178	ASN
11	7	292	ASN
11	7	311	GLN
11	7	336	ASN
11	7	518	ASN
11	7	568	ASN
6	a	528	ASN
6	a	641	GLN
6	a	651	ASN
7	b	29	GLN
7	b	239	ASN
7	b	259	GLN
7	b	417	GLN
7	b	424	ASN
7	b	530	HIS
8	c	563	ASN
8	c	685	ASN
9	d	223	ASN
9	d	259	GLN
9	d	499	GLN
9	d	576	HIS
9	d	579	ASN
9	d	581	ASN
10	e	633	ASN
10	e	814	ASN
11	f	112	HIS
11	f	178	ASN
11	f	336	ASN
11	f	412	ASN
11	f	568	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 18 ligands modelled in this entry, 14 are monoatomic - leaving 4 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
13	ADP	b	1001	14	27,29,29	1.36	4 (14%)	42,45,45	2.00	10 (23%)
13	ADP	d	801	14	27,29,29	1.37	4 (14%)	42,45,45	1.98	10 (23%)
13	ADP	5	801	14	27,29,29	1.37	4 (14%)	42,45,45	2.04	10 (23%)
13	ADP	3	1001	14	27,29,29	1.36	4 (14%)	42,45,45	1.99	10 (23%)

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
13	ADP	b	1001	14	-	5/16/32/32	0/3/3/3
13	ADP	d	801	14	-	3/16/32/32	0/3/3/3
13	ADP	5	801	14	-	0/16/32/32	0/3/3/3
13	ADP	3	1001	14	-	5/16/32/32	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
13	5	801	ADP	C5-C4	4.62	1.47	1.39
13	d	801	ADP	C5-C4	4.60	1.47	1.39
13	b	1001	ADP	C5-C4	4.54	1.47	1.39
13	3	1001	ADP	C5-C4	4.52	1.47	1.39
13	5	801	ADP	C5-C6	2.73	1.48	1.41
13	d	801	ADP	C5-C6	2.70	1.48	1.41
13	b	1001	ADP	C5-C6	2.67	1.48	1.41
13	3	1001	ADP	C5-C6	2.66	1.48	1.41
13	d	801	ADP	C8-N7	2.43	1.36	1.31
13	b	1001	ADP	C8-N7	2.39	1.36	1.31
13	5	801	ADP	C8-N7	2.37	1.36	1.31
13	3	1001	ADP	C8-N7	2.35	1.36	1.31
13	3	1001	ADP	C5-N7	-2.24	1.34	1.39
13	5	801	ADP	C5-N7	-2.21	1.34	1.39
13	d	801	ADP	C5-N7	-2.17	1.34	1.39
13	b	1001	ADP	C5-N7	-2.14	1.35	1.39

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	5	801	ADP	C5-C4-N3	-6.57	118.17	126.75
13	b	1001	ADP	C5-C4-N3	-6.37	118.44	126.75
13	3	1001	ADP	C5-C4-N3	-6.30	118.53	126.75
13	d	801	ADP	C5-C4-N3	-6.29	118.55	126.75
13	5	801	ADP	N3-C4-N9	5.18	135.62	127.08
13	3	1001	ADP	N3-C4-N9	4.96	135.25	127.08
13	b	1001	ADP	N3-C4-N9	4.95	135.25	127.08
13	d	801	ADP	N3-C4-N9	4.89	135.13	127.08
13	d	801	ADP	PA-O3A-PB	-4.00	119.09	132.83
13	5	801	ADP	C2-N3-C4	4.00	121.20	111.75
13	5	801	ADP	PA-O3A-PB	-3.98	119.17	132.83
13	b	1001	ADP	C2-N3-C4	3.89	120.94	111.75
13	d	801	ADP	C2-N3-C4	3.87	120.90	111.75
13	3	1001	ADP	C2-N3-C4	3.84	120.83	111.75
13	b	1001	ADP	PA-O3A-PB	-3.80	119.80	132.83
13	3	1001	ADP	PA-O3A-PB	-3.73	120.02	132.83
13	d	801	ADP	C4-C5-N7	-3.24	106.67	110.62
13	5	801	ADP	C4-C5-N7	-3.22	106.70	110.62
13	b	1001	ADP	C4-C5-N7	-3.20	106.72	110.62
13	3	1001	ADP	C4-C5-N7	-3.17	106.75	110.62
13	d	801	ADP	N3-C2-N1	-3.13	123.71	128.60
13	5	801	ADP	N3-C2-N1	-3.10	123.76	128.60

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	b	1001	ADP	N3-C2-N1	-3.09	123.77	128.60
13	3	1001	ADP	N3-C2-N1	-3.04	123.85	128.60
13	5	801	ADP	C5-N7-C8	2.75	107.41	103.51
13	d	801	ADP	C5-N7-C8	2.71	107.36	103.51
13	3	1001	ADP	C5-N7-C8	2.68	107.31	103.51
13	b	1001	ADP	C5-N7-C8	2.63	107.24	103.51
13	b	1001	ADP	C3'-C2'-C1'	2.60	106.36	101.43
13	5	801	ADP	C3'-C2'-C1'	2.51	106.20	101.43
13	3	1001	ADP	C4-N9-C8	2.49	108.43	105.73
13	3	1001	ADP	C3'-C2'-C1'	2.49	106.16	101.43
13	d	801	ADP	C4-N9-C8	2.41	108.33	105.73
13	b	1001	ADP	C4-N9-C8	2.38	108.31	105.73
13	5	801	ADP	C4-N9-C8	2.37	108.29	105.73
13	d	801	ADP	C3'-C2'-C1'	2.24	105.69	101.43
13	d	801	ADP	C6-C5-N7	2.09	135.92	132.02
13	b	1001	ADP	C6-C5-N7	2.03	135.79	132.02
13	3	1001	ADP	C6-C5-N7	2.02	135.79	132.02
13	5	801	ADP	C6-C5-N7	2.00	135.75	132.02

There are no chirality outliers.

All (13) torsion outliers are listed below:

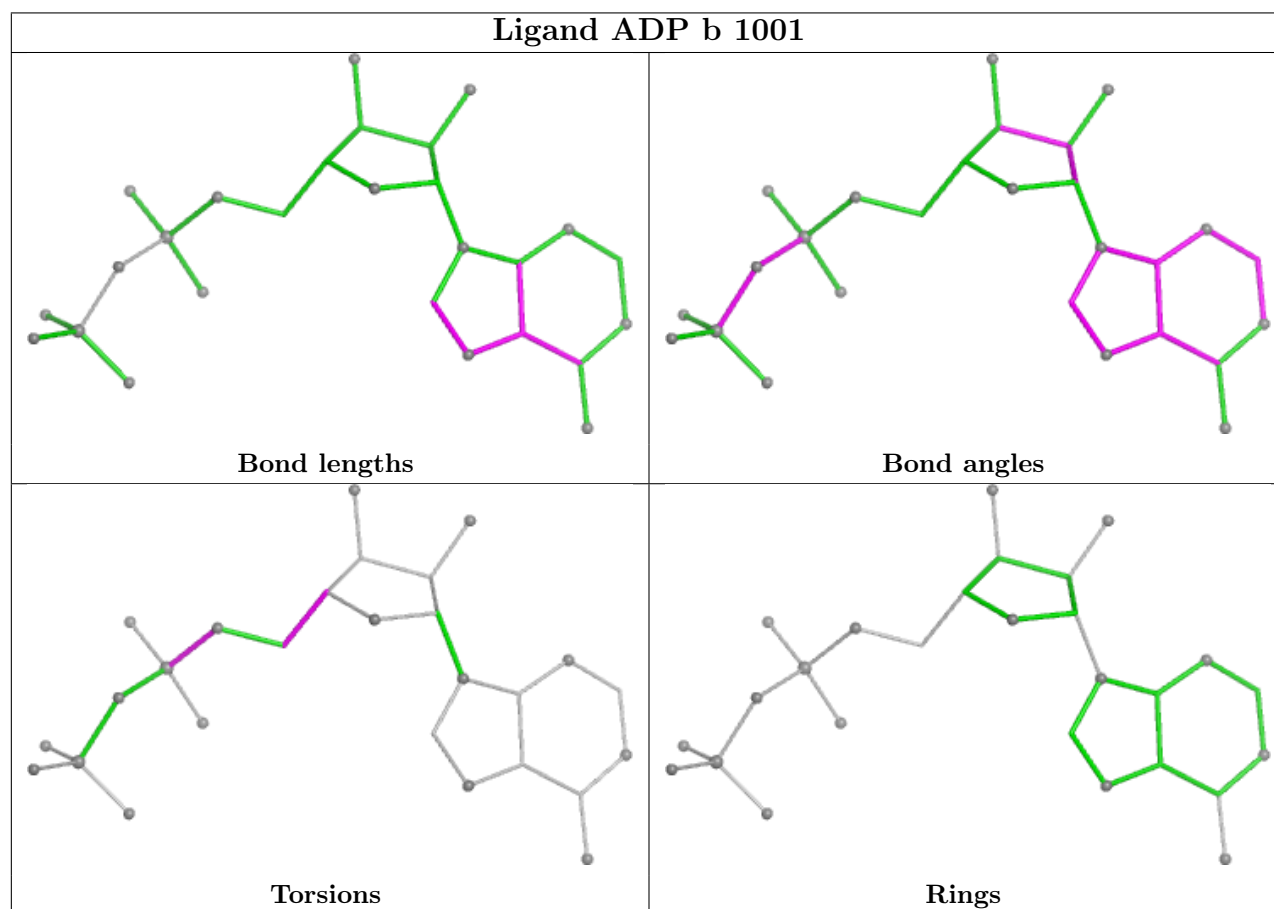
Mol	Chain	Res	Type	Atoms
13	3	1001	ADP	C5'-O5'-PA-O1A
13	3	1001	ADP	C5'-O5'-PA-O2A
13	3	1001	ADP	O4'-C4'-C5'-O5'
13	b	1001	ADP	C5'-O5'-PA-O1A
13	b	1001	ADP	C5'-O5'-PA-O2A
13	d	801	ADP	O4'-C4'-C5'-O5'
13	d	801	ADP	C3'-C4'-C5'-O5'
13	3	1001	ADP	C3'-C4'-C5'-O5'
13	b	1001	ADP	O4'-C4'-C5'-O5'
13	b	1001	ADP	C3'-C4'-C5'-O5'
13	d	801	ADP	PA-O3A-PB-O3B
13	3	1001	ADP	C5'-O5'-PA-O3A
13	b	1001	ADP	C5'-O5'-PA-O3A

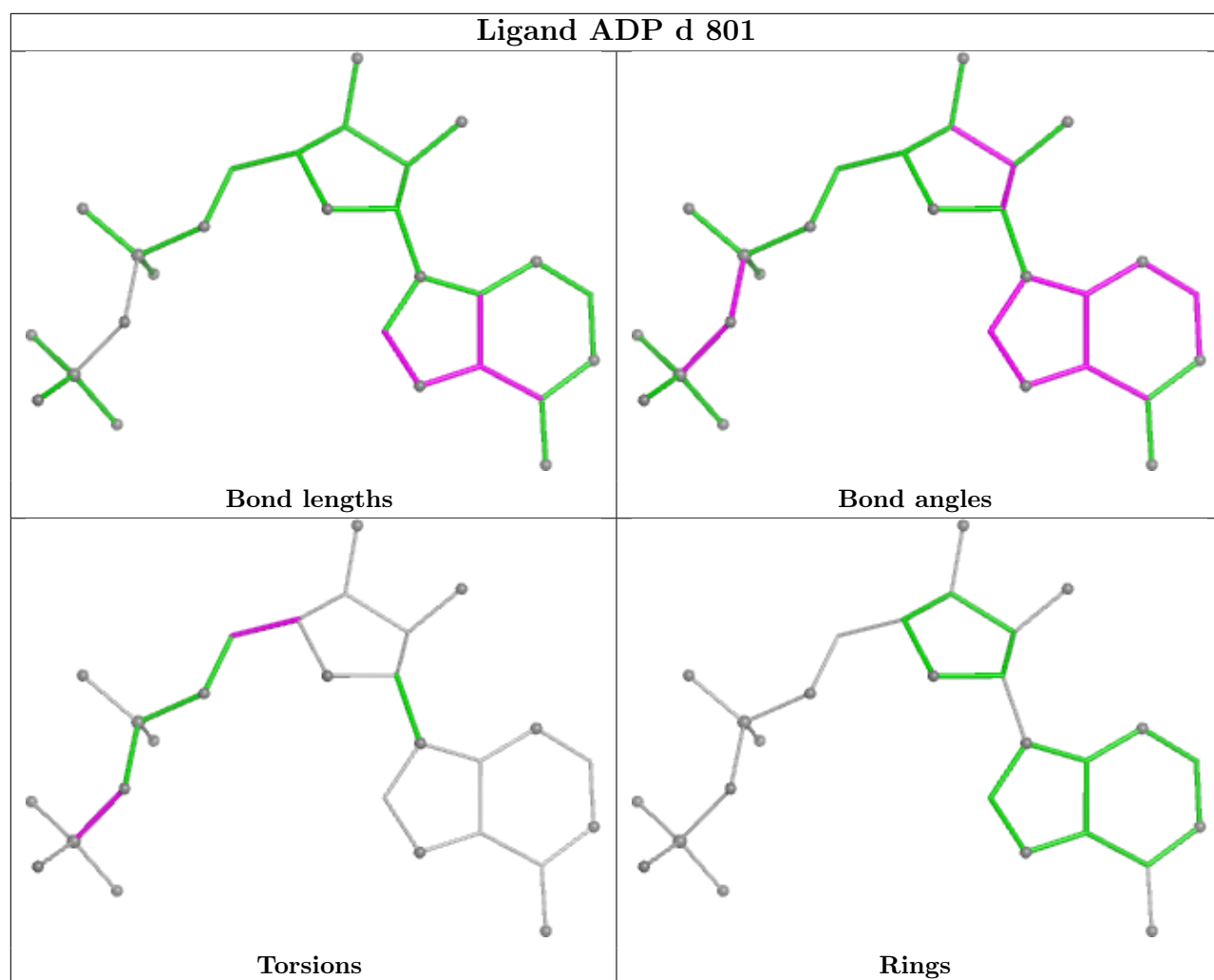
There are no ring outliers.

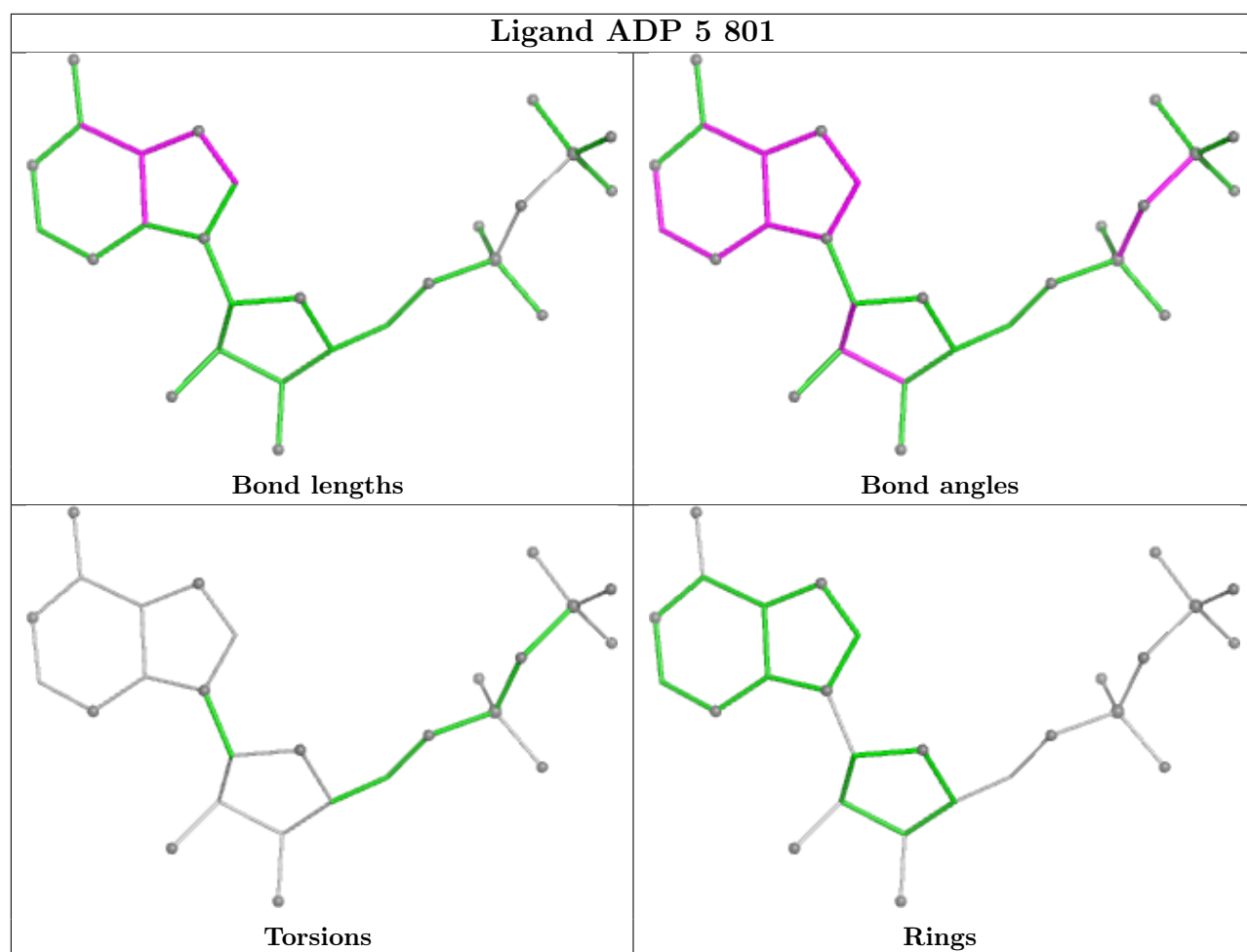
4 monomers are involved in 8 short contacts:

