



Full wwPDB EM Validation Report ⓘ

Mar 27, 2026 – 03:39 PM UTC

PDB ID : 6QYJ / pdb_00006qyj
EMDB ID : EMD-4678
Title : The cryo-EM structure of the connector of the mature bacteriophage phi29
Authors : Xu, J.; Wang, D.; Gui, M.; Xiang, Y.
Deposited on : 2019-03-09
Resolution : 3.40 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:

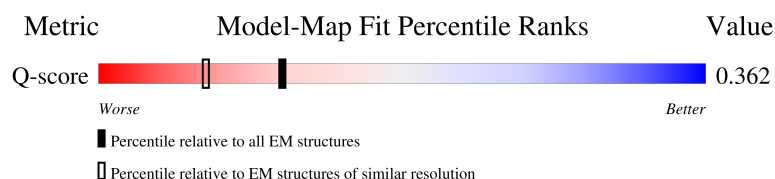
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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

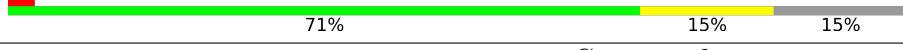
The reported resolution of this entry is 3.40 Å.

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.



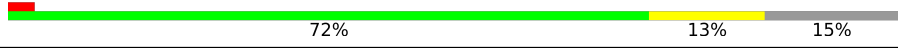
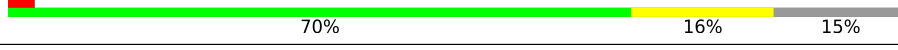
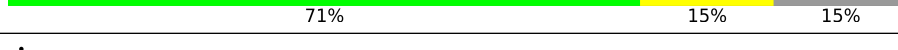
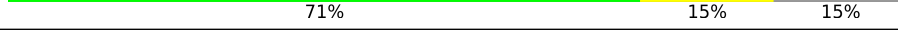
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14717 (2.90 - 3.90)

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	0a	309	
1	0b	309	
1	0c	309	
1	0d	309	

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Mol	Chain	Length	Quality of chain
1	0e	309	 71%14%15%
1	0f	309	 71%14%15%
1	0g	309	 72%13%15%
1	0h	309	 70%16%15%
1	0i	309	 70%16%15%
1	0j	309	 70%15%15%
1	0k	309	 71%15%15%
1	0l	309	 71%15%15%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 25956 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.

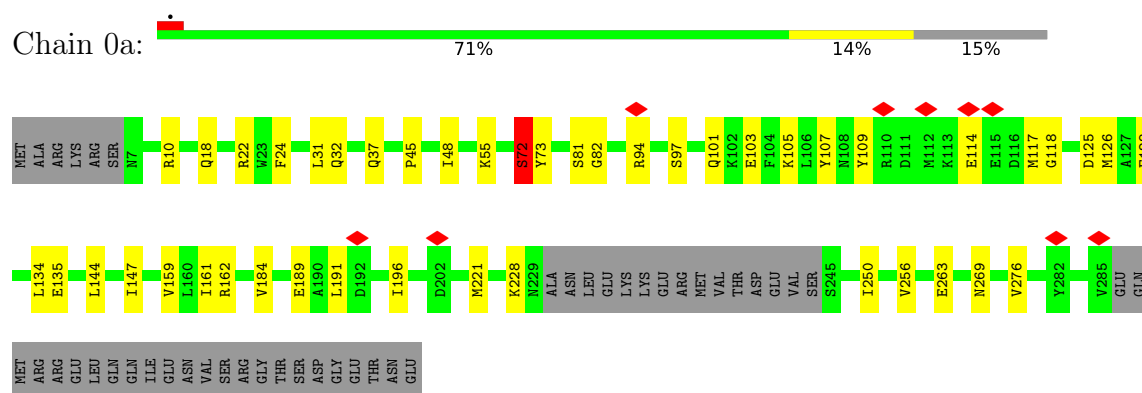
- Molecule 1 is a protein called Portal protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	0a	264	Total 2163	C 1384	N 360	O 412	S 7	0	0
1	0b	264	Total 2163	C 1384	N 360	O 412	S 7	0	0
1	0c	264	Total 2163	C 1384	N 360	O 412	S 7	0	0
1	0d	264	Total 2163	C 1384	N 360	O 412	S 7	0	0
1	0e	264	Total 2163	C 1384	N 360	O 412	S 7	0	0
1	0f	264	Total 2163	C 1384	N 360	O 412	S 7	0	0
1	0g	264	Total 2163	C 1384	N 360	O 412	S 7	0	0
1	0h	264	Total 2163	C 1384	N 360	O 412	S 7	0	0
1	0i	264	Total 2163	C 1384	N 360	O 412	S 7	0	0
1	0j	264	Total 2163	C 1384	N 360	O 412	S 7	0	0
1	0k	264	Total 2163	C 1384	N 360	O 412	S 7	0	0
1	0l	264	Total 2163	C 1384	N 360	O 412	S 7	0	0

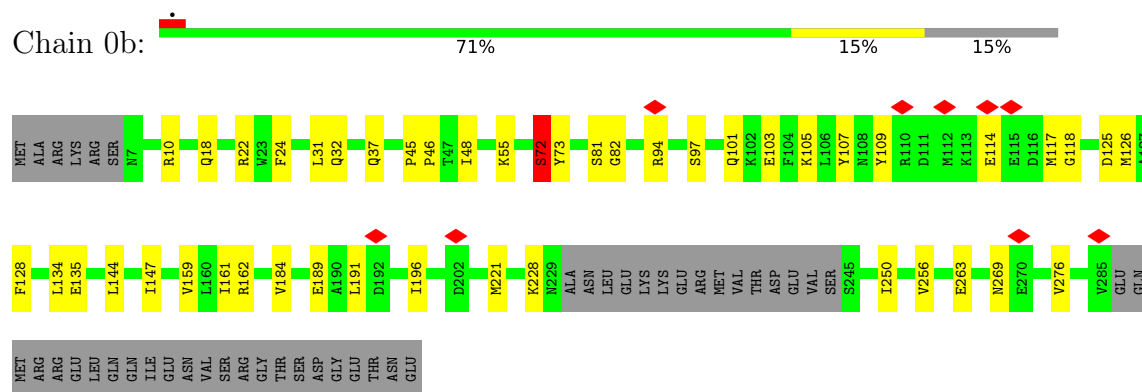
3 Residue-property plots

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.

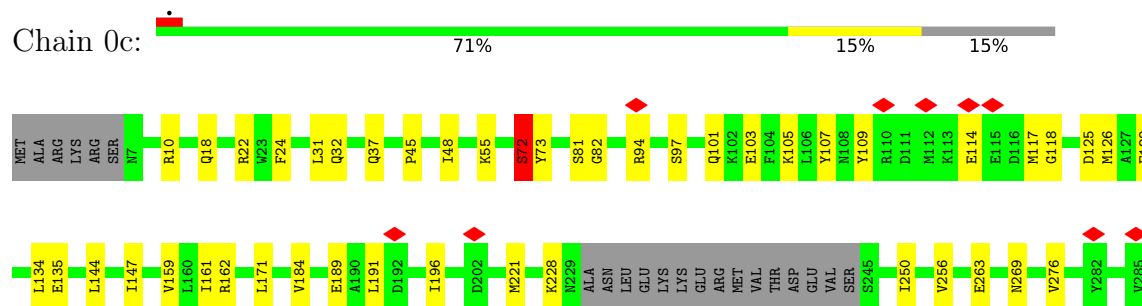
- Molecule 1: Portal protein



- Molecule 1: Portal protein



- Molecule 1: Portal protein



GLU
GLN
MET
ARG
ARG
ARG
GLU
LEU
GLN
GLN
ILE
GLU
ASN
VAL
SER
SER
GLY
THR
SER
ASP
GLY
GLU
THR
ASN
GLU

• Molecule 1: Portal protein

Chain 0d: 

MET
ALA
ARG
LYS
ARG
SER
N7
R10
Q18
R22
W23
F24
L31
Q32
Q37
P45
P46
T47
I48
K55
S72
Y73
R94
S97
Q101
K102
E103
F104
K105
L106
Y107
H108
Y109
R110
D111
M112
K113
E114
E115
D116
M117
G118
D125
M126
A127
F128
L134

E135
L144
I147
V159
L160
R162
I161
L171
V184
E189
A190
L191
D192
I196
D202
M221
K228
W229
ALA
ASN
LEU
GLU
LYS
LYS
GLU
ARG
MET
VAL
THR
ASP
GLU
VAL
SER
S245
E248
Q249
I250
V256
E263
N269
V276
Y282
V285

GLU
GLN
MET
ARG
ARG
GLU
LEU
GLN
GLN
ILE
GLU
ASN
VAL
SER
ARG
THR
SER
ASP
GLY
GLU
THR
ASN
GLU

• Molecule 1: Portal protein

Chain 0e: 

MET
ALA
ARG
LYS
ARG
SER
N7
R10
S11
I12
Q18
F24
Q32
Q37
P45
I48
K55
S72
Y73
S81
G82
D85
V86
R94
S97
Q101
K102
E103
F104
K105
L106
Y107
H108
Y109
R110
D111
M112
K113
E114
E115
D116
M117
G119
D125

