



Full wwPDB EM Validation Report ⓘ

Jul 28, 2026 – 06:23 PM EDT

PDB ID : 11YE / pdb_000011ye
EMDB ID : EMD-76184
Title : Curved structure of mPiezo1 in plasma membrane vesicles
Authors : Vaisey, G.V.; MacKinnon, R.M.
Deposited on : 2026-03-18
Resolution : 3.70 Å(reported)

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

EMDB validation analysis : 0.0.1.dev133
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) : 2.50

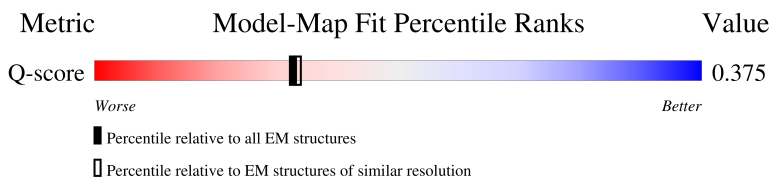
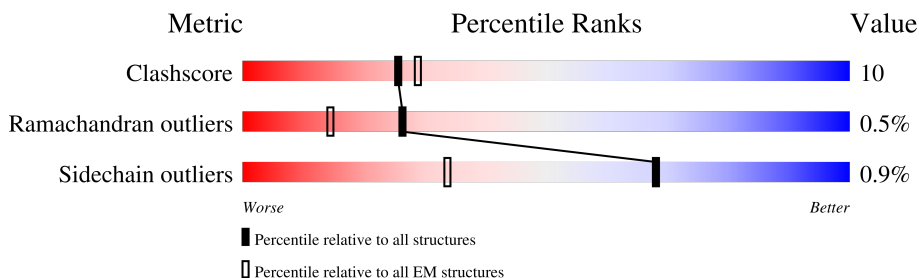
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.70 Å.

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



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

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	C	2561	
1	D	2561	
1	E	2561	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 29295 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 Piezo-type mechanosensitive ion channel component 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	C	1335	Total	C	N	O	S	0	0
			9765	6243	1715	1756	51		
1	D	1335	Total	C	N	O	S	0	0
			9765	6243	1715	1756	51		
1	E	1335	Total	C	N	O	S	0	0
			9765	6243	1715	1756	51		

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	2420A	SER	-	insertion	UNP E2JF22
C	2420B	ARG	-	insertion	UNP E2JF22
C	2420C	LEU	-	insertion	UNP E2JF22
C	2420D	GLU	-	insertion	UNP E2JF22
C	2420E	GLU	-	insertion	UNP E2JF22
C	2420F	GLU	-	insertion	UNP E2JF22
C	2420G	LEU	-	insertion	UNP E2JF22
C	2420H	ARG	-	insertion	UNP E2JF22
C	2420I	ARG	-	insertion	UNP E2JF22
C	2430J	ARG	-	insertion	UNP E2JF22
C	2420K	LEU	-	insertion	UNP E2JF22
C	2420L	THR	-	insertion	UNP E2JF22
C	2420M	GLU	-	insertion	UNP E2JF22
C	2548	SER	-	expression tag	UNP E2JF22
D	2420A	SER	-	insertion	UNP E2JF22
D	2420B	ARG	-	insertion	UNP E2JF22
D	2420C	LEU	-	insertion	UNP E2JF22
D	2420D	GLU	-	insertion	UNP E2JF22
D	2420E	GLU	-	insertion	UNP E2JF22
D	2420F	GLU	-	insertion	UNP E2JF22
D	2420G	LEU	-	insertion	UNP E2JF22
D	2420H	ARG	-	insertion	UNP E2JF22
D	2420I	ARG	-	insertion	UNP E2JF22
D	2420J	ARG	-	insertion	UNP E2JF22

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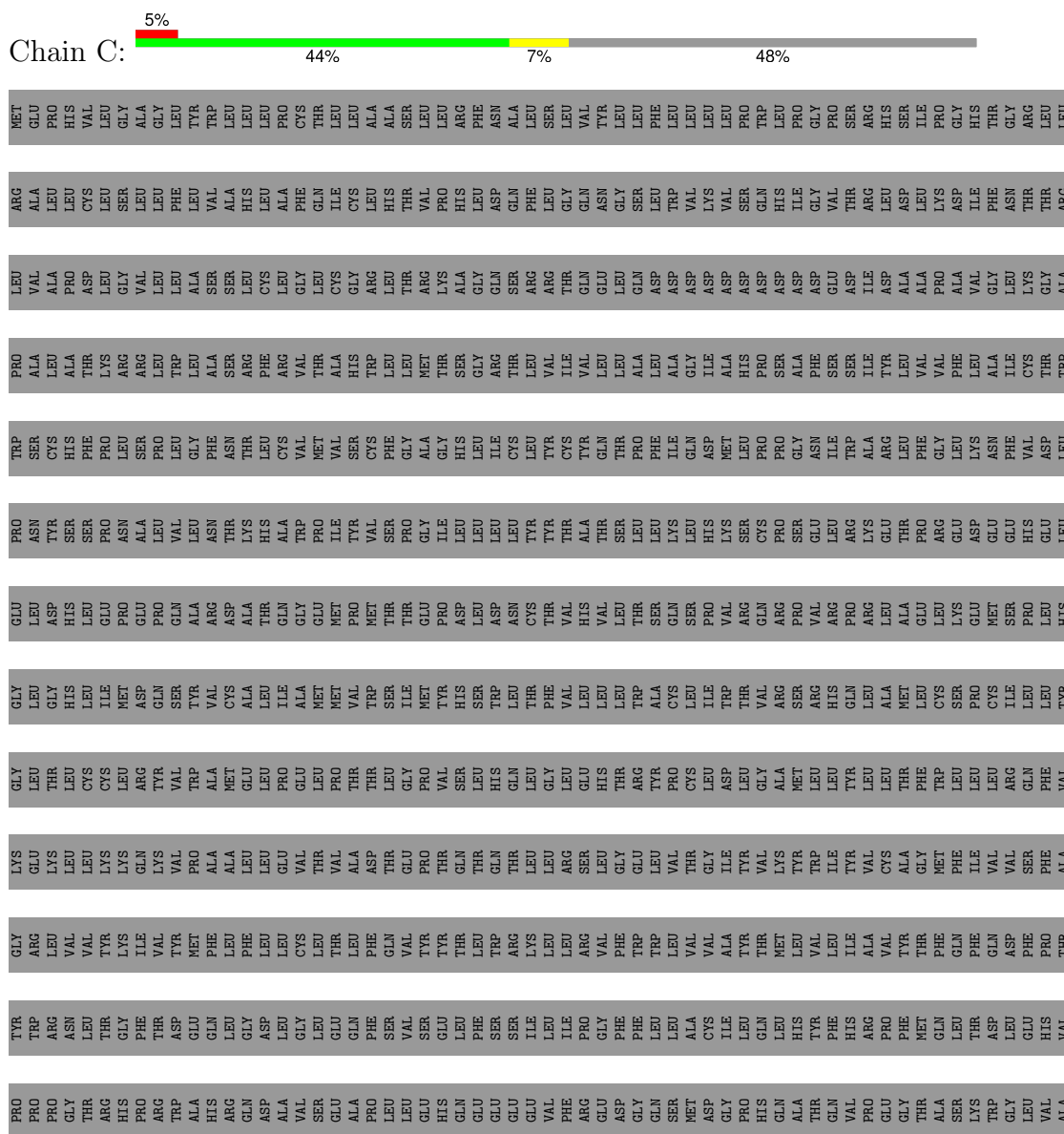
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Chain	Residue	Modelled	Actual	Comment	Reference
D	2420K	LEU	-	insertion	UNP E2JF22
D	2420L	THR	-	insertion	UNP E2JF22
D	2420M	GLU	-	insertion	UNP E2JF22
D	2548	SER	-	expression tag	UNP E2JF22
E	2420A	SER	-	insertion	UNP E2JF22
E	2420B	ARG	-	insertion	UNP E2JF22
E	2420C	LEU	-	insertion	UNP E2JF22
E	2420D	GLU	-	insertion	UNP E2JF22
E	2420E	GLU	-	insertion	UNP E2JF22
E	2420F	GLU	-	insertion	UNP E2JF22
E	2420G	LEU	-	insertion	UNP E2JF22
E	2420H	ARG	-	insertion	UNP E2JF22
E	2420I	ARG	-	insertion	UNP E2JF22
E	2420J	ARG	-	insertion	UNP E2JF22
E	2420K	LEU	-	insertion	UNP E2JF22
E	2420L	THR	-	insertion	UNP E2JF22
E	2420M	GLU	-	insertion	UNP E2JF22
E	2548	SER	-	expression tag	UNP E2JF22