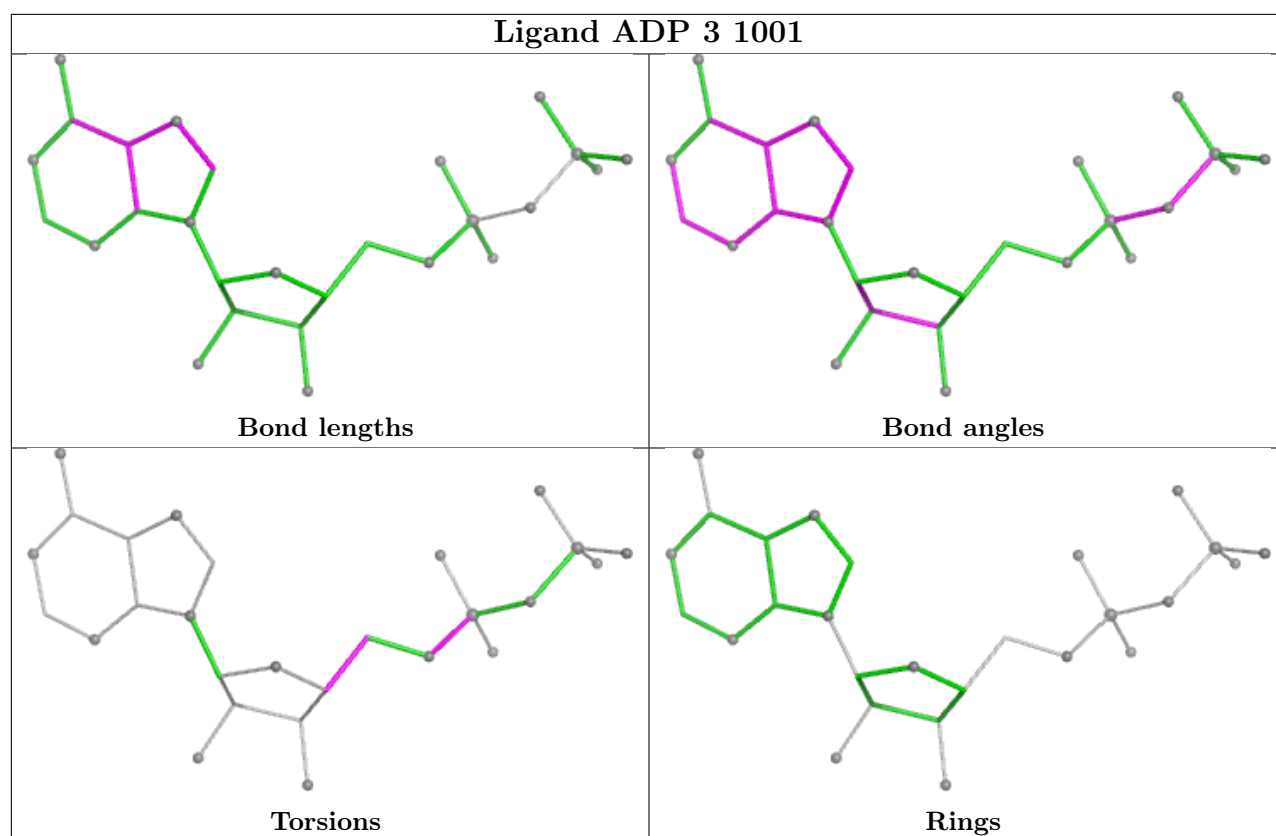
Mol	Chain	Res	Type	Clashes	Symm-Clashes
13	b	1001	ADP	1	0
13	d	801	ADP	3	0
13	5	801	ADP	3	0
13	3	1001	ADP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

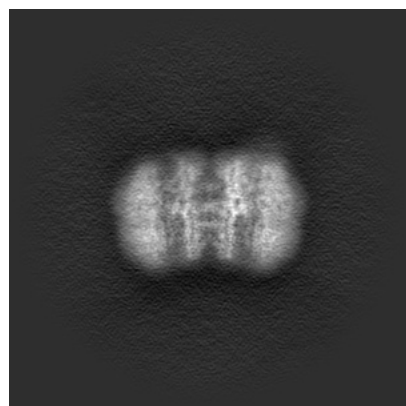
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-53973. These allow visual inspection of the internal detail of the map and identification of artifacts.

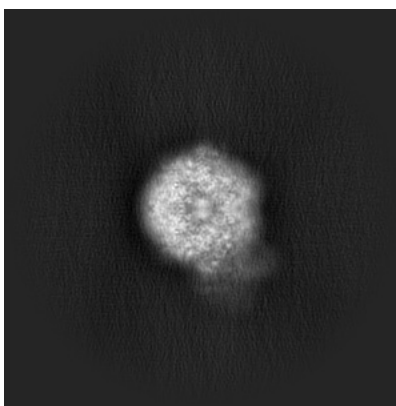
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