M126
A127
F128
L134
E135
L144
I147
V159
L160
I161
R162
A163
M164
V184
L191
D192
I196
D202
M221
K228
W229
ALA
ASN
LEU
GLU
LYS
LYS
GLU
ARG
MET
VAL
THR
ASP
GLU
VAL
SER
S245
A246
D247
I250
V256
E263
N269
V276

Y282
Y285
GLU
GLN
MET
ARG
ARG
GLU
LEU
GLN
GLN
ILE
GLU
ASN
VAL
SER
SER
ASP
GLY
GLU
THR
ASN
GLU

• Molecule 1: Portal protein

Chain 0f: 

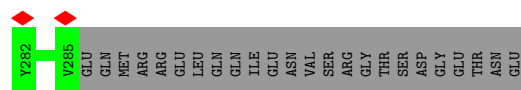
MET
ALA
ARG
LYS
ARG
SER
N7
R10
Q18
R22
W23
F24
L31
Q32
Q37
P45
I48
K55
S72
I74
S81
G82
R94
S97
Q101
K102
E103
F104
K105
Y109
R110
D111
M112
K113
E114
E115
D116
M117
D125
M126
A127
F128
L134

E135
L144
I147
V159
L160
R162
L171
V184
H188
E189
A190
L191
D192
I196
D202
M221
K228
W229
ALA
ASN
LEU
GLU
LYS
LYS
GLU
ARG
MET
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SER
S245
I250
V256
E263
N269
V276
Y282
V285
GLU

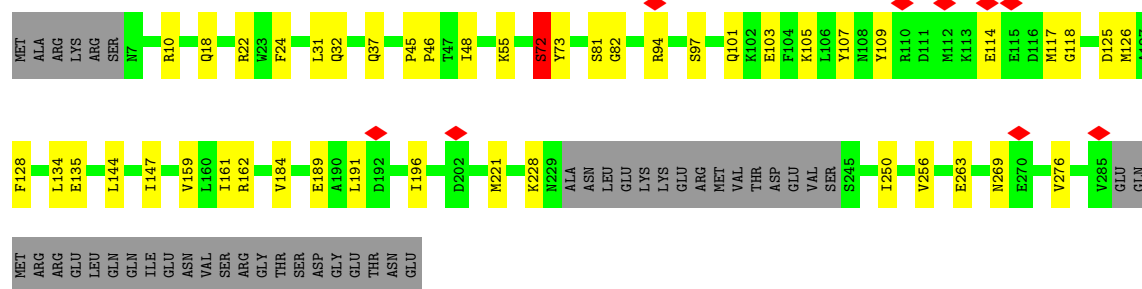
GLN
MET
ARG
ARG
GLU
LEU
GLN
GLN
ILE
GLU
ASN
VAL
SER
ARG
GLY
THR
SER
ASP
GLY
GLU
THR
ASN
GLU

• Molecule 1: Portal protein

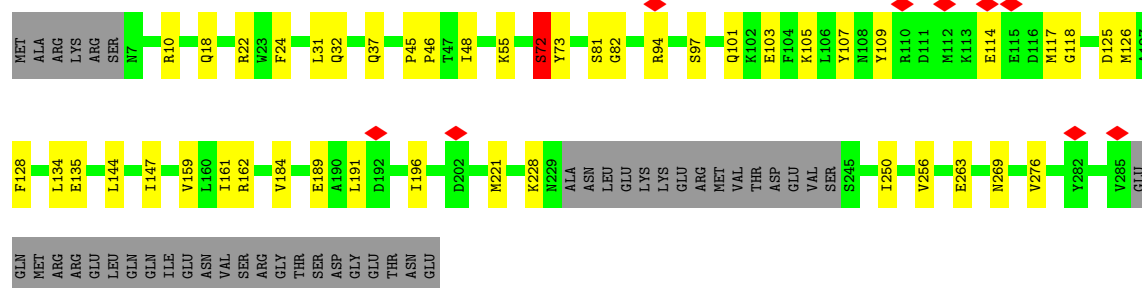
Chain 0g: 



• Molecule 1: Portal protein



• Molecule 1: Portal protein



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	36730	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	42.791	Depositor
Minimum map value	-0.417	Depositor
Average map value	0.005	Depositor
Map value standard deviation	0.271	Depositor
Recommended contour level	3.33	Depositor
Map size ($\text{\AA}$)	743.6, 743.6, 743.6	wwPDB
Map dimensions	572, 572, 572	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.3, 1.3, 1.3	Depositor

5 Model quality

5.1 Standard geometry

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	0a	0.22	0/2211	0.57	1/2996 (0.0%)
1	0b	0.23	0/2211	0.57	1/2996 (0.0%)
1	0c	0.23	0/2211	0.57	1/2996 (0.0%)
1	0d	0.23	0/2211	0.57	1/2996 (0.0%)
1	0e	0.23	0/2211	0.57	1/2996 (0.0%)
1	0f	0.23	0/2211	0.57	1/2996 (0.0%)
1	0g	0.23	0/2211	0.57	1/2996 (0.0%)
1	0h	0.23	0/2211	0.57	1/2996 (0.0%)
1	0i	0.23	0/2211	0.57	1/2996 (0.0%)
1	0j	0.23	0/2211	0.57	1/2996 (0.0%)
1	0k	0.23	0/2211	0.57	1/2996 (0.0%)
1	0l	0.23	0/2211	0.57	1/2996 (0.0%)
All	All	0.23	0/26532	0.57	12/35952 (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	0a	0	1
1	0b	0	1
1	0c	0	1
1	0d	0	1
1	0e	0	1
1	0f	0	1
1	0g	0	1
1	0h	0	1
1	0i	0	1
1	0j	0	1
1	0k	0	1
1	0l	0	1
All	All	0	12

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	0e	72	SER	N-CA-C	5.89	118.61	109.60
1	0d	72	SER	N-CA-C	5.89	118.61	109.60
1	0b	72	SER	N-CA-C	5.88	118.60	109.60
1	0k	72	SER	N-CA-C	5.88	118.60	109.60
1	0c	72	SER	N-CA-C	5.88	118.59	109.60
1	0f	72	SER	N-CA-C	5.88	118.59	109.60
1	0i	72	SER	N-CA-C	5.88	118.59	109.60
1	0a	72	SER	N-CA-C	5.87	118.59	109.60
1	0g	72	SER	N-CA-C	5.87	118.59	109.60
1	0j	72	SER	N-CA-C	5.87	118.59	109.60
1	0l	72	SER	N-CA-C	5.86	118.56	109.60
1	0h	72	SER	N-CA-C	5.85	118.54	109.60

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	0a	72	SER	Peptide
1	0b	72	SER	Peptide
1	0c	72	SER	Peptide
1	0d	72	SER	Peptide
1	0e	72	SER	Peptide
1	0f	72	SER	Peptide
1	0g	72	SER	Peptide
1	0h	72	SER	Peptide
1	0i	72	SER	Peptide
1	0j	72	SER	Peptide
1	0k	72	SER	Peptide
1	0l	72	SER	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	0a	2163	0	2104	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	0b	2163	0	2104	32	0
1	0c	2163	0	2104	31	0
1	0d	2163	0	2104	30	0
1	0e	2163	0	2104	32	0
1	0f	2163	0	2104	32	0
1	0g	2163	0	2104	29	0
1	0h	2163	0	2104	33	0
1	0i	2163	0	2104	33	0
1	0j	2163	0	2104	31	0
1	0k	2163	0	2104	31	0
1	0l	2163	0	2104	31	0
All	All	25956	0	25248	300	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 6.

