



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 02:50 AM UTC

PDB ID : 7BOZ / pdb_00007boz
EMDB ID : EMD-30137
Title : N-teminal of mature bacteriophage T7 tail fiber protein gp17
Authors : Chen, W.Y.; Xiao, H.
Deposited on : 2020-03-20
Resolution : 3.80 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:

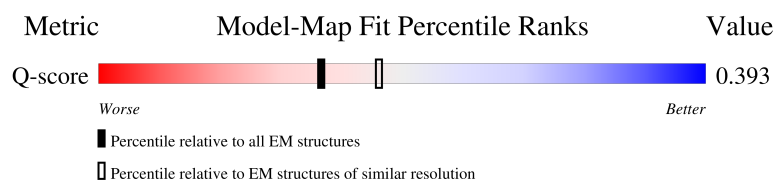
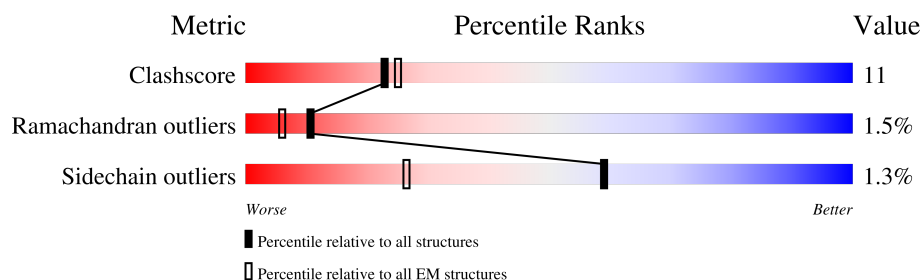
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 3.80 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.



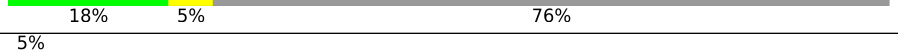
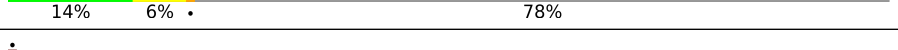
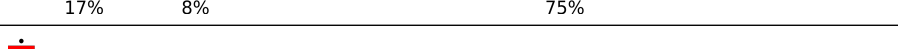
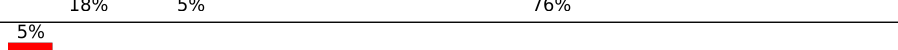
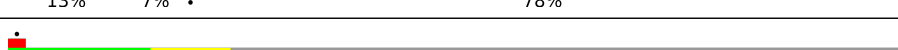
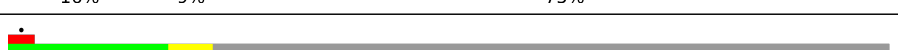
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	10198 (3.30 - 4.30)

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	553	
1	b	553	
1	c	553	
1	d	553	

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Mol	Chain	Length	Quality of chain
1	e	553	 17% 7% 76%
1	f	553	 14% 6% 78%
1	g	553	 17% 8% 75%
1	h	553	 18% 5% 76%
1	i	553	 5% 15% 5% 78%
1	j	553	 16% 8% 75%
1	k	553	 18% 5% 76%
1	l	553	 5% 14% 6% 78%
1	m	553	 17% 8% 75%
1	n	553	 18% 5% 76%
1	o	553	 5% 13% 7% 78%
1	p	553	 16% 9% 75%
1	q	553	 18% 5% 76%
1	r	553	 5% 13% 7% 78%



2 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 18516 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 N-terminal of mature bacteriophage T7 tail fiber protein gp17.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	a	138	Total	C	N	O	S	0	0
			1093	686	196	210	1		
1	b	131	Total	C	N	O	S	0	0
			1045	656	187	201	1		
1	c	119	Total	C	N	O	S	0	0
			948	598	164	185	1		
1	d	138	Total	C	N	O	S	0	0
			1093	686	196	210	1		
1	e	131	Total	C	N	O	S	0	0
			1045	656	187	201	1		
1	f	119	Total	C	N	O	S	0	0
			948	598	164	185	1		
1	g	138	Total	C	N	O	S	0	0
			1093	686	196	210	1		
1	h	131	Total	C	N	O	S	0	0
			1045	656	187	201	1		
1	i	119	Total	C	N	O	S	0	0
			948	598	164	185	1		
1	j	138	Total	C	N	O	S	0	0
			1093	686	196	210	1		
1	k	131	Total	C	N	O	S	0	0
			1045	656	187	201	1		
1	l	119	Total	C	N	O	S	0	0
			948	598	164	185	1		
1	m	138	Total	C	N	O	S	0	0
			1093	686	196	210	1		
1	n	131	Total	C	N	O	S	0	0
			1045	656	187	201	1		
1	o	119	Total	C	N	O	S	0	0
			948	598	164	185	1		
1	p	138	Total	C	N	O	S	0	0
			1093	686	196	210	1		
1	q	131	Total	C	N	O	S	0	0
			1045	656	187	201	1		

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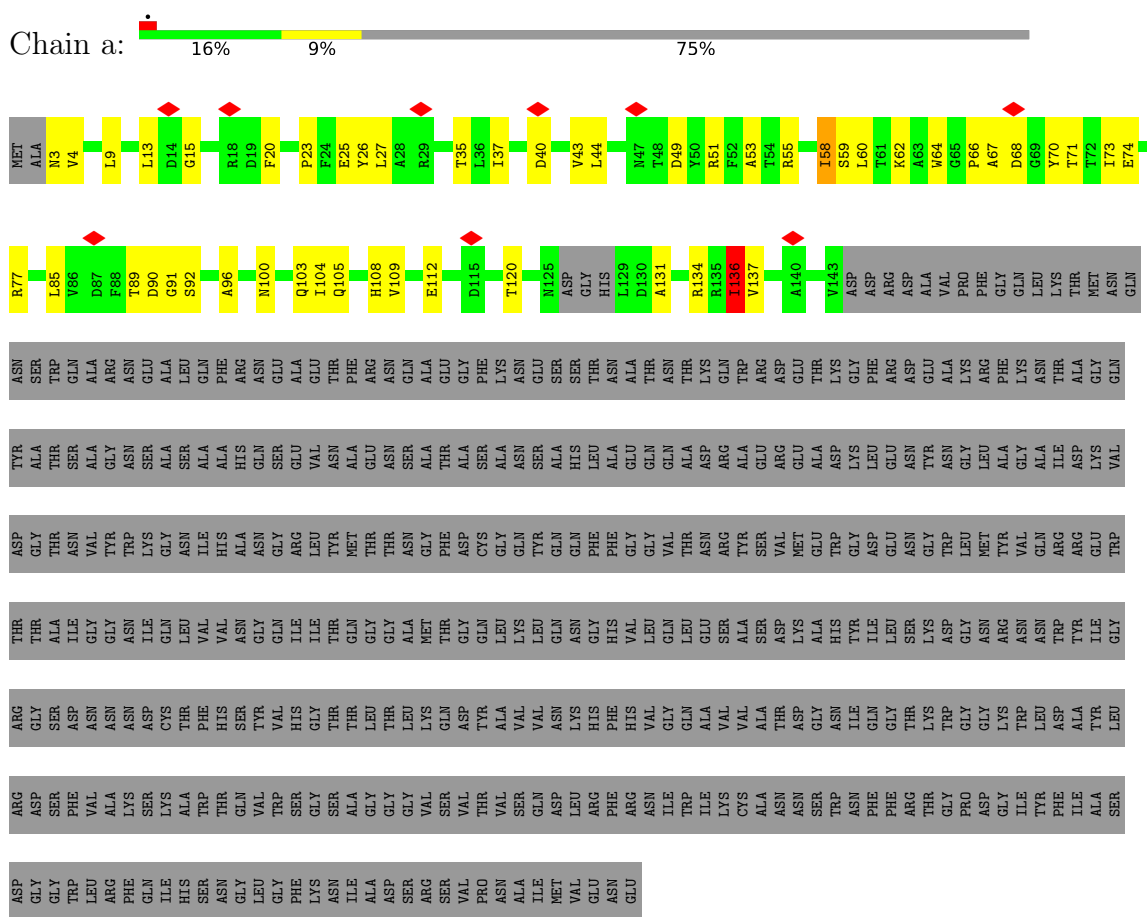
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Mol	Chain	Residues	Atoms					AltConf	Trace
1	r	119	Total	C	N	O	S	0	0
			948	598	164	185	1		

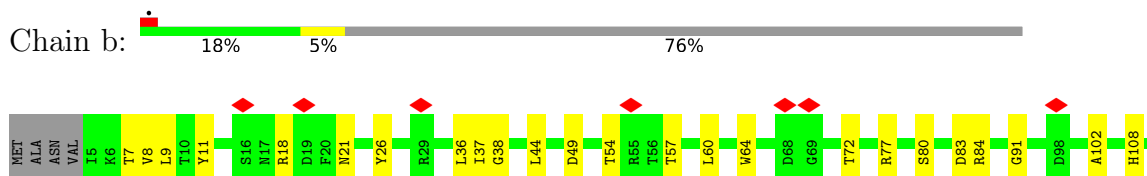
3 Residue-property plots

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.

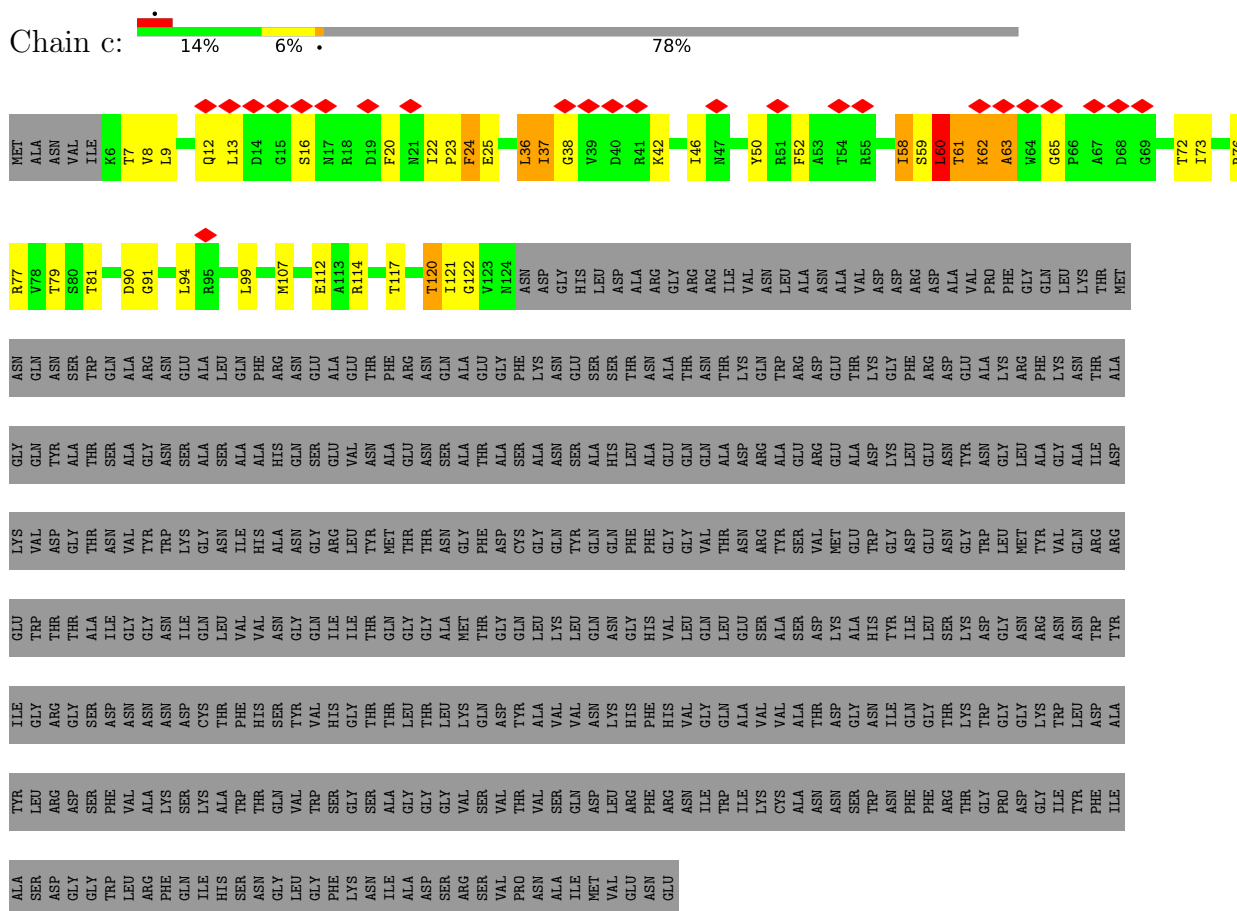
- Molecule 1: N-terminal of mature bacteriophage T7 tail fiber protein gp17



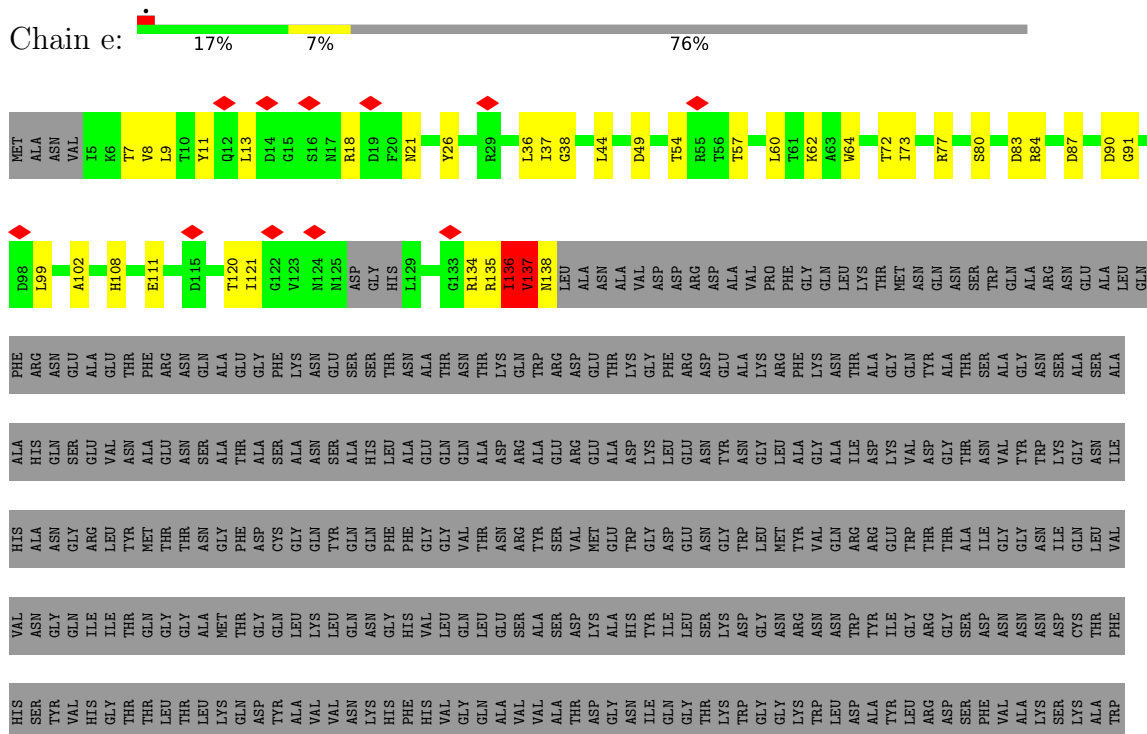
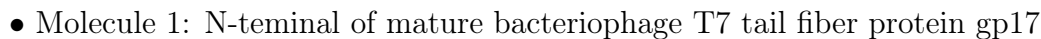
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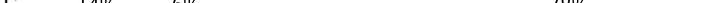


