

Mar 16, 2024 - 08:51 am GMT

	:	EMD-17516
Title	:	Cryo-EM structure of CAK in complex with inhibitor ICEC0829
Authors	:	Cushing, V.I.; Koh, A.F.; Feng, J.; Jurgaityte, K.; Bahl, A.K.; Ali, S.; Kotecha,
		A.; Greber, B.J.
Deposited on	:	2023-05-30
Resolution	:	2.00 Å(reported)
Based on initial model	:	80RM

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 (i) 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 (1)) were used in the production of this report:

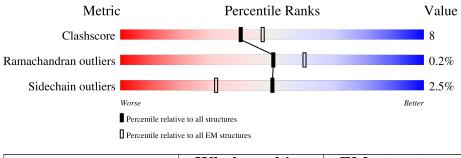
EMDB validation analysis Mogul		0.0.1.dev70 1.8.4, CSD as541be (2020)
MolProbity		
buster-report		
1		20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ		1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $ELECTRON\ MICROSCOPY$ 

The reported resolution of this entry is 2.00 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f EM} {f structures} \ (\#{f Entries})$
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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 <=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	Н	93	58% 119	•	30%
2	Ι	324	<u>6%</u> 69%	16%	• 15%
3	J	349	7%           64%	18%	18%



# 2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 5169 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 CDK-activating kinase assembly factor MAT1.

Mol	Chain	Residues		Atc	$\mathbf{ms}$			AltConf	Trace
1	Н	65	Total 512	C 329	N 87	0 94	$\begin{array}{c} \mathrm{S} \\ \mathrm{2} \end{array}$	0	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Н	217	SER	-	expression tag	UNP P51948
Н	218	ASN	-	expression tag	UNP P51948
Н	219	ALA	-	expression tag	UNP P51948

• Molecule 2 is a protein called Cyclin-H.

Mol	Chain	Residues		At	oms			AltConf	Trace
2	Ι	276	Total 2247	C 1437	N 385	O 407	S 18	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Ι	0	ACE	-	acetylation	UNP P51946

• Molecule 3 is a protein called Cyclin-dependent kinase 7.

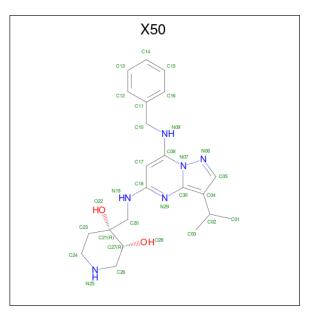
Mol	Chain	Residues		At	oms			AltConf	Trace
3	J	286	Total 2283	C 1482	N 386	0 404	S 11	0	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
J	-2	SER	-	expression tag	UNP P50613
J	-1	ASN	-	expression tag	UNP P50613
J	0	ALA	-	expression tag	UNP P50613

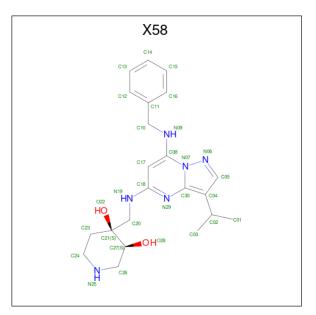


• Molecule 4 is (3R,4R)-4-[[[7-[(phenylmethyl)amino]-3-propan-2-yl-pyrazolo[1,5-a]pyrimi din-5-yl]amino]methyl]piperidine-3,4-diol (three-letter code: X50) (formula: C<sub>22</sub>H<sub>30</sub>N<sub>6</sub>O<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).



Ι	Mol	Chain	Residues	A	Aton	ns		AltConf
	4	J	1	Total 30	C 22	N 6	O 2	1

• Molecule 5 is (3S,4S)-4-[[[7-[(phenylmethyl)amino]-3-propan-2-yl-pyrazolo[1,5-a]pyrimid in-5-yl]amino]methyl]piperidine-3,4-diol (three-letter code: X58) (formula: C<sub>22</sub>H<sub>30</sub>N<sub>6</sub>O<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	A	Aton	ıs		AltConf
5	Т	1	Total	С	Ν	Ο	1
0	J		30	22	6	2	1

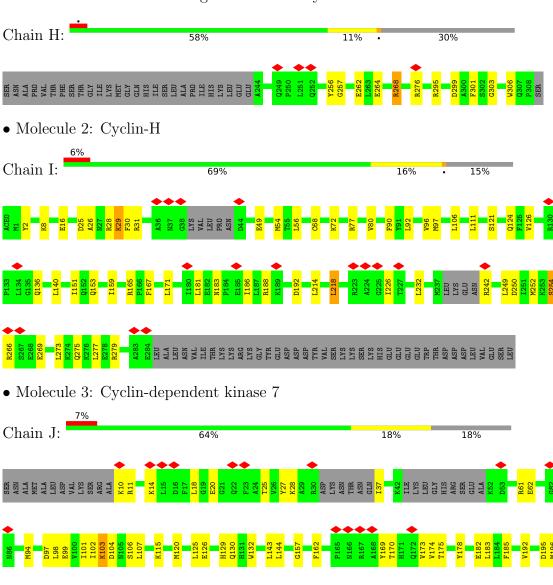
• Molecule 6 is water.

