



wwPDB EM Validation Summary Report ⓘ

Mar 9, 2026 – 09:17 AM UTC

PDB ID : 7C4J / pdb_00007c4j
EMDB ID : EMD-30285
Title : Cryo-EM structure of the yeast Swi/Snf complex in a nucleosome free state
Authors : Wang, C.C.; Guo, Z.Y.; Zhan, X.C.; Zhang, X.F.
Deposited on : 2020-05-18
Resolution : 2.89 Å(reported)

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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

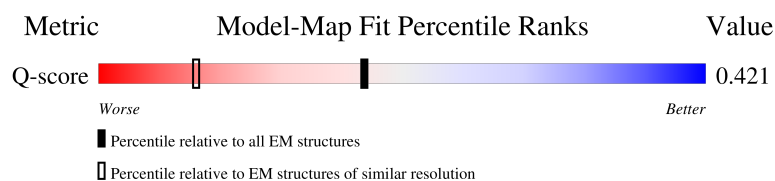
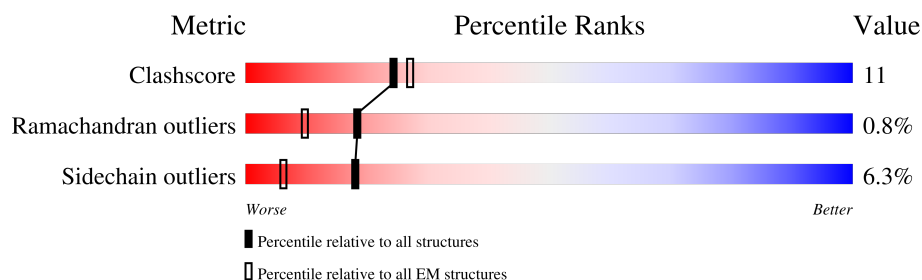
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.89 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.



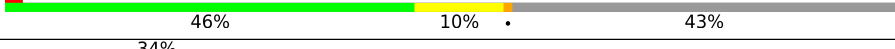
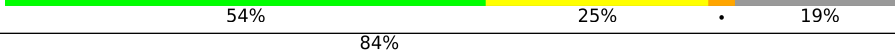
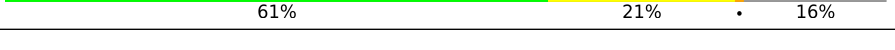
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	12148 (2.39 - 3.39)

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	566	
2	B	825	
2	D	825	
3	C	332	

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Mol	Chain	Length	Quality of chain
4	E	623	
5	F	905	
6	H	1703	
7	I	1314	
8	G	82	
9	J	157	
10	K	477	
11	L	467	

2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 29757 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 Transcription regulatory protein SNF12.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	374	Total	C	N	O	S	0	0
			3048	1928	529	581	10		

- Molecule 2 is a protein called SWI/SNF complex subunit SWI3.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	395	Total	C	N	O	S	0	0
			3250	2082	560	597	11		
2	D	421	Total	C	N	O	S	0	0
			3445	2203	592	638	12		

- Molecule 3 is a protein called Transcription regulatory protein SNF6.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	154	Total	C	N	O	S	0	0
			1266	793	232	236	5		

- Molecule 4 is a protein called SWI/SNF global transcription activator complex subunit SWP82.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	E	288	Total	C	N	O	S	0	0
			2355	1529	412	411	3		

- Molecule 5 is a protein called SWI/SNF chromatin-remodeling complex subunit SNF5.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	F	383	Total	C	N	O	S	0	0
			3134	1977	529	618	10		

- Molecule 6 is a protein called Transcription regulatory protein SNF2.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	H	242	Total	C	N	O	S	0	0
			1966	1233	354	373	6		

- Molecule 7 is a protein called SWI/SNF chromatin-remodeling complex subunit SWI1.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	I	524	Total	C	N	O	S	0	0
			4242	2754	703	774	11		

- Molecule 8 is a protein called Unkown.

Mol	Chain	Residues	Atoms				AltConf	Trace
8	G	47	Total	C	N	O	0	0
			271	166	56	49		

- Molecule 9 is a protein called Regulator of Ty1 transposition protein 102.

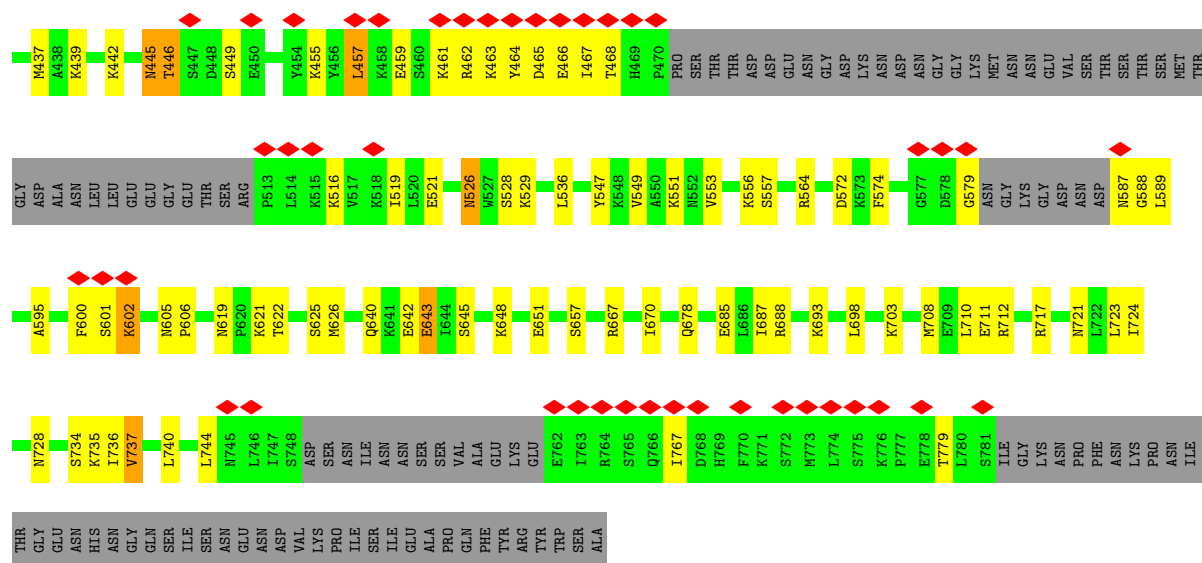
Mol	Chain	Residues	Atoms				AltConf	Trace
9	J	53	Total	C	N	O	0	0
			482	308	83	91		

- Molecule 10 is a protein called Actin-related protein 7.