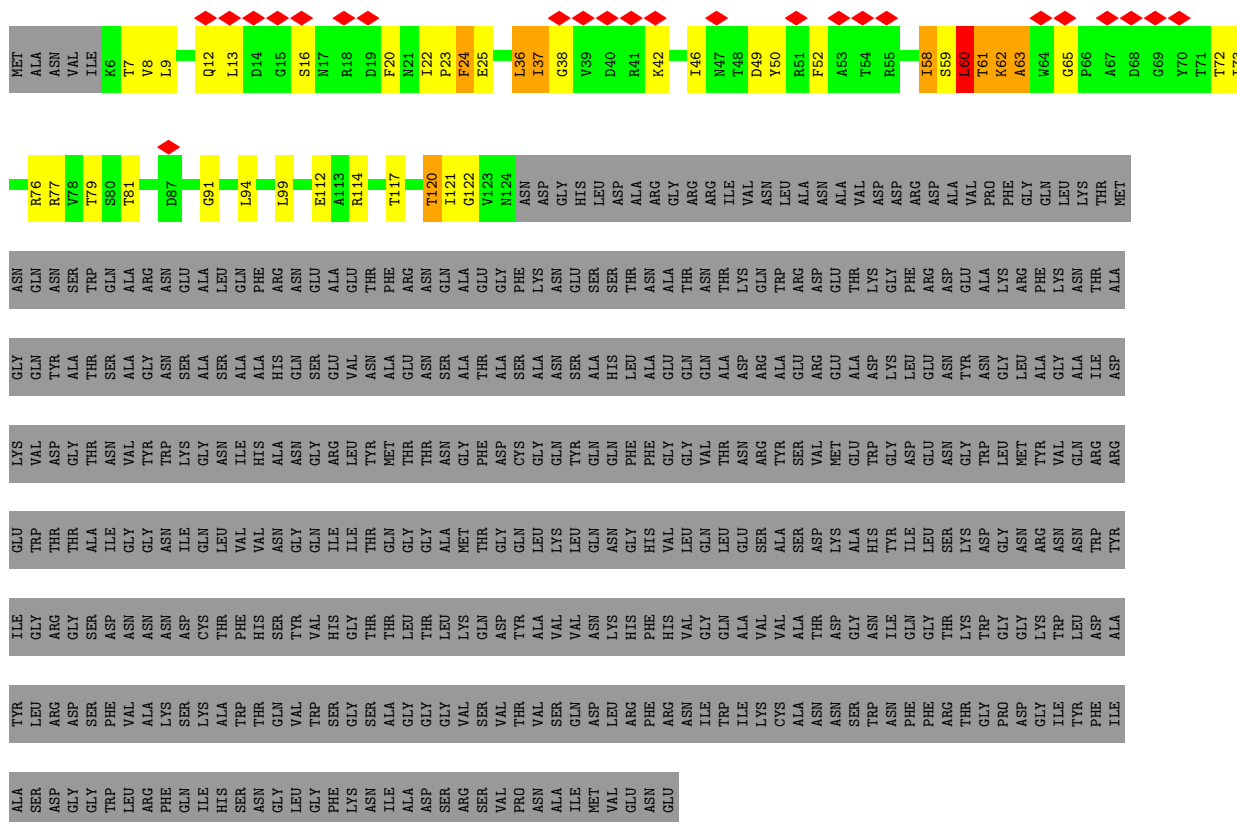
- Molecule 1: N-terminal of mature bacteriophage T7 tail fiber protein gp17



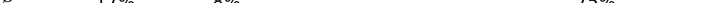
[illegible]

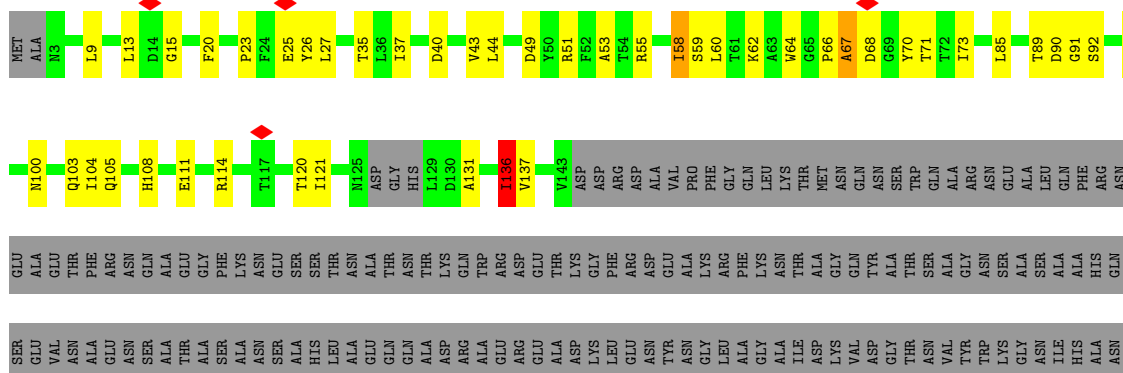
- Molecule 1: N-terminal of mature bacteriophage T7 tail fiber protein gp17

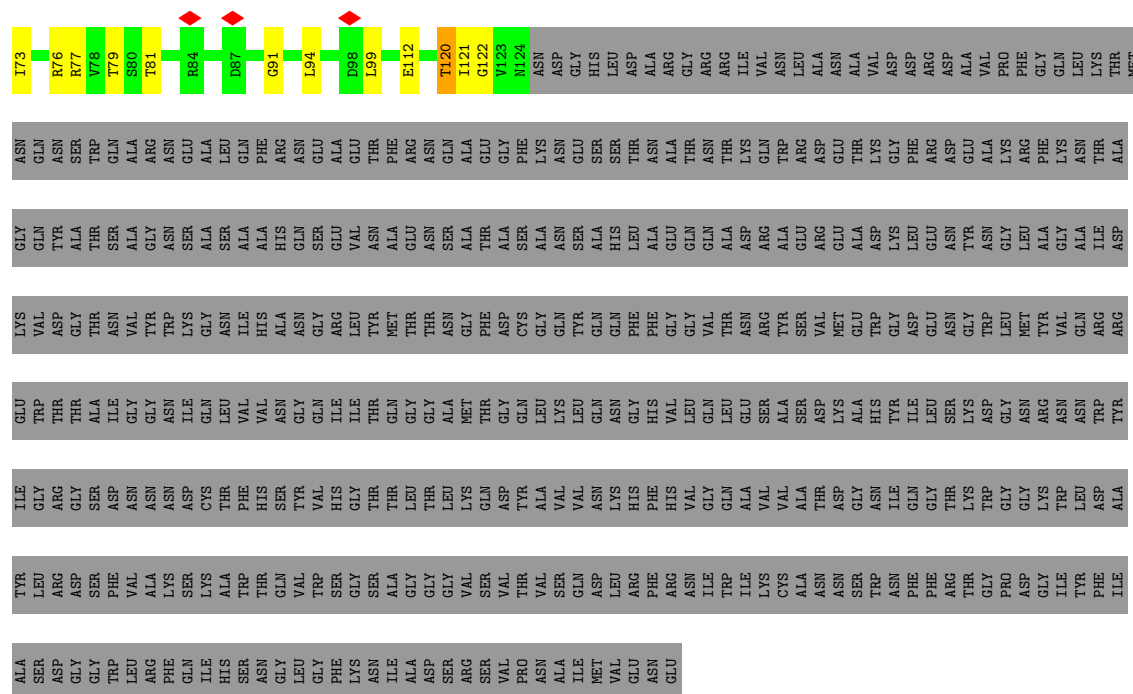
Chain f:  14% 6% 78%



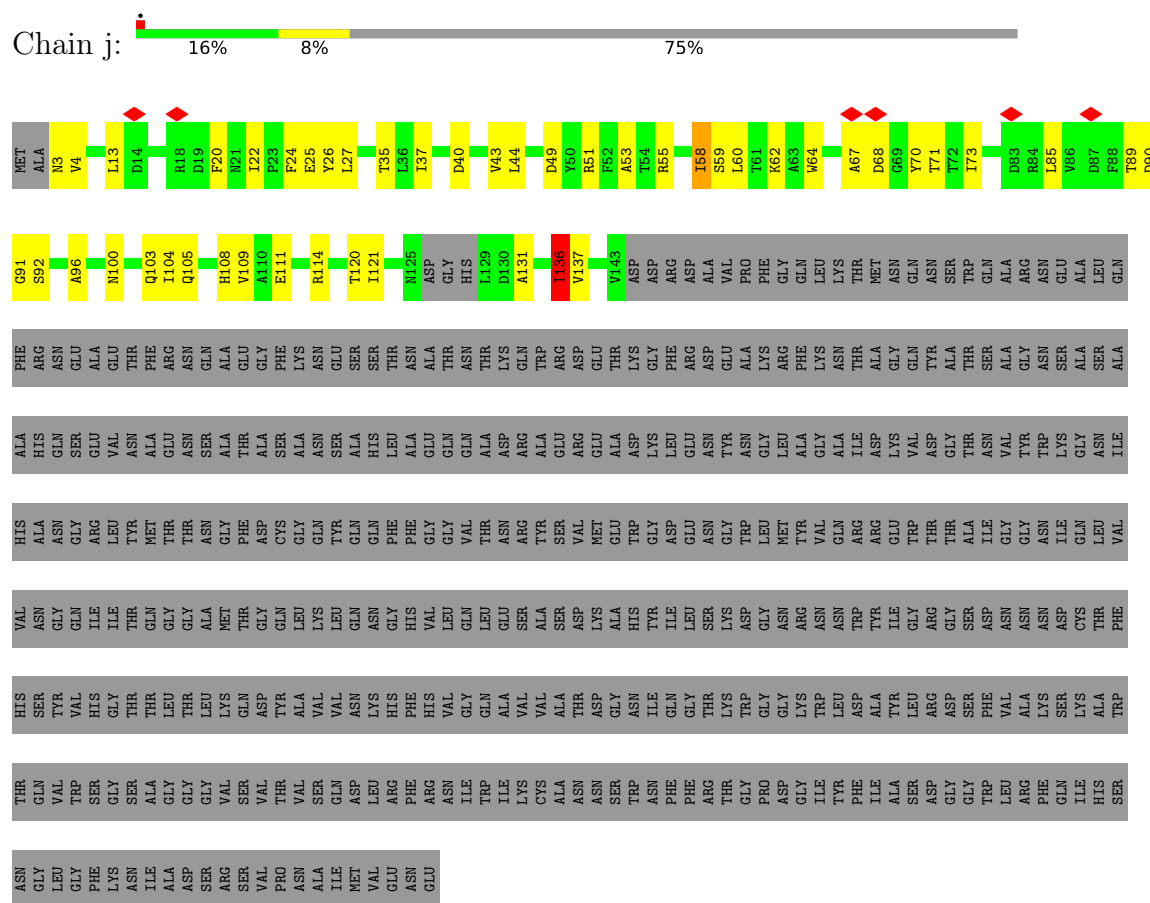
- Molecule 1: N-terminal of mature bacteriophage T7 tail fiber protein gp17

Chain g: 

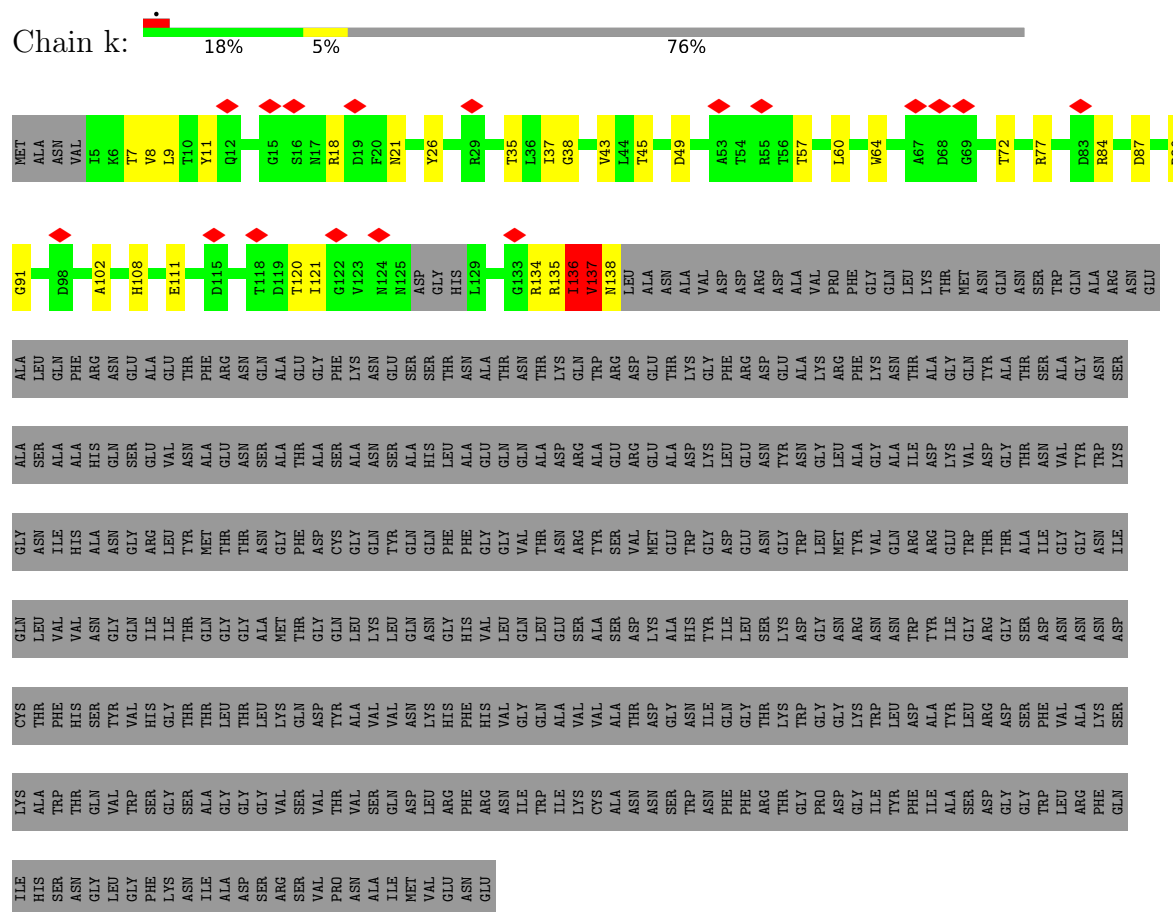




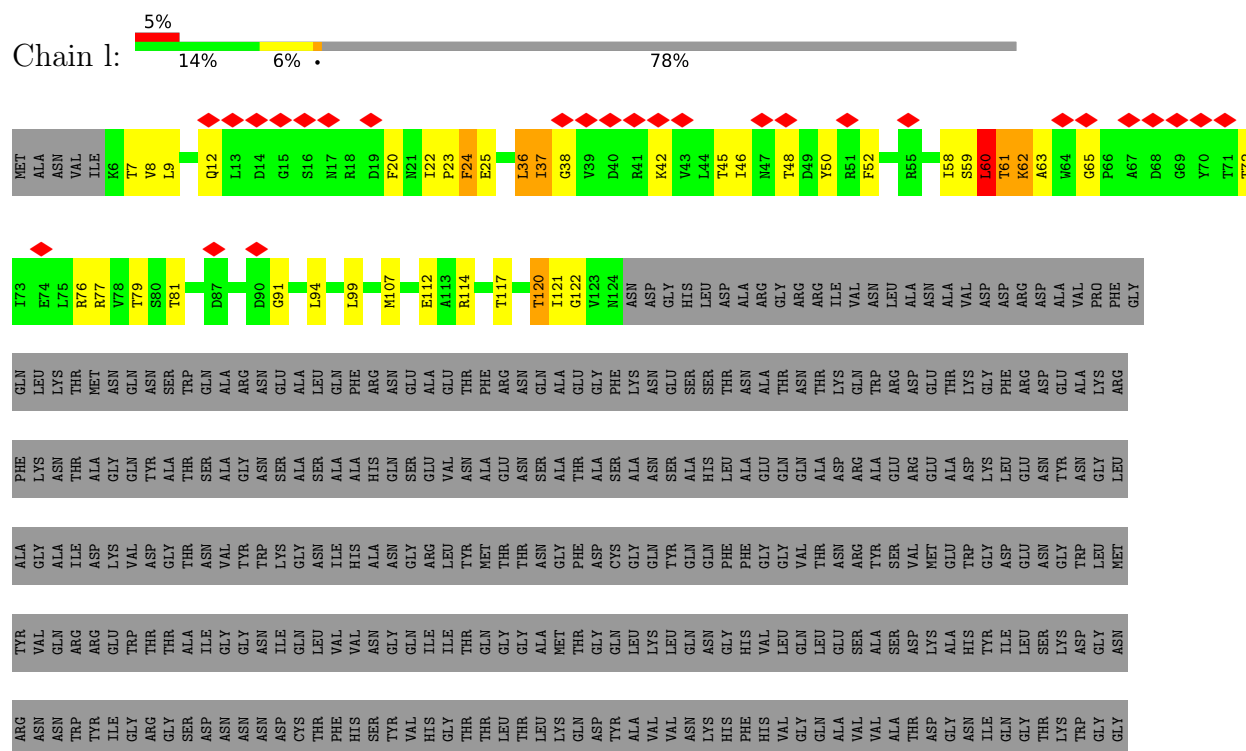
- Molecule 1: N-terminal of mature bacteriophage T7 tail fiber protein gp17