Mol	Chain	Residues	Atoms	AltConf
6	Н	6	Total O 6 6	0
6	Ι	28	TotalO2828	0
6	J	33	Total         O           33         33	0



# 3 Residue-property plots (i)

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: CDK-activating kinase assembly factor MAT1



#### ALA ILE LYS LYS LYS ARG THR ALA CLU GLU GLU GLU GLY CLU CLU PRO LYS LLU PHE



# 4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	253285	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose $(e^-/\text{\AA}^2)$	70	Depositor
Minimum defocus (nm)	400	Depositor
Maximum defocus (nm)	1900	Depositor
Magnification	245614	Depositor
Image detector	TFS FALCON 4i (4k x 4k)	Depositor
Maximum map value	0.054	Depositor
Minimum map value	-0.029	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.008	Depositor
Map size (Å)	273.59998, 273.59998, 273.59998	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles $(^{\circ})$	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.7124999, 0.7124999, 0.7124999	Depositor



# 5 Model quality (i)

## 5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: X50, X58, ACE

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	Bond lengths		angles
	Unam	RMSZ	# Z  > 5	RMSZ	# Z  > 5
1	Н	0.30	0/528	0.46	0/720
2	Ι	0.29	0/2292	0.51	0/3091
3	J	0.26	0/2340	0.49	0/3172
All	All	0.28	0/5160	0.49	0/6983

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

#### 5.2 Too-close contacts (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	Н	512	0	485	9	0
2	Ι	2247	0	2265	34	0
3	J	2283	0	2306	39	0
4	J	30	0	0	1	0
5	J	30	0	0	1	0
6	Н	6	0	0	0	0
6	Ι	28	0	0	0	0
6	J	33	0	0	1	0
All	All	5169	0	5056	78	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 8.

All (78) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
		distance (Å)	overlap (Å)
2:I:111:LEU:HD13	2:I:151:ILE:HD11	1.66	0.77
2:I:29:LYS:HD2	2:I:30:PHE:N	2.10	0.66
2:I:126:VAL:HG11	2:I:140:LEU:HB2	1.80	0.63
3:J:101:ILE:HG23	3:J:107:LEU:HD23	1.83	0.61
3:J:102:ILE:HD13	3:J:205:GLU:HG2	1.83	0.60
3:J:125:LEU:HD21	3:J:195:ASP:HB3	1.83	0.60
3:J:126:GLU:O	3:J:130:GLN:HG2	2.01	0.59
3:J:185:PHE:CD1	3:J:246:TYR:HA	2.36	0.59
2:I:31:ARG:NH2	2:I:49:GLU:OE2	2.37	0.57
3:J:98:LEU:HD12	3:J:143:LEU:HD12	1.85	0.57
2:I:68:CYS:HB3	2:I:80:VAL:HG22	1.86	0.57
2:I:249:LEU:HA	2:I:252:MET:HE2	1.86	0.57
3:J:213:LEU:HD13	3:J:225:ILE:HG12	1.87	0.56
1:H:295:ARG:HD2	3:J:132:TRP:CD2	2.42	0.54
1:H:256:TYR:HB2	1:H:303:GLY:HA2	1.90	0.53
3:J:28:LYS:HE3	3:J:37:ILE:HG21	1.90	0.52
1:H:264:GLU:OE1	1:H:264:GLU:N	2.39	0.52
3:J:62:GLU:OE1	6:J:501:HOH:O	2.18	0.51
3:J:99:GLU:HG3	3:J:103:LYS:HD2	1.93	0.51
3:J:231:THR:OG1	3:J:250:LYS:O	2.24	0.51
1:H:299:ASP:OD2	2:I:165:ARG:HD2	2.11	0.50
2:I:56:LEU:HD13	2:I:277:LEU:HD22	1.92	0.50
1:H:268:ARG:HB3	1:H:268:ARG:CZ	2.42	0.50
3:J:97:ASP:HB3	3:J:144:LEU:HD23	1.94	0.50
2:I:183:ASN:O	2:I:186:ILE:HG13	2.12	0.50
2:I:2:TYR:CZ	2:I:8:LYS:HB2	2.48	0.49
3:J:228:THR:HA	3:J:255:ILE:HD12	1.94	0.49
1:H:306:TRP:CE2	2:I:8:LYS:HE2	2.48	0.49
3:J:14:LYS:HD3	3:J:27:TYR:CD2	2.48	0.48
2:I:275:GLN:O	2:I:279:ARG:HD3	2.15	0.47
2:I:218:LEU:HD11	2:I:232:LEU:HD11	1.96	0.47
3:J:94:MET:HB2	5:J:402[B]:X58:N06	2.29	0.47
3:J:236:GLN:O	3:J:279:ASN:HB2	2.15	0.47
1:H:257:GLY:HA3	1:H:301:PHE:O	2.16	0.46
3:J:115:LYS:NZ	3:J:299:PRO:O	2.38	0.46
3:J:175:THR:HG23	3:J:178:TYR:HD1	1.81	0.46
3:J:192:VAL:O	3:J:196:MET:HG2	2.16	0.46
2:I:167:PHE:HE2	2:I:188:ARG:HG3	1.81	0.45