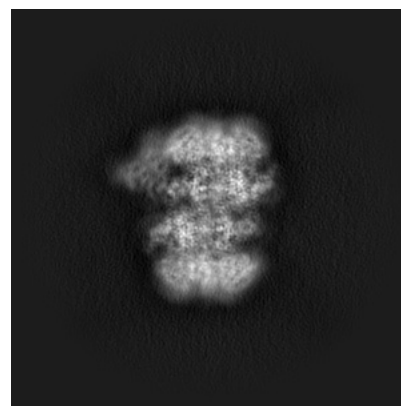
6.1.1 Primary map



X

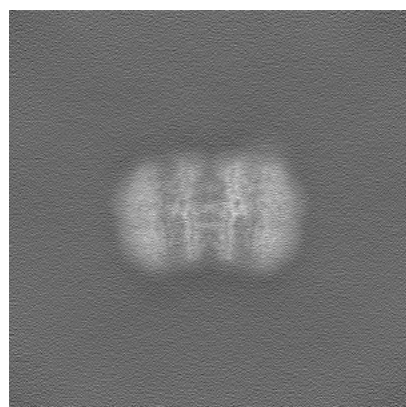


Y

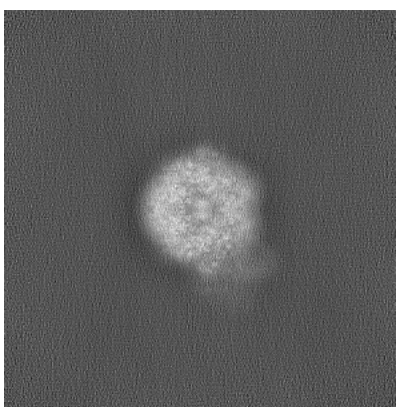


Z

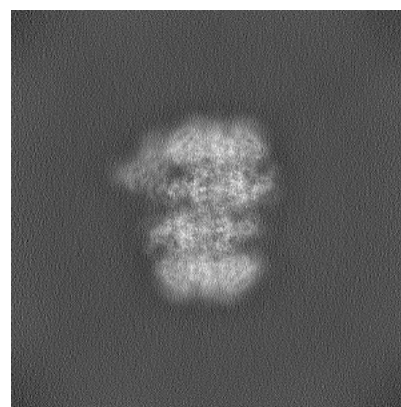
6.1.2 Raw map



X



Y

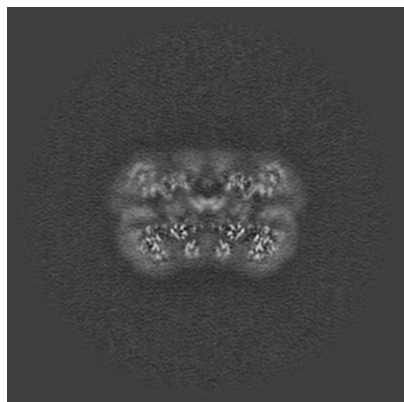


Z

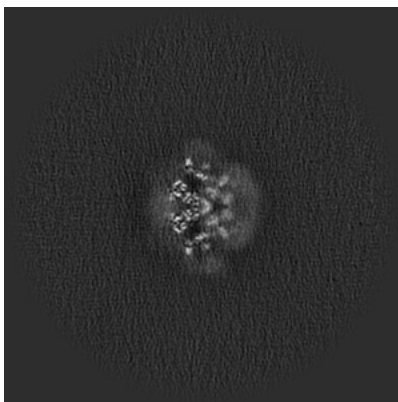
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

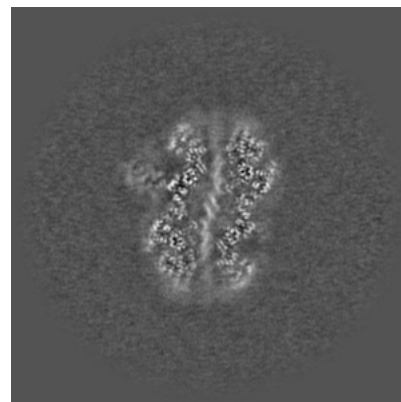
6.2.1 Primary map



X Index: 224

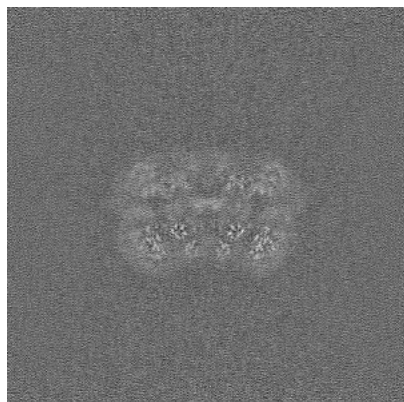


Y Index: 224

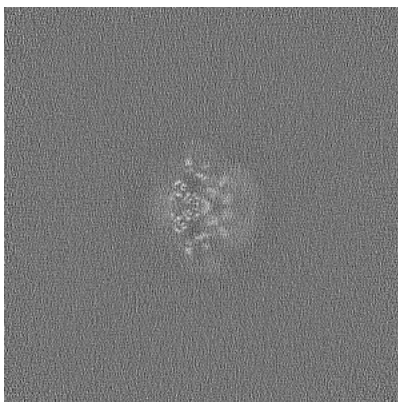


Z Index: 224

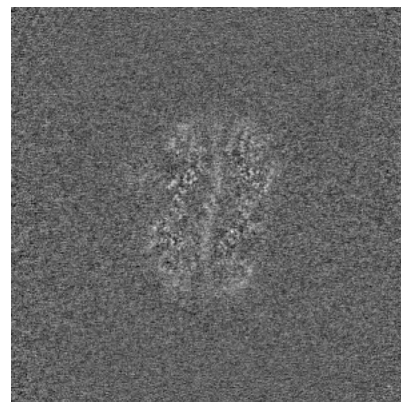
6.2.2 Raw map



X Index: 224



Y Index: 224

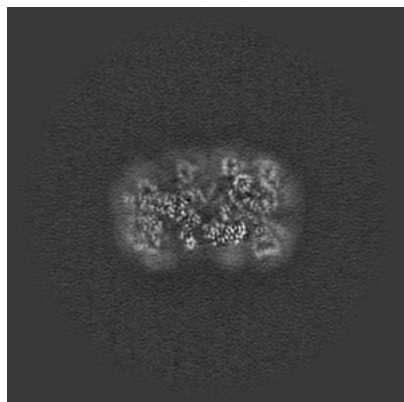


Z Index: 224

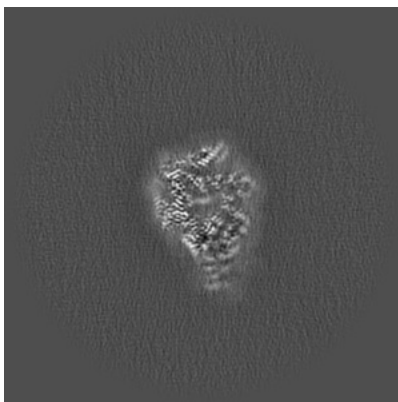
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

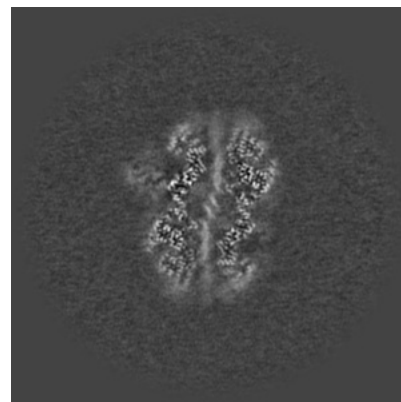
6.3.1 Primary map



X Index: 246

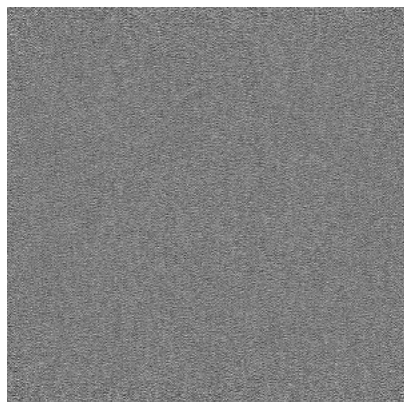


Y Index: 248

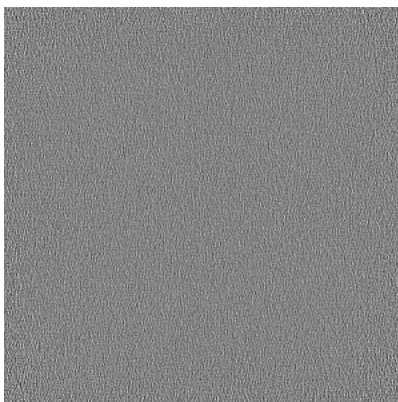


Z Index: 223

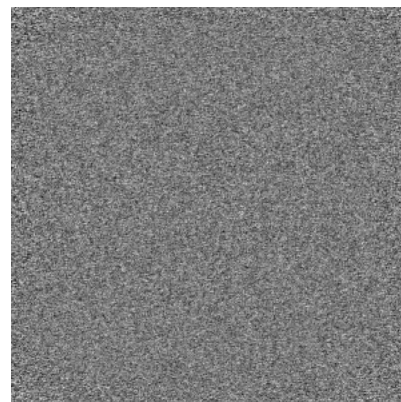
6.3.2 Raw map



X Index: 0



Y Index: 0

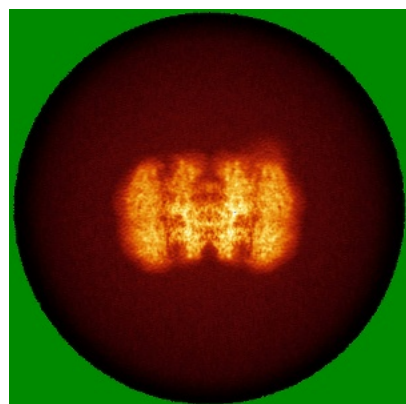


Z Index: 0

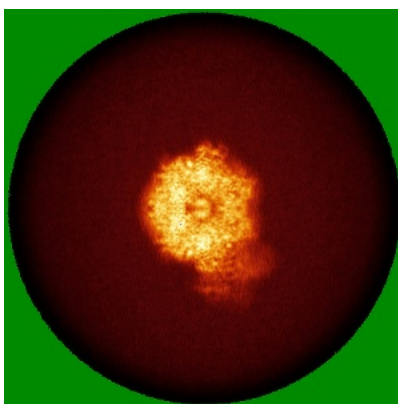
The images above show the largest variance slices of the map in three orthogonal directions.