- Molecule 1: N-terminal of mature bacteriophage T7 tail fiber protein gp17

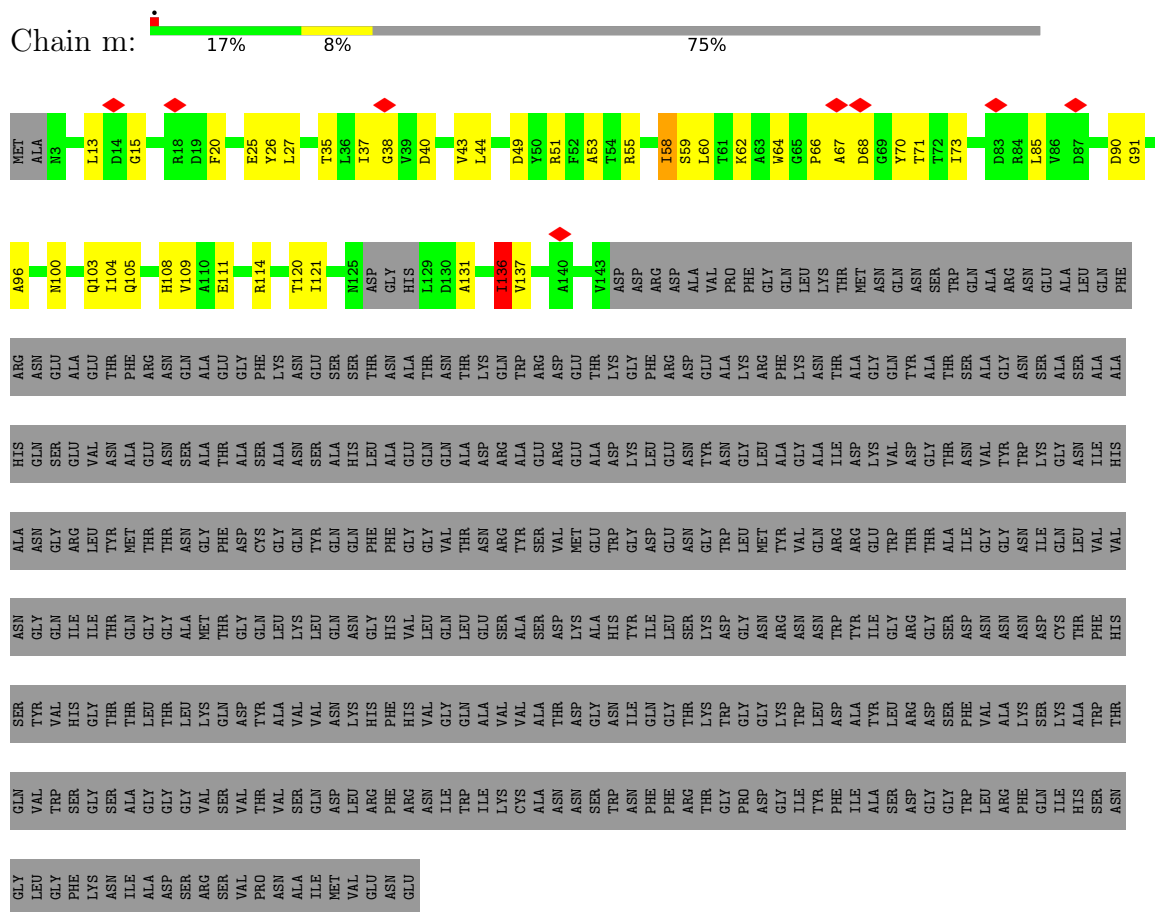


● Molecule 1: N-terminal of mature bacteriophage T7 tail fiber protein gp17

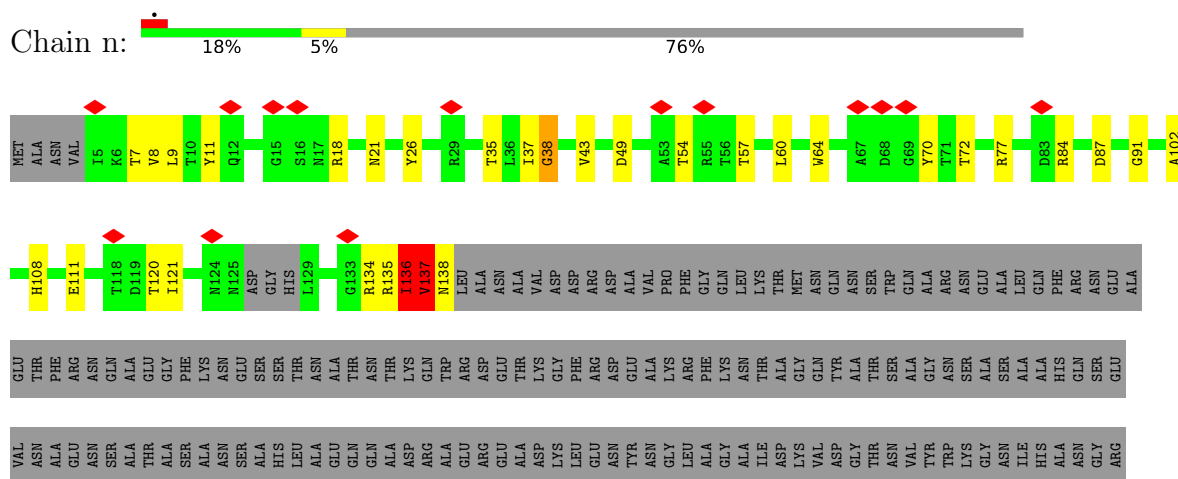


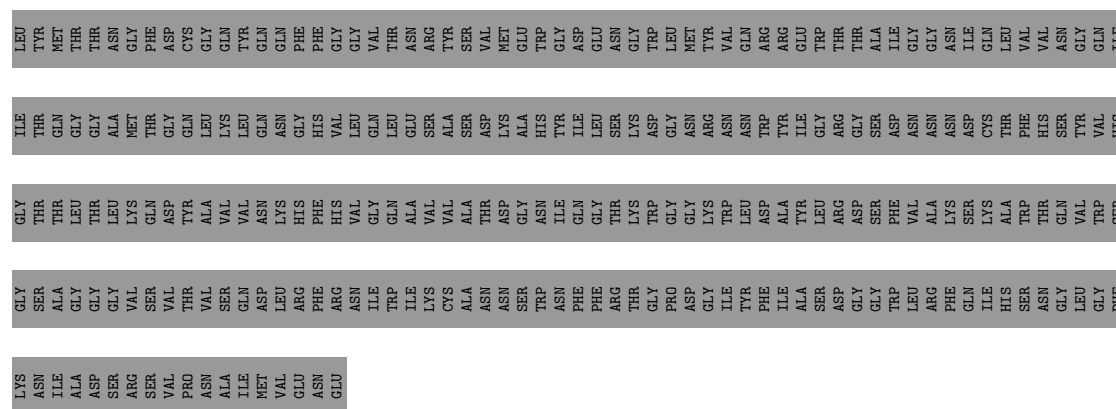
LYS	TRP	ASP	ALA	TYR	LEU	ARG	ASP	PHE	TRP	LEU	VAL	ALA	LYS	SER	GLN	THR	GLN	VAL	TRP	SER	GLY	SER	ALA	ALA	GLY	GLY	GLY	VAL	ARG	VAL	SER	VAL	THR	SER	ASN	VAL	ALA	SER	GLN	ASP	LEU	ARG	PHE	ASN	ARG	ASN	ASP	GLU	ASN	GLU
GLY	ILE	TYR	PHE	ILE	ALA	SER	ASP	GLY	TRP	LEU	ARG	PHE	GLN	ILE	HIS	SER	ASN	GLY	LEU	GLY	PHE	LYS	ASN	ILE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

● Molecule 1: N-terminal of mature bacteriophage T7 tail fiber protein gp17

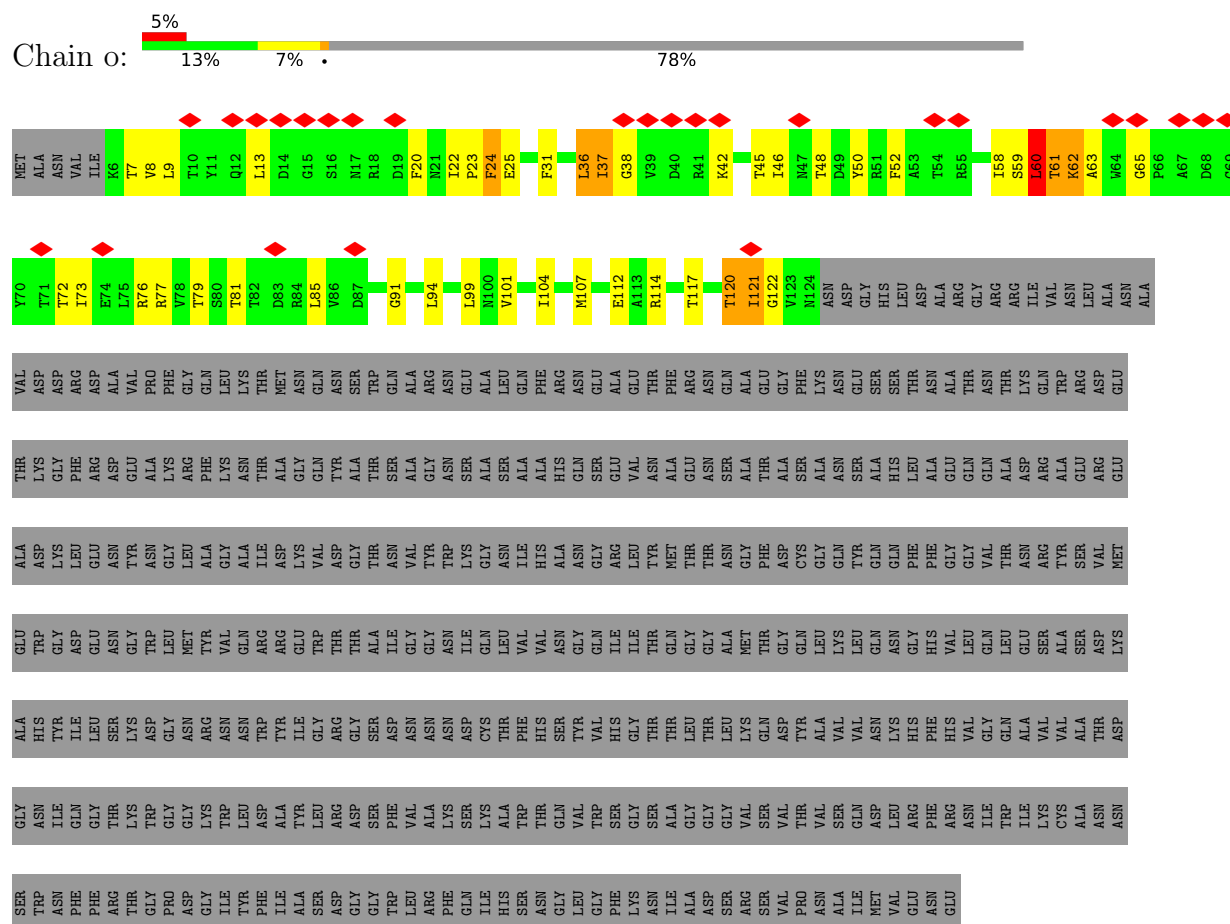


● Molecule 1: N-terminal of mature bacteriophage T7 tail fiber protein gp17

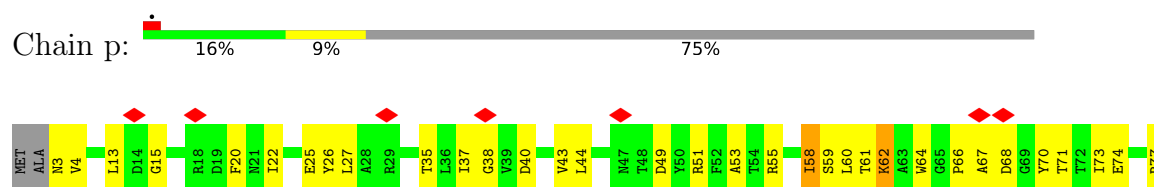




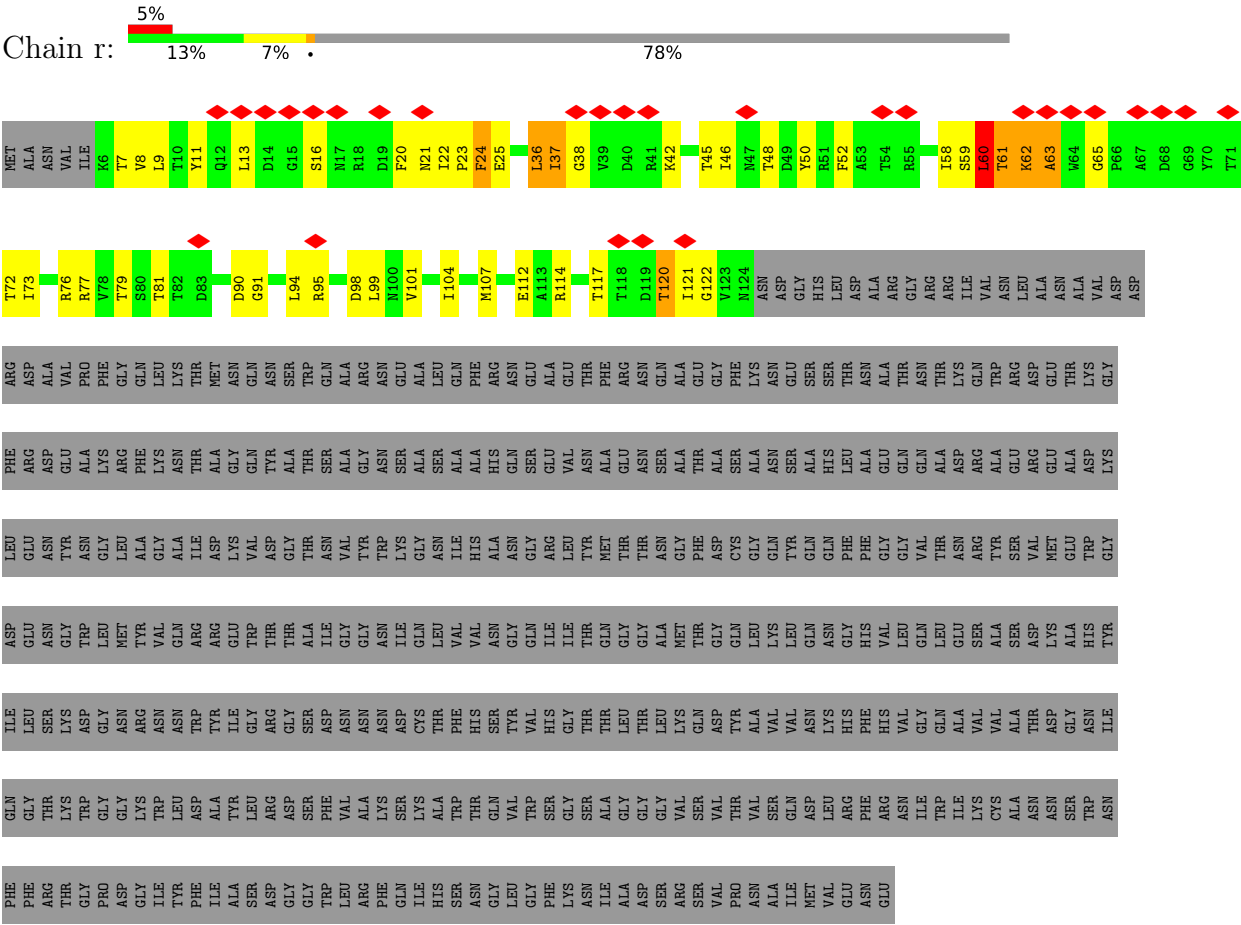
- Molecule 1: N-terminal of mature bacteriophage T7 tail fiber protein gp17



- Molecule 1: N-terminal of mature bacteriophage T7 tail fiber protein gp17







4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	55212	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TECNAI ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	24	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON II (4k x 4k)	Depositor
Maximum map value	27.532	Depositor
Minimum map value	-14.838	Depositor
Average map value	-0.319	Depositor
Map value standard deviation	0.994	Depositor
Recommended contour level	7	Depositor
Map size (Å)	447.03998, 447.03998, 447.03998	wwPDB
Map dimensions	352, 352, 352	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.27, 1.27, 1.27	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	a	0.47	0/1106	1.01	7/1504 (0.5%)
1	b	0.47	0/1058	1.09	10/1437 (0.7%)
1	c	0.47	0/962	1.14	9/1310 (0.7%)
1	d	0.48	0/1106	1.03	7/1504 (0.5%)
1	e	0.47	0/1058	1.09	10/1437 (0.7%)
1	f	0.48	0/962	1.14	9/1310 (0.7%)
1	g	0.47	0/1106	1.04	7/1504 (0.5%)
1	h	0.47	0/1058	1.06	8/1437 (0.6%)
1	i	0.48	0/962	1.14	9/1310 (0.7%)
1	j	0.45	0/1106	1.01	7/1504 (0.5%)
1	k	0.47	0/1058	1.04	8/1437 (0.6%)
1	l	0.46	0/962	1.13	9/1310 (0.7%)
1	m	0.45	0/1106	1.00	7/1504 (0.5%)
1	n	0.46	0/1058	1.04	9/1437 (0.6%)
1	o	0.46	0/962	1.13	9/1310 (0.7%)
1	p	0.46	0/1106	1.00	7/1504 (0.5%)
1	q	0.47	0/1058	1.07	9/1437 (0.6%)
1	r	0.46	0/962	1.13	9/1310 (0.7%)
All	All	0.47	0/18756	1.07	150/25506 (0.6%)

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	a	0	6
1	b	0	5
1	c	0	11
1	d	0	7
1	e	0	6
1	f	0	10
1	g	0	7

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Mol	Chain	#Chirality outliers	#Planarity outliers
1	h	0	6
1	i	0	10
1	j	0	6
1	k	0	7
1	l	0	10
1	m	0	7
1	n	0	5
1	o	0	10
1	p	0	7
1	q	0	5
1	r	0	11
All	All	0	136

There are no bond length outliers.