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Atom-1	Atom-2	Interatomic	Clash
		distance (Å)	overlap (Å)
2:I:250:ASP:O	2:I:254:SER:OG	2.32	0.45
2:I:26:ALA:HA	2:I:29:LYS:HE3	1.99	0.45
3:J:143:LEU:HD11	3:J:202:ILE:HD13	1.99	0.45
2:I:269:GLU:O	2:I:273:LEU:HG	2.17	0.45
3:J:18:LEU:HD11	3:J:28:LYS:HB2	1.99	0.45
1:H:262:GLU:HB3	1:H:264:GLU:OE1	2.17	0.44
2:I:126:VAL:O	2:I:136:GLN:HG2	2.18	0.44
2:I:121:SER:O	2:I:124:GLN:HG2	2.17	0.44
3:J:104:ASP:OD1	3:J:106:SER:OG	2.23	0.44
3:J:210:VAL:HG22	3:J:211:PRO:HD2	2.00	0.44
2:I:214:LEU:HD21	2:I:232:LEU:HD21	1.99	0.44
3:J:94:MET:HG3	3:J:144:LEU:HB2	1.99	0.44
3:J:120:MET:HE3	3:J:120:MET:HB2	1.80	0.43
3:J:232:PRO:HB3	3:J:278:PHE:CD2	2.53	0.43
2:I:26:ALA:HA	2:I:29:LYS:HG3	2.00	0.43
1:H:295:ARG:HD2	3:J:132:TRP:CG	2.53	0.43
2:I:159:ILE:HD13	2:I:159:ILE:HA	1.91	0.43
3:J:61:ARG:HD2	3:J:157:GLY:O	2.20	0.42
3:J:94:MET:HB2	4:J:401[A]:X50:N06	2.33	0.42
3:J:125:LEU:HG	3:J:129:HIS:NE2	2.34	0.42
3:J:232:PRO:HG2	3:J:240:MET:HE1	2.01	0.42
3:J:20:GLU:HG2	3:J:25:THR:HG22	2.00	0.42
3:J:170:THR:HG23	3:J:173:VAL:HG13	2.02	0.42
2:I:72:LYS:HB3	2:I:72:LYS:HE2	1.73	0.42
2:I:181:LEU:HD13	2:I:226:ILE:HG21	2.02	0.42
3:J:256:PRO:HB3	3:J:258:HIS:CE1	2.55	0.42
3:J:10:LYS:HB3	3:J:11:ARG:H	1.62	0.41
3:J:98:LEU:HD23	3:J:98:LEU:HA	1.89	0.41
2:I:266:ARG:HB2	2:I:269:GLU:HG2	2.02	0.41
3:J:162:PHE:HA	3:J:169:TYR:OH	2.20	0.41
2:I:90:PHE:CE1	2:I:153:GLN:HG3	2.56	0.41
2:I:92:LEU:HD23	2:I:92:LEU:HA	1.95	0.41
2:I:106:LEU:HD23	2:I:106:LEU:HA	1.94	0.41
2:I:54:MET:O	2:I:54:MET:HE3	2.21	0.40
2:I:96:VAL:O	2:I:277:LEU:HD11	2.22	0.40
2:I:275:GLN:HB3	2:I:279:ARG:HH11	1.86	0.40
2:I:28:ARG:HG3	2:I:29:LYS:N	2.36	0.40
2:I:30:PHE:CD2	2:I:97:MET:HE3	2.56	0.40
2:I:111:LEU:HD22	2:I:151:ILE:HD12	2.04	0.40
3:J:182:GLU:HG2	3:J:183:LEU:N	2.37	0.40