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

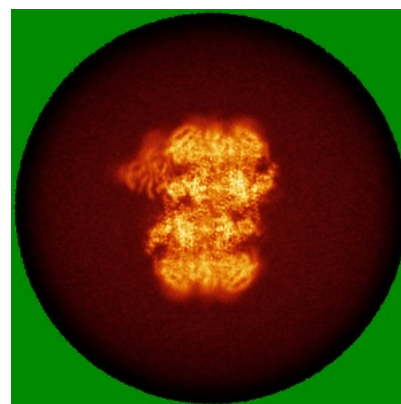
6.4.1 Primary map



X

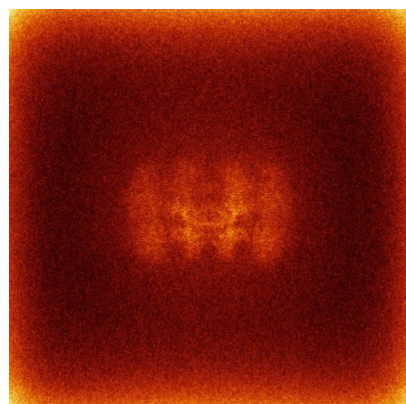


Y

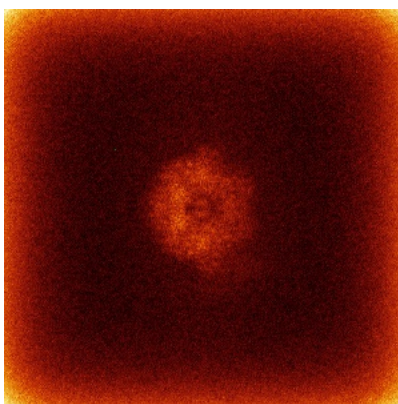


Z

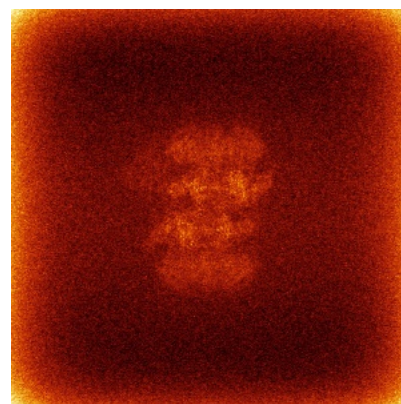
6.4.2 Raw map



X



Y

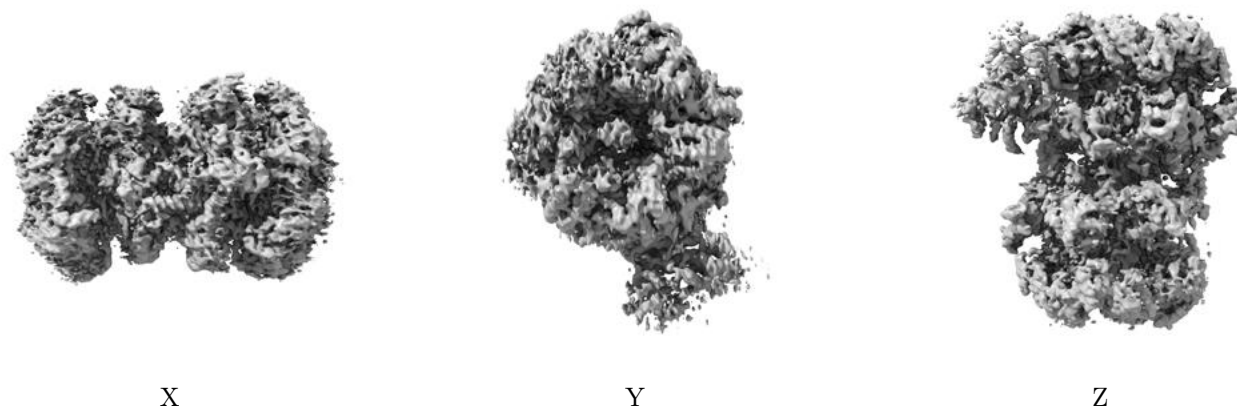


Z

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.

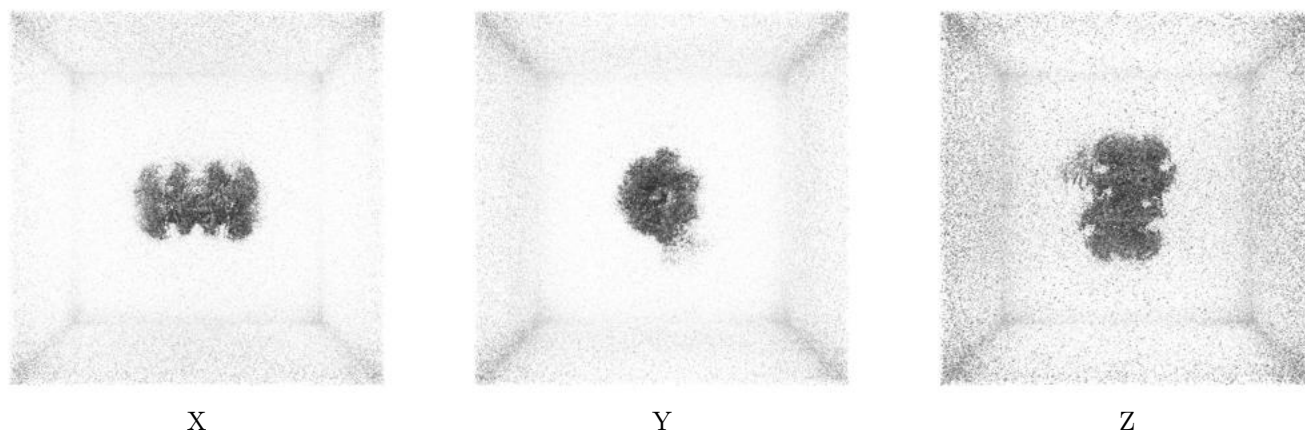
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.15. 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



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.