3 Residue-property plots

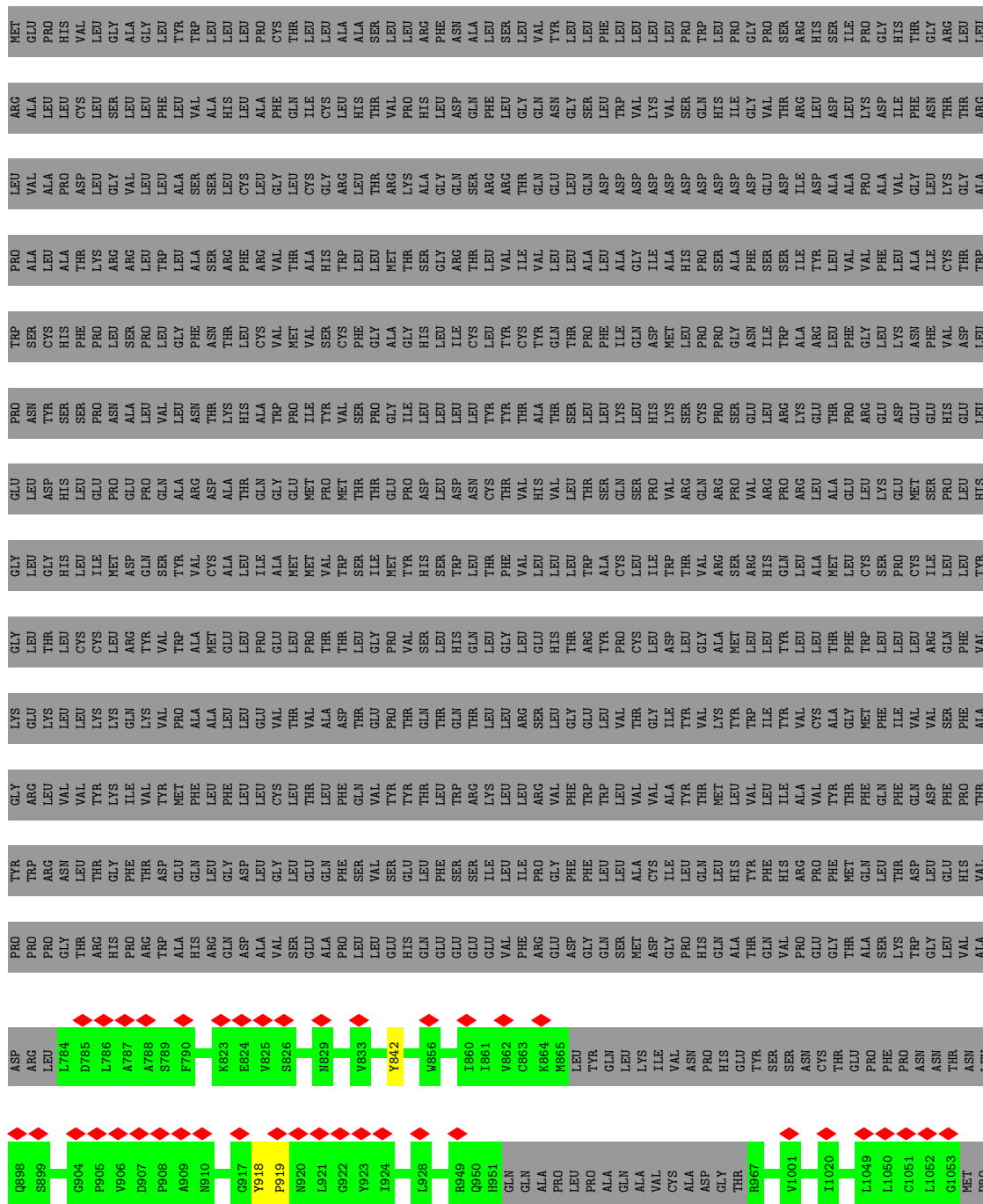
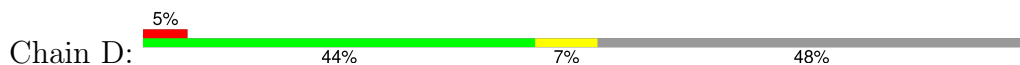
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Piezo-type mechanosensitive ion channel component 1



T2143	T2144	T2145	T2146	L2147	N2151	W2152	V2016	P2017	F2020	W2031	L2037	R2040	K2046	T2069	Q2075	N2076	A2077	V2078	A2079	Q2080	L2081	W2082	K2086	Y2089	S2093	L2097	G2100	T2103	R2104	N2108	H2116	L2117	Q2123	R2126	L2127	W2128	P2129	V2132	D2139																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																
SER	SER	LEU	SER	ASP	ASP	Q2015	Q2015	P2017	F2020	W2031	L2037	R2040	K2046	T2069	Q2075	N2076	A2077	V2078	A2079	Q2080	L2081	W2082	K2086	Y2089	S2093	L2097	G2100	T2103	R2104	N2108	H2116	L2117	Q2123	R2126	L2127	W2128	P2129	V2132	D2139																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																
GLY	THR	PRO	ARG	THR	THR	GLU	ARG	LYS	ARG	HIS	GLN	GLY	SER	ARG	GLU	ARG	MET	LYS	ALA	GLY	ARG	GLN	SER	GLY	Y2089	S2093	L2097	R1960	F1961	F1962	K1969	Y1970	R1971	I1988	F1998	GLY	LYS	HIS	THR	ALA																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																															
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ALA	ARG	THR	ARG	MET	ARG	THR	R1652	G1670	R1671	L1673	R1674	L1675	H1686	T1695	L1696	L1697	H1698	H1699	M1700	V1701	A1706	S1707	L1710	P1711	T1721	I1722	P1723	R1724	P1725	M1740	K1744	V1759	R1762	Y1768	F1769	P1770	P1771	HIS	PRO	GLY	THR	ALA																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																													
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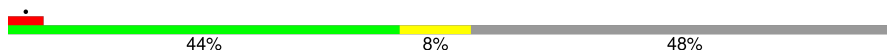
- Molecule 1: Piezo-type mechanosensitive ion channel component 1





- Molecule 1: Piezo-type mechanosensitive ion channel component 1

Chain E:

[illegible]



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	69370	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 ($e^-/\text{\AA}^2$)	60.00	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	34.226	Depositor
Minimum map value	-19.697	Depositor
Average map value	-0.010	Depositor
Map value standard deviation	0.997	Depositor
Recommended contour level	4.5	Depositor
Map size ($\text{\AA}$)	376.0, 376.0, 376.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.94, 0.94, 0.94	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	C	0.16	0/9982	0.34	0/13408
1	D	0.16	0/9982	0.34	0/13408
1	E	0.16	0/9982	0.34	0/13408
All	All	0.16	0/29946	0.34	0/40224

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	C	9765	0	9060	202	0
1	D	9765	0	9060	198	0
1	E	9765	0	9060	198	0
All	All	29295	0	27180	583	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 10.