All (300) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0e:162:ARG:NH1	1:0f:191:LEU:O	2.26	0.68
1:0d:256:VAL:HA	1:0e:37:GLN:HG3	1.77	0.65
1:0a:256:VAL:HA	1:0b:37:GLN:HG3	1.79	0.64
1:0k:256:VAL:HA	1:0l:37:GLN:HG3	1.79	0.64
1:0j:256:VAL:HA	1:0k:37:GLN:HG3	1.79	0.64
1:0a:37:GLN:HG3	1:0l:256:VAL:HA	1.79	0.64
1:0b:256:VAL:HA	1:0c:37:GLN:HG3	1.79	0.64
1:0g:256:VAL:HA	1:0h:37:GLN:HG3	1.79	0.63
1:0i:256:VAL:HA	1:0j:37:GLN:HG3	1.79	0.63
1:0f:256:VAL:HA	1:0g:37:GLN:HG3	1.79	0.63
1:0c:256:VAL:HA	1:0d:37:GLN:HG3	1.79	0.63
1:0h:256:VAL:HA	1:0i:37:GLN:HG3	1.79	0.63
1:0h:72:SER:OG	1:0h:73:TYR:N	2.33	0.62
1:0g:72:SER:OG	1:0g:73:TYR:N	2.33	0.61
1:0i:72:SER:OG	1:0i:73:TYR:N	2.33	0.61
1:0j:72:SER:OG	1:0j:73:TYR:N	2.33	0.61
1:0k:72:SER:OG	1:0k:73:TYR:N	2.33	0.60
1:0j:162:ARG:NH1	1:0k:191:LEU:O	2.36	0.59
1:0i:162:ARG:NH1	1:0j:191:LEU:O	2.36	0.59
1:0l:72:SER:OG	1:0l:73:TYR:N	2.33	0.59
1:0c:162:ARG:NH1	1:0d:191:LEU:O	2.36	0.59
1:0g:162:ARG:NH1	1:0h:191:LEU:O	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0k:162:ARG:NH1	1:0l:191:LEU:O	2.36	0.59
1:0f:162:ARG:NH1	1:0g:191:LEU:O	2.36	0.59
1:0a:72:SER:OG	1:0a:73:TYR:N	2.33	0.59
1:0b:72:SER:OG	1:0b:73:TYR:N	2.33	0.59
1:0b:162:ARG:NH1	1:0c:191:LEU:O	2.36	0.59
1:0a:191:LEU:O	1:0l:162:ARG:NH1	2.36	0.59
1:0a:162:ARG:NH1	1:0b:191:LEU:O	2.36	0.59
1:0h:162:ARG:NH1	1:0i:191:LEU:O	2.36	0.59
1:0e:256:VAL:HA	1:0f:37:GLN:HG3	1.85	0.57
1:0h:161:ILE:HG12	1:0h:196:ILE:HG12	1.87	0.57
1:0i:161:ILE:HG12	1:0i:196:ILE:HG12	1.87	0.57
1:0j:161:ILE:HG12	1:0j:196:ILE:HG12	1.87	0.57
1:0g:161:ILE:HG12	1:0g:196:ILE:HG12	1.87	0.57
1:0k:161:ILE:HG12	1:0k:196:ILE:HG12	1.87	0.57
1:0f:161:ILE:HG12	1:0f:196:ILE:HG12	1.87	0.56
1:0i:105:LYS:HB3	1:0i:117:MET:HE3	1.88	0.56
1:0l:105:LYS:HB3	1:0l:117:MET:HE3	1.88	0.56
1:0c:161:ILE:HG12	1:0c:196:ILE:HG12	1.87	0.56
1:0e:228:LYS:HA	1:0e:250:ILE:HD13	1.88	0.56
1:0j:105:LYS:HB3	1:0j:117:MET:HE3	1.88	0.56
1:0a:105:LYS:HB3	1:0a:117:MET:HE3	1.88	0.56
1:0b:161:ILE:HG12	1:0b:196:ILE:HG12	1.87	0.56
1:0d:228:LYS:HA	1:0d:250:ILE:HD13	1.88	0.56
1:0f:228:LYS:HA	1:0f:250:ILE:HD13	1.88	0.55
1:0d:161:ILE:HG12	1:0d:196:ILE:HG12	1.87	0.55
1:0l:161:ILE:HG12	1:0l:196:ILE:HG12	1.87	0.55
1:0a:161:ILE:HG12	1:0a:196:ILE:HG12	1.87	0.55
1:0c:228:LYS:HA	1:0c:250:ILE:HD13	1.88	0.55
1:0g:228:LYS:HA	1:0g:250:ILE:HD13	1.88	0.55
1:0h:228:LYS:HA	1:0h:250:ILE:HD13	1.88	0.55
1:0k:105:LYS:HB3	1:0k:117:MET:HE3	1.88	0.55
1:0k:228:LYS:HA	1:0k:250:ILE:HD13	1.88	0.55
1:0h:105:LYS:HB3	1:0h:117:MET:HE3	1.88	0.55
1:0l:228:LYS:HA	1:0l:250:ILE:HD13	1.88	0.55
1:0c:105:LYS:HB3	1:0c:117:MET:HE3	1.88	0.54
1:0e:161:ILE:HG12	1:0e:196:ILE:HG12	1.87	0.54
1:0f:105:LYS:HB3	1:0f:117:MET:HE3	1.88	0.54
1:0g:105:LYS:HB3	1:0g:117:MET:HE3	1.88	0.54
1:0i:228:LYS:HA	1:0i:250:ILE:HD13	1.88	0.54
1:0b:228:LYS:HA	1:0b:250:ILE:HD13	1.88	0.54
1:0f:97:SER:O	1:0f:101:GLN:NE2	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0d:105:LYS:HB3	1:0d:117:MET:HE3	1.88	0.54
1:0e:105:LYS:HB3	1:0e:117:MET:HE3	1.88	0.54
1:0j:228:LYS:HA	1:0j:250:ILE:HD13	1.88	0.54
1:0a:125:ASP:HB3	1:0b:55:LYS:HD2	1.89	0.54
1:0b:125:ASP:HB3	1:0c:55:LYS:HD2	1.89	0.54
1:0e:97:SER:O	1:0e:101:GLN:NE2	2.41	0.54
1:0g:97:SER:O	1:0g:101:GLN:NE2	2.41	0.54
1:0h:97:SER:O	1:0h:101:GLN:NE2	2.41	0.54
1:0k:125:ASP:HB3	1:0l:55:LYS:HD2	1.89	0.54
1:0a:55:LYS:HD2	1:0l:125:ASP:HB3	1.89	0.54
1:0a:228:LYS:HA	1:0a:250:ILE:HD13	1.88	0.54
1:0i:97:SER:O	1:0i:101:GLN:NE2	2.41	0.54
1:0a:97:SER:O	1:0a:101:GLN:NE2	2.41	0.54
1:0b:97:SER:O	1:0b:101:GLN:NE2	2.41	0.54
1:0c:97:SER:O	1:0c:101:GLN:NE2	2.41	0.54
1:0c:125:ASP:HB3	1:0d:55:LYS:HD2	1.89	0.54
1:0b:105:LYS:HB3	1:0b:117:MET:HE3	1.88	0.54
1:0d:97:SER:O	1:0d:101:GLN:NE2	2.41	0.54
1:0j:97:SER:O	1:0j:101:GLN:NE2	2.41	0.54
1:0j:125:ASP:HB3	1:0k:55:LYS:HD2	1.89	0.54
1:0l:97:SER:O	1:0l:101:GLN:NE2	2.41	0.54
1:0c:107:TYR:OH	1:0c:114:GLU:O	2.25	0.53
1:0f:125:ASP:HB3	1:0g:55:LYS:HD2	1.89	0.53
1:0k:97:SER:O	1:0k:101:GLN:NE2	2.41	0.53
1:0h:125:ASP:HB3	1:0i:55:LYS:HD2	1.89	0.53
1:0e:72:SER:OG	1:0e:73:TYR:N	2.33	0.53
1:0f:72:SER:OG	1:0f:73:TYR:N	2.33	0.53
1:0b:107:TYR:OH	1:0b:114:GLU:O	2.25	0.53
1:0i:125:ASP:HB3	1:0j:55:LYS:HD2	1.89	0.53
1:0d:72:SER:OG	1:0d:73:TYR:N	2.33	0.52
1:0g:125:ASP:HB3	1:0h:55:LYS:HD2	1.89	0.52
1:0d:125:ASP:HB3	1:0e:55:LYS:HD2	1.92	0.52
1:0d:162:ARG:NH1	1:0e:191:LEU:O	2.43	0.52
1:0e:135:GLU:OE1	1:0f:22:ARG:NH2	2.44	0.51
1:0a:107:TYR:OH	1:0a:114:GLU:O	2.25	0.51
1:0b:10:ARG:NH1	1:0b:18:GLN:OE1	2.44	0.51
1:0d:94:ARG:HA	1:0d:103:GLU:HA	1.93	0.51
1:0e:94:ARG:HA	1:0e:103:GLU:HA	1.93	0.51
1:0i:10:ARG:NH1	1:0i:18:GLN:OE1	2.44	0.51
1:0c:94:ARG:HA	1:0c:103:GLU:HA	1.93	0.51
1:0e:162:ARG:HB3	1:0f:189:GLU:HA	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0j:10:ARG:NH1	1:0j:18:GLN:OE1	2.44	0.51
1:0a:10:ARG:NH1	1:0a:18:GLN:OE1	2.44	0.51
1:0c:10:ARG:NH1	1:0c:18:GLN:OE1	2.44	0.51
1:0b:94:ARG:HA	1:0b:103:GLU:HA	1.93	0.51
1:0f:94:ARG:HA	1:0f:103:GLU:HA	1.93	0.51
1:0g:10:ARG:NH1	1:0g:18:GLN:OE1	2.44	0.51