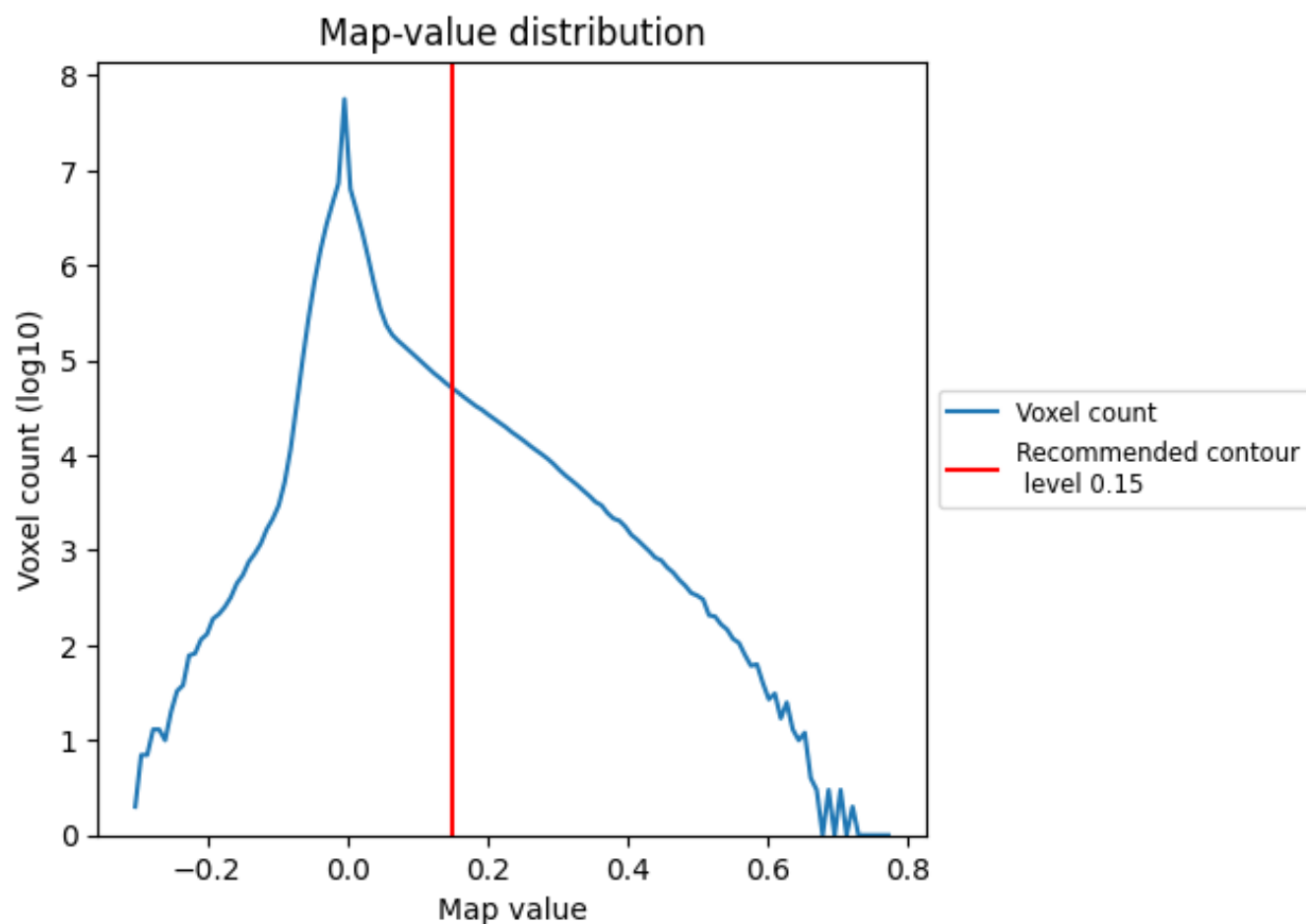
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

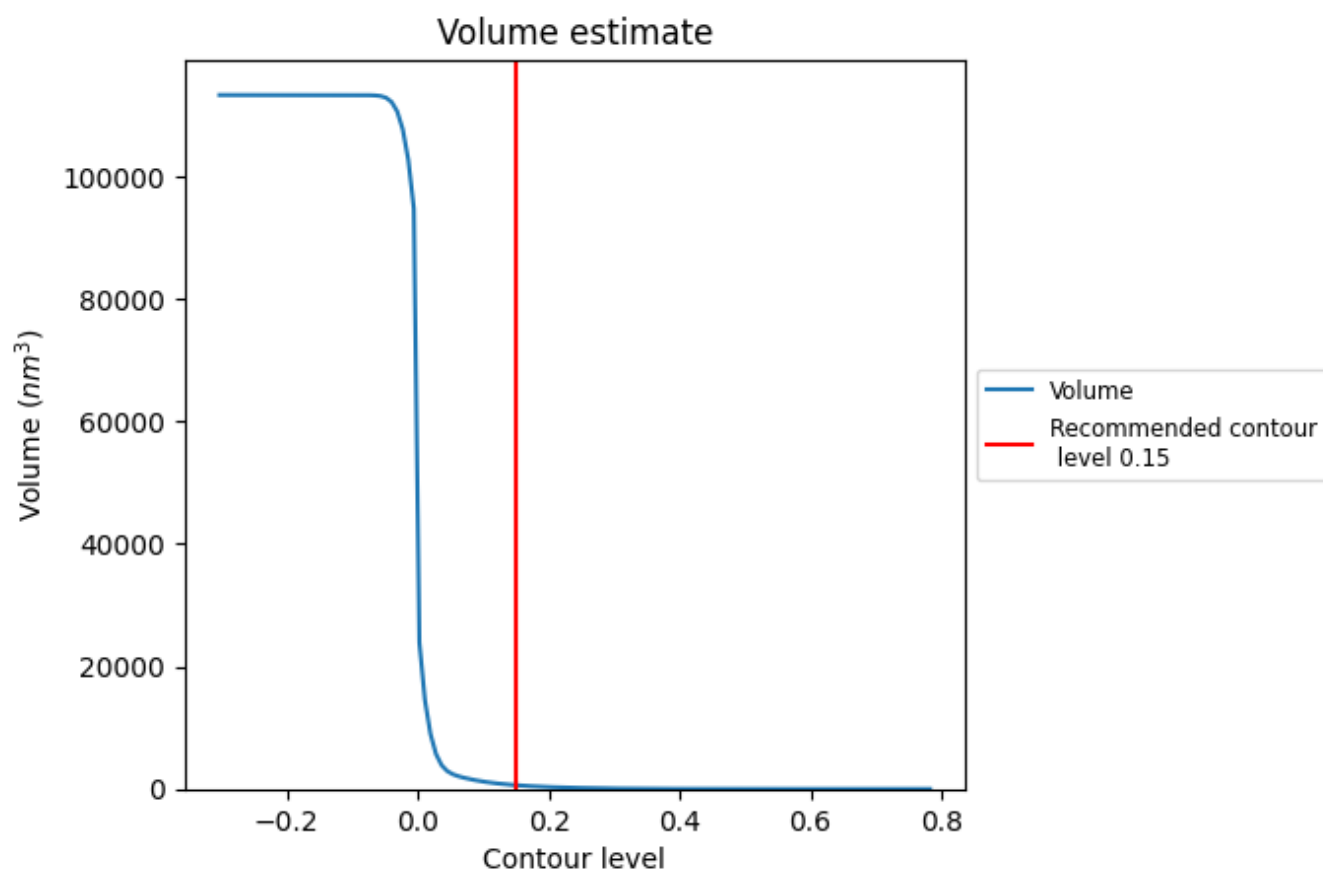
This section contains the results of statistical analysis of the map.

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.