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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	Н	63/93~(68%)	63 (100%)	0	0	100 100
2	Ι	270/324 (83%)	267~(99%)	3 (1%)	0	100 100
3	J	280/349~(80%)	273~(98%)	6(2%)	1 (0%)	34 30
All	All	613/766~(80%)	603 (98%)	9(2%)	1 (0%)	50 44

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	J	174	VAL

#### 5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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	Н	52/76~(68%)	50~(96%)	2~(4%)	33 31
2	Ι	250/296~(84%)	240 (96%)	10 (4%)	31 29
3	J	248/301~(82%)	246~(99%)	2(1%)	81 86
All	All	550/673~(82%)	536~(98%)	14 (2%)	50 49

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	Н	268	ARG
1	Н	276	ARG

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Mol	Chain	Res	Type
2	Ι	16	GLU
2	Ι	25	ASP
2	Ι	29	LYS
2	Ι	77	ARG
2	Ι	171	LEU
2	Ι	192	ASP
2	Ι	218	LEU
2	Ι	242	ARG
2	Ι	254	SER
2	Ι	261	LYS
3	J	103	LYS
3	J	242	SER

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Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
3	J	129	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 monosaccharides in this entry.

#### 5.6 Ligand geometry (i)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the



Mol	Type	Chain	Res	Link	Bo	ond leng	$\mathbf{ths}$	B	ond ang	les
10101	туре	Chain	nes	LIIIK	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	X58	J	402[B]	-	$28,\!33,\!33$	1.84	4 (14%)	29,47,47	1.41	5 (17%)
4	X50	J	401[A]	_	28,33,33	2.54	8 (28%)	29,47,47	1.37	5 (17%)

expected value. A bond length (or angle) with |Z| > 2 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Ν	Лоl	Type	Chain	Res	Link	Chirals	Torsions	Rings
	5	X58	J	402[B]	-	-	0/13/29/29	1/4/4/4
	4	X50	J	401[A]	-	-	2/13/29/29	1/4/4/4

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	J	401[A]	X50	C18-N19	9.29	1.50	1.36
5	J	402[B]	X58	C18-N19	6.89	1.46	1.36
4	J	401[A]	X50	C08-N09	4.82	1.48	1.37
4	J	401[A]	X50	C23-C21	-3.85	1.48	1.53
5	J	402[B]	X58	C23-C21	-3.77	1.48	1.53
5	J	402[B]	X58	C08-N09	3.69	1.46	1.37
4	J	401[A]	X50	C15-C16	2.69	1.44	1.38
5	J	402[B]	X58	C21-C27	-2.66	1.51	1.53
4	J	401[A]	X50	C30-N29	2.62	1.39	1.35
4	J	401[A]	X50	C20-C21	2.43	1.56	1.53
4	J	401[A]	X50	O28-C27	-2.27	1.39	1.43
4	J	401[A]	X50	C16-C11	-2.01	1.34	1.38

All (12) bond length outliers are listed below:

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
5	J	402[B]	X58	N09-C08-N07	3.96	119.42	116.65
5	J	402[B]	X58	C20-N19-C18	-3.53	117.87	123.42
4	J	401[A]	X50	N09-C08-N07	3.45	119.07	116.65
4	J	401[A]	X50	C20-N19-C18	-3.38	118.10	123.42
4	J	401[A]	X50	C05-N06-N07	3.13	108.88	103.94
5	J	402[B]	X58	C05-N06-N07	3.12	108.86	103.94
4	J	401[A]	X50	C18-N29-C30	3.10	120.42	117.59

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Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
5	J	402[B]	X58	C18-N29-C30	2.73	120.08	117.59
4	J	401[A]	X50	C17-C18-N29	-2.24	119.80	123.91
5	J	402[B]	X58	C17-C18-N29	-2.09	120.09	123.91

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There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	J	401[A]	X50	C11-C10-N09-C08
4	J	401[A]	X50	N19-C20-C21-C23

All (2) ring outliers are listed below:

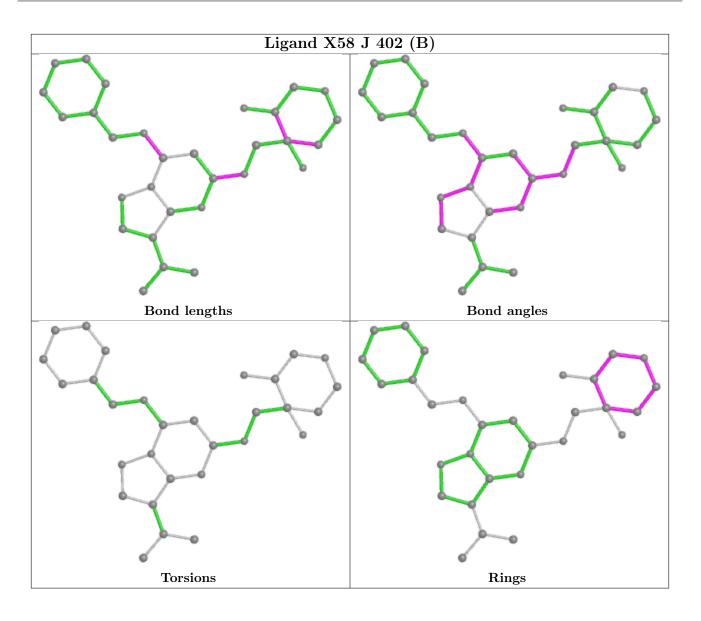
Mol	Chain	Res	Type	Atoms
5	J	402[B]	X58	C21-C23-C24-C26-C27-N25
4	J	401[A]	X50	C21-C23-C24-C26-C27-N25

2 monomers are involved in 2 short contacts:

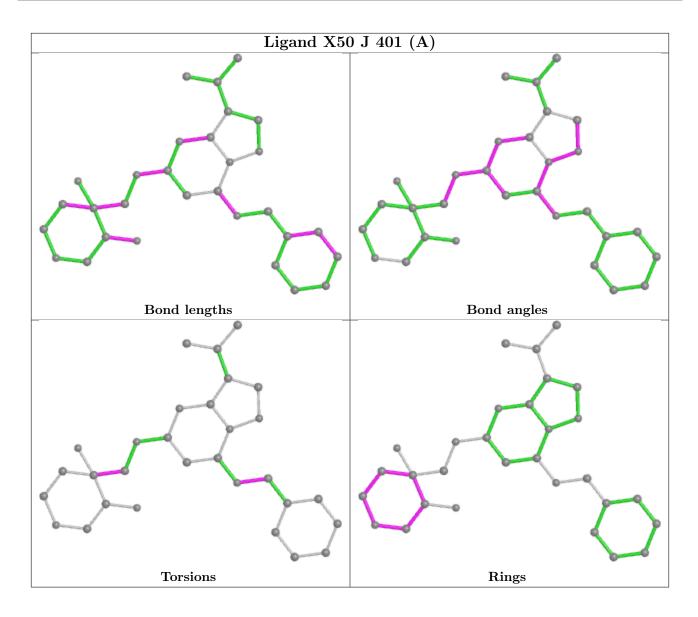
Ι	Mol	Chain	Res	Type	Clashes	Symm-Clashes
	5	J	402[B]	X58	1	0
	4	J	401[A]	X50	1	0

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









## 5.7 Other polymers (i)

There are no such residues in this entry.

## 5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



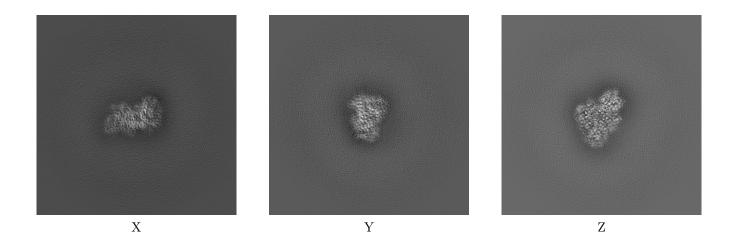
# 6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-17516. These allow visual inspection of the internal detail of the map and identification of artifacts.

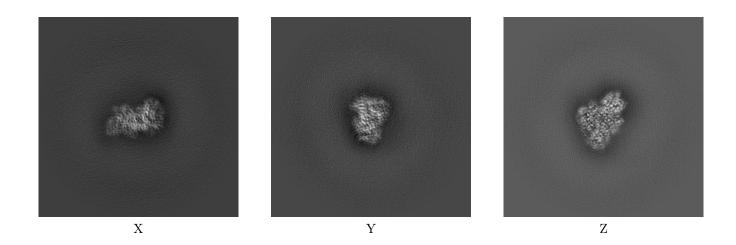
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

## 6.1 Orthogonal projections (i)

#### 6.1.1 Primary map



6.1.2 Raw map

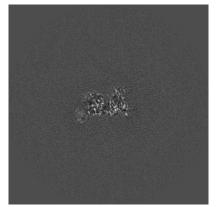


The images above show the map projected in three orthogonal directions.