1:0d:10:ARG:NH1	1:0d:18:GLN:OE1	2.44	0.50
1:0e:10:ARG:NH1	1:0e:18:GLN:OE1	2.44	0.50
1:0h:10:ARG:NH1	1:0h:18:GLN:OE1	2.44	0.50
1:0l:10:ARG:NH1	1:0l:18:GLN:OE1	2.44	0.50
1:0f:10:ARG:NH1	1:0f:18:GLN:OE1	2.44	0.50
1:0g:94:ARG:HA	1:0g:103:GLU:HA	1.93	0.50
1:0k:10:ARG:NH1	1:0k:18:GLN:OE1	2.44	0.50
1:0a:94:ARG:HA	1:0a:103:GLU:HA	1.93	0.50
1:0h:107:TYR:OH	1:0h:114:GLU:O	2.25	0.50
1:0h:94:ARG:HA	1:0h:103:GLU:HA	1.93	0.50
1:0e:125:ASP:HB3	1:0f:55:LYS:HD2	1.93	0.50
1:0l:94:ARG:HA	1:0l:103:GLU:HA	1.93	0.49
1:0i:94:ARG:HA	1:0i:103:GLU:HA	1.93	0.49
1:0j:94:ARG:HA	1:0j:103:GLU:HA	1.93	0.49
1:0k:94:ARG:HA	1:0k:103:GLU:HA	1.93	0.49
1:0l:107:TYR:OH	1:0l:114:GLU:O	2.25	0.49
1:0c:72:SER:OG	1:0c:73:TYR:N	2.33	0.48
1:0e:86:VAL:HG21	1:0f:74:ILE:HD11	1.95	0.48
1:0g:159:VAL:HB	1:0h:184:VAL:HG22	1.96	0.48
1:0h:159:VAL:HB	1:0i:184:VAL:HG22	1.96	0.48
1:0f:159:VAL:HB	1:0g:184:VAL:HG22	1.96	0.48
1:0i:159:VAL:HB	1:0j:184:VAL:HG22	1.96	0.48
1:0c:159:VAL:HB	1:0d:184:VAL:HG22	1.96	0.48
1:0b:159:VAL:HB	1:0c:184:VAL:HG22	1.96	0.48
1:0d:159:VAL:HB	1:0e:184:VAL:HG22	1.96	0.48
1:0a:184:VAL:HG22	1:0l:159:VAL:HB	1.96	0.47
1:0a:159:VAL:HB	1:0b:184:VAL:HG22	1.96	0.47
1:0j:159:VAL:HB	1:0k:184:VAL:HG22	1.96	0.47
1:0k:24:PHE:HZ	1:0k:135:GLU:HG2	1.80	0.47
1:0k:159:VAL:HB	1:0l:184:VAL:HG22	1.96	0.47
1:0b:24:PHE:HZ	1:0b:135:GLU:HG2	1.80	0.47
1:0c:24:PHE:HZ	1:0c:135:GLU:HG2	1.80	0.47
1:0b:72:SER:HG	1:0b:73:TYR:H	1.59	0.47
1:0d:24:PHE:HZ	1:0d:135:GLU:HG2	1.80	0.47
1:0h:24:PHE:HZ	1:0h:135:GLU:HG2	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0e:24:PHE:HZ	1:0e:135:GLU:HG2	1.80	0.47
1:0l:24:PHE:HZ	1:0l:135:GLU:HG2	1.80	0.47
1:0d:144:LEU:HD23	1:0d:147:ILE:HD12	1.97	0.47
1:0d:269:ASN:OD1	1:0d:276:VAL:N	2.48	0.47
1:0e:144:LEU:HD23	1:0e:147:ILE:HD12	1.97	0.47
1:0e:269:ASN:OD1	1:0e:276:VAL:N	2.48	0.47
1:0f:24:PHE:HZ	1:0f:135:GLU:HG2	1.80	0.47
1:0g:24:PHE:HZ	1:0g:135:GLU:HG2	1.80	0.47
1:0i:24:PHE:HZ	1:0i:135:GLU:HG2	1.80	0.47
1:0k:269:ASN:OD1	1:0k:276:VAL:N	2.48	0.47
1:0c:144:LEU:HD23	1:0c:147:ILE:HD12	1.97	0.46
1:0j:269:ASN:OD1	1:0j:276:VAL:N	2.48	0.46
1:0l:269:ASN:OD1	1:0l:276:VAL:N	2.48	0.46
1:0b:144:LEU:HD23	1:0b:147:ILE:HD12	1.97	0.46
1:0d:248:GLU:OE2	1:0e:228:LYS:NZ	2.40	0.46
1:0f:144:LEU:HD23	1:0f:147:ILE:HD12	1.97	0.46
1:0h:109:TYR:OH	1:0h:263:GLU:OE2	2.34	0.46
1:0i:109:TYR:OH	1:0i:263:GLU:OE2	2.34	0.46
1:0a:24:PHE:HZ	1:0a:135:GLU:HG2	1.80	0.46
1:0c:269:ASN:OD1	1:0c:276:VAL:N	2.48	0.46
1:0f:109:TYR:OH	1:0f:263:GLU:OE2	2.34	0.46
1:0j:24:PHE:HZ	1:0j:135:GLU:HG2	1.80	0.46
1:0b:31:LEU:HD22	1:0b:221:MET:HE2	1.98	0.46
1:0g:109:TYR:OH	1:0g:263:GLU:OE2	2.34	0.46
1:0i:269:ASN:OD1	1:0i:276:VAL:N	2.48	0.46
1:0j:109:TYR:OH	1:0j:263:GLU:OE2	2.34	0.46
1:0c:31:LEU:HD22	1:0c:221:MET:HE2	1.98	0.46
1:0e:109:TYR:OH	1:0e:263:GLU:OE2	2.34	0.46
1:0a:31:LEU:HD22	1:0a:221:MET:HE2	1.98	0.46
1:0k:31:LEU:HD22	1:0k:221:MET:HE2	1.98	0.46
1:0a:144:LEU:HD23	1:0a:147:ILE:HD12	1.97	0.46
1:0g:144:LEU:HD23	1:0g:147:ILE:HD12	1.98	0.46
1:0c:135:GLU:OE1	1:0d:22:ARG:NH2	2.49	0.46
1:0g:135:GLU:OE1	1:0h:22:ARG:NH2	2.49	0.46
1:0j:31:LEU:HD22	1:0j:221:MET:HE2	1.98	0.46
1:0k:135:GLU:OE1	1:0l:22:ARG:NH2	2.49	0.46
1:0l:31:LEU:HD22	1:0l:221:MET:HE2	1.98	0.46
1:0d:31:LEU:HD22	1:0d:221:MET:HE2	1.98	0.45
1:0d:109:TYR:OH	1:0d:263:GLU:OE2	2.34	0.45
1:0k:109:TYR:OH	1:0k:263:GLU:OE2	2.34	0.45
1:0i:135:GLU:OE1	1:0j:22:ARG:NH2	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0a:22:ARG:NH2	1:0l:135:GLU:OE1	2.49	0.45
1:0a:135:GLU:OE1	1:0b:22:ARG:NH2	2.49	0.45
1:0i:31:LEU:HD22	1:0i:221:MET:HE2	1.98	0.45
1:0j:135:GLU:OE1	1:0k:22:ARG:NH2	2.49	0.45
1:0b:269:ASN:OD1	1:0b:276:VAL:N	2.48	0.45
1:0d:107:TYR:OH	1:0d:114:GLU:O	2.25	0.45
1:0h:144:LEU:HD23	1:0h:147:ILE:HD12	1.97	0.45
1:0l:144:LEU:HD23	1:0l:147:ILE:HD12	1.97	0.45
1:0b:135:GLU:OE1	1:0c:22:ARG:NH2	2.49	0.45
1:0c:109:TYR:OH	1:0c:263:GLU:OE2	2.34	0.45
1:0h:31:LEU:HD22	1:0h:221:MET:HE2	1.98	0.45
1:0h:269:ASN:OD1	1:0h:276:VAL:N	2.48	0.45
1:0i:144:LEU:HD23	1:0i:147:ILE:HD12	1.97	0.45
1:0k:144:LEU:HD23	1:0k:147:ILE:HD12	1.97	0.45
1:0e:31:LEU:HD22	1:0e:221:MET:HE2	1.98	0.45
1:0f:45:PRO:HD2	1:0f:48:ILE:HD12	1.99	0.45
1:0h:135:GLU:OE1	1:0i:22:ARG:NH2	2.49	0.45
1:0l:109:TYR:OH	1:0l:263:GLU:OE2	2.34	0.45
1:0d:45:PRO:HD2	1:0d:48:ILE:HD12	1.99	0.45
1:0e:45:PRO:HD2	1:0e:48:ILE:HD12	1.99	0.45
1:0j:45:PRO:HD2	1:0j:48:ILE:HD12	1.99	0.45
1:0j:144:LEU:HD23	1:0j:147:ILE:HD12	1.97	0.45
1:0a:109:TYR:OH	1:0a:263:GLU:OE2	2.34	0.45
1:0b:109:TYR:OH	1:0b:263:GLU:OE2	2.34	0.45
1:0c:45:PRO:HD2	1:0c:48:ILE:HD12	1.99	0.45
1:0f:135:GLU:OE1	1:0g:22:ARG:NH2	2.49	0.45
1:0i:45:PRO:HD2	1:0i:48:ILE:HD12	1.99	0.45
1:0f:31:LEU:HD22	1:0f:221:MET:HE2	1.98	0.45
1:0f:269:ASN:OD1	1:0f:276:VAL:N	2.48	0.45
1:0b:45:PRO:HD2	1:0b:48:ILE:HD12	1.99	0.44
1:0g:31:LEU:HD22	1:0g:221:MET:HE2	1.98	0.44
1:0k:45:PRO:HD2	1:0k:48:ILE:HD12	1.99	0.44
1:0c:126:MET:HG3	1:0c:128:PHE:HB2	2.00	0.44
1:0b:126:MET:HG3	1:0b:128:PHE:HB2	2.00	0.44
1:0g:126:MET:HG3	1:0g:128:PHE:HB2	2.00	0.44
1:0j:45:PRO:HA	1:0j:46:PRO:HD3	1.88	0.44