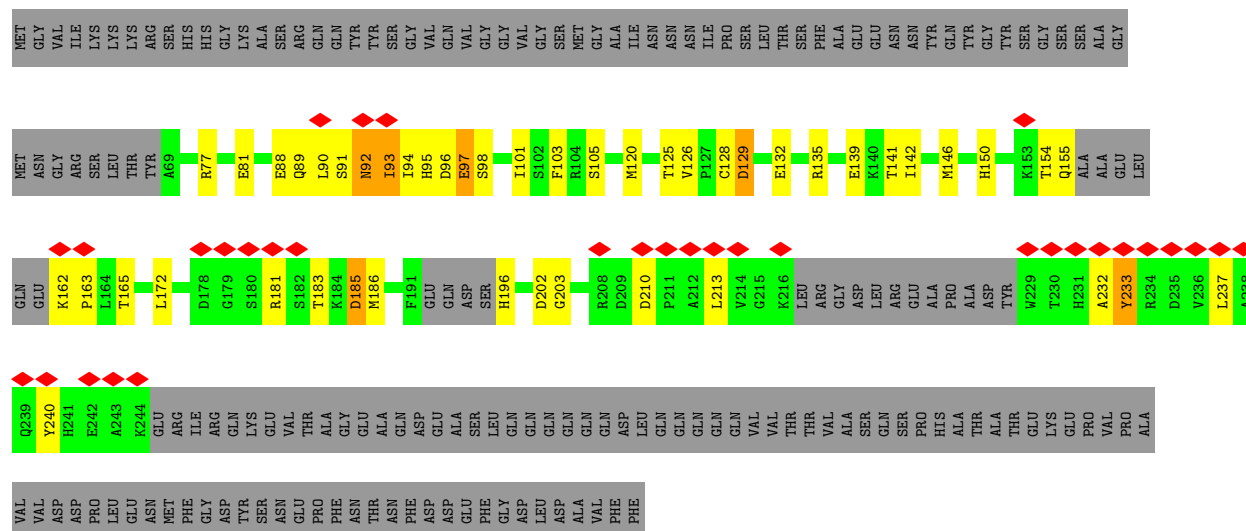
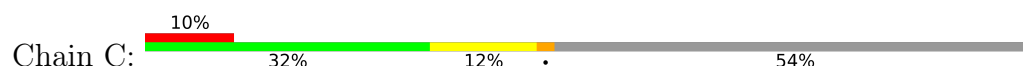
Mol	Chain	Residues	Atoms					AltConf	Trace
10	K	387	Total	C	N	O	S	3	0
			3131	2020	516	591	4		

- Molecule 11 is a protein called Actin-like protein ARP9.

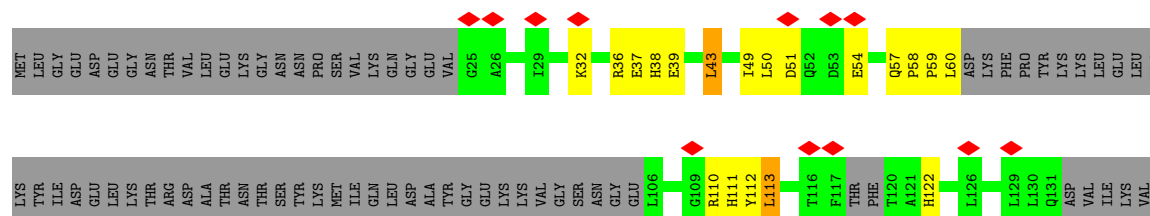
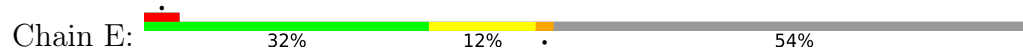
Mol	Chain	Residues	Atoms					AltConf	Trace
11	L	392	Total	C	N	O	S	1	0
			3167	2033	519	611	4		



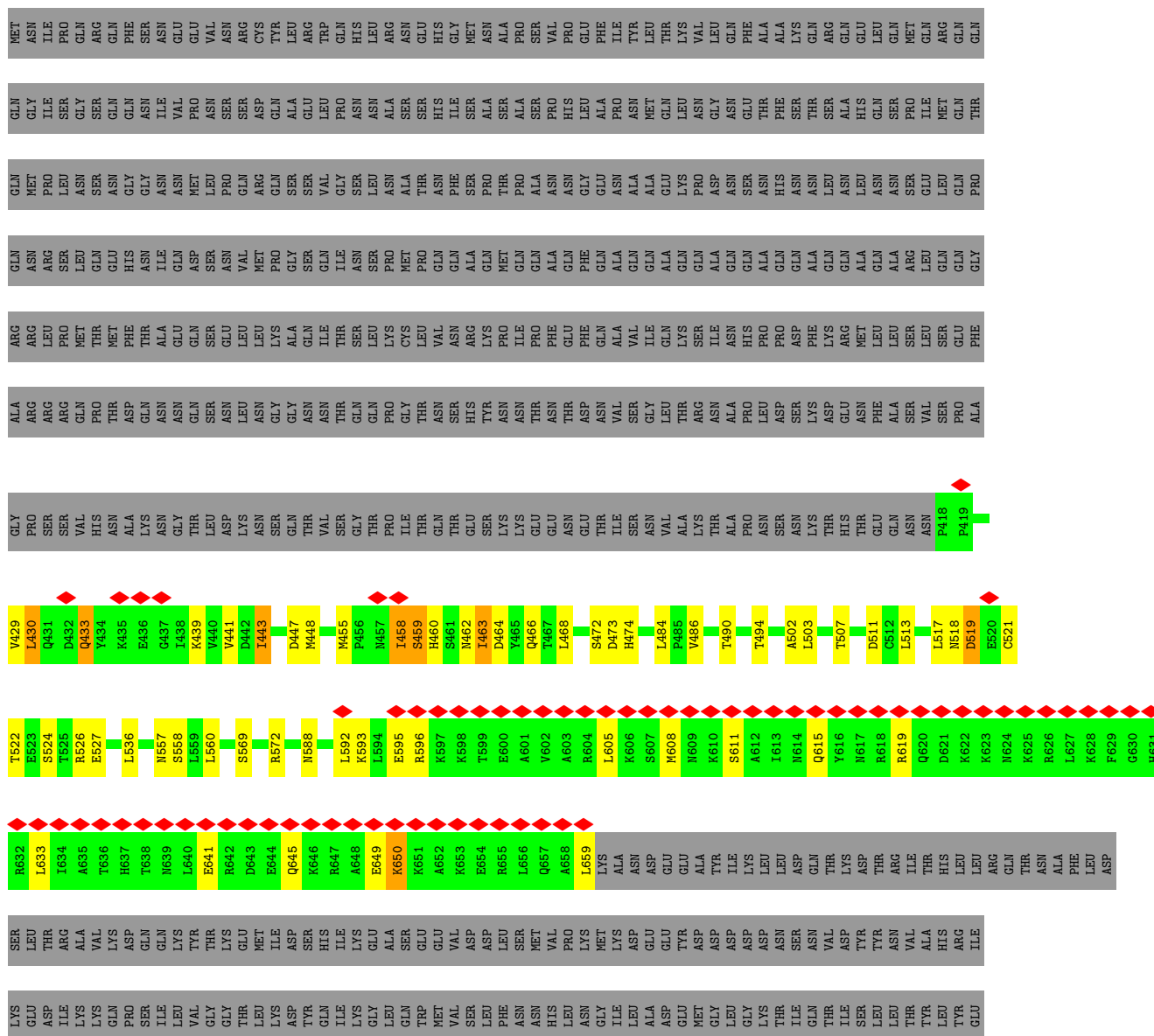
• Molecule 3: Transcription regulatory protein SNF6