All (150) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	k	136	ILE	CA-C-N	7.67	135.78	121.97
1	k	136	ILE	C-N-CA	7.67	135.78	121.97
1	h	136	ILE	CA-C-N	7.67	135.78	121.97
1	h	136	ILE	C-N-CA	7.67	135.78	121.97
1	n	136	ILE	CA-C-N	7.61	135.67	121.97
1	n	136	ILE	C-N-CA	7.61	135.67	121.97
1	e	136	ILE	CA-C-N	7.55	135.56	121.97
1	e	136	ILE	C-N-CA	7.55	135.56	121.97
1	q	136	ILE	CA-C-N	7.50	135.48	121.97
1	q	136	ILE	C-N-CA	7.50	135.48	121.97
1	b	136	ILE	CA-C-N	7.49	135.46	121.97
1	b	136	ILE	C-N-CA	7.49	135.46	121.97
1	k	91	GLY	CA-C-N	7.35	134.93	121.70
1	k	91	GLY	C-N-CA	7.35	134.93	121.70
1	h	91	GLY	CA-C-N	7.29	134.81	121.70
1	h	91	GLY	C-N-CA	7.29	134.81	121.70
1	f	91	GLY	CA-C-N	7.28	134.80	121.70
1	f	91	GLY	C-N-CA	7.28	134.80	121.70
1	c	91	GLY	CA-C-N	7.17	134.60	121.70
1	c	91	GLY	C-N-CA	7.17	134.60	121.70
1	m	91	GLY	CA-C-N	7.16	134.59	121.70
1	m	91	GLY	C-N-CA	7.16	134.59	121.70
1	g	91	GLY	CA-C-N	7.15	134.57	121.70
1	g	91	GLY	C-N-CA	7.15	134.57	121.70
1	j	91	GLY	CA-C-N	7.13	134.54	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	j	91	GLY	C-N-CA	7.13	134.54	121.70
1	r	24	PHE	CA-C-N	7.12	135.13	121.54
1	r	24	PHE	C-N-CA	7.12	135.13	121.54
1	e	91	GLY	CA-C-N	7.12	134.51	121.70
1	e	91	GLY	C-N-CA	7.12	134.51	121.70
1	d	91	GLY	CA-C-N	7.10	134.49	121.70
1	d	91	GLY	C-N-CA	7.10	134.49	121.70
1	r	91	GLY	CA-C-N	7.10	134.48	121.70
1	r	91	GLY	C-N-CA	7.10	134.48	121.70
1	p	91	GLY	CA-C-N	7.08	134.44	121.70
1	p	91	GLY	C-N-CA	7.08	134.44	121.70
1	f	60	LEU	CA-CB-CG	7.07	141.03	116.30
1	o	24	PHE	CA-C-N	7.06	135.02	121.54
1	o	24	PHE	C-N-CA	7.06	135.02	121.54
1	a	91	GLY	CA-C-N	7.05	134.38	121.70
1	a	91	GLY	C-N-CA	7.05	134.38	121.70
1	f	24	PHE	CA-C-N	6.97	134.86	121.54
1	f	24	PHE	C-N-CA	6.97	134.86	121.54
1	c	24	PHE	CA-C-N	6.96	134.84	121.54
1	c	24	PHE	C-N-CA	6.96	134.84	121.54
1	c	60	LEU	CA-CB-CG	6.86	140.31	116.30
1	i	91	GLY	CA-C-N	6.85	134.03	121.70
1	i	91	GLY	C-N-CA	6.85	134.03	121.70
1	l	24	PHE	CA-C-N	6.85	134.62	121.54
1	l	24	PHE	C-N-CA	6.85	134.62	121.54
1	i	24	PHE	CA-C-N	6.84	134.60	121.54
1	i	24	PHE	C-N-CA	6.84	134.60	121.54
1	o	91	GLY	CA-C-N	6.82	133.98	121.70
1	o	91	GLY	C-N-CA	6.82	133.98	121.70
1	l	91	GLY	CA-C-N	6.78	133.91	121.70
1	l	91	GLY	C-N-CA	6.78	133.91	121.70
1	i	60	LEU	CA-CB-CG	6.73	139.86	116.30
1	g	67	ALA	CA-C-N	6.68	134.22	123.93
1	g	67	ALA	C-N-CA	6.68	134.22	123.93
1	f	61	THR	CA-C-N	6.67	134.27	121.54
1	f	61	THR	C-N-CA	6.67	134.27	121.54
1	i	61	THR	CA-C-N	6.65	134.24	121.54
1	i	61	THR	C-N-CA	6.65	134.24	121.54
1	k	137	VAL	N-CA-CB	-6.58	100.36	111.23
1	h	137	VAL	N-CA-CB	-6.58	100.38	111.23
1	d	67	ALA	CA-C-N	6.57	134.05	123.93
1	d	67	ALA	C-N-CA	6.57	134.05	123.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	c	61	THR	CA-C-N	6.57	134.09	121.54
1	c	61	THR	C-N-CA	6.57	134.09	121.54
1	o	37	ILE	CA-C-N	6.55	134.26	121.41
1	o	37	ILE	C-N-CA	6.55	134.26	121.41
1	r	37	ILE	CA-C-N	6.54	134.24	121.41
1	r	37	ILE	C-N-CA	6.54	134.24	121.41
1	b	137	VAL	N-CA-CB	-6.54	100.43	111.23
1	l	61	THR	CA-C-N	6.52	134.00	121.54
1	l	61	THR	C-N-CA	6.52	134.00	121.54
1	o	61	THR	CA-C-N	6.48	133.92	121.54
1	o	61	THR	C-N-CA	6.48	133.92	121.54
1	i	37	ILE	CA-C-N	6.46	134.08	121.41
1	i	37	ILE	C-N-CA	6.46	134.08	121.41
1	r	61	THR	CA-C-N	6.46	133.88	121.54
1	r	61	THR	C-N-CA	6.46	133.88	121.54
1	e	137	VAL	N-CA-CB	-6.45	100.59	111.23
1	l	37	ILE	CA-C-N	6.39	133.94	121.41
1	l	37	ILE	C-N-CA	6.39	133.94	121.41
1	n	91	GLY	CA-C-N	6.38	133.18	121.70
1	n	91	GLY	C-N-CA	6.38	133.18	121.70
1	q	137	VAL	N-CA-CB	-6.26	100.89	111.23
1	c	37	ILE	CA-C-N	6.26	133.68	121.41
1	c	37	ILE	C-N-CA	6.26	133.68	121.41
1	j	137	VAL	N-CA-C	6.22	116.70	110.23
1	r	60	LEU	CA-CB-CG	6.17	137.91	116.30
1	n	137	VAL	N-CA-CB	-6.17	101.05	111.23
1	f	37	ILE	CA-C-N	6.17	133.50	121.41
1	f	37	ILE	C-N-CA	6.17	133.50	121.41
1	q	91	GLY	CA-C-N	6.17	132.80	121.70
1	q	91	GLY	C-N-CA	6.17	132.80	121.70
1	m	137	VAL	N-CA-C	6.10	116.58	110.23
1	b	91	GLY	CA-C-N	6.08	132.63	121.70
1	b	91	GLY	C-N-CA	6.08	132.63	121.70
1	p	137	VAL	N-CA-C	6.04	116.51	110.23
1	l	60	LEU	CA-CB-CG	5.99	137.25	116.30
1	g	137	VAL	N-CA-C	5.93	116.40	110.23
1	a	137	VAL	N-CA-C	5.83	116.29	110.23
1	o	60	LEU	CA-CB-CG	5.80	136.61	116.30
1	d	137	VAL	N-CA-C	5.66	116.11	110.23
1	a	136	ILE	CA-C-N	5.60	128.11	120.77
1	a	136	ILE	C-N-CA	5.60	128.11	120.77
1	g	136	ILE	CA-C-N	5.53	128.01	120.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	g	136	ILE	C-N-CA	5.53	128.01	120.77
1	d	136	ILE	CA-C-N	5.47	127.94	120.77
1	d	136	ILE	C-N-CA	5.47	127.94	120.77
1	j	136	ILE	CA-C-N	5.41	127.86	120.77
1	j	136	ILE	C-N-CA	5.41	127.86	120.77
1	m	67	ALA	CA-C-N	5.39	134.41	124.82
1	m	67	ALA	C-N-CA	5.39	134.41	124.82
1	a	67	ALA	CA-C-N	5.36	134.36	124.82
1	a	67	ALA	C-N-CA	5.36	134.36	124.82
1	p	67	ALA	CA-C-N	5.35	134.34	124.82
1	p	67	ALA	C-N-CA	5.35	134.34	124.82
1	j	67	ALA	CA-C-N	5.34	134.32	124.82
1	j	67	ALA	C-N-CA	5.34	134.32	124.82
1	h	137	VAL	N-CA-C	5.30	120.37	109.34
1	p	136	ILE	CA-C-N	5.27	127.67	120.77
1	p	136	ILE	C-N-CA	5.27	127.67	120.77
1	e	26	TYR	CA-C-N	5.26	131.59	121.54
1	e	26	TYR	C-N-CA	5.26	131.59	121.54
1	m	136	ILE	CA-C-N	5.25	127.64	120.77
1	m	136	ILE	C-N-CA	5.25	127.64	120.77
1	h	26	TYR	CA-C-N	5.23	131.53	121.54
1	h	26	TYR	C-N-CA	5.23	131.53	121.54
1	q	54	THR	CA-C-N	5.18	131.43	121.54
1	q	54	THR	C-N-CA	5.18	131.43	121.54
1	b	26	TYR	CA-C-N	5.17	131.42	121.54
1	b	26	TYR	C-N-CA	5.17	131.42	121.54
1	n	54	THR	CA-C-N	5.16	131.40	121.54
1	n	54	THR	C-N-CA	5.16	131.40	121.54
1	e	137	VAL	N-CA-C	5.14	120.04	109.34
1	b	54	THR	CA-C-N	5.11	131.30	121.54
1	b	54	THR	C-N-CA	5.11	131.30	121.54
1	e	54	THR	CA-C-N	5.11	131.30	121.54
1	e	54	THR	C-N-CA	5.11	131.30	121.54
1	q	26	TYR	CA-C-N	5.11	131.29	121.54
1	q	26	TYR	C-N-CA	5.11	131.29	121.54
1	n	26	TYR	CA-C-N	5.04	131.18	121.54
1	n	26	TYR	C-N-CA	5.04	131.18	121.54
1	k	137	VAL	N-CA-C	5.02	119.78	109.34
1	b	137	VAL	N-CA-C	5.01	119.76	109.34
1	k	26	TYR	CA-C-N	5.00	131.09	121.54
1	k	26	TYR	C-N-CA	5.00	131.09	121.54

There are no chirality outliers.

All (136) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	a	131	ALA	Peptide
1	a	136	ILE	Peptide
1	a	51	ARG	Peptide
1	a	62	LYS	Peptide
1	a	68	ASP	Peptide
1	a	90	ASP	Peptide
1	b	120	THR	Peptide
1	b	134	ARG	Peptide
1	b	136	ILE	Peptide
1	b	137	VAL	Peptide
1	b	38	GLY	Peptide
1	c	120	THR	Peptide
1	c	22	ILE	Peptide
1	c	23	PRO	Peptide
1	c	24	PHE	Peptide
1	c	36	LEU	Peptide
1	c	37	ILE	Peptide
1	c	59	SER	Peptide
1	c	62	LYS	Peptide
1	c	63	ALA	Peptide
1	c	65	GLY	Peptide
1	c	90	ASP	Peptide
1	d	131	ALA	Peptide
1	d	136	ILE	Peptide
1	d	51	ARG	Peptide
1	d	62	LYS	Peptide
1	d	67	ALA	Peptide
1	d	68	ASP	Peptide
1	d	90	ASP	Peptide
1	e	120	THR	Peptide
1	e	134	ARG	Peptide
1	e	136	ILE	Peptide
1	e	137	VAL	Peptide
1	e	38	GLY	Peptide
1	e	90	ASP	Peptide
1	f	120	THR	Peptide
1	f	22	ILE	Peptide
1	f	23	PRO	Peptide
1	f	24	PHE	Peptide
1	f	36	LEU	Peptide
1	f	37	ILE	Peptide
1	f	59	SER	Peptide

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Mol	Chain	Res	Type	Group
1	f	62	LYS	Peptide
1	f	63	ALA	Peptide
1	f	65	GLY	Peptide
1	g	131	ALA	Peptide
1	g	136	ILE	Peptide
1	g	51	ARG	Peptide
1	g	62	LYS	Peptide
1	g	67	ALA	Peptide
1	g	68	ASP	Peptide
1	g	90	ASP	Peptide
1	h	120	THR	Peptide
1	h	134	ARG	Peptide
1	h	136	ILE	Peptide
1	h	137	VAL	Peptide
1	h	38	GLY	Peptide
1	h	90	ASP	Peptide
1	i	120	THR	Peptide
1	i	22	ILE	Peptide
1	i	23	PRO	Peptide
1	i	24	PHE	Peptide
1	i	36	LEU	Peptide
1	i	37	ILE	Peptide
1	i	59	SER	Peptide
1	i	62	LYS	Peptide
1	i	63	ALA	Peptide
1	i	65	GLY	Peptide
1	j	131	ALA	Peptide
1	j	136	ILE	Peptide
1	j	51	ARG	Peptide
1	j	62	LYS	Peptide
1	j	68	ASP	Peptide
1	j	90	ASP	Peptide
1	k	120	THR	Peptide
1	k	134	ARG	Peptide
1	k	136	ILE	Peptide
1	k	137	VAL	Peptide
1	k	38	GLY	Peptide
1	k	45	THR	Peptide
1	k	90	ASP	Peptide
1	l	120	THR	Peptide
1	l	22	ILE	Peptide
1	l	23	PRO	Peptide

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Mol	Chain	Res	Type	Group
1	l	24	PHE	Peptide
1	l	36	LEU	Peptide
1	l	37	ILE	Peptide
1	l	59	SER	Peptide
1	l	62	LYS	Peptide
1	l	63	ALA	Peptide
1	l	65	GLY	Peptide
1	m	131	ALA	Peptide
1	m	136	ILE	Peptide
1	m	38	GLY	Peptide
1	m	51	ARG	Peptide
1	m	62	LYS	Peptide
1	m	68	ASP	Peptide
1	m	90	ASP	Peptide
1	n	120	THR	Peptide
1	n	134	ARG	Peptide
1	n	136	ILE	Peptide
1	n	137	VAL	Peptide
1	n	38	GLY	Peptide
1	o	120	THR	Peptide
1	o	22	ILE	Peptide
1	o	23	PRO	Peptide
1	o	24	PHE	Peptide
1	o	36	LEU	Peptide
1	o	37	ILE	Peptide
1	o	59	SER	Peptide
1	o	62	LYS	Peptide
1	o	63	ALA	Peptide
1	o	65	GLY	Peptide
1	p	131	ALA	Peptide
1	p	136	ILE	Peptide
1	p	38	GLY	Peptide
1	p	51	ARG	Peptide
1	p	62	LYS	Peptide
1	p	68	ASP	Peptide
1	p	90	ASP	Peptide
1	q	120	THR	Peptide
1	q	134	ARG	Peptide
1	q	136	ILE	Peptide
1	q	137	VAL	Peptide
1	q	38	GLY	Peptide
1	r	120	THR	Peptide

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Mol	Chain	Res	Type	Group
1	r	22	ILE	Peptide
1	r	23	PRO	Peptide
1	r	24	PHE	Peptide
1	r	36	LEU	Peptide
1	r	37	ILE	Peptide
1	r	59	SER	Peptide
1	r	62	LYS	Peptide
1	r	63	ALA	Peptide
1	r	65	GLY	Peptide
1	r	90	ASP	Peptide

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	1093	0	1107	31	0
1	b	1045	0	1056	19	0
1	c	948	0	952	23	0
1	d	1093	0	1107	28	0
1	e	1045	0	1056	25	0
1	f	948	0	952	25	0
1	g	1093	0	1107	30	0
1	h	1045	0	1056	23	0
1	i	948	0	952	22	0
1	j	1093	0	1107	32	0
1	k	1045	0	1056	22	0
1	l	948	0	952	23	0
1	m	1093	0	1107	29	0
1	n	1045	0	1056	23	0
1	o	948	0	952	26	0
1	p	1093	0	1107	35	0
1	q	1045	0	1056	23	0
1	r	948	0	952	28	0
All	All	18516	0	18690	418	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 11.