1:0d:126:MET:HG3	1:0d:128:PHE:HB2	2.00	0.44
1:0h:126:MET:HG3	1:0h:128:PHE:HB2	2.00	0.44
1:0a:45:PRO:HD2	1:0a:48:ILE:HD12	1.99	0.44
1:0a:126:MET:HG3	1:0a:128:PHE:HB2	2.00	0.44
1:0f:126:MET:HG3	1:0f:128:PHE:HB2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0g:45:PRO:HD2	1:0g:48:ILE:HD12	1.99	0.44
1:0g:269:ASN:OD1	1:0g:276:VAL:N	2.48	0.44
1:0i:45:PRO:HA	1:0i:46:PRO:HD3	1.88	0.44
1:0j:126:MET:HG3	1:0j:128:PHE:HB2	2.00	0.44
1:0a:269:ASN:OD1	1:0a:276:VAL:N	2.48	0.44
1:0e:126:MET:HG3	1:0e:128:PHE:HB2	2.00	0.44
1:0i:126:MET:HG3	1:0i:128:PHE:HB2	2.00	0.44
1:0h:45:PRO:HD2	1:0h:48:ILE:HD12	1.99	0.44
1:0k:45:PRO:HA	1:0k:46:PRO:HD3	1.88	0.44
1:0k:126:MET:HG3	1:0k:128:PHE:HB2	2.00	0.44
1:0l:45:PRO:HD2	1:0l:48:ILE:HD12	1.99	0.44
1:0l:126:MET:HG3	1:0l:128:PHE:HB2	2.00	0.44
1:0k:107:TYR:OH	1:0k:114:GLU:O	2.25	0.44
1:0b:45:PRO:HA	1:0b:46:PRO:HD3	1.88	0.43
1:0a:189:GLU:HA	1:0l:162:ARG:HB3	2.01	0.43
1:0i:107:TYR:OH	1:0i:114:GLU:O	2.25	0.43
1:0e:159:VAL:HB	1:0f:184:VAL:HG22	2.01	0.43
1:0f:162:ARG:HB3	1:0g:189:GLU:HA	2.01	0.43
1:0a:162:ARG:HB3	1:0b:189:GLU:HA	2.01	0.43
1:0k:162:ARG:HB3	1:0l:189:GLU:HA	2.01	0.43
1:0l:45:PRO:HA	1:0l:46:PRO:HD3	1.88	0.43
1:0g:162:ARG:HB3	1:0h:189:GLU:HA	2.01	0.43
1:0b:162:ARG:HB3	1:0c:189:GLU:HA	2.01	0.42
1:0e:12:ILE:HD11	1:0g:183:PRO:HD2	2.01	0.42
1:0h:45:PRO:HA	1:0h:46:PRO:HD3	1.88	0.42
1:0b:107:TYR:N	1:0b:118:GLY:O	2.52	0.42
1:0h:162:ARG:HB3	1:0i:189:GLU:HA	2.01	0.42
1:0j:162:ARG:HB3	1:0k:189:GLU:HA	2.01	0.42
1:0c:162:ARG:HB3	1:0d:189:GLU:HA	2.01	0.42
1:0c:171:LEU:HD23	1:0c:171:LEU:HA	1.90	0.42
1:0c:107:TYR:N	1:0c:118:GLY:O	2.52	0.42
1:0d:45:PRO:HA	1:0d:46:PRO:HD3	1.88	0.42
1:0j:81:SER:OG	1:0j:82:GLY:N	2.53	0.42
1:0a:107:TYR:N	1:0a:118:GLY:O	2.52	0.42
1:0i:81:SER:OG	1:0i:82:GLY:N	2.53	0.42
1:0j:107:TYR:N	1:0j:118:GLY:O	2.52	0.42
1:0k:107:TYR:N	1:0k:118:GLY:O	2.52	0.42
1:0b:81:SER:OG	1:0b:82:GLY:N	2.53	0.42
1:0e:164:ASN:HA	1:0f:188:HIS:NE2	2.35	0.42
1:0k:81:SER:OG	1:0k:82:GLY:N	2.53	0.42
1:0a:32:GLN:HA	1:0a:134:LEU:HD13	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0d:107:TYR:N	1:0d:118:GLY:O	2.52	0.41
1:0h:107:TYR:N	1:0h:118:GLY:O	2.52	0.41
1:0c:81:SER:OG	1:0c:82:GLY:N	2.53	0.41
1:0i:32:GLN:HA	1:0i:134:LEU:HD13	2.03	0.41
1:0i:107:TYR:N	1:0i:118:GLY:O	2.52	0.41
1:0j:32:GLN:HA	1:0j:134:LEU:HD13	2.03	0.41
1:0l:32:GLN:HA	1:0l:134:LEU:HD13	2.03	0.41
1:0b:32:GLN:HA	1:0b:134:LEU:HD13	2.03	0.41
1:0h:32:GLN:HA	1:0h:134:LEU:HD13	2.02	0.41
1:0a:81:SER:OG	1:0a:82:GLY:N	2.53	0.41
1:0k:32:GLN:HA	1:0k:134:LEU:HD13	2.03	0.41
1:0l:107:TYR:N	1:0l:118:GLY:O	2.52	0.41
1:0f:81:SER:OG	1:0f:82:GLY:N	2.53	0.41
1:0h:81:SER:OG	1:0h:82:GLY:N	2.53	0.41
1:0h:171:LEU:HD23	1:0h:171:LEU:HA	1.90	0.41
1:0l:81:SER:OG	1:0l:82:GLY:N	2.53	0.41
1:0c:32:GLN:HA	1:0c:134:LEU:HD13	2.03	0.41
1:0i:162:ARG:HB3	1:0j:189:GLU:HA	2.01	0.41
1:0d:32:GLN:HA	1:0d:134:LEU:HD13	2.03	0.41
1:0d:171:LEU:HD23	1:0d:171:LEU:HA	1.90	0.41
1:0f:32:GLN:HA	1:0f:134:LEU:HD13	2.02	0.41
1:0g:32:GLN:HA	1:0g:134:LEU:HD13	2.03	0.41
1:0g:171:LEU:HD23	1:0g:171:LEU:HA	1.90	0.41
1:0i:171:LEU:HD23	1:0i:171:LEU:HA	1.90	0.41
1:0e:32:GLN:HA	1:0e:134:LEU:HD13	2.03	0.41
1:0e:81:SER:OG	1:0e:82:GLY:N	2.53	0.41
1:0e:107:TYR:N	1:0e:118:GLY:O	2.52	0.40
1:0e:125:ASP:OD2	1:0f:37:GLN:NE2	2.43	0.40
1:0f:171:LEU:HD23	1:0f:171:LEU:HA	1.90	0.40
1:0i:174:VAL:O	1:0i:178:TYR:N	2.51	0.40
1:0h:174:VAL:O	1:0h:178:TYR:N	2.51	0.40
1:0j:174:VAL:O	1:0j:178:TYR:N	2.51	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	0a	260/309 (84%)	243 (94%)	17 (6%)	0	100	100
1	0b	260/309 (84%)	243 (94%)	17 (6%)	0	100	100
1	0c	260/309 (84%)	243 (94%)	17 (6%)	0	100	100
1	0d	260/309 (84%)	243 (94%)	17 (6%)	0	100	100
1	0e	260/309 (84%)	243 (94%)	17 (6%)	0	100	100
1	0f	260/309 (84%)	243 (94%)	17 (6%)	0	100	100
1	0g	260/309 (84%)	243 (94%)	17 (6%)	0	100	100
1	0h	260/309 (84%)	243 (94%)	17 (6%)	0	100	100
1	0i	260/309 (84%)	243 (94%)	17 (6%)	0	100	100
1	0j	260/309 (84%)	243 (94%)	17 (6%)	0	100	100
1	0k	260/309 (84%)	243 (94%)	17 (6%)	0	100	100
1	0l	260/309 (84%)	243 (94%)	17 (6%)	0	100	100
All	All	3120/3708 (84%)	2916 (94%)	204 (6%)	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	0a	236/277 (85%)	236 (100%)	0	100	100
1	0b	236/277 (85%)	236 (100%)	0	100	100
1	0c	236/277 (85%)	236 (100%)	0	100	100
1	0d	236/277 (85%)	236 (100%)	0	100	100
1	0e	236/277 (85%)	236 (100%)	0	100	100
1	0f	236/277 (85%)	236 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	0g	236/277 (85%)	236 (100%)	0	100	100
1	0h	236/277 (85%)	236 (100%)	0	100	100
1	0i	236/277 (85%)	236 (100%)	0	100	100
1	0j	236/277 (85%)	236 (100%)	0	100	100
1	0k	236/277 (85%)	236 (100%)	0	100	100
1	0l	236/277 (85%)	236 (100%)	0	100	100
All	All	2832/3324 (85%)	2832 (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 (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	0a	16	GLN
1	0b	16	GLN
1	0c	16	GLN
1	0d	16	GLN
1	0e	16	GLN
1	0f	16	GLN
1	0g	16	GLN
1	0h	16	GLN
1	0h	21	ASN
1	0i	16	GLN
1	0j	16	GLN
1	0k	16	GLN
1	0l	16	GLN