• Molecule 4: SWI/SNF global transcription activator complex subunit SWP82



- Molecule 6: Transcription regulatory protein SNF2

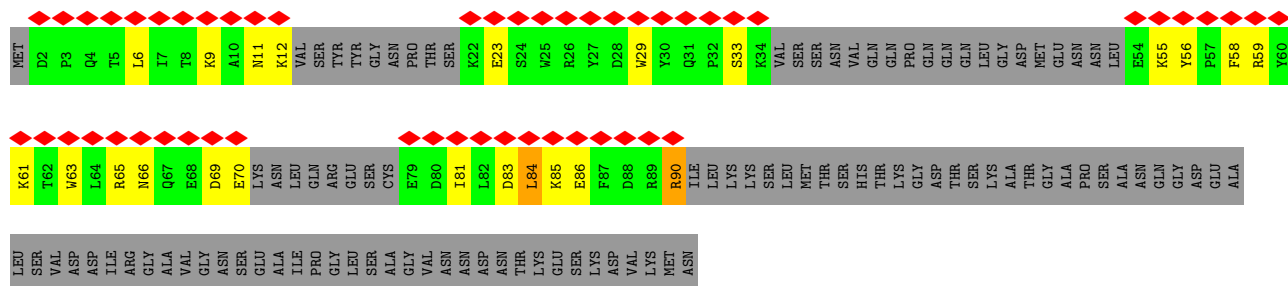


TYR	CYS	GLU	TYR	VAL	ASN	ARG	GLU	GLU	ASP	ARG	GLU	GLY	ASP	GLY	GLN	THR	MET	THR	GLY	ARG	LEU	PRO	THR	MET
	ILE	GLU	TYR	THR	ASP	GLY	ARG	MET	LYS	ARG	GLY	LYS	GLY	GLY	GLN	GLU	THR	GLY	GLY	ARG	GLN	GLY	LYS	
	PRO	MET	VAL	THR	PRO	ARG	THR	VAL	ILE	THR	VAL	ILE	ILE	VAL	ASP	ASP	ILE	ILE	LYS	VAL	LEU	ILE	ILE	
	ALA	GLY	ALA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

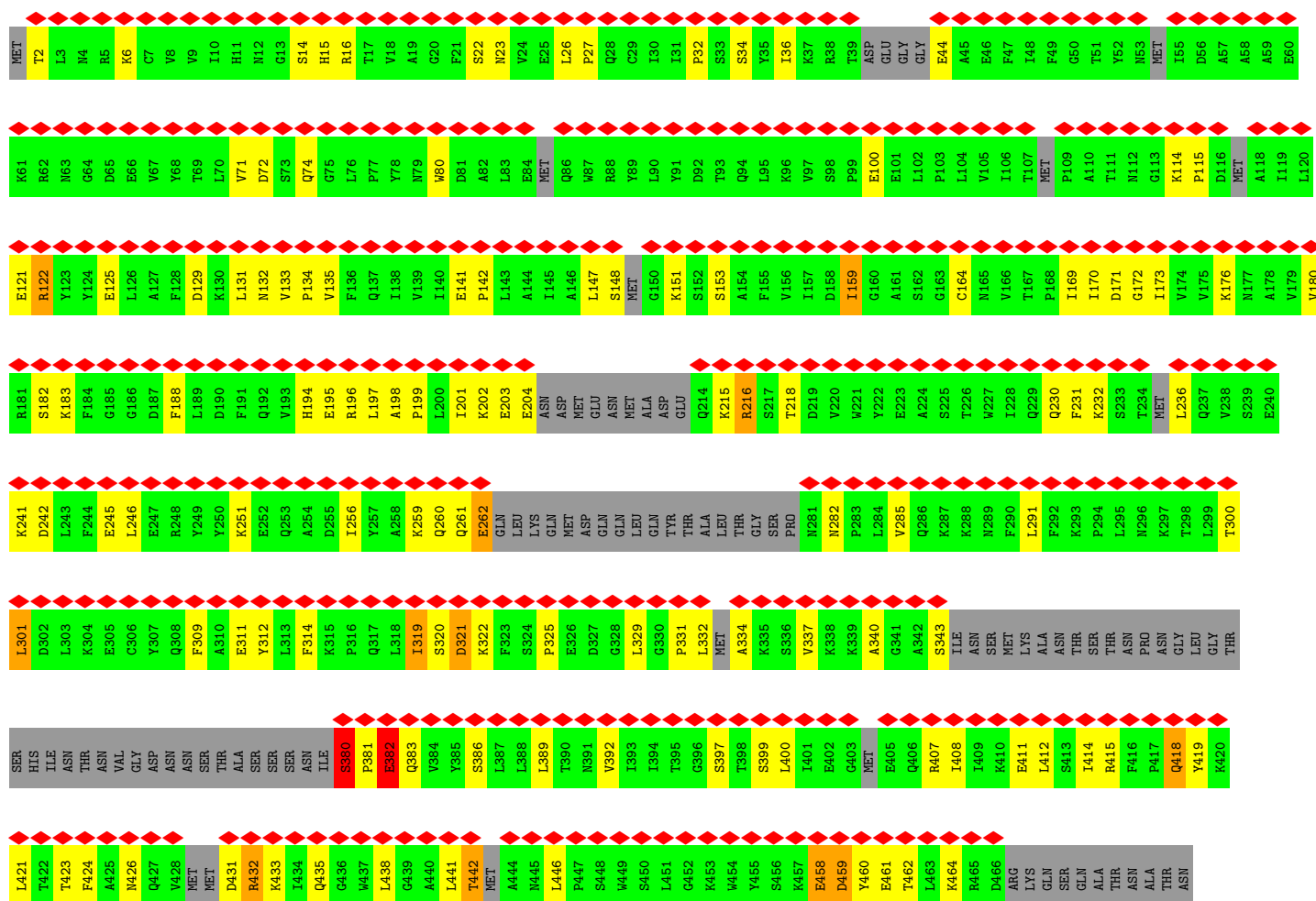
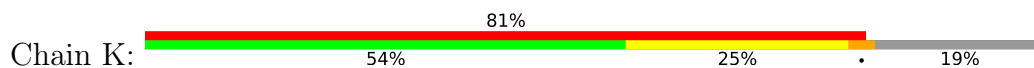
- Molecule 7: SWI/SNF chromatin-remodeling complex subunit SWI1

Chain I: 32% 7% 60%

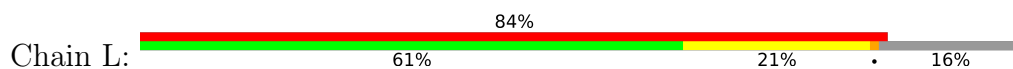
[illegible]