All (418) 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:j:136:ILE:HD12	1:k:135:ARG:NH1	1.85	0.91
1:p:136:ILE:HD12	1:q:135:ARG:NH1	1.86	0.91
1:a:136:ILE:HD12	1:b:135:ARG:NH1	1.88	0.89
1:g:136:ILE:HD12	1:h:135:ARG:NH1	1.89	0.88
1:j:136:ILE:CD1	1:k:135:ARG:NH1	2.38	0.86
1:m:136:ILE:HD12	1:n:135:ARG:NH1	1.89	0.86
1:p:136:ILE:CD1	1:q:135:ARG:NH1	2.40	0.84
1:m:136:ILE:CD1	1:n:135:ARG:NH1	2.43	0.81
1:d:136:ILE:HD12	1:e:135:ARG:NH1	1.97	0.79
1:j:136:ILE:O	1:j:136:ILE:HG22	1.82	0.79
1:m:136:ILE:HG22	1:m:136:ILE:O	1.82	0.79
1:a:136:ILE:HG22	1:a:136:ILE:O	1.82	0.79
1:a:136:ILE:CD1	1:b:135:ARG:NH1	2.47	0.78
1:g:136:ILE:CD1	1:h:135:ARG:NH1	2.47	0.78
1:p:136:ILE:HG22	1:p:136:ILE:O	1.81	0.77
1:g:136:ILE:HG22	1:g:136:ILE:O	1.84	0.76
1:d:136:ILE:HG22	1:d:136:ILE:O	1.86	0.75
1:q:111:GLU:HG3	1:r:81:THR:HG21	1.71	0.71
1:b:111:GLU:HG3	1:c:81:THR:HG21	1.73	0.70
1:k:37:ILE:HG22	1:k:72:THR:H	1.56	0.69
1:q:37:ILE:HG22	1:q:72:THR:H	1.56	0.69
1:n:37:ILE:HG22	1:n:72:THR:H	1.57	0.68
1:n:111:GLU:HG3	1:o:81:THR:HG21	1.75	0.68
1:e:37:ILE:HG22	1:e:72:THR:H	1.57	0.67
1:h:37:ILE:HG22	1:h:72:THR:H	1.57	0.67
1:e:111:GLU:HG3	1:f:81:THR:HG21	1.77	0.66
1:l:61:THR:OG1	1:l:62:LYS:N	2.29	0.66
1:k:111:GLU:HG3	1:l:81:THR:HG21	1.77	0.66
1:f:36:LEU:HB2	1:f:42:LYS:HB3	1.78	0.65
1:b:37:ILE:HG22	1:b:72:THR:H	1.57	0.65
1:d:136:ILE:CD1	1:e:135:ARG:NH1	2.60	0.65
1:c:36:LEU:HB2	1:c:42:LYS:HB3	1.79	0.65
1:h:111:GLU:HG3	1:i:81:THR:HG21	1.79	0.65
1:r:36:LEU:HB2	1:r:42:LYS:HB3	1.80	0.64
1:i:36:LEU:HB2	1:i:42:LYS:HB3	1.79	0.64
1:o:36:LEU:HB2	1:o:42:LYS:HB3	1.80	0.64
1:o:61:THR:OG1	1:o:62:LYS:N	2.29	0.64
1:m:26:TYR:OH	1:m:55:ARG:O	2.16	0.63
1:r:60:LEU:HD23	1:r:61:THR:HG23	1.80	0.63
1:c:60:LEU:HD23	1:c:61:THR:HG23	1.79	0.63
1:f:20:PHE:HD2	1:f:58:ILE:HD11	1.64	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:l:36:LEU:HB2	1:l:42:LYS:HB3	1.81	0.61
1:l:120:THR:O	1:l:122:GLY:N	2.34	0.61
1:o:120:THR:O	1:o:122:GLY:N	2.33	0.61
1:a:26:TYR:OH	1:a:55:ARG:O	2.15	0.61
1:d:60:LEU:HB3	1:d:64:TRP:HZ3	1.66	0.61
1:r:120:THR:O	1:r:122:GLY:N	2.34	0.61
1:c:61:THR:OG1	1:c:62:LYS:N	2.30	0.61
1:i:20:PHE:HD2	1:i:58:ILE:HD11	1.66	0.60
1:i:60:LEU:HD23	1:i:61:THR:HG23	1.83	0.60
1:i:120:THR:O	1:i:122:GLY:N	2.34	0.60
1:j:136:ILE:HD12	1:k:135:ARG:HH12	1.64	0.60
1:p:26:TYR:OH	1:p:55:ARG:O	2.14	0.60
1:f:60:LEU:HD23	1:f:61:THR:HG23	1.83	0.60
1:l:60:LEU:HD23	1:l:61:THR:HG23	1.82	0.60
1:a:60:LEU:HB3	1:a:64:TRP:HZ3	1.66	0.60
1:c:120:THR:O	1:c:122:GLY:N	2.35	0.60
1:f:120:THR:O	1:f:122:GLY:N	2.35	0.59
1:o:60:LEU:HD23	1:o:61:THR:HG23	1.84	0.59
1:o:94:LEU:HD13	1:o:99:LEU:HD21	1.84	0.59
1:b:21:ASN:HA	1:b:57:THR:HA	1.85	0.59
1:j:26:TYR:OH	1:j:55:ARG:O	2.19	0.59
1:f:61:THR:OG1	1:f:62:LYS:N	2.30	0.59
1:p:136:ILE:HD12	1:q:135:ARG:HH12	1.65	0.59
1:i:61:THR:OG1	1:i:62:LYS:N	2.30	0.58
1:r:36:LEU:HD22	1:r:42:LYS:HD3	1.85	0.58
1:n:11:TYR:OH	1:n:21:ASN:O	2.18	0.58
1:d:26:TYR:OH	1:d:55:ARG:O	2.19	0.58
1:e:21:ASN:HA	1:e:57:THR:HA	1.86	0.58
1:g:60:LEU:HB3	1:g:64:TRP:HZ3	1.69	0.58
1:l:94:LEU:HD13	1:l:99:LEU:HD21	1.86	0.58
1:c:36:LEU:HD22	1:c:42:LYS:HD3	1.86	0.58
1:p:60:LEU:HB3	1:p:64:TRP:HZ3	1.69	0.58
1:f:36:LEU:HD22	1:f:42:LYS:HD3	1.85	0.57
1:g:89:THR:O	1:g:92:SER:OG	2.22	0.57
1:f:94:LEU:HD13	1:f:99:LEU:HD21	1.86	0.57
1:m:53:ALA:HB2	1:m:59:SER:HB3	1.85	0.57
1:m:136:ILE:HD12	1:n:135:ARG:HH12	1.68	0.57
1:j:53:ALA:HB2	1:j:59:SER:HB3	1.87	0.57
1:m:109:VAL:HG21	1:o:107:MET:HE3	1.87	0.57
1:p:37:ILE:O	1:p:71:THR:N	2.36	0.57
1:i:94:LEU:HD13	1:i:99:LEU:HD21	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:i:36:LEU:HD22	1:i:42:LYS:HD3	1.87	0.56
1:j:89:THR:O	1:j:92:SER:OG	2.23	0.56
1:h:21:ASN:HA	1:h:57:THR:HA	1.88	0.56
1:m:60:LEU:HB3	1:m:64:TRP:HZ3	1.68	0.56
1:p:53:ALA:HB2	1:p:59:SER:HB3	1.87	0.56
1:a:25:GLU:HG3	1:a:108:HIS:HE1	1.70	0.56
1:g:26:TYR:OH	1:g:55:ARG:O	2.21	0.56
1:r:61:THR:OG1	1:r:62:LYS:N	2.30	0.56
1:o:36:LEU:HD22	1:o:42:LYS:HD3	1.87	0.55
1:d:25:GLU:HG3	1:d:108:HIS:HE1	1.71	0.55
1:j:60:LEU:HB3	1:j:64:TRP:HZ3	1.71	0.55
1:h:7:THR:HB	1:h:77:ARG:HG2	1.88	0.55
1:q:21:ASN:HA	1:q:57:THR:HA	1.88	0.55
1:n:21:ASN:HA	1:n:57:THR:HA	1.89	0.55
1:p:109:VAL:HG21	1:r:107:MET:HE3	1.89	0.54
1:l:36:LEU:HD22	1:l:42:LYS:HD3	1.89	0.54
1:a:37:ILE:O	1:a:71:THR:N	2.37	0.54
1:a:53:ALA:HB2	1:a:59:SER:HB3	1.89	0.54
1:k:11:TYR:OH	1:k:21:ASN:O	2.17	0.54
1:g:53:ALA:HB2	1:g:59:SER:HB3	1.90	0.54
1:h:49:ASP:HB2	1:h:64:TRP:HH2	1.72	0.54
1:j:37:ILE:O	1:j:71:THR:N	2.41	0.54
1:p:25:GLU:HG3	1:p:108:HIS:HE1	1.72	0.54
1:b:49:ASP:HB2	1:b:64:TRP:HH2	1.73	0.54
1:q:49:ASP:HB2	1:q:64:TRP:HH2	1.72	0.54
1:e:7:THR:HB	1:e:77:ARG:HG2	1.90	0.54
1:k:21:ASN:HA	1:k:57:THR:HA	1.91	0.53
1:p:103:GLN:HG3	1:q:102:ALA:HB1	1.90	0.53
1:e:11:TYR:OH	1:e:21:ASN:O	2.20	0.53
1:q:7:THR:OG1	1:q:8:VAL:N	2.42	0.53
1:m:136:ILE:O	1:m:136:ILE:CG2	2.54	0.53
1:e:49:ASP:HB2	1:e:64:TRP:HH2	1.74	0.53
1:k:7:THR:HB	1:k:77:ARG:HG2	1.90	0.52
1:r:94:LEU:HD13	1:r:99:LEU:HD21	1.90	0.52
1:c:20:PHE:HD2	1:c:58:ILE:HD11	1.73	0.52
1:j:103:GLN:HE22	1:k:84:ARG:HD2	1.74	0.52
1:h:7:THR:OG1	1:h:8:VAL:N	2.41	0.52
1:j:40:ASP:HB2	1:j:70:TYR:HE1	1.75	0.52
1:d:103:GLN:HG3	1:e:102:ALA:HB1	1.92	0.52
1:b:7:THR:HB	1:b:77:ARG:HG2	1.91	0.52
1:b:11:TYR:OH	1:b:21:ASN:O	2.20	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:a:27:LEU:HD22	1:a:85:LEU:HD22	1.92	0.52
1:j:111:GLU:OE1	1:j:114:ARG:NH2	2.43	0.52
1:g:40:ASP:HB2	1:g:70:TYR:HE1	1.74	0.51
1:g:103:GLN:HG3	1:h:102:ALA:HB1	1.92	0.51
1:p:13:LEU:HD12	1:p:73:ILE:HG13	1.91	0.51
1:p:27:LEU:HD22	1:p:85:LEU:HD22	1.93	0.51
1:a:40:ASP:HB2	1:a:70:TYR:HE1	1.75	0.51
1:b:7:THR:OG1	1:b:8:VAL:N	2.41	0.51
1:q:77:ARG:HH12	1:q:108:HIS:HB3	1.76	0.51
1:a:103:GLN:HG3	1:b:102:ALA:HB1	1.93	0.51
1:m:13:LEU:HD12	1:m:73:ILE:HG13	1.93	0.51
1:m:37:ILE:O	1:m:71:THR:N	2.37	0.51
1:c:58:ILE:HD12	1:c:60:LEU:HA	1.93	0.51
1:h:11:TYR:OH	1:h:21:ASN:O	2.19	0.51
1:k:7:THR:OG1	1:k:8:VAL:N	2.42	0.51
1:m:103:GLN:HG3	1:n:102:ALA:HB1	1.91	0.51
1:n:7:THR:HB	1:n:77:ARG:HG2	1.92	0.51
1:d:40:ASP:HB2	1:d:70:TYR:HE1	1.75	0.51
1:g:103:GLN:HE22	1:h:84:ARG:HD2	1.75	0.50
1:r:46:ILE:O	1:r:50:TYR:N	2.44	0.50
1:e:7:THR:OG1	1:e:8:VAL:N	2.42	0.50
1:m:103:GLN:HE22	1:n:84:ARG:HD2	1.76	0.50
1:d:53:ALA:HB2	1:d:59:SER:HB3	1.93	0.50
1:f:52:PHE:CE2	1:f:58:ILE:HG22	2.47	0.50
1:r:45:THR:HG1	1:r:48:THR:HG1	1.58	0.50
1:c:46:ILE:O	1:c:50:TYR:N	2.45	0.50
1:l:77:ARG:HD2	1:l:79:THR:HG23	1.94	0.50
1:o:46:ILE:O	1:o:50:TYR:N	2.45	0.50
1:c:94:LEU:HD13	1:c:99:LEU:HD21	1.92	0.50
1:l:7:THR:OG1	1:l:112:GLU:OE1	2.29	0.50
1:p:103:GLN:HE22	1:q:84:ARG:HD2	1.76	0.50
1:n:77:ARG:HH12	1:n:108:HIS:HB3	1.76	0.50
1:o:20:PHE:HD2	1:o:58:ILE:HD11	1.76	0.49
1:a:136:ILE:O	1:a:136:ILE:CG2	2.55	0.49
1:m:111:GLU:OE1	1:m:114:ARG:NH2	2.45	0.49
1:p:40:ASP:HB2	1:p:70:TYR:HE1	1.77	0.49
1:i:52:PHE:CE2	1:i:58:ILE:HG22	2.48	0.49
1:d:103:GLN:HE22	1:e:84:ARG:HD2	1.76	0.49
1:q:7:THR:HB	1:q:77:ARG:HG2	1.93	0.49
1:r:7:THR:HG23	1:r:8:VAL:HG23	1.94	0.49
1:m:27:LEU:HD22	1:m:85:LEU:HD22	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:m:40:ASP:HB2	1:m:70:TYR:HE1	1.77	0.49
1:r:8:VAL:HG13	1:r:76:ARG:HB3	1.94	0.49
1:r:20:PHE:HD2	1:r:58:ILE:HD11	1.76	0.49
1:g:96:ALA:O	1:g:100:ASN:ND2	2.46	0.49
1:o:77:ARG:HD2	1:o:79:THR:HG23	1.95	0.49
1:q:11:TYR:OH	1:q:21:ASN:O	2.19	0.49
1:l:20:PHE:HD2	1:l:58:ILE:HD11	1.78	0.48
1:c:7:THR:HG23	1:c:8:VAL:HG23	1.94	0.48
1:i:7:THR:HG23	1:i:8:VAL:HG23	1.95	0.48
1:a:103:GLN:HE22	1:b:84:ARG:HD2	1.79	0.48
1:b:77:ARG:HH12	1:b:108:HIS:HB3	1.77	0.48
1:g:25:GLU:HG3	1:g:108:HIS:HE1	1.78	0.48
1:a:13:LEU:HD12	1:a:73:ILE:HG13	1.96	0.48
1:g:111:GLU:OE1	1:g:114:ARG:NH2	2.46	0.48
1:l:7:THR:HG23	1:l:8:VAL:HG23	1.96	0.48
1:o:7:THR:HG23	1:o:8:VAL:HG23	1.96	0.48
1:d:89:THR:O	1:d:92:SER:OG	2.31	0.48
1:n:35:THR:HG22	1:n:43:VAL:HA	1.96	0.48
1:o:52:PHE:CE1	1:o:58:ILE:HG22	2.49	0.48
1:c:52:PHE:CE2	1:c:58:ILE:HG22	2.49	0.47
1:h:136:ILE:HD12	1:h:137:VAL:H	1.79	0.47
1:n:7:THR:OG1	1:n:8:VAL:N	2.46	0.47
1:l:46:ILE:O	1:l:50:TYR:N	2.47	0.47
1:a:20:PHE:HB2	1:a:58:ILE:HD11	1.95	0.47
1:d:27:LEU:HD22	1:d:85:LEU:HD22	1.96	0.47
1:e:136:ILE:HD12	1:e:137:VAL:H	1.79	0.47
1:m:100:ASN:O	1:m:104:ILE:N	2.38	0.47
1:e:77:ARG:HH12	1:e:108:HIS:HB3	1.79	0.47
1:f:7:THR:HG23	1:f:8:VAL:HG23	1.96	0.47
1:p:20:PHE:HB2	1:p:58:ILE:HD11	1.95	0.47
1:r:52:PHE:CE1	1:r:58:ILE:HG22	2.49	0.47
1:n:121:ILE:HD11	1:o:120:THR:HB	1.97	0.47
1:d:136:ILE:O	1:d:136:ILE:CG2	2.58	0.47
1:f:16:SER:HA	1:f:63:ALA:HA	1.97	0.47
1:j:103:GLN:HG3	1:k:102:ALA:HB1	1.95	0.47
1:k:84:ARG:NH1	1:k:87:ASP:OD1	2.47	0.47
1:o:8:VAL:HG13	1:o:76:ARG:HB3	1.96	0.47
1:r:77:ARG:HD2	1:r:79:THR:HG23	1.97	0.47
1:a:120:THR:HA	1:a:134:ARG:HH12	1.79	0.47
1:d:37:ILE:O	1:d:71:THR:N	2.44	0.47
1:e:49:ASP:HB2	1:e:64:TRP:CH2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:j:27:LEU:HD22	1:j:85:LEU:HD22	1.96	0.47
1:a:136:ILE:CD1	1:b:135:ARG:HH11	2.27	0.46
1:b:136:ILE:HD12	1:b:137:VAL:H	1.81	0.46
1:h:77:ARG:HH22	1:h:108:HIS:HB3	1.80	0.46
1:j:13:LEU:HD12	1:j:73:ILE:HG13	1.96	0.46
1:k:21:ASN:HB3	1:k:57:THR:HG22	1.97	0.46
1:p:120:THR:HA	1:p:134:ARG:HH12	1.80	0.46
1:c:114:ARG:O	1:c:117:THR:OG1	2.33	0.46
1:d:13:LEU:HD12	1:d:73:ILE:HG13	1.98	0.46
1:d:96:ALA:O	1:d:100:ASN:ND2	2.48	0.46
1:j:3:ASN:HA	1:j:4:VAL:HA	1.79	0.46
1:k:35:THR:HG22	1:k:43:VAL:HA	1.98	0.46
1:j:96:ALA:O	1:j:100:ASN:ND2	2.48	0.46
1:p:136:ILE:CD1	1:q:135:ARG:HH11	2.28	0.46
1:d:9:LEU:HD21	1:d:23:PRO:HD2	1.98	0.46
1:i:16:SER:HA	1:i:63:ALA:HA	1.97	0.46