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 [i](#)

There are no ligands in this entry.

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.

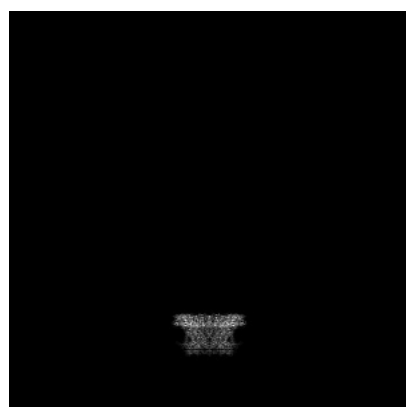
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-4678. These allow visual inspection of the internal detail of the map and identification of artifacts.

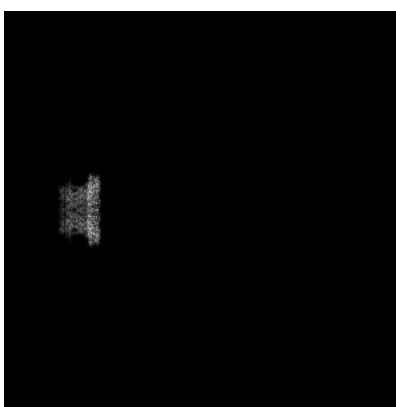
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

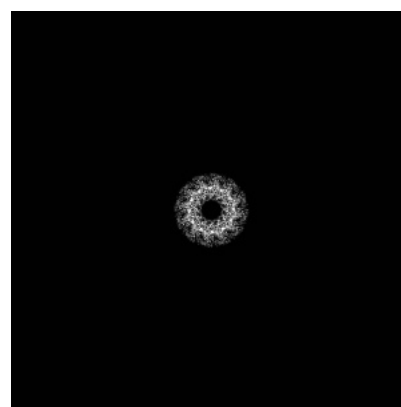
6.1.1 Primary map



X



Y

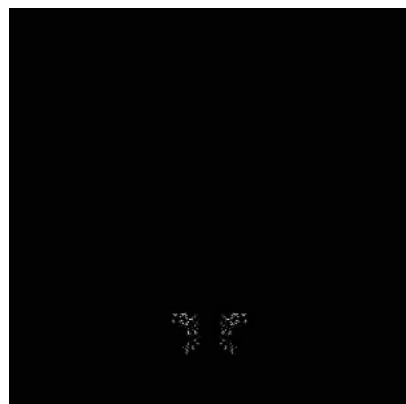


Z

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

6.2 Central slices [i](#)

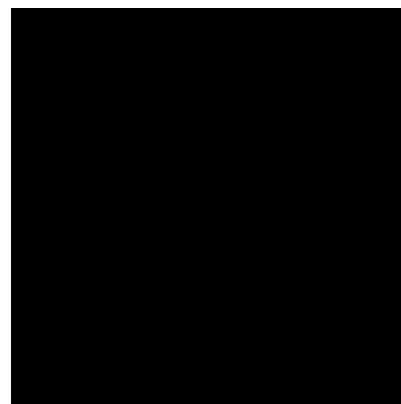
6.2.1 Primary map



X Index: 286



Y Index: 286



Z Index: 286

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

6.3 Largest variance slices [i](#)

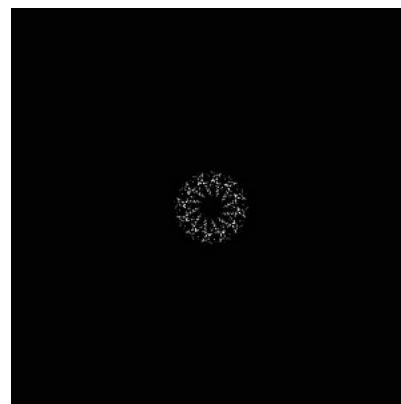
6.3.1 Primary map



X Index: 260



Y Index: 260

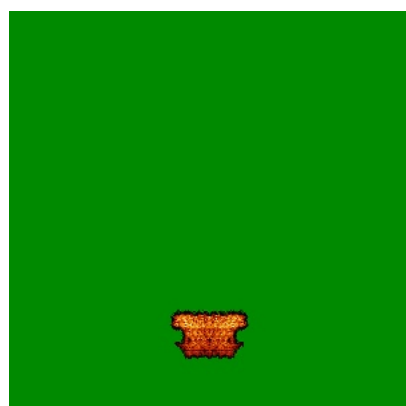


Z Index: 121

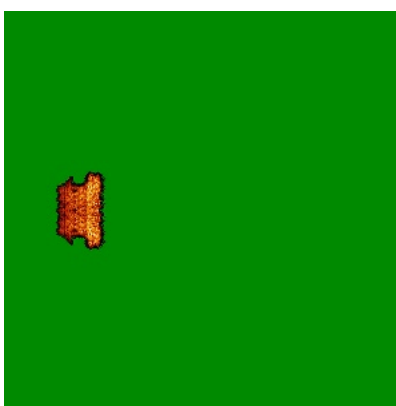
The images above show the largest variance slices of the map in three orthogonal directions.

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

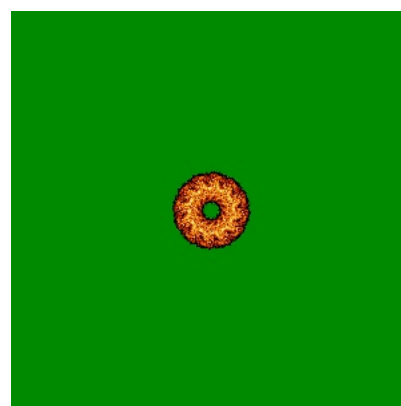
6.4.1 Primary map



X



Y



Z

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

This section was not generated.