• Molecule 10: Actin-related protein 7



• Molecule 11: Actin-like protein ARP9



K421	S422	G423	Y424	S425	E426	I427	I428	F429	L430	G431	A432	Q433	I434	V435	S436	K437	Q438	I439	F440	T441	H442	P443	K444	D445	T446	F447	Y448	I449	T450	R451	E452	K453	Y454	M455	MET	K457	G458	P459	A460	A461	L462	V463	D464	V465	Q466	F467													
E361	P362	E363	E364	E365	K366	S367	K368	R369	E370	E371	E372	A373	K374	S375	VAL	LEU	PRO	ALA	ALA	THR	LYS	LYS	SER	LYS	PHE	MET	THR	ASN	SER	THR	ALA	V394	V395	P396	T397	I398	E399	Y400	V401	Q402	C403	P404	T405	V406	I407	K408	L409	A410	K411	Y412	P413	D414	Y415	F416	P417	E418	W419	K420	
G301	C302	N303	N304	L305	I306	K307	N308	I309	S310	N311	R312	V313	G314	L315	T316	L317	D318	N319	I320	D321	D322	I323	N324	K325	A326	K327	A328	V329	W330	E331	N332	I333	I334	I335	V336	G337	G338	T339	T340	S341	I342	S343	G344	F345	K346	E347	A348	L349	L350	G351	Q352	L353	L354	K355	D356	H357	L358	I359	I360
GLU	ASP	GLU	GLY	THR	LEU	ASN	VAL	ALA	GLU	ILE	ILE	THR	SER	GLY	ARG	ASP	THR	ARG	GLU	VAL	LEU	GLU	GLY	GLN	LYS	VAL	LYS	ASN	VAL	K275	N276	S277	D278	L279	E280	F281	N282	T283	F284	W285	D286	E287	K288	G289	N290	E291	I292	K293	V294	G295	K296	Q297	R298	F299	Q300				
D181	H182	L183	V184	S185	S186	I187	P188	MET	G190	G191	Q192	S193	I194	N195	D196	S197	L198	K199	K200	L201	L202	P203	Q204	W205	D206	D207	D208	Q209	I210	E211	S212	L213	K214	K215	S216	P217	I218	F219	E220	V221	L222	S223	ASP	ASP	ALA	LYS	LYS	LEU	SER	SER	PHE	ASP	PHE	GLY	ASN	GLU	ASN	GLU	ASP
Q121	S122	D123	L124	E125	I126	I127	T128	Q129	Y130	V131	F132	E133	I134	L135	E136	I137	N138	N139	L140	I141	Q142	L143	P144	A145	S146	L147	A148	A149	T150	Y151	S152	MET	I154	S155	L156	Q157	N158	C159	C160	I161	I162	D163	V164	G165	T166	H167	H168	T169	D170	I171	I172	P173	I174	V175	D176	Y177	A178	Q179	L180
K61	A62	E63	L64	I65	K66	P67	I68	Q69	N70	G71	E72	I73	I74	D75	I76	S77	A78	F79	T80	Q81	F82	L83	R84	L85	I86	F87	V88	S89	I90	L91	S92	D93	R94	A95	N96	K97	N98	Q99	D100	A101	F102	E103	A104	E105	L106	S107	N108	I109	P110	L111	L112	L113	I114	T115	H116	H117	S118	W119	S120

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	386469	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.094	Depositor
Minimum map value	-0.042	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.02	Depositor
Map size ($\text{\AA}$)	347.84, 347.84, 347.84	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.087, 1.087, 1.087	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.30	0/3104	0.47	0/4188
2	B	0.33	0/3318	0.49	0/4465
2	D	0.32	0/3517	0.46	0/4735
3	C	0.34	0/1287	0.56	1/1729 (0.1%)
4	E	0.35	0/2404	0.56	0/3237
5	F	0.35	0/3201	0.55	3/4336 (0.1%)
6	H	0.43	0/2000	0.67	0/2706
7	I	0.40	0/4316	0.54	2/5843 (0.0%)
8	G	0.25	0/101	0.60	0/123
9	J	0.45	0/493	0.83	0/659
10	K	0.55	1/3197 (0.0%)	0.88	4/4313 (0.1%)
11	L	0.59	0/3234	0.89	2/4382 (0.0%)
All	All	0.41	1/30172 (0.0%)	0.63	12/40716 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	K	380	SER	C-N	5.29	1.40	1.34

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	468	ARG	N-CA-C	-6.59	100.91	108.49
5	F	357	THR	CA-C-N	6.54	134.02	121.54
5	F	357	THR	C-N-CA	6.54	134.02	121.54
10	K	380	SER	CA-C-N	-5.65	113.10	118.97
10	K	380	SER	C-N-CA	-5.65	113.10	118.97

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	3048	0	3012	60	0
2	B	3250	0	3302	78	0
2	D	3445	0	3505	86	0
3	C	1266	0	1264	38	0
4	E	2355	0	2416	69	0
5	F	3134	0	3045	56	0
6	H	1966	0	1991	64	0
7	I	4242	0	4401	74	0
8	G	271	0	129	10	0
9	J	482	0	455	19	0
10	K	3131	0	3132	112	0
11	L	3167	0	3149	90	0
All	All	29757	0	29801	637	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.

The worst 5 of 637 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)
7:I:1051:LYS:HB3	7:I:1052:PRO:CD	1.70	1.19
11:L:403:CYS:SG	11:L:404:PRO:HD3	1.89	1.11
2:B:437:MET:HE3	6:H:502:ALA:HB1	1.33	1.08
7:I:1051:LYS:HE2	7:I:1051:LYS:HA	1.17	1.08
10:K:216:ARG:HG2	10:K:216:ARG:HH11	1.17	1.07

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	360/566 (64%)	344 (96%)	15 (4%)	1 (0%)	36	65
2	B	385/825 (47%)	361 (94%)	21 (6%)	3 (1%)	16	44
2	D	413/825 (50%)	383 (93%)	27 (6%)	3 (1%)	18	48
3	C	146/332 (44%)	134 (92%)	10 (7%)	2 (1%)	9	30
4	E	272/623 (44%)	245 (90%)	24 (9%)	3 (1%)	11	36
5	F	379/905 (42%)	333 (88%)	43 (11%)	3 (1%)	16	44
6	H	240/1703 (14%)	230 (96%)	9 (4%)	1 (0%)	30	59
7	I	506/1314 (38%)	486 (96%)	15 (3%)	5 (1%)	12	39
8	G	13/82 (16%)	12 (92%)	1 (8%)	0	100	100
9	J	45/157 (29%)	42 (93%)	2 (4%)	1 (2%)	5	20
10	K	360/477 (76%)	340 (94%)	17 (5%)	3 (1%)	16	44
11	L	381/467 (82%)	360 (94%)	19 (5%)	2 (0%)	24	54
All	All	3500/8276 (42%)	3270 (93%)	203 (6%)	27 (1%)	18	44

5 of 27 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	431	SER
2	B	432	VAL
4	E	208	ASP
5	F	358	ASN
7	I	867	PRO

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	345/517 (67%)	323 (94%)	22 (6%)	16	44
2	B	368/751 (49%)	343 (93%)	25 (7%)	14	42

Continued on next page...

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	392/751 (52%)	372 (95%)	20 (5%)	21	53
3	C	141/288 (49%)	131 (93%)	10 (7%)	13	40
4	E	257/558 (46%)	235 (91%)	22 (9%)	10	30
5	F	351/823 (43%)	335 (95%)	16 (5%)	24	57
6	H	222/1520 (15%)	204 (92%)	18 (8%)	11	33
7	I	500/1218 (41%)	471 (94%)	29 (6%)	18	49
8	G	7/15 (47%)	6 (86%)	1 (14%)	3	11
9	J	52/140 (37%)	50 (96%)	2 (4%)	29	64
10	K	345/420 (82%)	320 (93%)	25 (7%)	13	39
11	L	359/423 (85%)	338 (94%)	21 (6%)	18	49
All	All	3339/7424 (45%)	3128 (94%)	211 (6%)	18	45

5 of 211 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
6	H	429	VAL
7	I	932	ASP
11	L	94	ARG
6	H	447	ASP
6	H	649	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 74 such sidechains are listed below:

Mol	Chain	Res	Type
7	I	915	ASN
11	L	303	ASN
7	I	981	ASN
7	I	1284	GLN
2	D	587	ASN

5.3.3 RNA ⓘ

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.