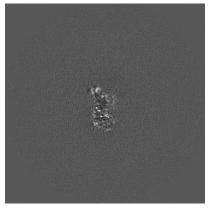


#### 6.2 Central slices (i)

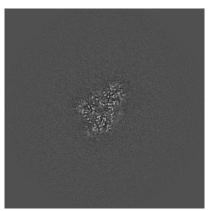
#### 6.2.1 Primary map



X Index: 192

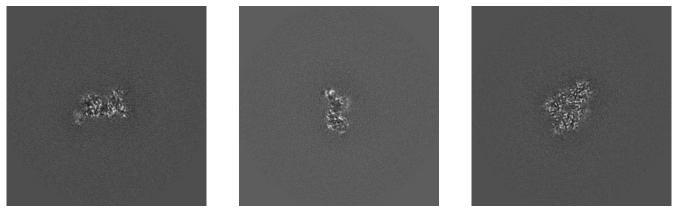


Y Index: 192



Z Index: 192

#### 6.2.2 Raw map



X Index: 192

Y Index: 192

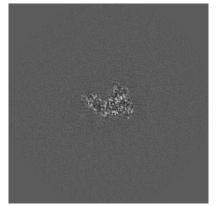
Z Index: 192

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



#### 6.3 Largest variance slices (i)

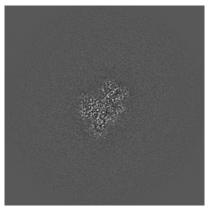
#### 6.3.1 Primary map



X Index: 202



Y Index: 208



Z Index: 179

#### 6.3.2 Raw map



X Index: 202

Y Index: 208

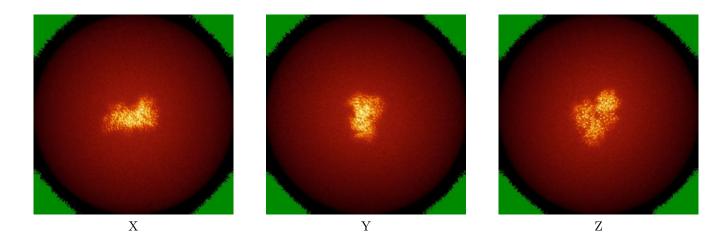


The images above show the largest variance slices of the map in three orthogonal directions.

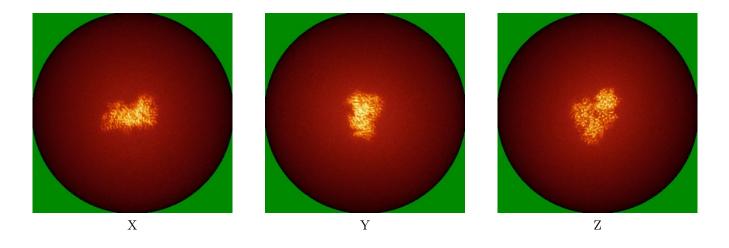


### 6.4 Orthogonal standard-deviation projections (False-color) (i)

#### 6.4.1 Primary map



6.4.2 Raw map

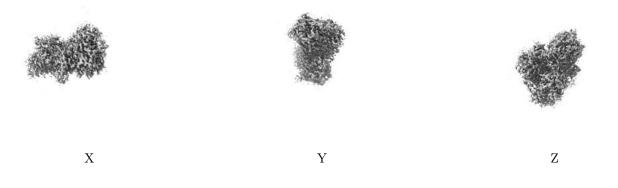


The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.



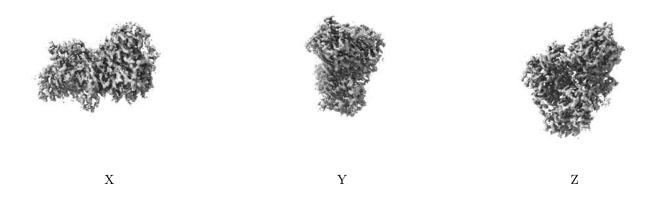
#### 6.5 Orthogonal surface views (i)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.008. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

#### 6.5.2 Raw map



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



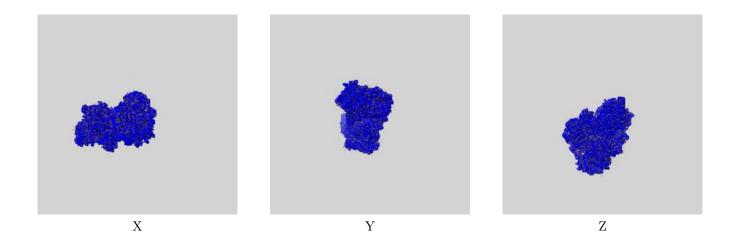
### 6.6 Mask visualisation (i)