All (583) 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:C:1744:LYS:CG	1:C:1776:LEU:HD22	1.59	1.31

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:1744:LYS:CG	1:E:1776:LEU:HD22	1.59	1.30
1:D:1744:LYS:CG	1:D:1776:LEU:HD22	1.59	1.30
1:C:1523:PHE:HZ	1:C:1686:HIS:NE2	1.37	1.21
1:D:1744:LYS:HG2	1:D:1776:LEU:CD2	1.71	1.21
1:C:1744:LYS:HG2	1:C:1776:LEU:CD2	1.71	1.20
1:D:1523:PHE:HZ	1:D:1686:HIS:NE2	1.37	1.20
1:E:1523:PHE:HZ	1:E:1686:HIS:NE2	1.37	1.20
1:E:1744:LYS:HG2	1:E:1776:LEU:CD2	1.71	1.20
1:E:2139:ASP:OD1	1:E:2143:THR:HG21	1.41	1.18
1:C:2139:ASP:OD1	1:C:2143:THR:HG21	1.41	1.17
1:D:2139:ASP:OD1	1:D:2143:THR:HG21	1.41	1.15
1:C:1191:LEU:CD1	1:C:1198:LEU:CD2	2.32	1.08
1:E:1285:ILE:HG13	1:E:1289:PHE:CE2	1.89	1.08
1:D:1191:LEU:CD1	1:D:1198:LEU:CD2	2.32	1.08
1:E:1191:LEU:CD1	1:E:1198:LEU:CD2	2.32	1.06
1:C:1285:ILE:HG13	1:C:1289:PHE:CE2	1.89	1.06
1:C:1744:LYS:HE3	1:C:1776:LEU:HD13	1.38	1.06
1:D:1285:ILE:HG13	1:D:1289:PHE:CE2	1.89	1.06
1:D:2238:LEU:HD13	1:D:2430:TRP:CD2	1.91	1.05
1:E:2238:LEU:HD13	1:E:2430:TRP:CD2	1.91	1.05
1:D:1744:LYS:HE3	1:D:1776:LEU:HD13	1.38	1.05
1:E:1701:VAL:CG1	1:E:1784:LYS:HD3	1.87	1.05
1:C:2238:LEU:HD13	1:C:2430:TRP:CD2	1.91	1.04
1:D:1701:VAL:CG1	1:D:1784:LYS:HD3	1.87	1.04
1:E:1744:LYS:HE3	1:E:1776:LEU:HD13	1.38	1.04
1:C:1701:VAL:CG1	1:C:1784:LYS:HD3	1.87	1.03
1:D:1523:PHE:CZ	1:D:1686:HIS:NE2	2.26	1.03
1:D:1701:VAL:HG12	1:D:1784:LYS:CD	1.89	1.03
1:C:1523:PHE:CZ	1:C:1686:HIS:NE2	2.26	1.02
1:E:1523:PHE:CZ	1:E:1686:HIS:NE2	2.26	1.02
1:C:1282:TRP:O	1:C:1285:ILE:HG22	1.60	1.02
1:E:1701:VAL:HG12	1:E:1784:LYS:CD	1.89	1.02
1:D:1282:TRP:O	1:D:1285:ILE:HG22	1.60	1.01
1:C:1701:VAL:HG12	1:C:1784:LYS:CD	1.89	1.01
1:D:1695:ILE:HG12	1:D:1789:GLN:NE2	1.77	0.99
1:E:1282:TRP:O	1:E:1285:ILE:HG22	1.60	0.99
1:C:1191:LEU:CD1	1:C:1198:LEU:HD22	1.93	0.99
1:D:1701:VAL:HG12	1:D:1784:LYS:HD3	1.44	0.98
1:C:1695:ILE:HG12	1:C:1789:GLN:NE2	1.77	0.98
1:E:1695:ILE:HG12	1:E:1789:GLN:NE2	1.77	0.98
1:E:1191:LEU:CD1	1:E:1198:LEU:HD22	1.93	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1191:LEU:CD1	1:D:1198:LEU:HD22	1.93	0.97
1:E:1701:VAL:HG12	1:E:1784:LYS:HD3	1.44	0.97
1:C:1701:VAL:HG12	1:C:1784:LYS:HD3	1.44	0.96
1:E:2401:VAL:HG11	1:E:2432:ILE:HD11	1.53	0.91
1:C:1701:VAL:CG1	1:C:1784:LYS:HE2	2.01	0.90
1:E:1701:VAL:CG1	1:E:1784:LYS:HE2	2.01	0.90
1:D:2401:VAL:HG11	1:D:2432:ILE:HD11	1.53	0.90
1:D:1701:VAL:CG1	1:D:1784:LYS:HE2	2.01	0.89
1:C:2401:VAL:HG11	1:C:2432:ILE:HD11	1.53	0.88
1:C:2139:ASP:OD1	1:C:2143:THR:CG2	2.22	0.88
1:E:2139:ASP:OD1	1:E:2143:THR:CG2	2.22	0.88
1:C:1523:PHE:HZ	1:C:1686:HIS:HE2	0.87	0.87
1:C:1701:VAL:CG1	1:C:1784:LYS:CD	2.50	0.86
1:E:1523:PHE:HZ	1:E:1686:HIS:HE2	0.87	0.86
1:E:1701:VAL:CG1	1:E:1784:LYS:CD	2.50	0.86
1:D:2139:ASP:OD1	1:D:2143:THR:CG2	2.22	0.85
1:D:1523:PHE:HZ	1:D:1686:HIS:HE2	0.87	0.85
1:C:1217:ASN:O	1:C:1221:ILE:HG12	1.76	0.85
1:D:1191:LEU:HD13	1:D:1198:LEU:HD22	1.57	0.85
1:E:1191:LEU:HD13	1:E:1198:LEU:HD22	1.58	0.85
1:D:1282:TRP:O	1:D:1285:ILE:CG2	2.25	0.85
1:E:1217:ASN:O	1:E:1221:ILE:HG12	1.76	0.85
1:C:1282:TRP:O	1:C:1285:ILE:CG2	2.25	0.85
1:C:1701:VAL:CG1	1:C:1784:LYS:CE	2.56	0.84
1:E:1282:TRP:O	1:E:1285:ILE:CG2	2.25	0.84
1:D:1701:VAL:CG1	1:D:1784:LYS:CE	2.56	0.84
1:C:1191:LEU:HD13	1:C:1198:LEU:HD22	1.57	0.83
1:D:1217:ASN:O	1:D:1221:ILE:HG12	1.76	0.83
1:E:1701:VAL:CG1	1:E:1784:LYS:CE	2.56	0.83
1:D:1701:VAL:CG1	1:D:1784:LYS:CD	2.50	0.83
1:C:2370:HIS:ND1	1:C:2400:GLY:N	2.27	0.82
1:E:2370:HIS:ND1	1:E:2400:GLY:N	2.27	0.82
1:D:2370:HIS:ND1	1:D:2400:GLY:N	2.27	0.82
1:C:2348:SER:O	1:C:2352:ARG:HB3	1.80	0.81
1:D:2348:SER:O	1:D:2352:ARG:HB3	1.80	0.81
1:D:1721:THR:HG22	1:D:1962:PHE:HE1	1.46	0.81
1:C:1721:THR:HG22	1:C:1962:PHE:HE1	1.46	0.81
1:D:1315:GLN:HG3	1:D:1541:LEU:CD2	2.12	0.80