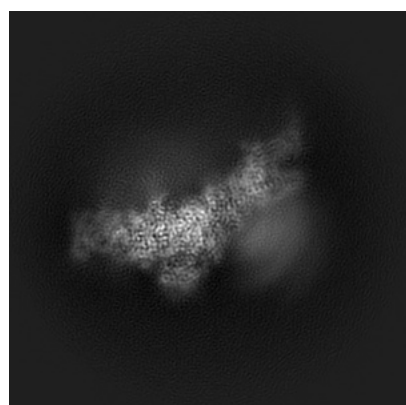
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-30285. These allow visual inspection of the internal detail of the map and identification of artifacts.

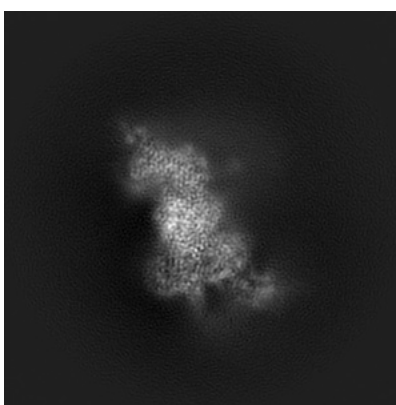
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

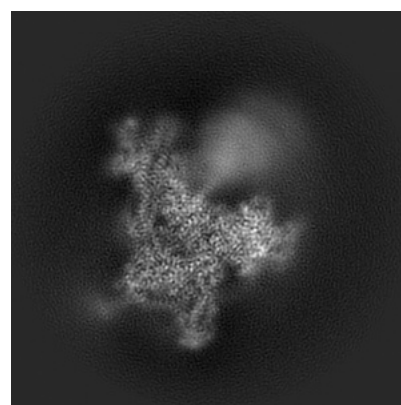
6.1.1 Primary map



X



Y

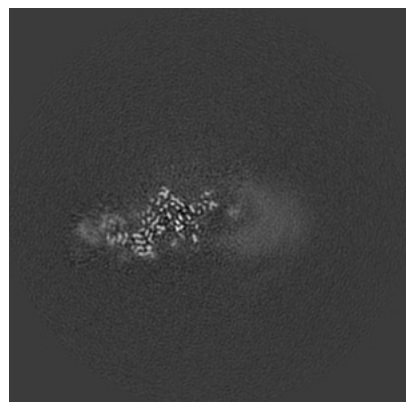


Z

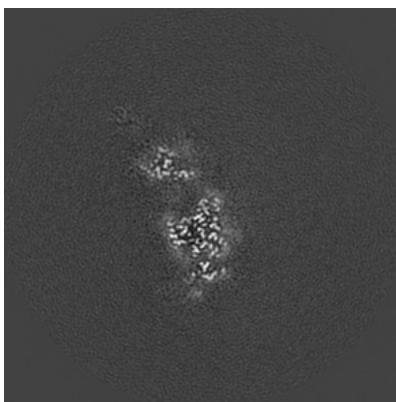
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

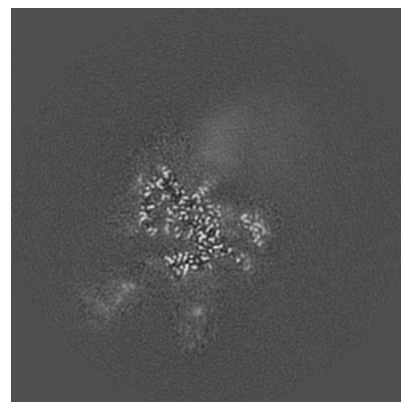
6.2.1 Primary map



X Index: 160



Y Index: 160

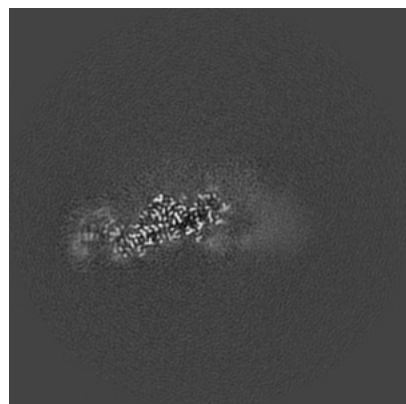


Z Index: 160

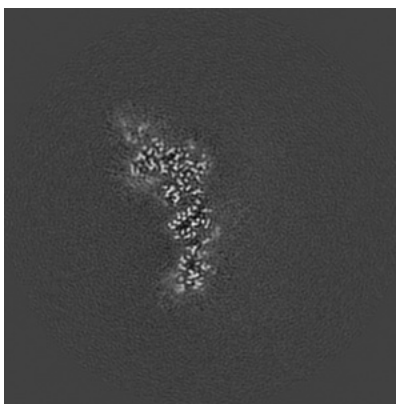
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

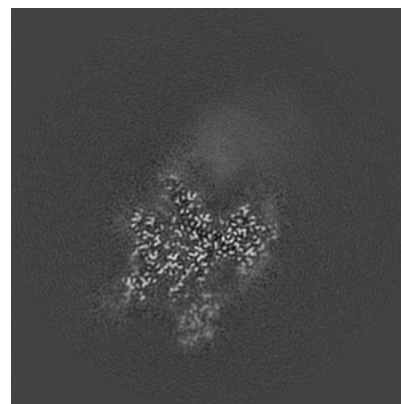
6.3.1 Primary map



X Index: 152



Y Index: 143

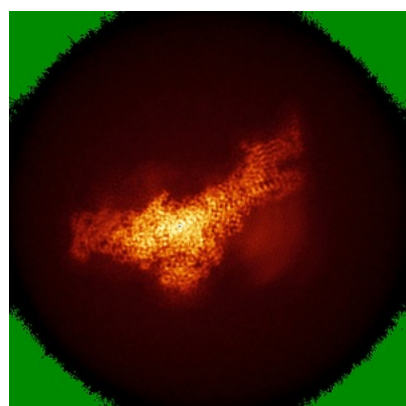


Z Index: 146

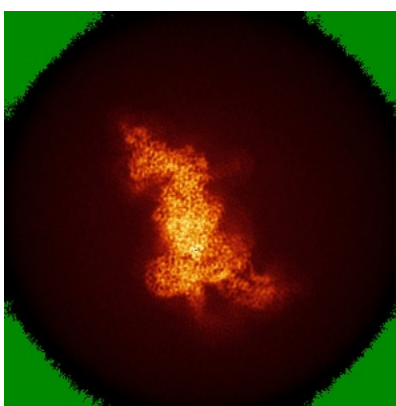
The images above show the largest variance slices of the map in three orthogonal directions.