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

A mask typically either:

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

#### 6.6.1 emd\_17516\_msk\_1.map (i)

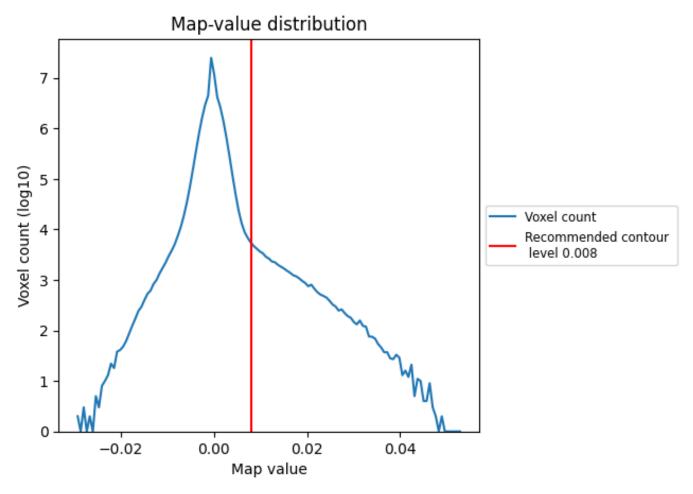




# 7 Map analysis (i)

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

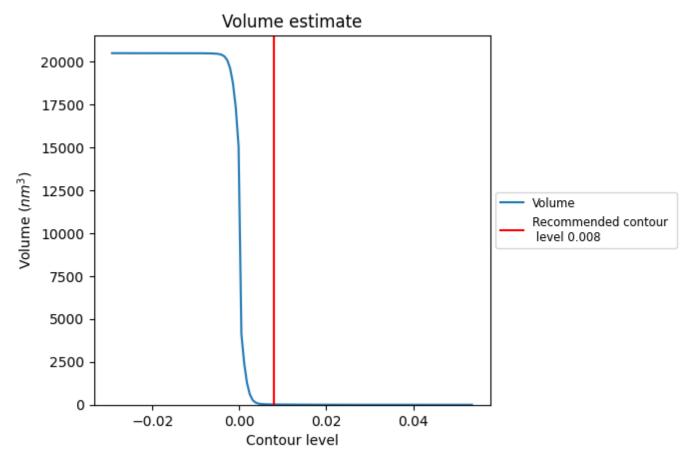
## 7.1 Map-value distribution (i)



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)

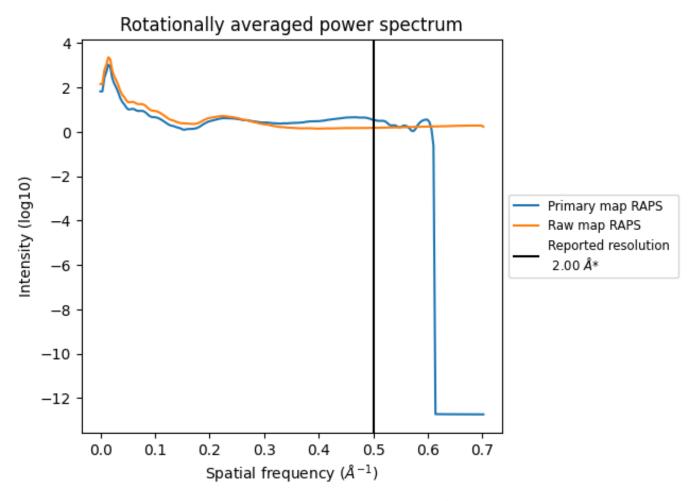


The volume at the recommended contour level is  $19~\mathrm{nm^3};$  this corresponds to an approximate mass of 17 kDa.

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)



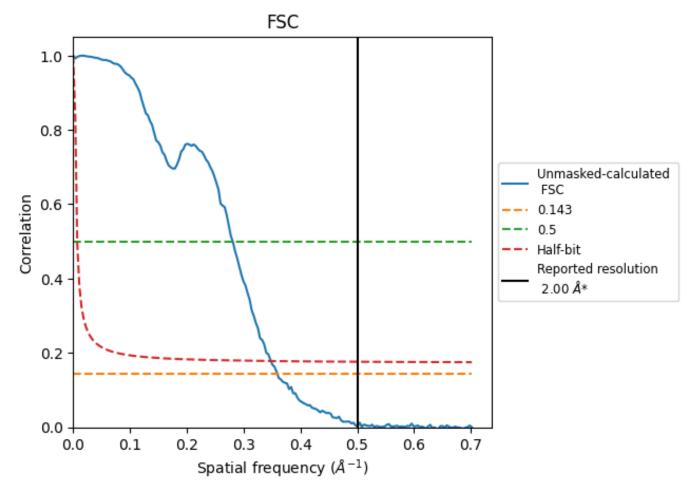
\*Reported resolution corresponds to spatial frequency of 0.500  ${\rm \AA^{-1}}$ 