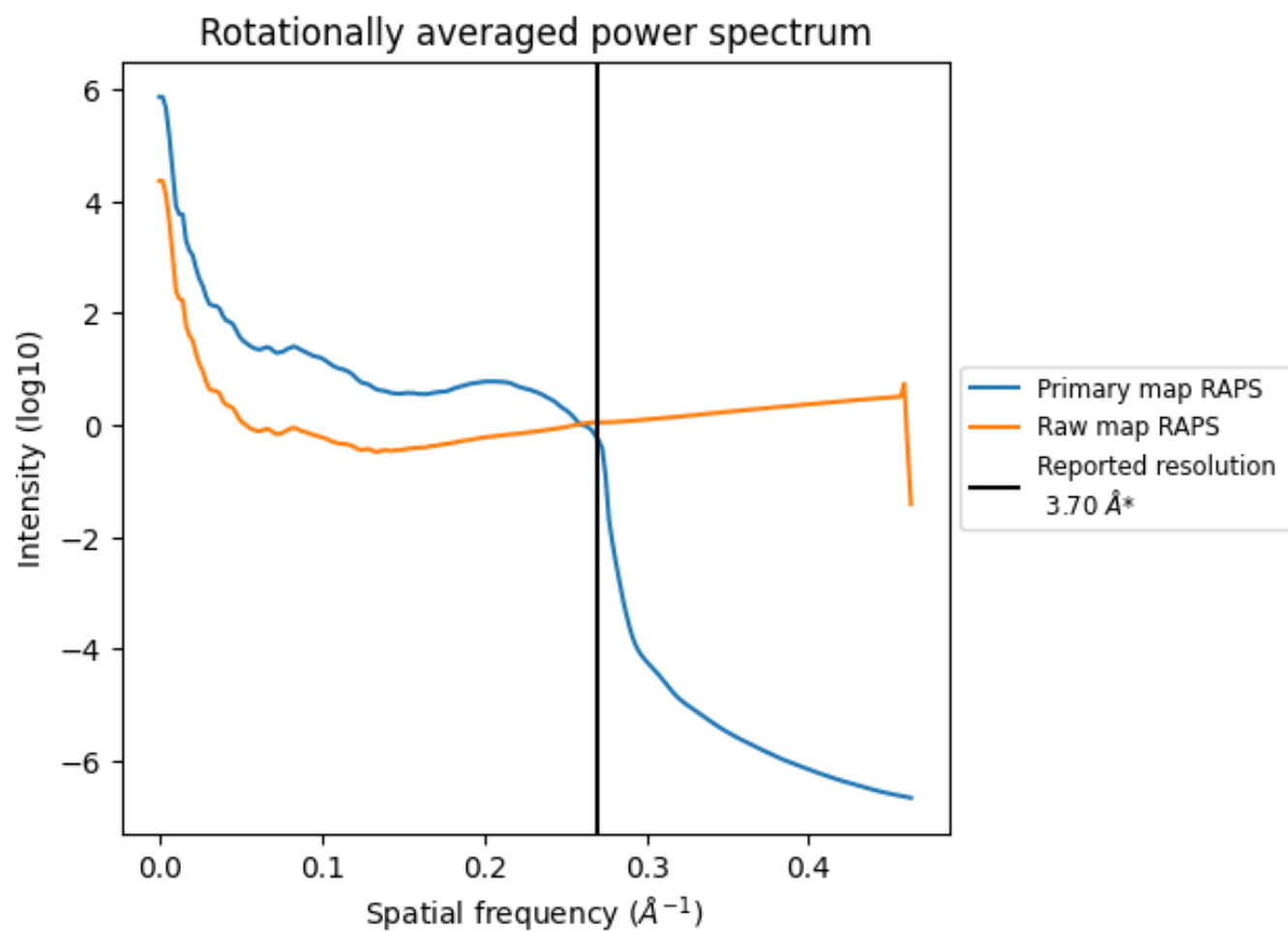
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 617 nm^3 ; this corresponds to an approximate mass of 557 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 ⓘ

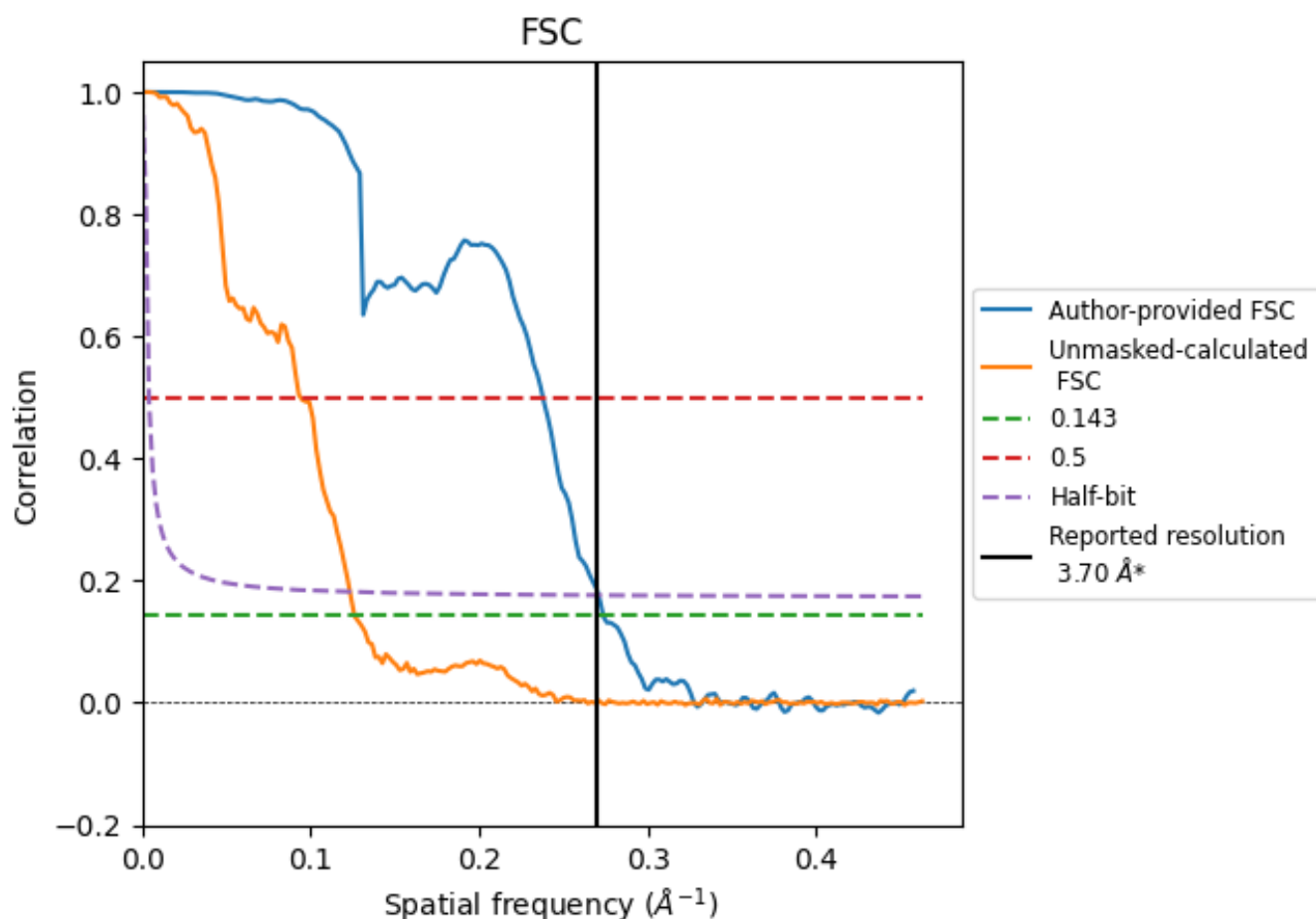


*Reported resolution corresponds to spatial frequency of 0.270 $\text{\AA}^{-1}$

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 0.270 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

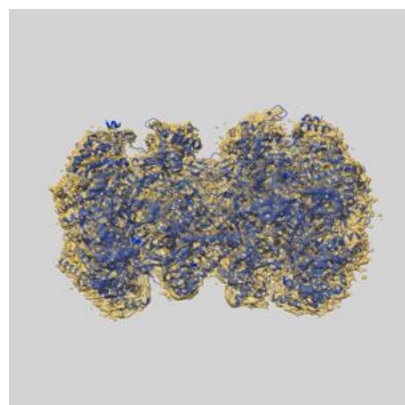
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.70	-	-
Author-provided FSC curve	3.65	4.20	3.69
Unmasked-calculated*	7.94	10.64	8.11

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 7.94 differs from the reported value 3.7 by more than 10 %

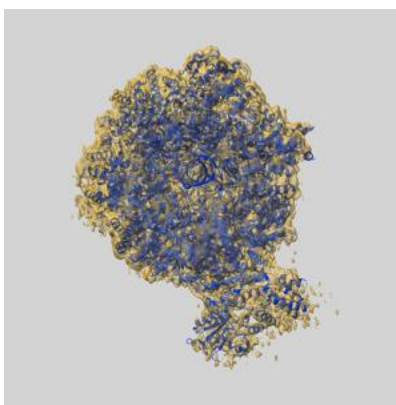
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-53973 and PDB model 9RHM. Per-residue inclusion information can be found in section [3](#) on page [12](#).