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

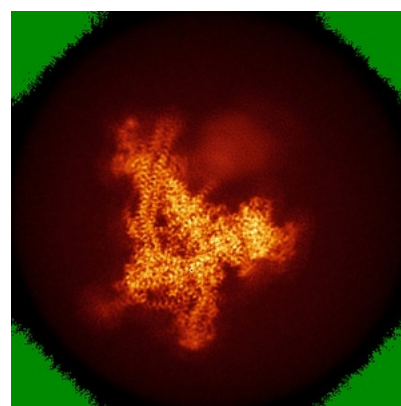
6.4.1 Primary map



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 0.02. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

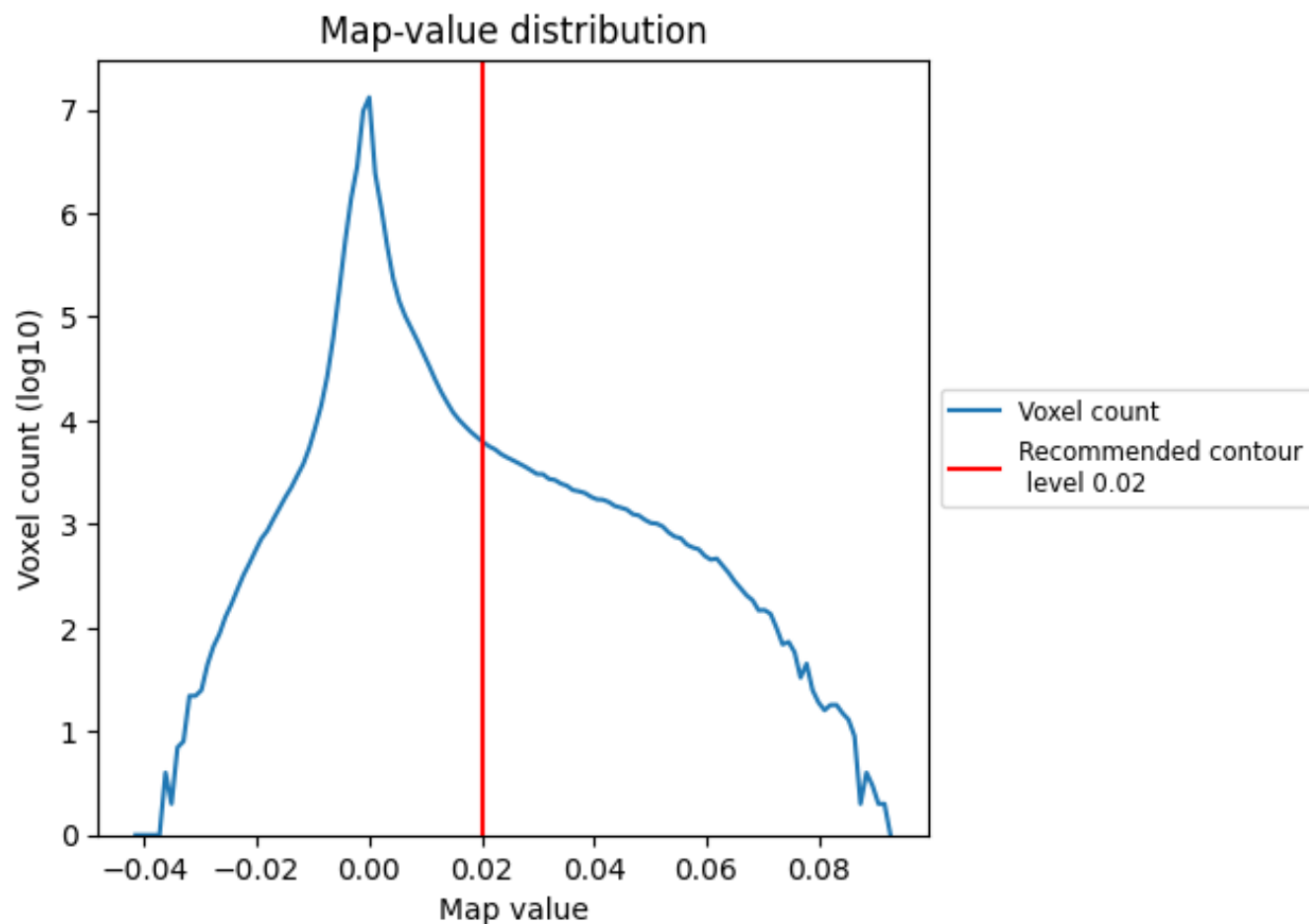
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

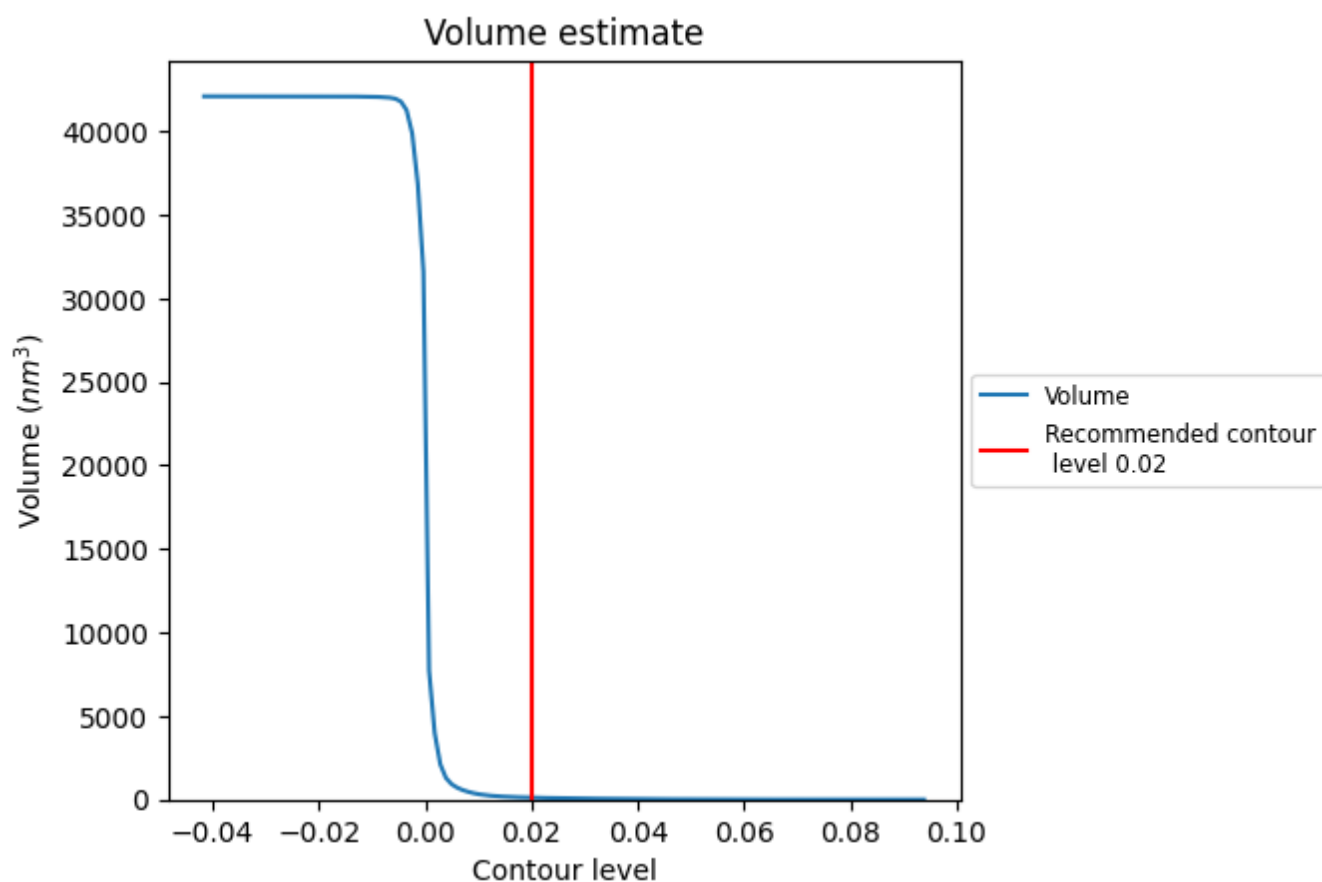
This section contains the results of statistical analysis of the map.