## 8 Fourier-Shell correlation (i)

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

#### 8.1 FSC (i)



\*Reported resolution corresponds to spatial frequency of 0.500  ${\rm \AA^{-1}}$ 



#### 8.2 Resolution estimates (i)

Resolution estimate (Å)	Estim	Estimation criterion (FSC cut-off)			
Resolution estimate (A)	0.143	0.5	Half-bit		
Reported by author	2.00	-	-		
Author-provided FSC curve	-	-	-		
Unmasked-calculated*	2.78	3.56	2.87		

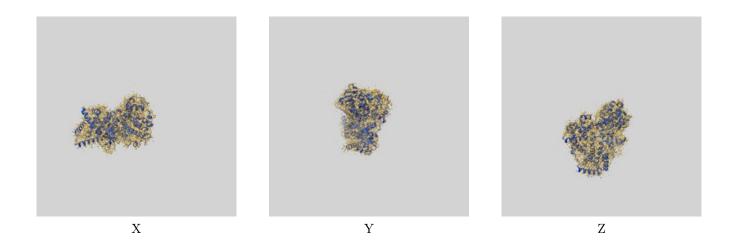
\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 2.78 differs from the reported value 2.0 by more than 10 %



# 9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-17516 and PDB model 8P73. Per-residue inclusion information can be found in section 3 on page 6.

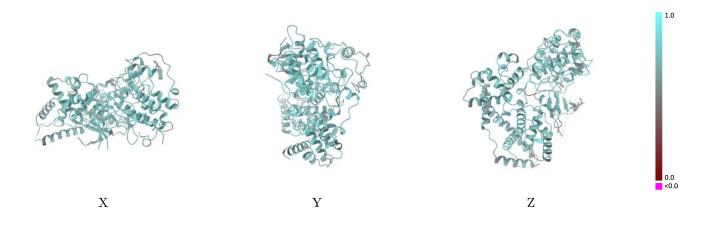
## 9.1 Map-model overlay (i)



The images above show the 3D surface view of the map at the recommended contour level 0.008 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.

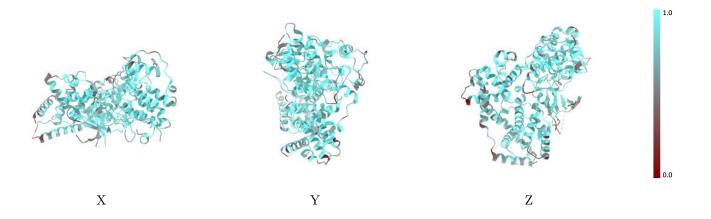


#### 9.2 Q-score mapped to coordinate model (i)



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.

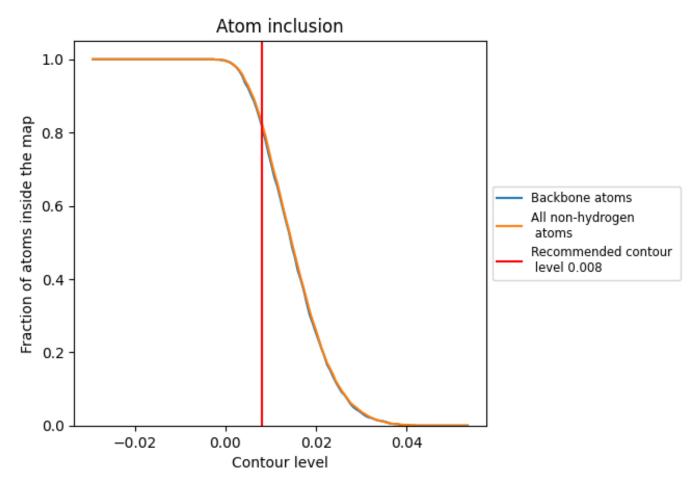
#### 9.3 Atom inclusion mapped to coordinate model (i)



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 (0.008).



#### 9.4 Atom inclusion (i)



At the recommended contour level, 82% of all backbone atoms, 83% of all non-hydrogen atoms, are inside the map.



1.0

0.0 <0.0

### 9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.008) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8280	0.6830
Н	0.8580	0.6940
Ι	0.8380	0.6850
J	0.8220	0.6770