1:E:1315:GLN:HG3	1:E:1541:LEU:CD2	2.12	0.80
1:E:1721:THR:HG22	1:E:1962:PHE:HE1	1.46	0.80
1:E:2348:SER:O	1:E:2352:ARG:HB3	1.80	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:2139:ASP:O	1:C:2143:THR:HG23	1.81	0.79
1:D:2238:LEU:HD22	1:D:2430:TRP:CZ2	2.18	0.79
1:E:2139:ASP:O	1:E:2143:THR:HG23	1.81	0.79
1:C:1315:GLN:HG3	1:C:1541:LEU:CD2	2.12	0.79
1:D:2139:ASP:O	1:D:2143:THR:HG23	1.81	0.79
1:E:2238:LEU:HD13	1:E:2430:TRP:CG	2.18	0.79
1:C:1191:LEU:HD12	1:C:1198:LEU:CD2	2.12	0.79
1:C:2496:GLU:OE1	1:E:2534:ARG:HD2	1.83	0.79
1:D:2238:LEU:HD13	1:D:2430:TRP:CG	2.18	0.79
1:C:2280:ASP:OD2	1:C:2282:VAL:HG23	1.83	0.79
1:D:1191:LEU:HD12	1:D:1198:LEU:CD2	2.12	0.79
1:E:1191:LEU:HD12	1:E:1198:LEU:CD2	2.12	0.78
1:E:2280:ASP:OD2	1:E:2282:VAL:HG23	1.83	0.78
1:C:2238:LEU:HD22	1:C:2430:TRP:CZ2	2.18	0.78
1:E:2238:LEU:HD22	1:E:2430:TRP:CZ2	2.18	0.78
1:E:2225:ASP:HB3	1:E:2324:GLN:HB3	1.65	0.77
1:C:2238:LEU:HD13	1:C:2430:TRP:CG	2.18	0.77
1:D:2225:ASP:HB3	1:D:2324:GLN:HB3	1.65	0.77
1:D:2147:LEU:HB3	1:D:2151:ASN:HB2	1.66	0.77
1:D:2210:MET:SD	1:D:2466:ILE:HG21	2.25	0.77
1:E:2147:LEU:HB3	1:E:2151:ASN:HB2	1.66	0.77
1:C:2210:MET:SD	1:C:2466:ILE:HG21	2.25	0.77
1:D:2280:ASP:OD2	1:D:2282:VAL:HG23	1.83	0.77
1:E:2210:MET:SD	1:E:2466:ILE:HG21	2.25	0.77
1:C:1191:LEU:HD11	1:C:1198:LEU:CD2	2.16	0.76
1:C:1285:ILE:HG13	1:C:1289:PHE:HE2	1.51	0.76
1:C:2225:ASP:HB3	1:C:2324:GLN:HB3	1.65	0.76
1:C:2147:LEU:HB3	1:C:2151:ASN:HB2	1.66	0.76
1:E:1282:TRP:C	1:E:1285:ILE:HG22	2.10	0.76
1:E:1191:LEU:HD11	1:E:1198:LEU:CD2	2.16	0.75
1:D:1282:TRP:C	1:D:1285:ILE:HG22	2.10	0.75
1:D:2016:VAL:HG22	1:D:2069:THR:HG21	1.68	0.75
1:C:1282:TRP:C	1:C:1285:ILE:HG22	2.10	0.74
1:C:1744:LYS:HG2	1:C:1776:LEU:HD22	0.78	0.74
1:C:2016:VAL:HG22	1:C:2069:THR:HG21	1.68	0.74
1:D:1191:LEU:HD11	1:D:1198:LEU:CD2	2.16	0.74
1:D:1744:LYS:HG2	1:D:1776:LEU:HD22	0.78	0.74
1:E:2016:VAL:HG22	1:E:2069:THR:HG21	1.68	0.74
1:D:1701:VAL:HG11	1:D:1784:LYS:HD3	1.70	0.74
1:C:2268:LEU:HB3	1:C:2389:GLN:OE1	1.88	0.74
1:E:2268:LEU:HB3	1:E:2389:GLN:OE1	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1701:VAL:HG11	1:C:1784:LYS:HD3	1.70	0.73
1:D:2268:LEU:HB3	1:D:2389:GLN:OE1	1.88	0.73
1:E:1744:LYS:HG2	1:E:1776:LEU:HD22	0.78	0.73
1:E:1744:LYS:CG	1:E:1776:LEU:CD2	2.47	0.73
1:D:1695:ILE:HG12	1:D:1789:GLN:HE21	1.54	0.72
1:E:2145:THR:HG21	1:E:2152:TRP:HZ3	1.55	0.72
1:E:1285:ILE:HG13	1:E:1289:PHE:HE2	1.51	0.72
1:E:1695:ILE:HG12	1:E:1789:GLN:HE21	1.54	0.72
1:D:1744:LYS:CG	1:D:1776:LEU:CD2	2.47	0.71
1:C:2145:THR:HG21	1:C:2152:TRP:HZ3	1.55	0.71
1:E:1701:VAL:HG11	1:E:1784:LYS:HD3	1.70	0.71
1:D:2348:SER:HB3	1:D:2352:ARG:HE	1.55	0.71
1:E:1191:LEU:HD11	1:E:1198:LEU:HD21	1.72	0.70
1:C:1744:LYS:CG	1:C:1776:LEU:CD2	2.47	0.70
1:C:1695:ILE:HG12	1:C:1789:GLN:HE21	1.54	0.70
1:D:2145:THR:HG21	1:D:2152:TRP:HZ3	1.55	0.70
1:E:2348:SER:HB3	1:E:2352:ARG:HE	1.55	0.70
1:E:1191:LEU:CD1	1:E:1198:LEU:HD21	2.20	0.70
1:C:2348:SER:HB3	1:C:2352:ARG:HE	1.55	0.69
1:C:2129:PRO:O	1:D:2482:ARG:NH2	2.25	0.69
1:D:2129:PRO:O	1:E:2482:ARG:NH2	2.25	0.69
1:D:1191:LEU:CD1	1:D:1198:LEU:HD21	2.20	0.69
1:C:1191:LEU:CD1	1:C:1198:LEU:HD21	2.20	0.69
1:D:2238:LEU:HD22	1:D:2430:TRP:CE2	2.28	0.69
1:D:1191:LEU:HD11	1:D:1198:LEU:HD21	1.72	0.69
1:C:1191:LEU:HD11	1:C:1198:LEU:HD21	1.72	0.68
1:E:2238:LEU:HD22	1:E:2430:TRP:CE2	2.28	0.68
1:C:2238:LEU:HD22	1:C:2430:TRP:CE2	2.28	0.68
1:D:1285:ILE:HG13	1:D:1289:PHE:HE2	1.51	0.68
1:D:2238:LEU:HD12	1:D:2238:LEU:O	1.94	0.68
1:D:2367:VAL:HG22	1:D:2402:ARG:HD3	1.76	0.68
1:E:2367:VAL:HG22	1:E:2402:ARG:HD3	1.76	0.68
1:C:2238:LEU:HD12	1:C:2238:LEU:O	1.94	0.67
1:D:2534:ARG:HD2	1:E:2496:GLU:OE1	1.94	0.67
1:D:2326:ASP:HB3	1:D:2329:LYS:HE2	1.76	0.67
1:E:2238:LEU:HD12	1:E:2238:LEU:O	1.94	0.67
1:E:1213:LEU:HD23	1:E:1290:LEU:HD11	1.77	0.67