1:m:25:GLU:HG3	1:m:108:HIS:HE1	1.80	0.46
1:b:136:ILE:O	1:b:138:ASN:N	2.49	0.46
1:i:46:ILE:O	1:i:50:TYR:N	2.49	0.46
1:i:58:ILE:HD12	1:i:60:LEU:HA	1.97	0.46
1:a:20:PHE:HE2	1:a:73:ILE:HB	1.81	0.46
1:a:77:ARG:HH21	1:a:112:GLU:HB2	1.81	0.46
1:b:49:ASP:HB2	1:b:64:TRP:CH2	2.51	0.46
1:g:37:ILE:O	1:g:71:THR:N	2.46	0.46
1:g:136:ILE:O	1:g:136:ILE:CG2	2.56	0.46
1:m:20:PHE:HB2	1:m:58:ILE:HD11	1.97	0.46
1:i:8:VAL:HG13	1:i:76:ARG:HB3	1.97	0.46
1:l:52:PHE:CE1	1:l:58:ILE:HG22	2.51	0.46
1:f:58:ILE:HD12	1:f:60:LEU:HA	1.97	0.45
1:a:9:LEU:HD21	1:a:23:PRO:HD2	1.99	0.45
1:c:7:THR:OG1	1:c:112:GLU:OE1	2.28	0.45
1:g:44:LEU:HD12	1:g:49:ASP:OD2	2.17	0.45
1:i:77:ARG:HD2	1:i:79:THR:HG23	1.98	0.45
1:h:18:ARG:HB2	1:h:60:LEU:HB2	1.99	0.45
1:k:121:ILE:HD11	1:l:120:THR:HB	1.99	0.45
1:l:58:ILE:HD12	1:l:60:LEU:HA	1.98	0.45
1:c:16:SER:HA	1:c:63:ALA:HA	1.99	0.45
1:k:77:ARG:HH12	1:k:108:HIS:HB3	1.81	0.45
1:p:100:ASN:O	1:p:104:ILE:N	2.38	0.45
1:h:84:ARG:NH1	1:h:87:ASP:OD1	2.50	0.45
1:j:44:LEU:HD12	1:j:49:ASP:OD2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:n:49:ASP:HB3	1:n:64:TRP:HH2	1.81	0.45
1:a:35:THR:HG22	1:a:43:VAL:HG22	1.98	0.45
1:g:9:LEU:HD21	1:g:23:PRO:HD2	1.99	0.45
1:g:27:LEU:HD22	1:g:85:LEU:HD22	1.99	0.45
1:l:8:VAL:HG13	1:l:76:ARG:HB3	1.98	0.45
1:n:21:ASN:HB3	1:n:57:THR:HG22	1.98	0.45
1:p:20:PHE:HE2	1:p:73:ILE:HB	1.81	0.45
1:r:72:THR:HG22	1:r:73:ILE:HG13	1.98	0.45
1:k:136:ILE:HD12	1:k:137:VAL:H	1.82	0.45
1:r:114:ARG:O	1:r:117:THR:OG1	2.34	0.45
1:g:136:ILE:HD12	1:h:135:ARG:HH12	1.79	0.45
1:j:35:THR:HG22	1:j:43:VAL:HG22	1.99	0.45
1:n:136:ILE:O	1:n:138:ASN:N	2.50	0.45
1:o:45:THR:HG1	1:o:48:THR:HG1	1.59	0.45
1:q:49:ASP:HB2	1:q:64:TRP:CH2	2.51	0.45
1:d:15:GLY:H	1:d:66:PRO:HD3	1.82	0.44
1:g:136:ILE:CD1	1:h:135:ARG:HH11	2.29	0.44
1:k:49:ASP:HB3	1:k:64:TRP:HH2	1.83	0.44
1:g:100:ASN:O	1:g:104:ILE:N	2.43	0.44
1:m:44:LEU:HD12	1:m:49:ASP:OD2	2.17	0.44
1:m:96:ALA:O	1:m:100:ASN:ND2	2.51	0.44
1:a:15:GLY:H	1:a:66:PRO:HD3	1.82	0.44
1:g:13:LEU:HD12	1:g:73:ILE:HG13	1.98	0.44
1:h:36:LEU:HD13	1:h:44:LEU:HD21	2.00	0.44
1:h:136:ILE:O	1:h:138:ASN:N	2.50	0.44
1:q:36:LEU:HD13	1:q:44:LEU:HD21	1.99	0.44
1:d:20:PHE:HB2	1:d:58:ILE:HD11	1.98	0.44
1:f:77:ARG:HH21	1:f:112:GLU:N	2.16	0.44
1:g:35:THR:HG22	1:g:43:VAL:HG22	2.00	0.44
1:k:18:ARG:HB2	1:k:60:LEU:HB2	1.98	0.44
1:q:136:ILE:O	1:q:138:ASN:N	2.50	0.44
1:a:109:VAL:HG21	1:c:107:MET:HE3	1.99	0.44
1:l:77:ARG:HE	1:l:112:GLU:HB2	1.82	0.44
1:e:136:ILE:O	1:e:138:ASN:N	2.49	0.44
1:i:77:ARG:HH21	1:i:112:GLU:N	2.15	0.44
1:h:49:ASP:HB2	1:h:64:TRP:CH2	2.51	0.44
1:i:72:THR:HG22	1:i:73:ILE:HG13	2.00	0.44
1:r:58:ILE:HD12	1:r:60:LEU:HA	1.99	0.44
1:c:72:THR:HG22	1:c:73:ILE:HG13	2.00	0.44
1:d:20:PHE:HE2	1:d:73:ILE:HB	1.83	0.44
1:p:35:THR:OG1	1:p:74:GLU:OE2	2.35	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:d:44:LEU:HD12	1:d:49:ASP:OD2	2.18	0.43
1:c:8:VAL:HG13	1:c:76:ARG:HB3	1.99	0.43
1:h:77:ARG:HH12	1:h:108:HIS:HB3	1.83	0.43
1:k:77:ARG:HH22	1:k:108:HIS:HB3	1.82	0.43
1:m:35:THR:HG22	1:m:43:VAL:HG22	1.99	0.43
1:p:111:GLU:OE1	1:p:114:ARG:NH2	2.51	0.43
1:c:77:ARG:HH21	1:c:112:GLU:N	2.17	0.43
1:g:27:LEU:HD13	1:g:105:GLN:NE2	2.33	0.43
1:h:21:ASN:HB3	1:h:57:THR:HG22	2.00	0.43
1:q:136:ILE:HD12	1:q:137:VAL:H	1.83	0.43
1:d:99:LEU:HD13	1:e:99:LEU:HD21	2.01	0.43
1:j:20:PHE:HB2	1:j:58:ILE:HD11	1.99	0.43
1:c:77:ARG:HD2	1:c:79:THR:HG23	2.01	0.43
1:l:77:ARG:HH21	1:l:112:GLU:N	2.15	0.43
1:q:121:ILE:HD11	1:r:120:THR:HB	2.01	0.43
1:d:100:ASN:O	1:d:104:ILE:HG12	2.18	0.43
1:n:136:ILE:HD12	1:n:137:VAL:H	1.84	0.43
1:d:27:LEU:HD13	1:d:105:GLN:NE2	2.34	0.43
1:f:72:THR:HG22	1:f:73:ILE:HG13	2.01	0.43
1:i:77:ARG:HE	1:i:112:GLU:HB2	1.84	0.43
1:m:20:PHE:HE2	1:m:73:ILE:HB	1.83	0.43
1:g:100:ASN:O	1:g:104:ILE:HG12	2.18	0.43
1:j:100:ASN:O	1:j:104:ILE:HG12	2.19	0.43
1:p:3:ASN:HA	1:p:4:VAL:HA	1.77	0.43
1:p:35:THR:HG22	1:p:43:VAL:HG22	2.00	0.43
1:r:95:ARG:HG3	1:r:98:ASP:H	1.83	0.43
1:a:96:ALA:O	1:a:100:ASN:ND2	2.52	0.42
1:p:136:ILE:O	1:p:136:ILE:CG2	2.54	0.42
1:e:77:ARG:HH22	1:e:108:HIS:HB3	1.84	0.42
1:f:61:THR:OG1	1:f:62:LYS:O	2.37	0.42
1:h:121:ILE:HD11	1:i:120:THR:HB	2.01	0.42
1:o:77:ARG:HH21	1:o:112:GLU:N	2.17	0.42
1:a:35:THR:OG1	1:a:74:GLU:OE2	2.36	0.42
1:a:100:ASN:O	1:a:104:ILE:HG12	2.19	0.42
1:d:35:THR:HG22	1:d:43:VAL:HG22	2.01	0.42
1:e:18:ARG:HB2	1:e:60:LEU:HB2	2.00	0.42
1:j:27:LEU:HD13	1:j:105:GLN:NE2	2.34	0.42
1:o:77:ARG:HE	1:o:112:GLU:HB2	1.84	0.42
1:e:121:ILE:HD11	1:f:120:THR:HB	2.02	0.42
1:r:77:ARG:HH21	1:r:112:GLU:N	2.18	0.42
1:e:84:ARG:NH1	1:e:87:ASP:OD1	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:j:20:PHE:HE2	1:j:73:ILE:HB	1.85	0.42
1:j:109:VAL:HG21	1:l:107:MET:HE3	2.01	0.42
1:g:120:THR:OG1	1:g:121:ILE:N	2.53	0.42
1:f:77:ARG:HD2	1:f:79:THR:HG23	2.01	0.42
1:g:20:PHE:HE2	1:g:73:ILE:HB	1.85	0.42
1:n:84:ARG:NH1	1:n:87:ASP:OD1	2.52	0.42
1:r:7:THR:OG1	1:r:112:GLU:OE1	2.29	0.42
1:g:15:GLY:H	1:g:66:PRO:HD3	1.85	0.42
1:i:61:THR:OG1	1:i:62:LYS:O	2.38	0.42
1:p:22:ILE:HG21	1:p:22:ILE:HD13	1.82	0.42
1:c:12:GLN:HA	1:c:72:THR:HB	2.02	0.42
1:n:18:ARG:HB2	1:n:60:LEU:HB2	2.02	0.42
1:p:15:GLY:H	1:p:66:PRO:HD3	1.85	0.42
1:j:136:ILE:CD1	1:k:135:ARG:HH11	2.28	0.42
1:m:60:LEU:HB3	1:m:64:TRP:CZ3	2.52	0.42
1:r:11:TYR:OH	1:r:21:ASN:O	2.32	0.42
1:b:18:ARG:HB2	1:b:60:LEU:HB2	2.02	0.41
1:j:22:ILE:HD13	1:j:22:ILE:HG21	1.85	0.41
1:q:62:LYS:HB3	1:q:64:TRP:CZ2	2.55	0.41
1:q:21:ASN:HB3	1:q:57:THR:HG22	2.01	0.41
1:a:27:LEU:HD13	1:a:105:GLN:NE2	2.35	0.41
1:b:36:LEU:HD13	1:b:44:LEU:HD21	2.01	0.41
1:j:120:THR:OG1	1:j:121:ILE:N	2.54	0.41
1:k:136:ILE:O	1:k:138:ASN:N	2.51	0.41
1:l:45:THR:HG1	1:l:48:THR:HG1	1.61	0.41
1:m:100:ASN:O	1:m:104:ILE:HG12	2.20	0.41
1:p:40:ASP:HB2	1:p:70:TYR:CE1	2.55	0.41
1:p:60:LEU:HB3	1:p:64:TRP:CZ3	2.52	0.41
1:q:38:GLY:HA2	1:q:70:TYR:HA	2.02	0.41
1:f:13:LEU:HD22	1:f:20:PHE:HZ	1.85	0.41
1:f:114:ARG:O	1:f:117:THR:OG1	2.32	0.41
1:n:38:GLY:HA2	1:n:70:TYR:HA	2.02	0.41
1:r:16:SER:HA	1:r:63:ALA:HA	2.02	0.41
1:a:60:LEU:HB3	1:a:64:TRP:CZ3	2.52	0.41
1:d:100:ASN:O	1:d:104:ILE:N	2.44	0.41
1:j:25:GLU:HG3	1:j:108:HIS:HE1	1.85	0.41
1:p:100:ASN:O	1:p:104:ILE:HG12	2.20	0.41
1:m:15:GLY:H	1:m:66:PRO:HD3	1.86	0.41
1:o:121:ILE:HD12	1:o:121:ILE:HG23	1.89	0.41
1:c:13:LEU:HD22	1:c:20:PHE:HZ	1.86	0.41
1:e:13:LEU:HD23	1:e:73:ILE:H	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:j:60:LEU:HB3	1:j:64:TRP:CZ3	2.53	0.41
1:l:12:GLN:HA	1:l:72:THR:HB	2.02	0.41
1:o:13:LEU:HD22	1:o:20:PHE:HZ	1.85	0.41
1:p:89:THR:O	1:p:92:SER:OG	2.38	0.41
1:f:12:GLN:HA	1:f:72:THR:HB	2.03	0.41
1:l:36:LEU:N	1:l:42:LYS:O	2.54	0.41
1:l:114:ARG:O	1:l:117:THR:OG1	2.32	0.41
1:o:58:ILE:HD12	1:o:60:LEU:HA	2.03	0.41
1:p:77:ARG:HH21	1:p:112:GLU:HB2	1.84	0.41
1:a:3:ASN:HA	1:a:4:VAL:HA	1.77	0.41
1:e:62:LYS:HB3	1:e:64:TRP:CZ2	2.55	0.41
1:f:8:VAL:HG22	1:f:76:ARG:HB2	2.02	0.41
1:f:46:ILE:O	1:f:50:TYR:N	2.54	0.41
1:o:72:THR:HG22	1:o:73:ILE:HG13	2.03	0.41
1:o:101:VAL:HA	1:o:104:ILE:HB	2.02	0.41
1:r:20:PHE:CD2	1:r:58:ILE:HD11	2.55	0.41
1:a:44:LEU:HD12	1:a:49:ASP:OD2	2.20	0.41
1:i:12:GLN:HA	1:i:72:THR:HB	2.03	0.41
1:m:136:ILE:CD1	1:n:135:ARG:HH11	2.32	0.41
1:o:31:PHE:CZ	1:o:85:LEU:HD11	2.56	0.41
1:b:80:SER:OG	1:b:83:ASP:HB3	2.22	0.40
1:f:8:VAL:HG13	1:f:76:ARG:HB3	2.02	0.40
1:g:60:LEU:HB3	1:g:64:TRP:CZ3	2.53	0.40
1:j:40:ASP:HB2	1:j:70:TYR:CE1	2.56	0.40
1:m:120:THR:OG1	1:m:121:ILE:N	2.54	0.40
1:n:11:TYR:HD1	1:n:11:TYR:HA	1.78	0.40
1:d:120:THR:HA	1:d:134:ARG:HH12	1.85	0.40
1:e:21:ASN:HB3	1:e:57:THR:HG22	2.03	0.40
1:g:20:PHE:HB2	1:g:58:ILE:HD11	2.02	0.40
1:m:27:LEU:HD13	1:m:105:GLN:NE2	2.36	0.40
1:o:114:ARG:O	1:o:117:THR:OG1	2.32	0.40
1:p:44:LEU:HD12	1:p:49:ASP:OD2	2.19	0.40
1:p:61:THR:HA	1:p:62:LYS:HA	1.85	0.40
1:d:40:ASP:HB2	1:d:70:TYR:CE1	2.55	0.40
1:q:11:TYR:HD1	1:q:11:TYR:HA	1.79	0.40
1:r:77:ARG:HE	1:r:112:GLU:HB2	1.87	0.40
1:a:89:THR:O	1:a:92:SER:OG	2.39	0.40
1:e:36:LEU:HD13	1:e:44:LEU:HD21	2.02	0.40
1:f:49:ASP:HB3	1:f:61:THR:HG21	2.03	0.40
1:f:77:ARG:HE	1:f:112:GLU:HB2	1.86	0.40
1:i:49:ASP:HB3	1:i:61:THR:HG21	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:c:61:THR:OG1	1:c:62:LYS:O	2.39	0.40
1:e:80:SER:OG	1:e:83:ASP:HB3	2.22	0.40
1:j:24:PHE:CE2	1:j:26:TYR:HB3	2.55	0.40
1:j:104:ILE:HD12	1:j:104:ILE:HG23	1.88	0.40
1:o:61:THR:OG1	1:o:62:LYS:O	2.40	0.40
1:p:27:LEU:HD13	1:p:105:GLN:NE2	2.36	0.40
1:r:13:LEU:HD22	1:r:20:PHE:HZ	1.86	0.40
1:r:101:VAL:HA	1:r:104:ILE:HB	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	134/553 (24%)	115 (86%)	19 (14%)	0	100	100
1	b	127/553 (23%)	111 (87%)	14 (11%)	2 (2%)	7	35
1	c	117/553 (21%)	97 (83%)	16 (14%)	4 (3%)	3	23
1	d	134/553 (24%)	115 (86%)	19 (14%)	0	100	100
1	e	127/553 (23%)	110 (87%)	15 (12%)	2 (2%)	7	35
1	f	117/553 (21%)	96 (82%)	17 (14%)	4 (3%)	3	23
1	g	134/553 (24%)	114 (85%)	20 (15%)	0	100	100
1	h	127/553 (23%)	111 (87%)	14 (11%)	2 (2%)	7	35
1	i	117/553 (21%)	97 (83%)	16 (14%)	4 (3%)	3	23
1	j	134/553 (24%)	115 (86%)	19 (14%)	0	100	100
1	k	127/553 (23%)	112 (88%)	13 (10%)	2 (2%)	7	35
1	l	117/553 (21%)	97 (83%)	16 (14%)	4 (3%)	3	23
1	m	134/553 (24%)	115 (86%)	19 (14%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	n	127/553 (23%)	111 (87%)	14 (11%)	2 (2%)	7	35
1	o	117/553 (21%)	97 (83%)	16 (14%)	4 (3%)	3	23
1	p	134/553 (24%)	115 (86%)	19 (14%)	0	100	100
1	q	127/553 (23%)	110 (87%)	16 (13%)	1 (1%)	16	48
1	r	117/553 (21%)	97 (83%)	16 (14%)	4 (3%)	3	23
All	All	2268/9954 (23%)	1935 (85%)	298 (13%)	35 (2%)	11	36