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.

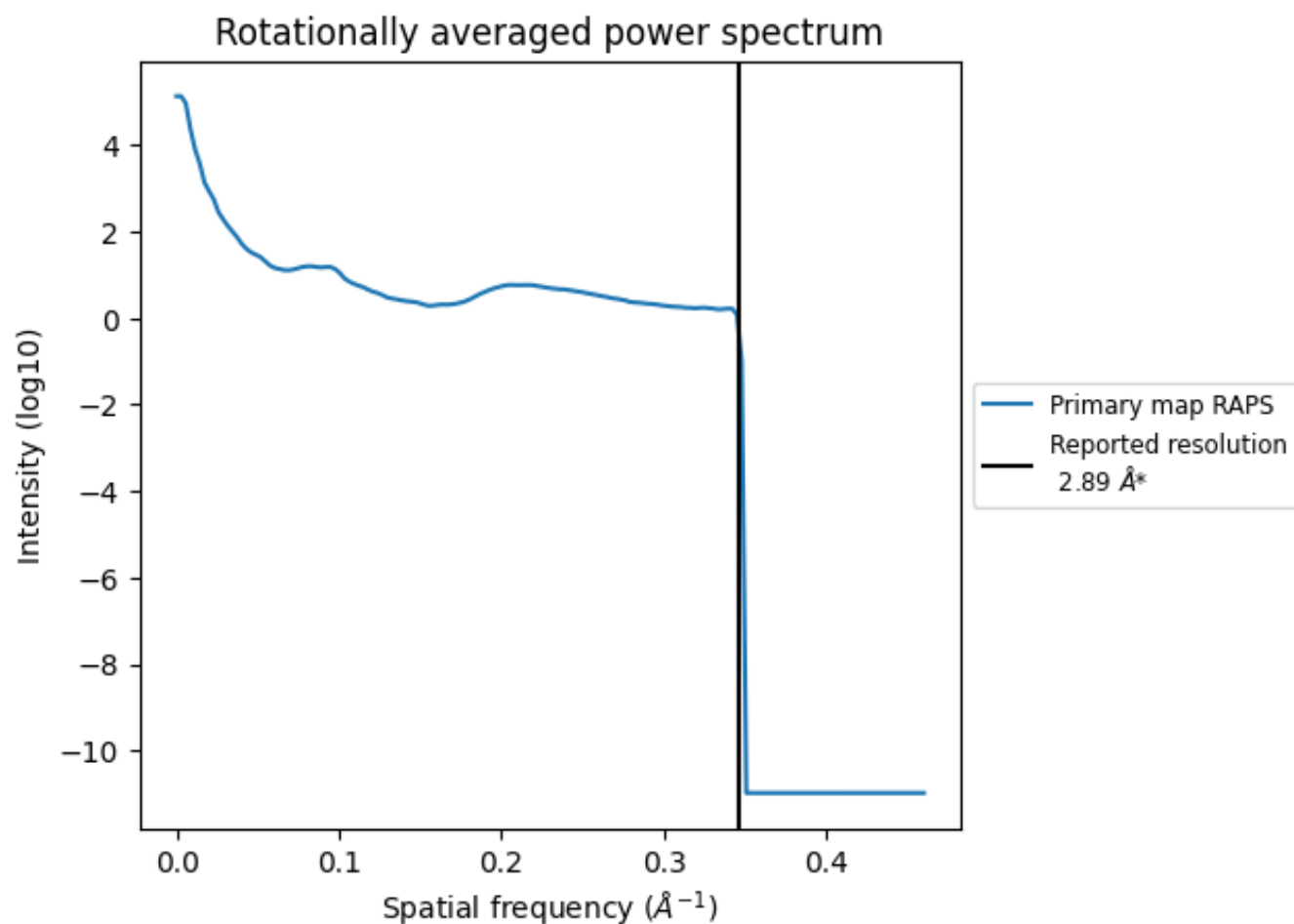
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 117 nm³; this corresponds to an approximate mass of 106 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 0.346 Å⁻¹

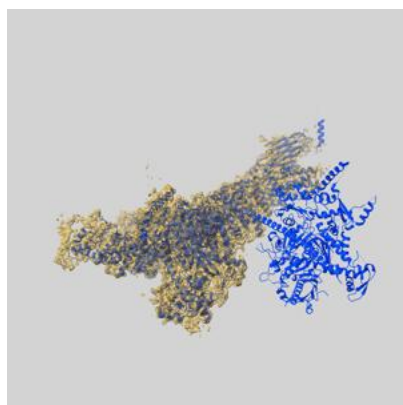
8 Fourier-Shell correlation

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

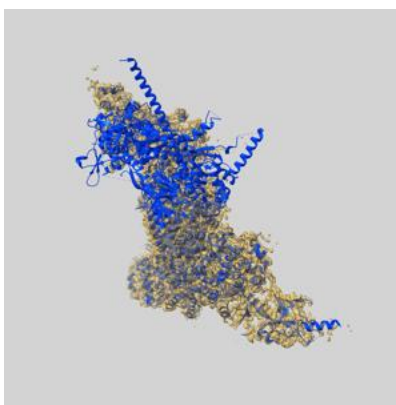
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-30285 and PDB model 7C4J. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

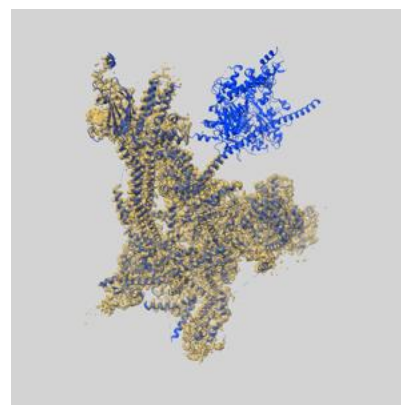
9.1 Map-model overlay [i](#)