1:E:2326:ASP:HB3	1:E:2329:LYS:HE2	1.76	0.67
1:C:1315:GLN:CG	1:C:1541:LEU:HD22	2.25	0.66
1:C:1213:LEU:HD23	1:C:1290:LEU:HD11	1.77	0.66
1:C:2367:VAL:HG22	1:C:2402:ARG:HD3	1.76	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1148:SER:O	1:D:1152:MET:CB	2.44	0.66
1:E:1744:LYS:HE3	1:E:1776:LEU:CD1	2.22	0.66
1:C:2326:ASP:HB3	1:C:2329:LYS:HE2	1.76	0.66
1:E:1315:GLN:CG	1:E:1541:LEU:HD22	2.25	0.66
1:C:1148:SER:O	1:C:1152:MET:CB	2.44	0.66
1:D:2222:GLN:HG3	1:D:2282:VAL:HG21	1.78	0.66
1:E:1148:SER:O	1:E:1152:MET:CB	2.44	0.66
1:D:1744:LYS:HE3	1:D:1776:LEU:CD1	2.22	0.66
1:E:2222:GLN:HG3	1:E:2282:VAL:HG21	1.78	0.66
1:C:1315:GLN:HG3	1:C:1541:LEU:HD22	1.78	0.66
1:D:1315:GLN:CG	1:D:1541:LEU:HD22	2.25	0.65
1:E:1315:GLN:HG3	1:E:1541:LEU:HD22	1.78	0.65
1:E:1523:PHE:HZ	1:E:1686:HIS:CE1	2.13	0.65
1:E:1701:VAL:HG12	1:E:1784:LYS:CE	2.25	0.65
1:C:1701:VAL:HG12	1:C:1784:LYS:CE	2.25	0.65
1:D:1701:VAL:HG11	1:D:1784:LYS:CE	2.27	0.65
1:C:2534:ARG:HD2	1:D:2496:GLU:OE1	1.97	0.65
1:D:1213:LEU:HD23	1:D:1290:LEU:HD11	1.77	0.64
1:D:1315:GLN:HG3	1:D:1541:LEU:HD22	1.78	0.64
1:C:1523:PHE:HZ	1:C:1686:HIS:CE1	2.14	0.64
1:C:2222:GLN:HG3	1:C:2282:VAL:HG21	1.78	0.64
1:E:1282:TRP:HA	1:E:1285:ILE:HG22	1.79	0.64
1:D:1744:LYS:CD	1:D:1776:LEU:HD22	2.27	0.64
1:D:1523:PHE:HZ	1:D:1686:HIS:CE1	2.14	0.64
1:D:1701:VAL:HG12	1:D:1784:LYS:CE	2.25	0.64
1:C:1701:VAL:HG11	1:C:1784:LYS:CE	2.27	0.63
1:C:1282:TRP:HA	1:C:1285:ILE:HG22	1.79	0.63
1:C:2482:ARG:NH2	1:E:2129:PRO:O	2.32	0.63
1:E:1701:VAL:HG11	1:E:1784:LYS:CE	2.27	0.63
1:C:1744:LYS:HE3	1:C:1776:LEU:CD1	2.22	0.63
1:D:1282:TRP:HA	1:D:1285:ILE:HG22	1.79	0.62
1:E:2401:VAL:HG11	1:E:2432:ILE:CD1	2.28	0.62
1:E:1744:LYS:CD	1:E:1776:LEU:HD22	2.27	0.62
1:D:2401:VAL:HG11	1:D:2432:ILE:CD1	2.28	0.62
1:C:1744:LYS:CD	1:C:1776:LEU:HD22	2.27	0.62
1:D:1699:HIS:HA	1:D:1707:SER:HB2	1.82	0.62
1:D:2534:ARG:HH21	1:E:2540:ILE:HG12	1.64	0.62
1:E:1699:HIS:HA	1:E:1707:SER:HB2	1.82	0.61
1:C:1699:HIS:HA	1:C:1707:SER:HB2	1.82	0.61
1:E:2145:THR:HG21	1:E:2152:TRP:CZ3	2.36	0.61
1:C:2145:THR:HG21	1:C:2152:TRP:CZ3	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:2534:ARG:HH21	1:D:2540:ILE:HG12	1.64	0.61
1:C:2405:LEU:HD11	1:C:2428:GLU:HB3	1.83	0.61
1:D:2348:SER:O	1:D:2352:ARG:CB	2.49	0.61
1:E:2405:LEU:HD11	1:E:2428:GLU:HB3	1.83	0.61
1:D:2405:LEU:HD11	1:D:2428:GLU:HB3	1.83	0.61
1:C:2144:ASP:O	1:C:2145:THR:HG23	2.01	0.60
1:D:2144:ASP:O	1:D:2145:THR:HG23	2.01	0.60
1:E:2144:ASP:O	1:E:2145:THR:HG23	2.01	0.60
1:C:1148:SER:O	1:C:1152:MET:HB3	2.01	0.60
1:D:1148:SER:O	1:D:1152:MET:HB3	2.01	0.60
1:E:1148:SER:O	1:E:1152:MET:HB3	2.01	0.59
1:C:1285:ILE:HG13	1:C:1289:PHE:CD2	2.36	0.59
1:C:1672:THR:O	1:C:1675:LEU:HB3	2.02	0.59
1:D:1672:THR:O	1:D:1675:LEU:HB3	2.02	0.59
1:D:1285:ILE:HG13	1:D:1289:PHE:CD2	2.36	0.59
1:D:2145:THR:HG21	1:D:2152:TRP:CZ3	2.36	0.59
1:E:1672:THR:O	1:E:1675:LEU:HB3	2.02	0.59
1:E:2348:SER:O	1:E:2352:ARG:CB	2.49	0.59
1:E:1285:ILE:HG13	1:E:1289:PHE:CD2	2.36	0.59
1:C:1988:ILE:HD11	1:C:2031:MET:SD	2.44	0.58
1:E:1523:PHE:CZ	1:E:1686:HIS:CE1	2.91	0.58
1:E:1988:ILE:HD11	1:E:2031:MET:SD	2.44	0.58
1:C:2348:SER:O	1:C:2352:ARG:CB	2.49	0.58
1:C:2370:HIS:CE1	1:C:2435:GLN:OE1	2.57	0.58
1:C:2401:VAL:HG11	1:C:2432:ILE:CD1	2.28	0.58
1:D:1523:PHE:CZ	1:D:1686:HIS:CE1	2.91	0.58
1:E:2139:ASP:O	1:E:2143:THR:CG2	2.51	0.58
1:E:2370:HIS:CE1	1:E:2435:GLN:OE1	2.57	0.57
1:C:1282:TRP:CA	1:C:1285:ILE:HG22	2.34	0.57
1:D:1282:TRP:CA	1:D:1285:ILE:HG22	2.34	0.57
1:E:2280:ASP:OD2	1:E:2282:VAL:CG2	2.53	0.57
1:D:2370:HIS:CE1	1:D:2435:GLN:OE1	2.57	0.57
1:E:1494:MET:HG3	1:E:1497:ARG:NH1	2.20	0.57
1:D:1988:ILE:HD11	1:D:2031:MET:SD	2.44	0.57
1:E:1273:LEU:HD12	1:E:1274:PRO:HD2	1.87	0.57
1:D:1701:VAL:HG11	1:D:1784:LYS:CD	2.31	0.57
1:E:1282:TRP:CA	1:E:1285:ILE:HG22	2.34	0.57
1:C:1523:PHE:CZ	1:C:1686:HIS:CE1	2.91	0.56