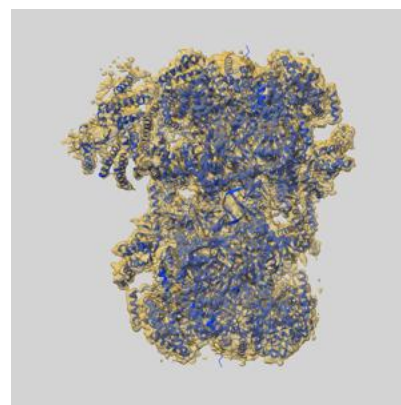
9.1 Map-model overlay [i](#)



X



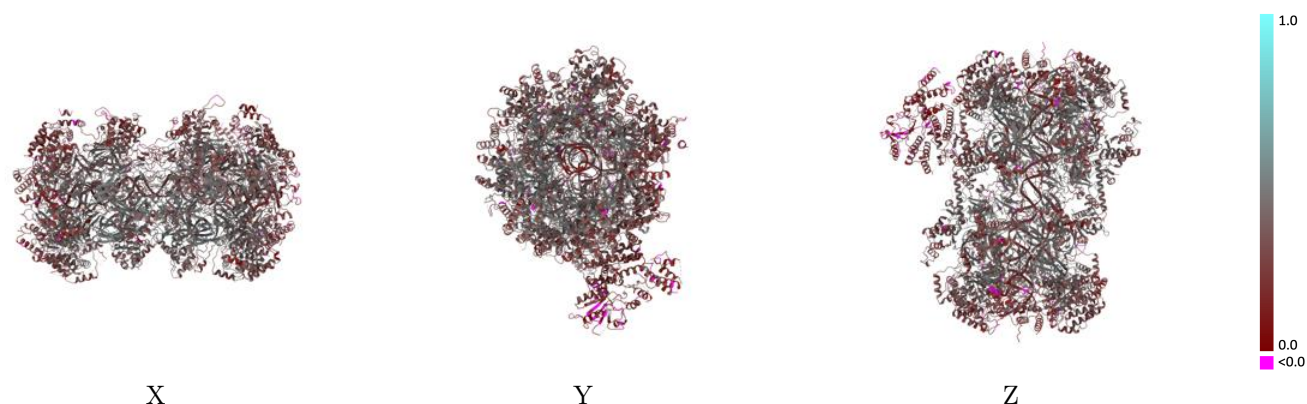
Y



Z

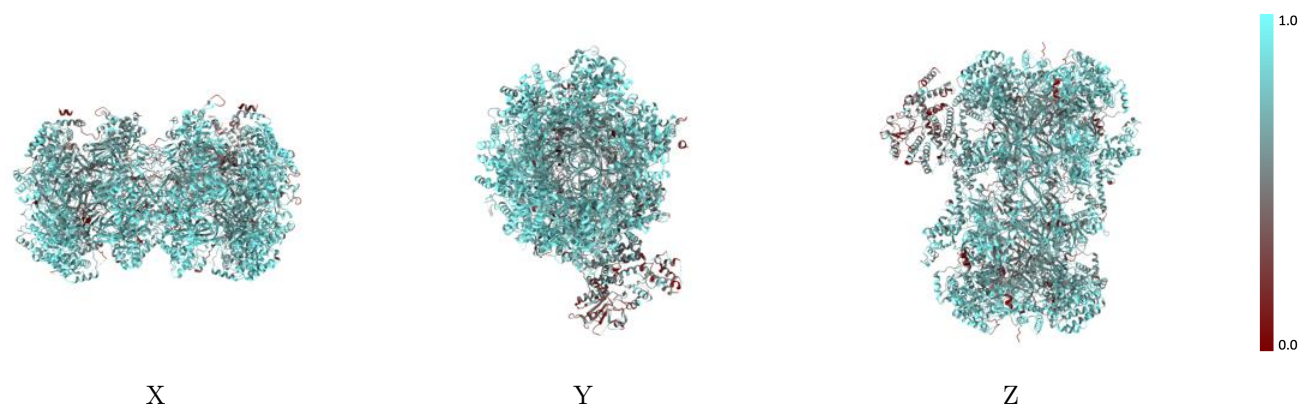
The images above show the 3D surface view of the map at the recommended contour level 0.15 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



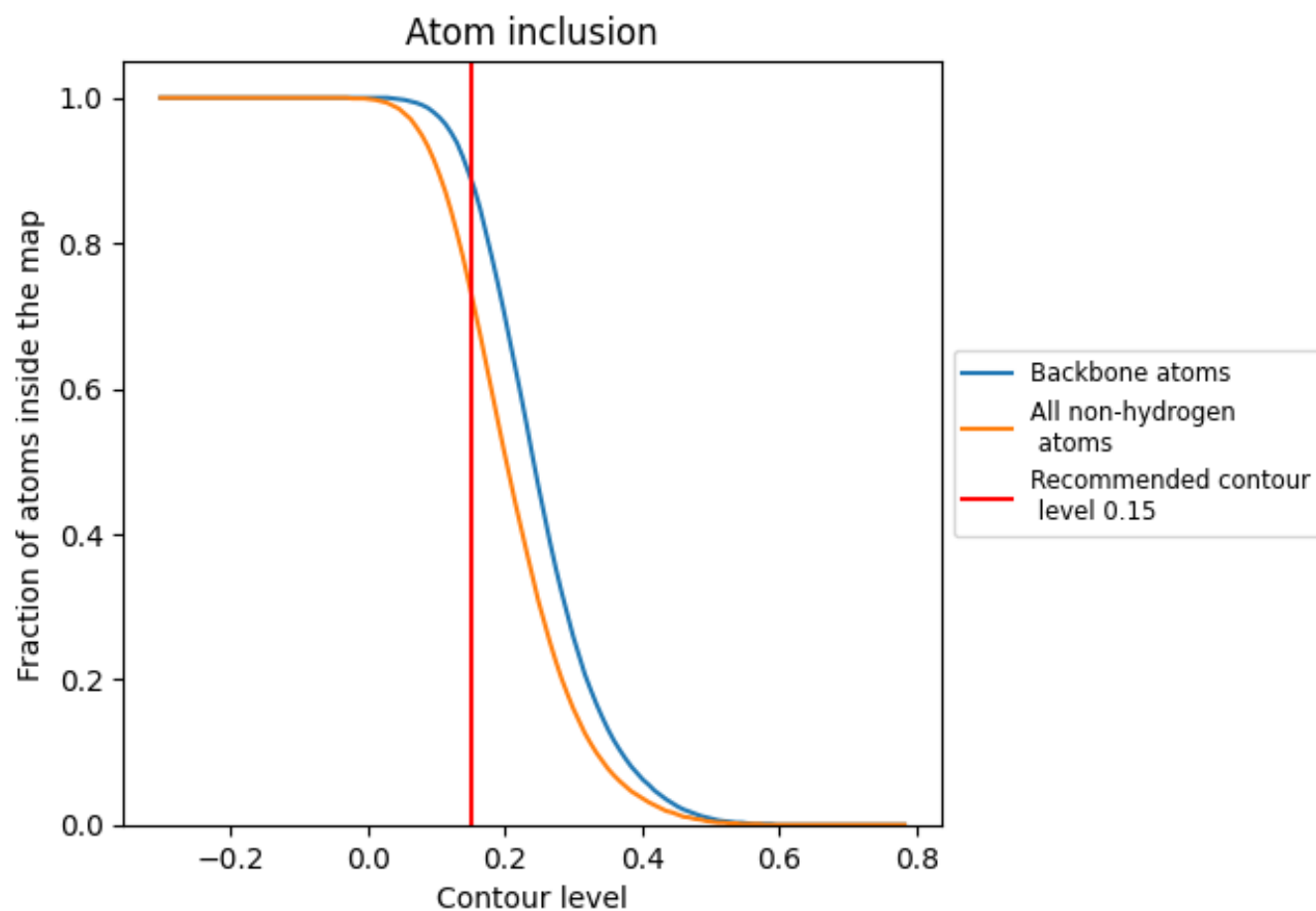
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.15).

9.4 Atom inclusion [i](#)



At the recommended contour level, 89% of all backbone atoms, 73% 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 (0.15) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7330	 0.3510
2	 0.7280	 0.3570
3	 0.8250	 0.4160
4	 0.7550	 0.3690
5	 0.7790	 0.3900
6	 0.7100	 0.3300
7	 0.7970	 0.3860
A	 0.1400	 0.2120
E	 0.4760	 0.2060
H	 0.5980	 0.3550
I	 0.4350	 0.1890
R	 0.1890	 0.1060
X	 0.7040	 0.2320
Y	 0.6960	 0.2220
a	 0.6800	 0.3310
b	 0.8110	 0.4010
c	 0.7280	 0.3520
d	 0.7640	 0.3780
e	 0.6780	 0.3020
f	 0.7750	 0.3730