All (35) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	b	137	VAL
1	c	60	LEU
1	c	121	ILE
1	e	137	VAL
1	f	60	LEU
1	f	121	ILE
1	h	137	VAL
1	i	60	LEU
1	i	121	ILE
1	k	137	VAL
1	l	60	LEU
1	l	121	ILE
1	n	137	VAL
1	o	60	LEU
1	o	121	ILE
1	q	137	VAL
1	r	60	LEU
1	r	121	ILE
1	c	38	GLY
1	f	38	GLY
1	i	38	GLY
1	l	38	GLY
1	o	38	GLY
1	r	38	GLY
1	c	25	GLU
1	l	25	GLU
1	o	25	GLU
1	r	25	GLU
1	f	25	GLU
1	i	25	GLU

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Mol	Chain	Res	Type
1	b	136	ILE
1	e	136	ILE
1	h	136	ILE
1	k	136	ILE
1	n	136	ILE

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	120/451 (27%)	119 (99%)	1 (1%)	73	76
1	b	115/451 (26%)	113 (98%)	2 (2%)	53	67
1	c	105/451 (23%)	103 (98%)	2 (2%)	50	66
1	d	120/451 (27%)	119 (99%)	1 (1%)	73	76
1	e	115/451 (26%)	113 (98%)	2 (2%)	53	67
1	f	105/451 (23%)	103 (98%)	2 (2%)	50	66
1	g	120/451 (27%)	119 (99%)	1 (1%)	73	76
1	h	115/451 (26%)	113 (98%)	2 (2%)	53	67
1	i	105/451 (23%)	103 (98%)	2 (2%)	50	66
1	j	120/451 (27%)	119 (99%)	1 (1%)	73	76
1	k	115/451 (26%)	113 (98%)	2 (2%)	53	67
1	l	105/451 (23%)	104 (99%)	1 (1%)	68	74
1	m	120/451 (27%)	119 (99%)	1 (1%)	73	76
1	n	115/451 (26%)	113 (98%)	2 (2%)	53	67
1	o	105/451 (23%)	104 (99%)	1 (1%)	68	74
1	p	120/451 (27%)	119 (99%)	1 (1%)	73	76
1	q	115/451 (26%)	113 (98%)	2 (2%)	53	67
1	r	105/451 (23%)	104 (99%)	1 (1%)	68	74
All	All	2040/8118 (25%)	2013 (99%)	27 (1%)	59	71

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	a	58	ILE
1	b	9	LEU
1	b	136	ILE
1	c	9	LEU
1	c	58	ILE
1	d	58	ILE
1	e	9	LEU
1	e	136	ILE
1	f	9	LEU
1	f	58	ILE
1	g	58	ILE
1	h	9	LEU
1	h	136	ILE
1	i	9	LEU
1	i	58	ILE
1	j	58	ILE
1	k	9	LEU
1	k	136	ILE
1	l	9	LEU
1	m	58	ILE
1	n	9	LEU
1	n	136	ILE
1	o	9	LEU
1	p	58	ILE
1	q	9	LEU
1	q	136	ILE
1	r	9	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	a	103	GLN
1	a	105	GLN
1	a	108	HIS
1	b	108	HIS
1	d	103	GLN
1	d	105	GLN
1	d	108	HIS
1	e	108	HIS
1	g	103	GLN
1	g	105	GLN

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Mol	Chain	Res	Type
1	h	108	HIS
1	j	103	GLN
1	j	105	GLN
1	k	108	HIS
1	m	103	GLN
1	m	105	GLN
1	m	108	HIS
1	n	108	HIS
1	p	103	GLN
1	p	105	GLN
1	p	108	HIS
1	q	108	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

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.

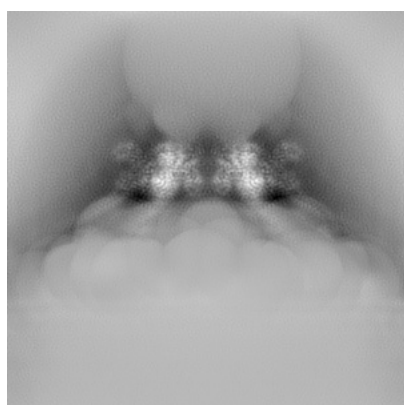
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-30137. These allow visual inspection of the internal detail of the map and identification of artifacts.

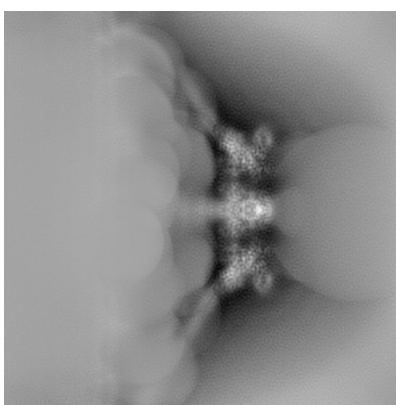
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

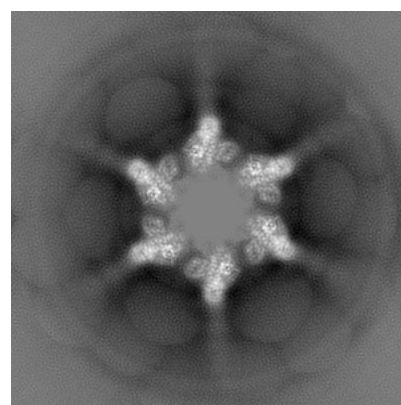
6.1.1 Primary map



X



Y

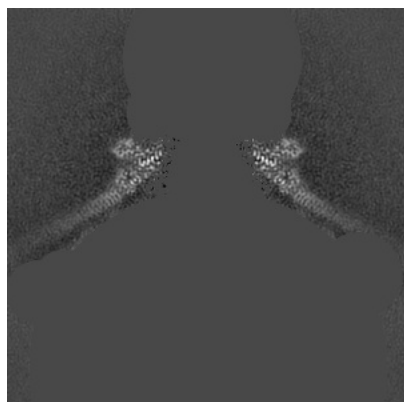


Z

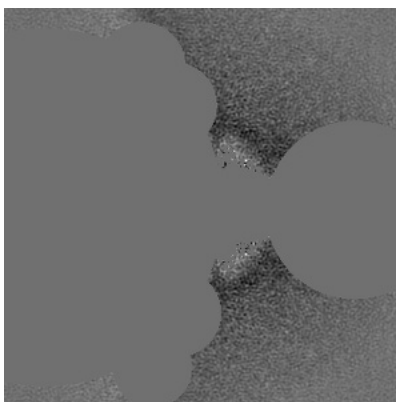
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

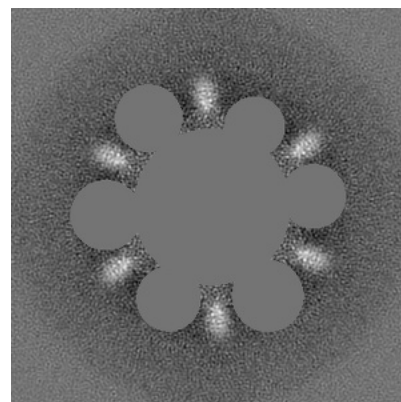
6.2.1 Primary map



X Index: 176



Y Index: 176

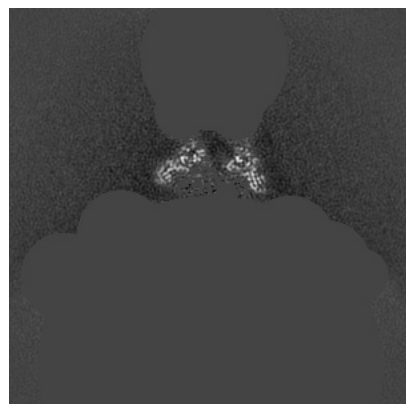


Z Index: 176

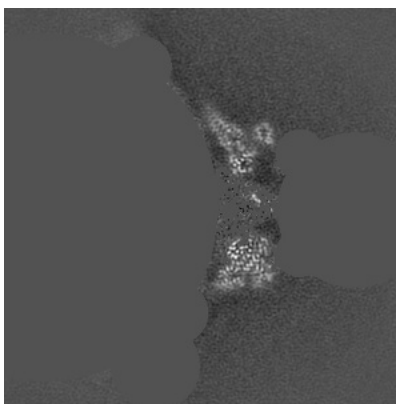
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

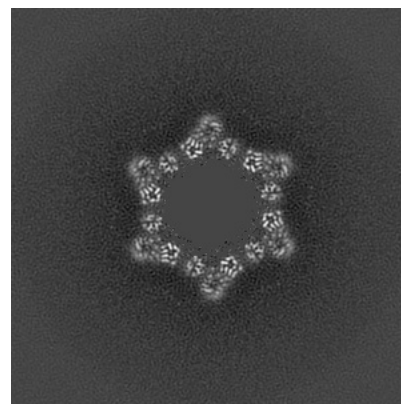
6.3.1 Primary map



X Index: 218



Y Index: 141

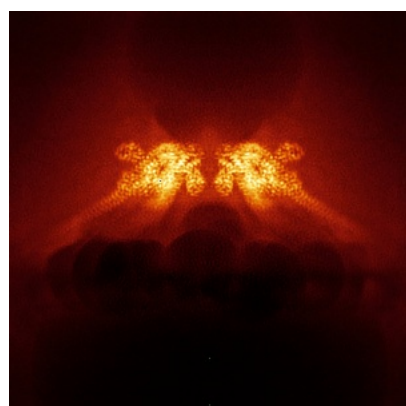


Z Index: 202

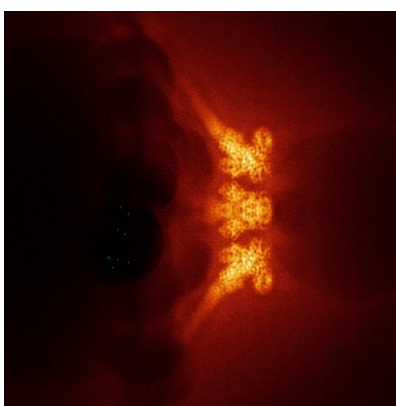
The images above show the largest variance slices of the map in three orthogonal directions.

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

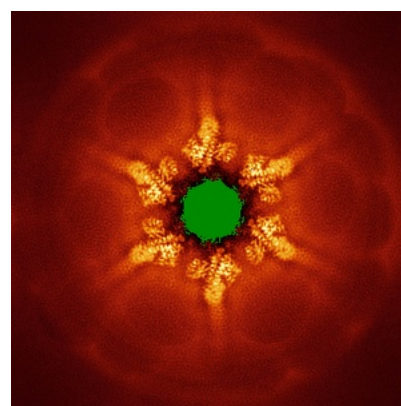
6.4.1 Primary map



X



Y

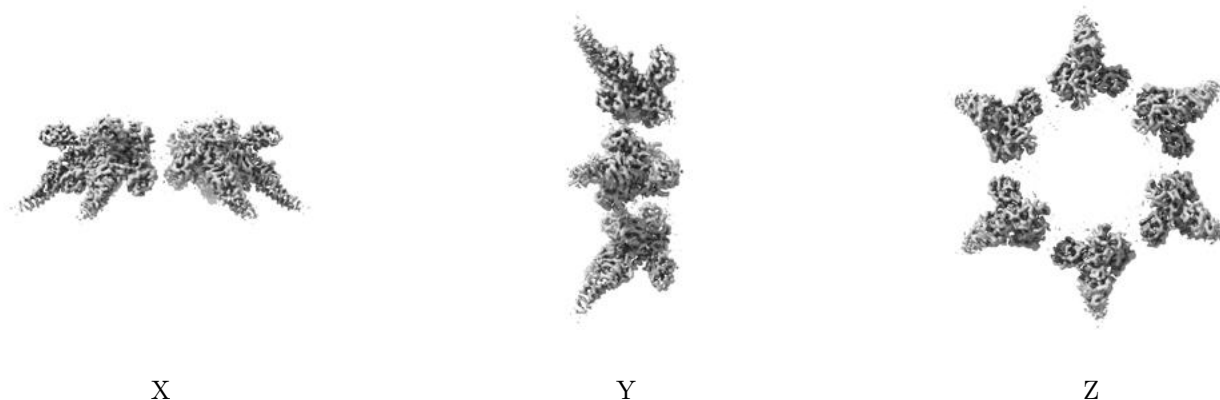


Z

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 [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 7.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