1:D:2139:ASP:O	1:D:2143:THR:CG2	2.51	0.56
1:C:1494:MET:HG3	1:C:1497:ARG:NH1	2.20	0.56
1:E:2016:VAL:HG22	1:E:2069:THR:CG2	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:2040:ARG:HH22	1:C:2144:ASP:CG	2.14	0.56
1:D:1721:THR:HG22	1:D:1962:PHE:CE1	2.36	0.56
1:D:2280:ASP:OD2	1:D:2282:VAL:CG2	2.53	0.56
1:E:1721:THR:HG22	1:E:1962:PHE:CE1	2.36	0.56
1:D:2040:ARG:HH22	1:D:2144:ASP:CG	2.14	0.56
1:E:1349:MET:HE3	1:E:2518:GLU:HG3	1.88	0.56
1:C:2280:ASP:OD2	1:C:2282:VAL:CG2	2.53	0.56
1:D:2016:VAL:HG22	1:D:2069:THR:CG2	2.36	0.56
1:D:1494:MET:HG3	1:D:1497:ARG:NH1	2.20	0.55
1:D:1805:LEU:HD23	1:D:1806:TRP:N	2.21	0.55
1:C:1273:LEU:HD12	1:C:1274:PRO:HD2	1.87	0.55
1:E:1805:LEU:HD23	1:E:1806:TRP:N	2.21	0.55
1:C:2139:ASP:O	1:C:2143:THR:CG2	2.51	0.55
1:C:2077:ALA:O	1:C:2080:GLN:HB3	2.06	0.55
1:E:2077:ALA:O	1:E:2080:GLN:HB3	2.07	0.55
1:C:2016:VAL:HG22	1:C:2069:THR:CG2	2.36	0.55
1:E:1969:LYS:HG2	1:E:1970:TYR:N	2.22	0.55
1:C:1148:SER:O	1:C:1152:MET:HB2	2.07	0.55
1:E:2040:ARG:HH22	1:E:2144:ASP:CG	2.14	0.55
1:C:1349:MET:HE3	1:C:2518:GLU:HG3	1.88	0.55
1:D:1148:SER:O	1:D:1152:MET:HB2	2.07	0.55
1:C:2016:VAL:HG11	1:C:2076:ASN:ND2	2.22	0.55
1:D:1349:MET:HE3	1:D:2518:GLU:HG3	1.88	0.55
1:C:1158:PHE:HB3	1:C:1302:PHE:CD1	2.42	0.55
1:E:1158:PHE:HB3	1:E:1302:PHE:CD1	2.42	0.54
1:D:2016:VAL:HG11	1:D:2076:ASN:ND2	2.22	0.54
1:D:1158:PHE:HB3	1:D:1302:PHE:CD1	2.42	0.54
1:D:1273:LEU:HD12	1:D:1274:PRO:HD2	1.87	0.54
1:E:1148:SER:O	1:E:1152:MET:HB2	2.07	0.54
1:C:1805:LEU:HD23	1:C:1806:TRP:N	2.21	0.54
1:D:1285:ILE:CG1	1:D:1289:PHE:CE2	2.80	0.54
1:D:2077:ALA:O	1:D:2080:GLN:HB3	2.07	0.54
1:C:2144:ASP:O	1:C:2145:THR:CG2	2.56	0.54
1:E:2016:VAL:HG11	1:E:2076:ASN:ND2	2.22	0.54
1:D:2401:VAL:CG1	1:D:2432:ILE:HG13	2.38	0.54
1:D:1969:LYS:HG2	1:D:1970:TYR:N	2.22	0.53
1:C:1320:PHE:HB2	1:C:1804:GLY:HA2	1.90	0.53
1:C:1969:LYS:HG2	1:C:1970:TYR:N	2.22	0.53
1:C:2519:LEU:O	1:C:2522:GLU:HB3	2.08	0.53
1:E:2401:VAL:CG1	1:E:2432:ILE:HG13	2.38	0.53
1:E:2519:LEU:O	1:E:2522:GLU:HB3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:2425:ASP:HB2	1:E:2410:VAL:HG13	1.91	0.53
1:E:1320:PHE:HB2	1:E:1804:GLY:HA2	1.90	0.53
1:E:1805:LEU:HD23	1:E:1806:TRP:C	2.34	0.53
1:E:2144:ASP:C	1:E:2145:THR:HG23	2.34	0.53
1:C:1805:LEU:HD23	1:C:1806:TRP:C	2.34	0.53
1:D:2519:LEU:O	1:D:2522:GLU:HB3	2.08	0.53
1:E:2370:HIS:HE1	1:E:2435:GLN:OE1	1.92	0.53
1:C:2370:HIS:HE1	1:C:2435:GLN:OE1	1.92	0.53
1:C:2401:VAL:CG1	1:C:2432:ILE:HG13	2.38	0.53
1:D:1320:PHE:HB2	1:D:1804:GLY:HA2	1.90	0.52
1:D:1805:LEU:HD23	1:D:1806:TRP:C	2.34	0.52
1:D:2144:ASP:O	1:D:2145:THR:CG2	2.56	0.52
1:E:2144:ASP:O	1:E:2145:THR:CG2	2.56	0.52
1:C:1282:TRP:HA	1:C:1285:ILE:CG2	2.40	0.52
1:C:2144:ASP:C	1:C:2145:THR:HG23	2.34	0.52
1:E:1147:ARG:O	1:E:1151:ASP:HB2	2.10	0.52
1:C:2238:LEU:HB2	1:C:2430:TRP:CD1	2.45	0.52
1:D:1147:ARG:O	1:D:1151:ASP:HB2	2.10	0.52
1:D:1176:ARG:HH22	1:D:1221:ILE:CD1	2.22	0.52
1:D:2144:ASP:C	1:D:2145:THR:HG23	2.34	0.52
1:D:2238:LEU:HB2	1:D:2430:TRP:CD1	2.45	0.52
1:C:1285:ILE:CG1	1:C:1289:PHE:CE2	2.80	0.52
1:C:1147:ARG:O	1:C:1151:ASP:HB2	2.10	0.52
1:E:1701:VAL:HG12	1:E:1784:LYS:HE2	1.84	0.52
1:E:1805:LEU:HD23	1:E:1806:TRP:O	2.10	0.52
1:C:1176:ARG:HH22	1:C:1221:ILE:CD1	2.22	0.52
1:C:2496:GLU:OE1	1:E:2534:ARG:CD	2.57	0.51
1:E:1176:ARG:HH22	1:E:1221:ILE:CD1	2.22	0.51
1:C:1805:LEU:HD23	1:C:1806:TRP:O	2.10	0.51
1:D:1805:LEU:HD23	1:D:1806:TRP:O	2.10	0.51
1:D:1239:GLN:HB3	1:D:1269:ARG:HH22	1.75	0.51
1:E:1239:GLN:HB3	1:E:1269:ARG:HH22	1.76	0.51
1:E:1282:TRP:HA	1:E:1285:ILE:CG2	2.40	0.51
1:D:2370:HIS:HE1	1:D:2435:GLN:OE1	1.92	0.51
1:D:1282:TRP:HA	1:D:1285:ILE:CG2	2.40	0.51
1:E:1701:VAL:HG11	1:E:1784:LYS:CD	2.31	0.51
1:D:1264:MET:SD	1:D:1273:LEU:HD22	2.51	0.50
1:D:2123:GLN:O	1:D:2127:LEU:HB2	2.12	0.50
1:E:2238:LEU:HB2	1:E:2430:TRP:CD1	2.45	0.50