X



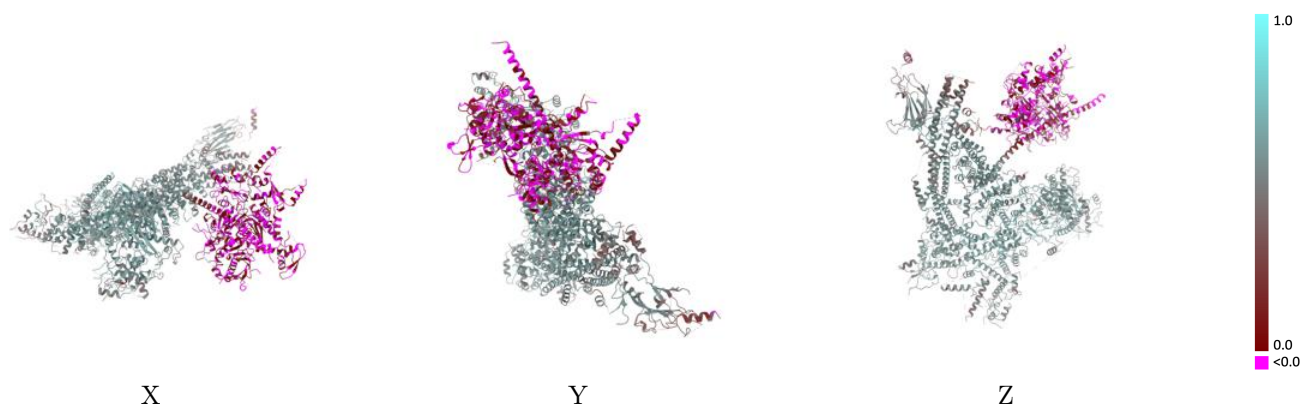
Y



Z

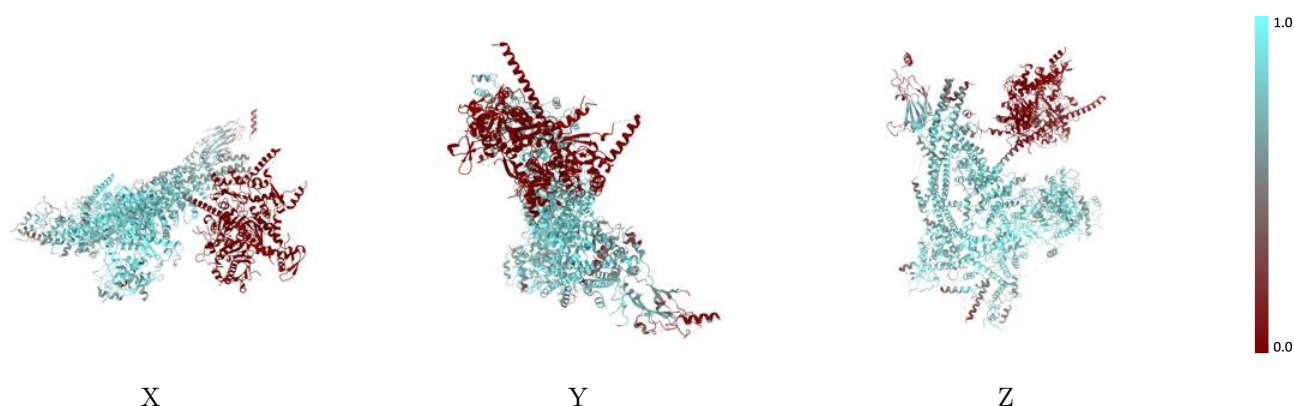
The images above show the 3D surface view of the map at the recommended contour level 0.02 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



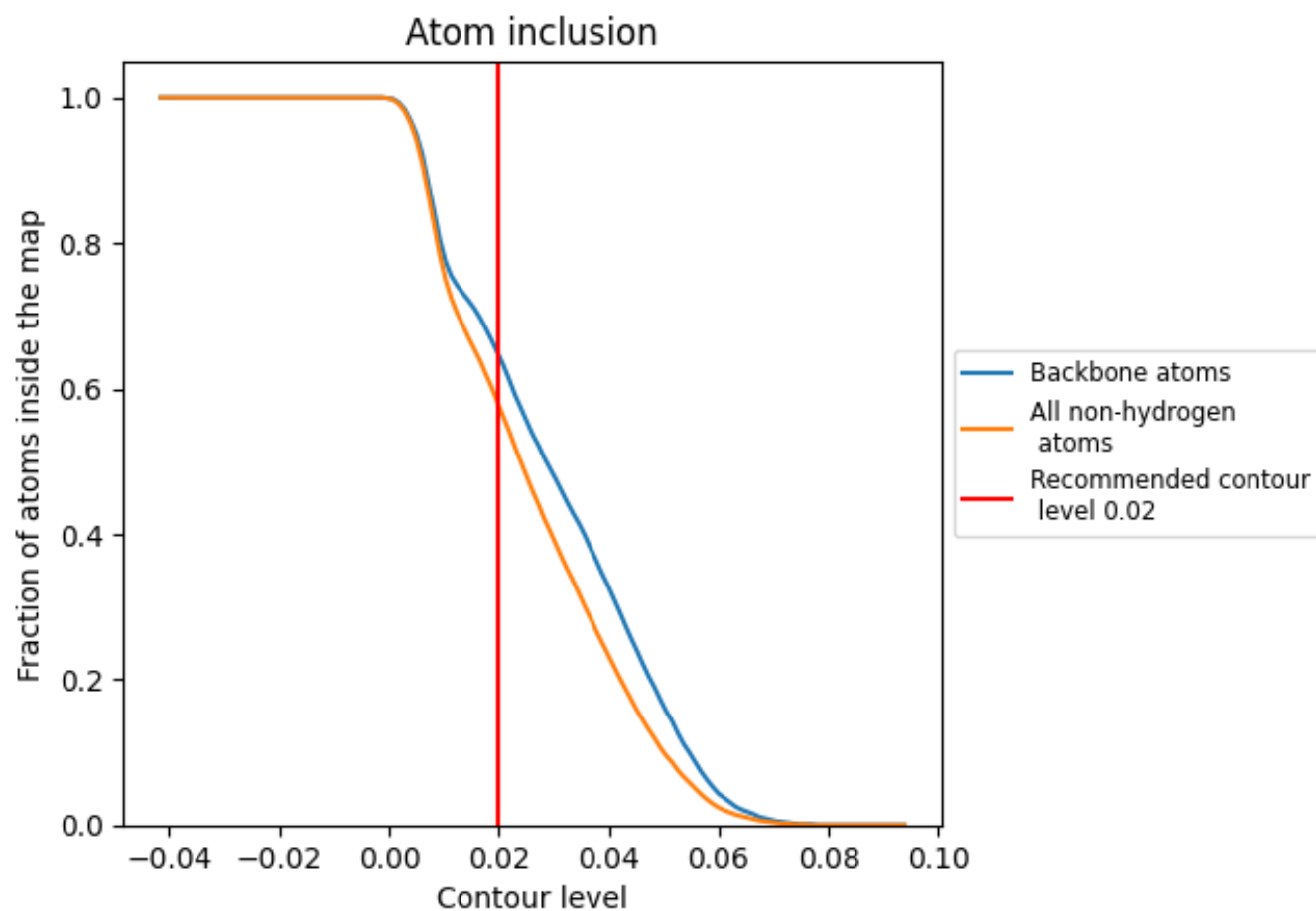
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.02).

9.4 Atom inclusion [i](#)



At the recommended contour level, 65% of all backbone atoms, 58% 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 (0.02) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.5770	 0.4210
A	 0.6750	 0.5210
B	 0.7750	 0.5570
C	 0.6390	 0.5020
D	 0.7580	 0.5560
E	 0.7080	 0.5270
F	 0.8220	 0.5720
G	 0.7980	 0.5180
H	 0.5780	 0.4090
I	 0.8420	 0.5850
J	 0.0000	 0.0470
K	 0.0000	 0.0130
L	 0.0000	 0.0160