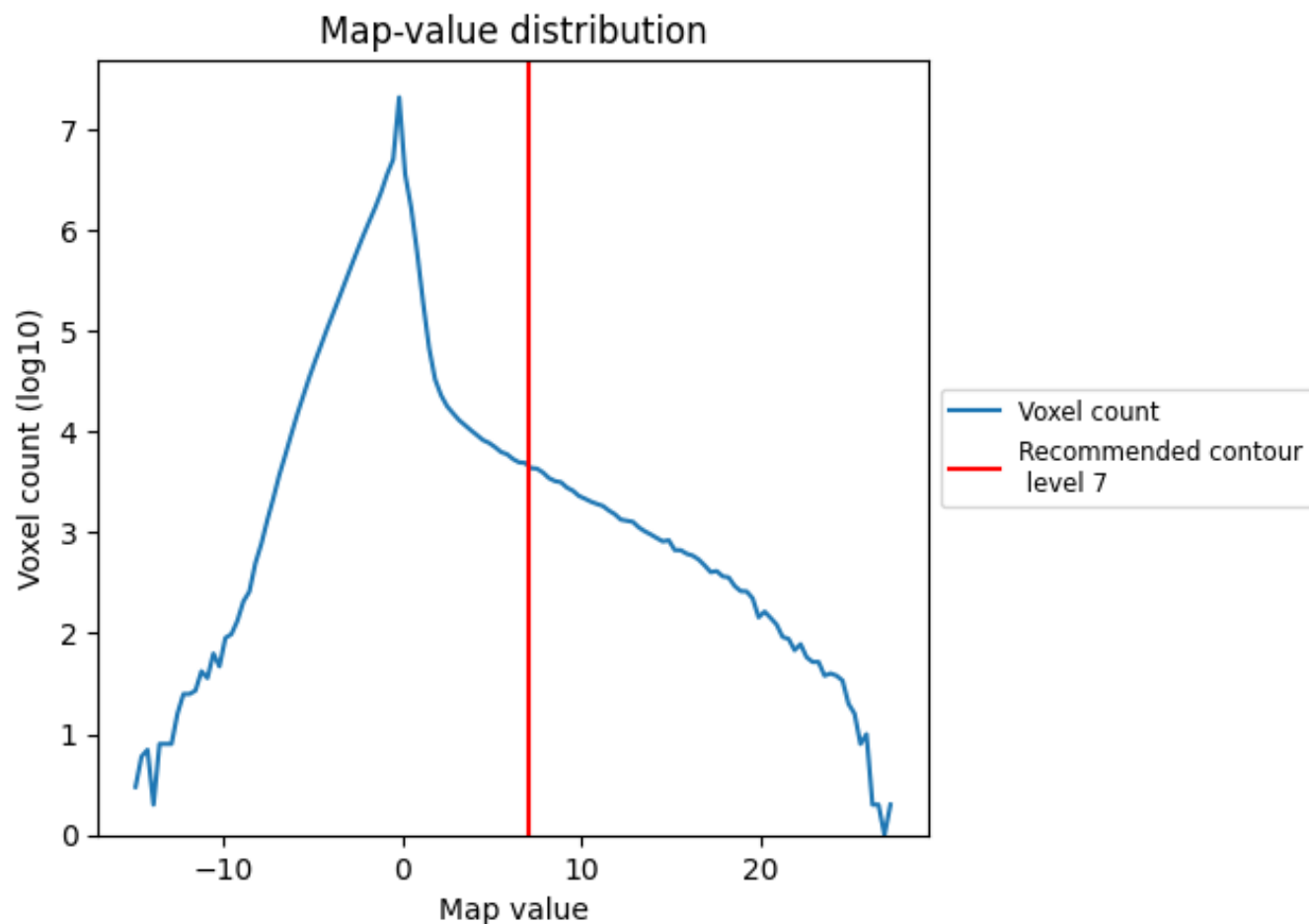
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

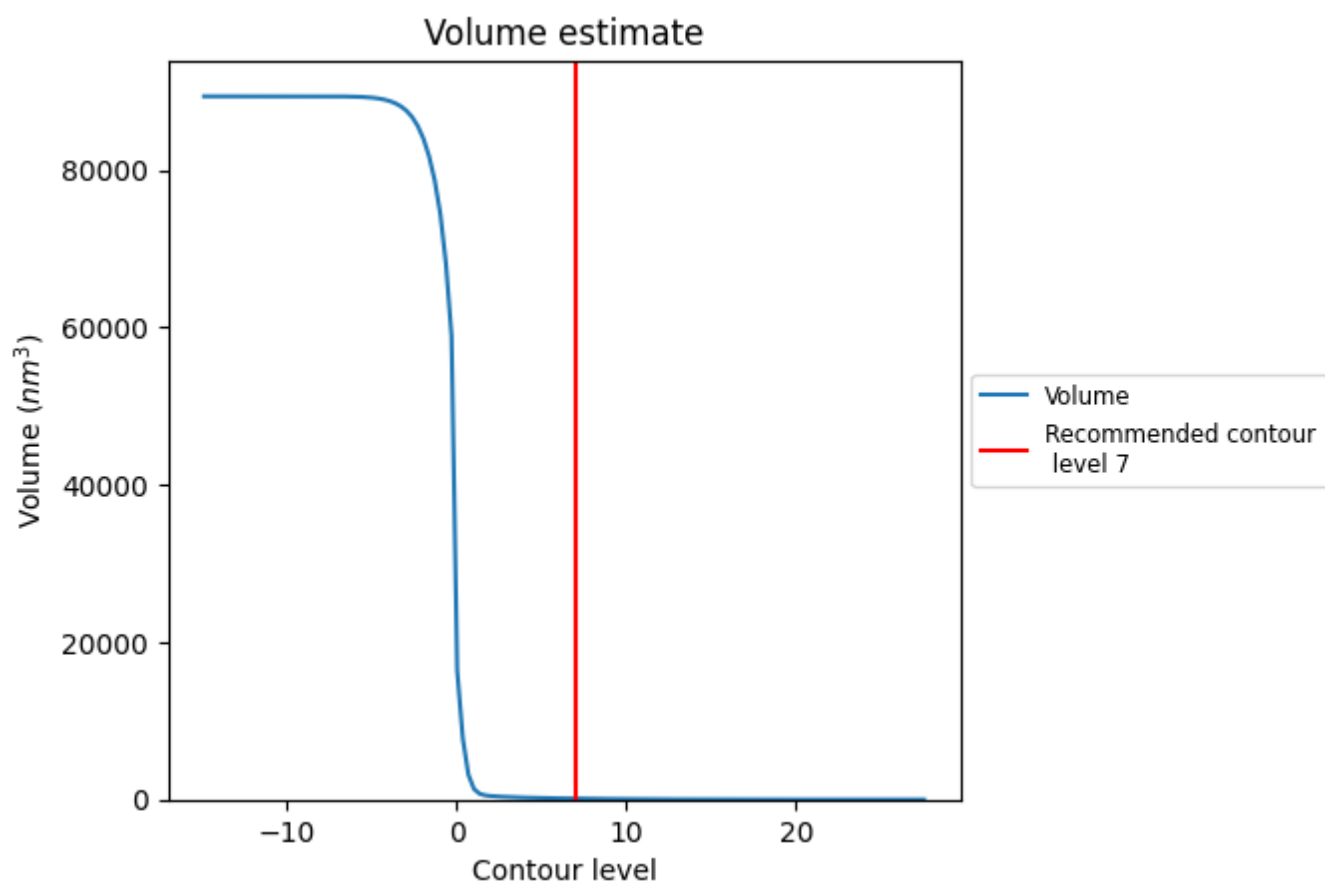
This section contains the results of statistical analysis of the map.

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.

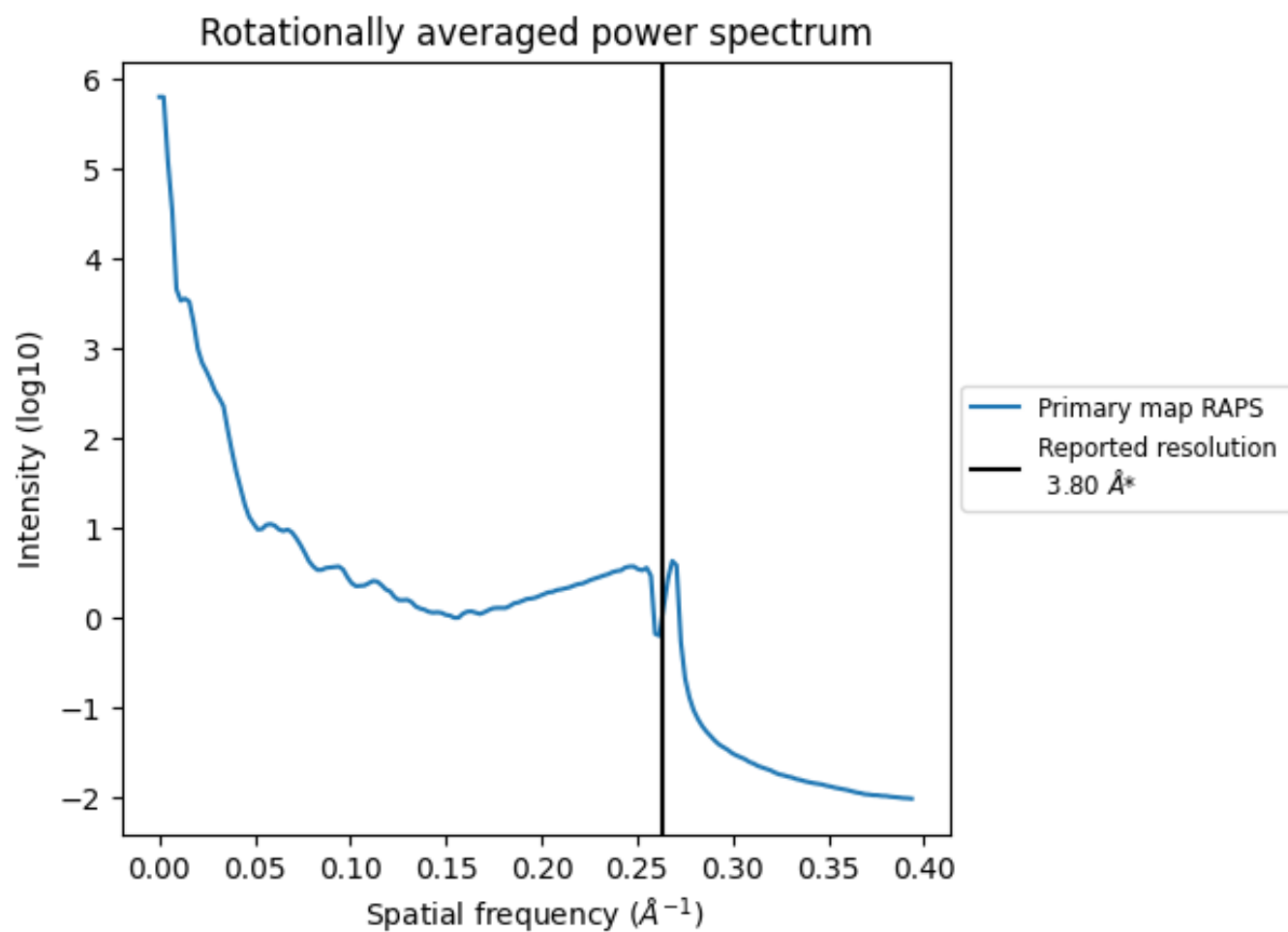
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 124 nm³; this corresponds to an approximate mass of 112 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 [i](#)



*Reported resolution corresponds to spatial frequency of 0.263 Å⁻¹

8 Fourier-Shell correlation ⓘ

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

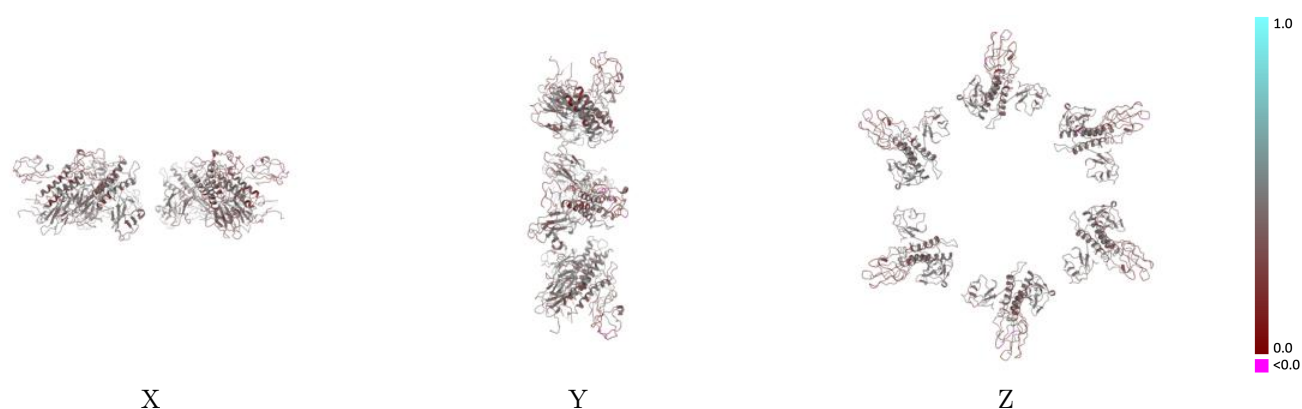
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-30137 and PDB model 7BOZ. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlay [i](#)

This section was not generated.

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

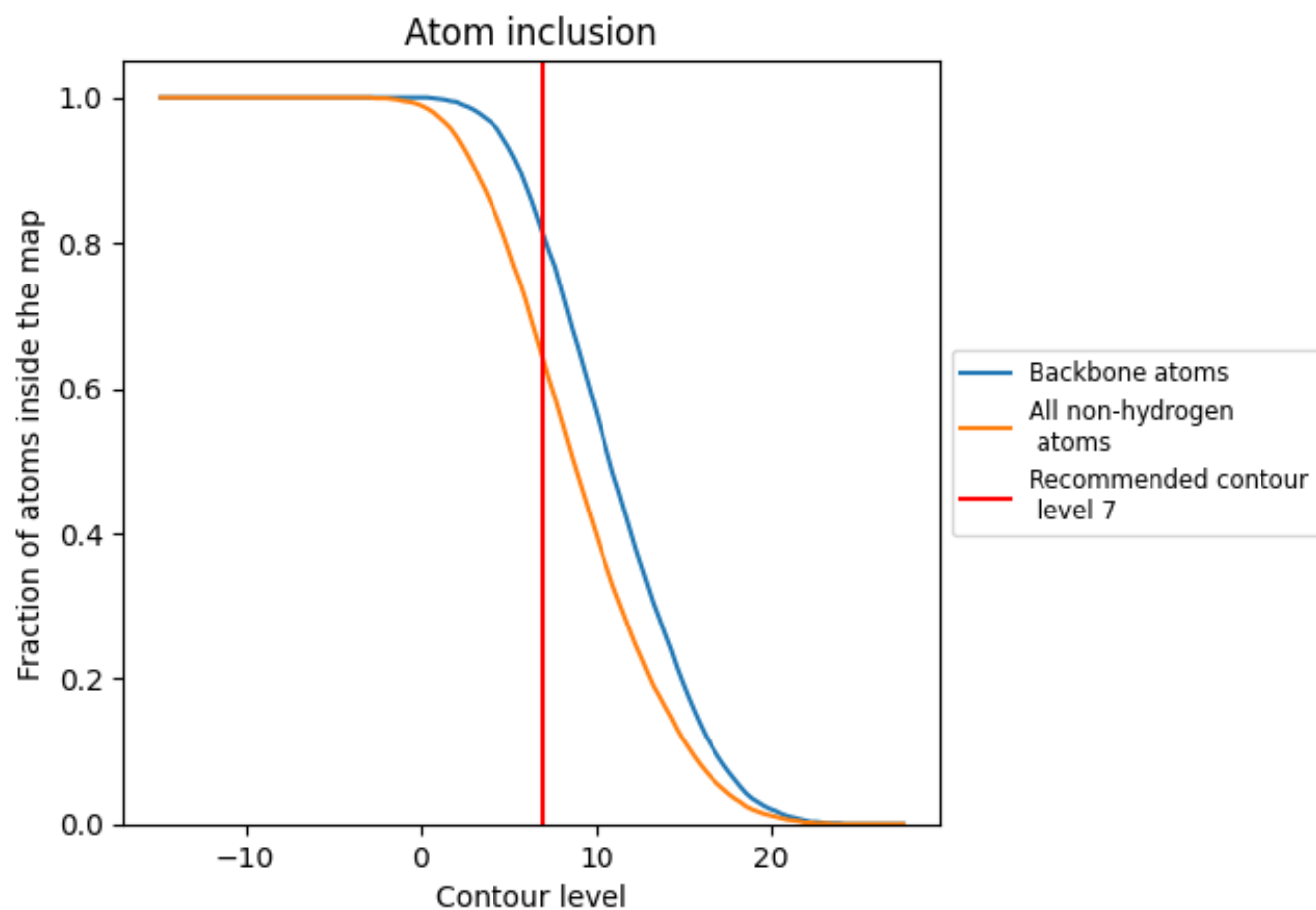


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 81% of all backbone atoms, 64% 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 (7) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6400	 0.3930
a	 0.6930	 0.4230
b	 0.6500	 0.4190
c	 0.5730	 0.3560
d	 0.6990	 0.4230
e	 0.6530	 0.4120
f	 0.5610	 0.3450
g	 0.7170	 0.4320
h	 0.6550	 0.4210
i	 0.5630	 0.3510
j	 0.7120	 0.4370
k	 0.6480	 0.4150
l	 0.5710	 0.3580
m	 0.6930	 0.4080
n	 0.6340	 0.3900
o	 0.5520	 0.3130
p	 0.6890	 0.4010
q	 0.6390	 0.3990
r	 0.5540	 0.3260