1:C:1239:GLN:HB3	1:C:1269:ARG:HH22	1.75	0.50
1:C:2123:GLN:O	1:C:2127:LEU:HB2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:1274:PRO:HB2	1:E:1277:GLU:HB2	1.94	0.50
1:E:1191:LEU:HD12	1:E:1198:LEU:HD23	1.91	0.50
1:D:1191:LEU:HD12	1:D:1198:LEU:HD23	1.91	0.50
1:C:1264:MET:SD	1:C:1273:LEU:HD22	2.51	0.49
1:D:2238:LEU:HD13	1:D:2430:TRP:CE3	2.45	0.49
1:C:1207:LEU:H	1:C:1207:LEU:HD23	1.77	0.49
1:C:2466:ILE:HG22	1:C:2466:ILE:O	2.12	0.49
1:D:1701:VAL:HG12	1:D:1784:LYS:HE2	1.84	0.49
1:C:1701:VAL:HG12	1:C:1784:LYS:HE2	1.84	0.49
1:D:2282:VAL:HG12	1:D:2283:THR:N	2.28	0.49
1:D:2401:VAL:CG1	1:D:2432:ILE:CG1	2.91	0.49
1:E:1264:MET:SD	1:E:1273:LEU:HD22	2.51	0.49
1:E:2401:VAL:CG1	1:E:2432:ILE:CG1	2.91	0.49
1:E:1207:LEU:H	1:E:1207:LEU:HD23	1.77	0.49
1:C:1143:PHE:HD1	1:C:1144:ILE:HG23	1.78	0.49
1:C:1721:THR:HG22	1:C:1962:PHE:CE1	2.36	0.49
1:C:2425:ASP:CB	1:E:2410:VAL:HG13	2.43	0.49
1:E:2282:VAL:HG12	1:E:2283:THR:N	2.27	0.49
1:C:1274:PRO:HB2	1:C:1277:GLU:HB2	1.94	0.49
1:C:2282:VAL:HG12	1:C:2283:THR:N	2.28	0.49
1:E:2123:GLN:O	1:E:2127:LEU:HB2	2.12	0.49
1:E:2370:HIS:HA	1:E:2400:GLY:HA2	1.95	0.49
1:E:1143:PHE:HD1	1:E:1144:ILE:HG23	1.78	0.48
1:C:1356:GLN:HG3	1:C:2521:LEU:HD11	1.95	0.48
1:D:1207:LEU:HD23	1:D:1207:LEU:H	1.77	0.48
1:D:2466:ILE:O	1:D:2466:ILE:HG22	2.12	0.48
1:D:1697:LEU:HD21	1:D:1785:TYR:CE1	2.49	0.48
1:E:1356:GLN:HG3	1:E:2521:LEU:HD11	1.95	0.48
1:E:1170:PHE:HB2	1:E:1188:CYS:SG	2.54	0.48
1:C:1170:PHE:HB2	1:C:1188:CYS:SG	2.54	0.48
1:C:1697:LEU:HD21	1:C:1785:TYR:CE1	2.49	0.48
1:D:1274:PRO:HB2	1:D:1277:GLU:HB2	1.94	0.48
1:E:2466:ILE:HG22	1:E:2466:ILE:O	2.12	0.48
1:E:1697:LEU:HD21	1:E:1785:TYR:CE1	2.49	0.48
1:E:2519:LEU:H	1:E:2519:LEU:HD22	1.79	0.48
1:D:1143:PHE:HD1	1:D:1144:ILE:HG23	1.78	0.48
1:C:2401:VAL:CG1	1:C:2432:ILE:CG1	2.91	0.48
1:D:1170:PHE:HB2	1:D:1188:CYS:SG	2.53	0.48
1:D:1356:GLN:HG3	1:D:2521:LEU:HD11	1.95	0.48
1:C:2370:HIS:HA	1:C:2400:GLY:HA2	1.95	0.47
1:D:2434:LEU:HG	1:D:2444:LEU:HD22	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:2373:PRO:HG2	1:E:2376:ILE:HD11	1.96	0.47
1:D:2370:HIS:HA	1:D:2400:GLY:HA2	1.95	0.47
1:E:1721:THR:OG1	1:E:1725:PRO:HB3	2.14	0.47
1:C:2339:LYS:HB3	1:C:2339:LYS:HE2	1.36	0.47
1:C:2373:PRO:HG2	1:C:2376:ILE:HD11	1.96	0.47
1:D:1150:LEU:HD23	1:D:1150:LEU:H	1.79	0.47
1:E:2401:VAL:HG13	1:E:2432:ILE:CG1	2.44	0.47
1:D:2519:LEU:HD22	1:D:2519:LEU:H	1.79	0.47
1:C:1721:THR:OG1	1:C:1725:PRO:HB3	2.14	0.47
1:C:2434:LEU:HG	1:C:2444:LEU:HD22	1.96	0.47
1:C:1229:LEU:HD22	1:C:1258:TYR:HD1	1.80	0.47
1:C:1701:VAL:HG11	1:C:1784:LYS:CD	2.32	0.47
1:C:2401:VAL:HG13	1:C:2432:ILE:CG1	2.44	0.47
1:C:2482:ARG:CG	1:C:2486:SER:OG	2.63	0.47
1:D:1797:ARG:HG3	1:D:1806:TRP:CZ3	2.50	0.47
1:E:2116:HIS:CD2	1:E:2117:LEU:HD22	2.50	0.47
1:E:2434:LEU:HG	1:E:2444:LEU:HD22	1.96	0.47
1:D:1721:THR:OG1	1:D:1725:PRO:HB3	2.14	0.47
1:D:2401:VAL:HG13	1:D:2432:ILE:CG1	2.44	0.47
1:E:1150:LEU:H	1:E:1150:LEU:HD23	1.79	0.47
1:C:1701:VAL:HG13	1:C:1784:LYS:HE2	1.93	0.47
1:D:2482:ARG:CG	1:D:2486:SER:OG	2.63	0.47
1:E:2482:ARG:CG	1:E:2486:SER:OG	2.63	0.47
1:C:1797:ARG:HG3	1:C:1806:TRP:CZ3	2.50	0.47
1:D:1229:LEU:HD22	1:D:1258:TYR:HD1	1.80	0.47
1:E:1797:ARG:HG3	1:E:1806:TRP:CZ3	2.50	0.47
1:D:2116:HIS:CD2	1:D:2117:LEU:HD22	2.50	0.46
1:E:2079:ALA:O	1:E:2082:TRP:HB3	2.15	0.46
1:C:1150:LEU:H	1:C:1150:LEU:HD23	1.79	0.46
1:D:2373:PRO:HG2	1:D:2376:ILE:HD11	1.97	0.46
1:E:1229:LEU:HD22	1:E:1258:TYR:HD1	1.80	0.46
1:C:1191:LEU:HD12	1:C:1198:LEU:HD23	1.91	0.46
1:C:2519:LEU:HD22	1:C:2519:LEU:H	1.79	0.46
1:E:2408:GLU:HG3	1:E:2429:TRP:HH2	1.80	0.46
1:D:1744:LYS:CB	1:D:1776:LEU:CD2	2.94	0.46
1:D:2079:ALA:O	1:D:2082:TRP:HB3	2.15	0.46
1:C:2116:HIS:CD2	1:C:2117:LEU:HD22	2.50	0.46
1:D:1264:MET:HG2	1:D:1267:ARG:HH21	1.81	0.46
1:E:2093:SER:O	1:E:2097:ILE:HG12	2.16	0.46