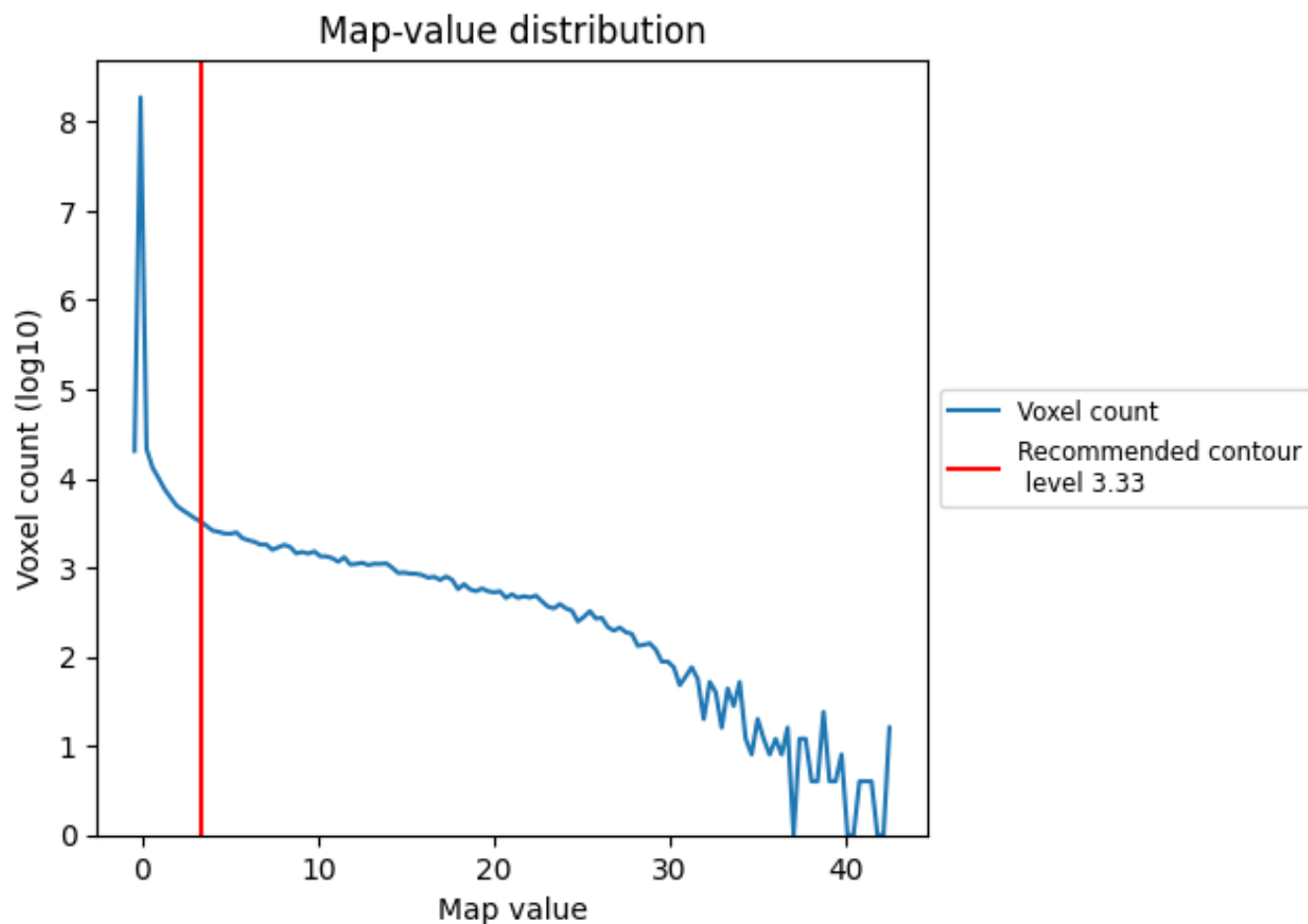
6.6 Mask visualisation

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

7 Map analysis [i](#)

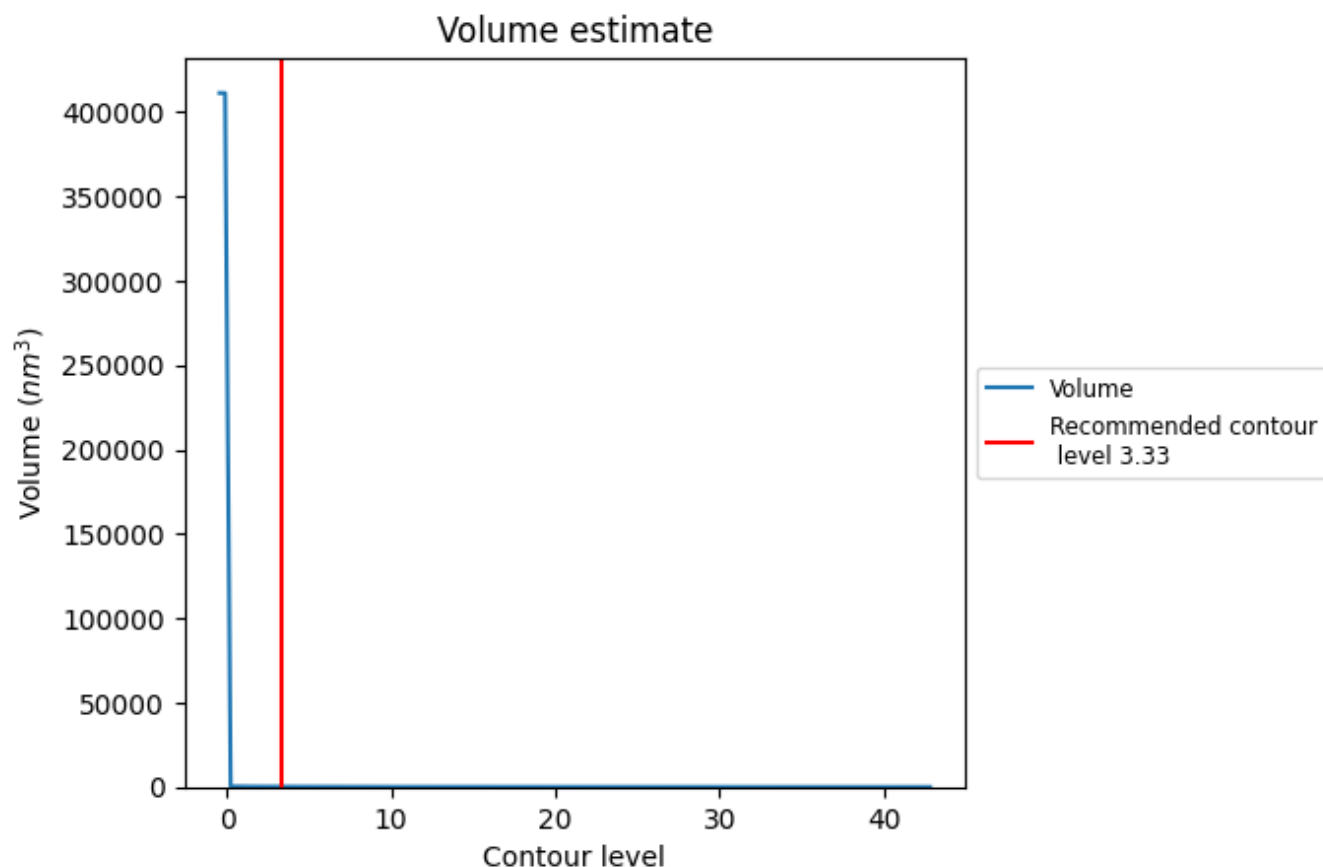
This section contains the results of statistical analysis of the map.

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.

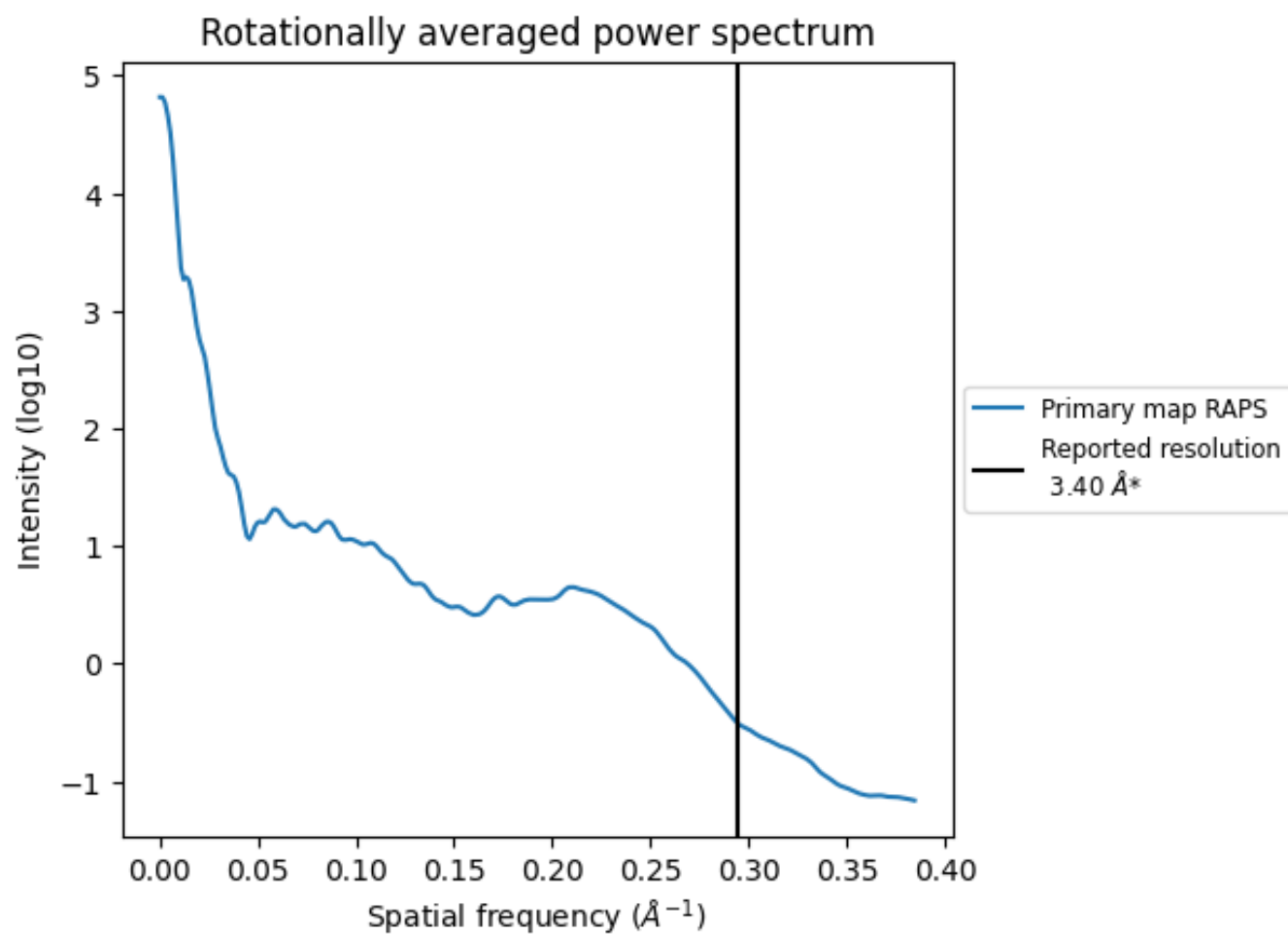
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 170 nm³; this corresponds to an approximate mass of 154 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 0.294 $\text{\AA}^{-1}$

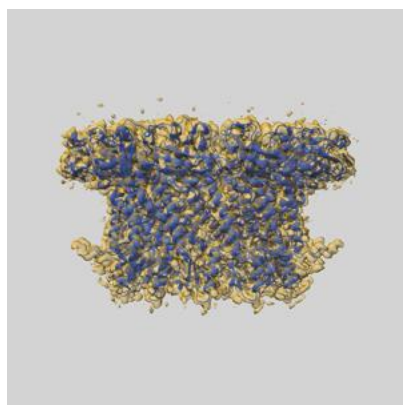
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

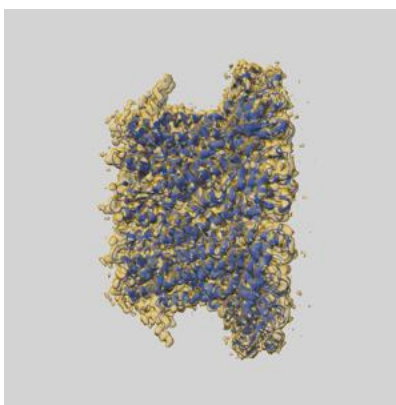
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-4678 and PDB model 6QYJ. Per-residue inclusion information can be found in section 3 on page 5.

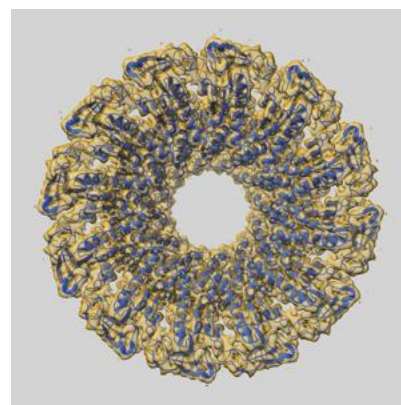
9.1 Map-model overlay [i](#)



X



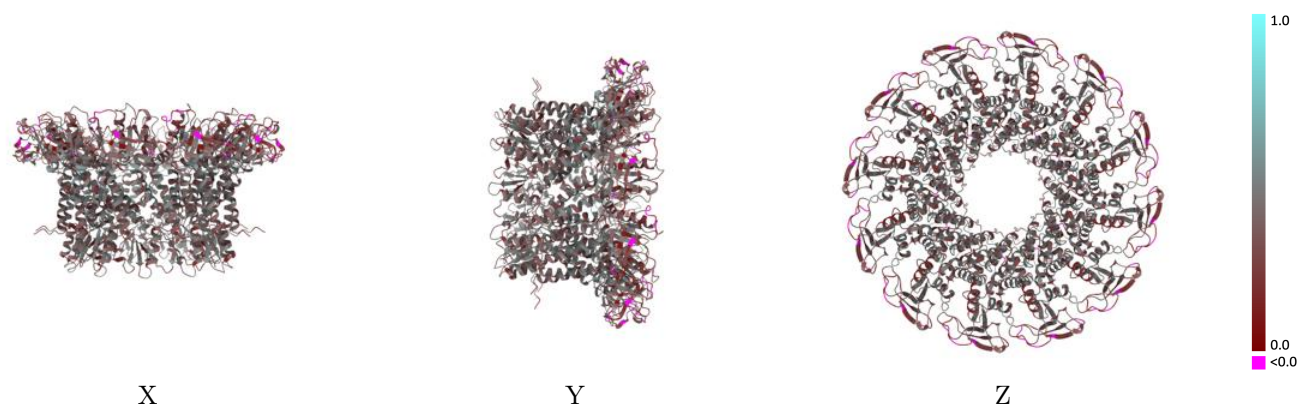
Y



Z

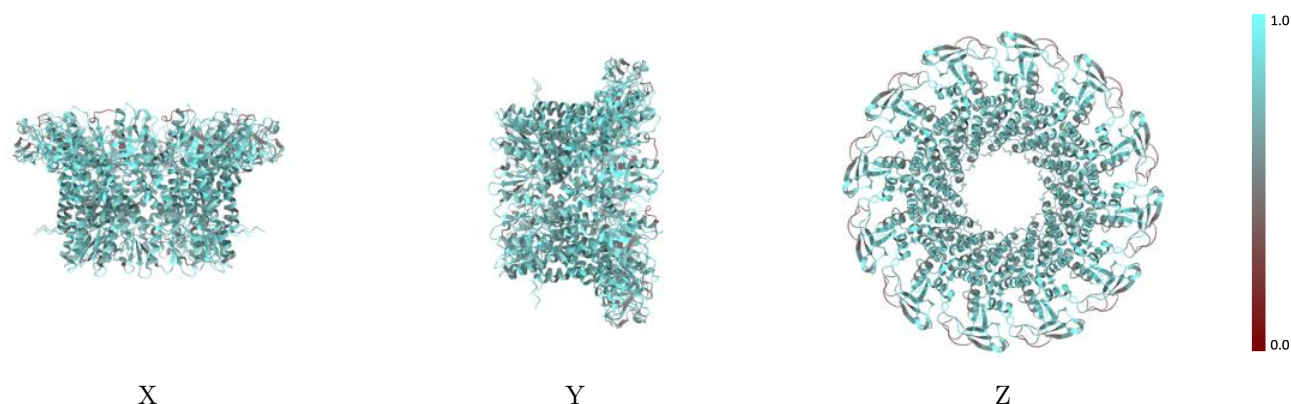
The images above show the 3D surface view of the map at the recommended contour level 3.33 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



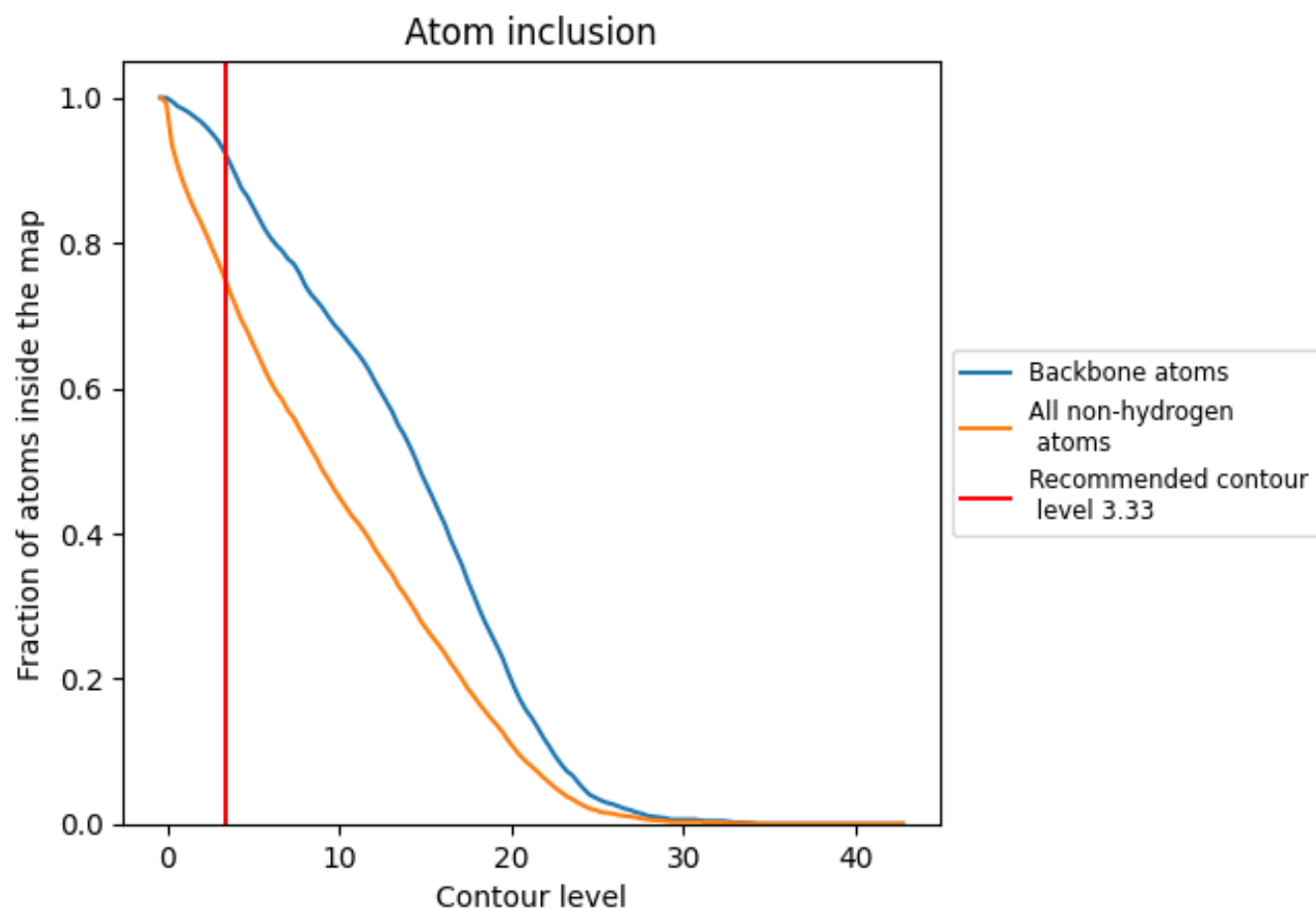
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (3.33).

9.4 Atom inclusion [i](#)



At the recommended contour level, 92% of all backbone atoms, 75% 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 (3.33) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7510	0.3620
0a	0.7520	0.3620
0b	0.7510	0.3620
0c	0.7500	0.3610
0d	0.7520	0.3620
0e	0.7510	0.3620
0f	0.7490	0.3630
0g	0.7520	0.3620
0h	0.7520	0.3630
0i	0.7490	0.3640
0j	0.7520	0.3610
0k	0.7500	0.3630
0l	0.7490	0.3630

1.0

0.0

<0.0