1:C:1264:MET:HG2	1:C:1267:ARG:HH21	1.81	0.46
1:C:2079:ALA:O	1:C:2082:TRP:HB3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1706:ALA:HB2	1:D:1740:MET:HG2	1.98	0.46
1:D:2408:GLU:HG3	1:D:2429:TRP:HH2	1.80	0.46
1:D:2401:VAL:HG13	1:D:2432:ILE:HG13	1.98	0.46
1:E:1285:ILE:CG1	1:E:1289:PHE:CE2	2.80	0.46
1:C:2093:SER:O	1:C:2097:ILE:HG12	2.16	0.46
1:C:1722:ILE:HG13	1:C:1724:ARG:H	1.81	0.45
1:D:1722:ILE:HG13	1:D:1724:ARG:H	1.82	0.45
1:E:1744:LYS:CB	1:E:1776:LEU:CD2	2.94	0.45
1:E:2401:VAL:HG13	1:E:2432:ILE:HG13	1.98	0.45
1:C:1176:ARG:HH22	1:C:1221:ILE:HD11	1.81	0.45
1:E:1259:TYR:OH	1:E:1273:LEU:HD11	2.17	0.45
1:E:1264:MET:HG2	1:E:1267:ARG:HH21	1.81	0.45
1:C:1706:ALA:HB2	1:C:1740:MET:HG2	1.98	0.45
1:E:1176:ARG:HH22	1:E:1221:ILE:HD11	1.81	0.45
1:C:1744:LYS:CB	1:C:1776:LEU:CD2	2.94	0.45
1:C:2126:ARG:HB2	1:C:2132:VAL:HG21	1.99	0.45
1:C:2238:LEU:HD13	1:C:2430:TRP:CE3	2.45	0.45
1:E:2297:SER:HB2	1:E:2300:SER:HB3	1.99	0.45
1:D:2297:SER:HB2	1:D:2300:SER:HB3	1.99	0.45
1:C:2401:VAL:HG13	1:C:2432:ILE:HG13	1.98	0.45
1:D:1176:ARG:HH22	1:D:1221:ILE:HD11	1.81	0.45
1:D:2093:SER:O	1:D:2097:ILE:HG12	2.16	0.45
1:E:1285:ILE:HG23	1:E:1286:CYS:N	2.32	0.45
1:E:1706:ALA:HB2	1:E:1740:MET:HG2	1.98	0.45
1:C:1315:GLN:CG	1:C:1541:LEU:CD2	2.86	0.45
1:D:1285:ILE:HG23	1:D:1286:CYS:N	2.32	0.45
1:C:2296:ILE:O	1:E:2293:LEU:HB2	2.16	0.45
1:C:2540:ILE:HG12	1:E:2534:ARG:HH21	1.81	0.45
1:D:2126:ARG:HB2	1:D:2132:VAL:HG21	1.99	0.45
1:D:2185:LYS:O	1:D:2189:TYR:HB2	2.17	0.45
1:E:2126:ARG:HB2	1:E:2132:VAL:HG21	1.99	0.45
1:E:2185:LYS:O	1:E:2189:TYR:HB2	2.17	0.45
1:C:2408:GLU:HG3	1:C:2429:TRP:HH2	1.80	0.45
1:C:2506:LEU:O	1:C:2509:ASP:HB3	2.17	0.45
1:D:2020:PHE:HD1	1:D:2020:PHE:HA	1.70	0.45
1:D:2248:ILE:HG23	1:D:2282:VAL:HG11	1.99	0.44
1:D:2506:LEU:O	1:D:2509:ASP:HB3	2.17	0.44
1:E:1722:ILE:HG13	1:E:1724:ARG:H	1.81	0.44
1:D:2503:ILE:HG12	1:D:2532:LEU:HD21	1.99	0.44
1:C:2297:SER:HB2	1:C:2300:SER:HB3	1.99	0.44
1:D:1238:MET:HA	1:D:1242:PHE:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:2339:LYS:HE2	1:D:2339:LYS:HB3	1.36	0.44
1:D:2401:VAL:CG1	1:D:2432:ILE:HD11	2.37	0.44
1:E:2232:LEU:HD23	1:E:2304:MET:HE3	1.99	0.44
1:C:2185:LYS:O	1:C:2189:TYR:HB2	2.17	0.44
1:C:2248:ILE:HG23	1:C:2282:VAL:HG11	1.99	0.44
1:E:2238:LEU:HD13	1:E:2430:TRP:CE3	2.45	0.44
1:C:1251:LEU:HD12	1:C:1251:LEU:HA	1.88	0.44
1:D:1259:TYR:OH	1:D:1273:LEU:HD11	2.17	0.44
1:E:2248:ILE:HG23	1:E:2282:VAL:HG11	1.99	0.44
1:E:2482:ARG:HG3	1:E:2486:SER:OG	2.17	0.44
1:C:2503:ILE:HG12	1:C:2532:LEU:HD21	1.99	0.44
1:E:2506:LEU:O	1:E:2509:ASP:HB3	2.17	0.44
1:D:2086:LYS:O	1:D:2089:TYR:HB3	2.18	0.44
1:C:2232:LEU:HD23	1:C:2304:MET:HE3	1.99	0.44
1:D:1184:TYR:HE1	1:D:1213:LEU:HD21	1.82	0.44
1:C:1285:ILE:HG23	1:C:1286:CYS:N	2.32	0.43
1:C:1769:PHE:HB3	1:C:1772:ARG:HB3	2.00	0.43
1:C:2297:SER:O	1:C:2298:PRO:C	2.61	0.43
1:C:2323:PHE:HB2	1:C:2336:THR:HG23	2.00	0.43
1:D:2482:ARG:HG3	1:D:2486:SER:OG	2.17	0.43
1:E:1710:LEU:HB2	1:E:1711:PRO:HD3	2.00	0.43
1:E:2503:ILE:HG12	1:E:2532:LEU:HD21	1.99	0.43
1:C:2466:ILE:O	1:C:2466:ILE:CG2	2.67	0.43
1:E:1184:TYR:HE1	1:E:1213:LEU:HD21	1.82	0.43
1:C:1259:TYR:OH	1:C:1273:LEU:HD11	2.17	0.43
1:D:1710:LEU:HB2	1:D:1711:PRO:HD3	2.00	0.43
1:C:1238:MET:HA	1:C:1242:PHE:HB2	1.99	0.43
1:C:2496:GLU:OE1	1:C:2496:GLU:O	2.37	0.43
1:E:1238:MET:HA	1:E:1242:PHE:HB2	1.99	0.43
1:E:2496:GLU:OE1	1:E:2496:GLU:O	2.37	0.43
1:C:1184:TYR:HE1	1:C:1213:LEU:HD21	1.82	0.43
1:C:1793:LEU:HD23	1:C:1793:LEU:HA	1.91	0.43
1:D:2323:PHE:HB2	1:D:2336:THR:HG23	2.00	0.43
1:D:2297:SER:O	1:D:2298:PRO:C	2.61	0.43
1:C:2086:LYS:O	1:C:2089:TYR:HB3	2.18	0.43
1:C:2482:ARG:HG3	1:C:2486:SER:OG	2.17	0.43
1:D:2496:GLU:OE1	1:D:2496:GLU:O	2.37	0.43
1:D:2466:ILE:O	1:D:2466:ILE:CG2	2.67	0.43
1:C:2315:ILE:HD12	1:C:2315:ILE:HA	1.95	0.43
1:D:2321:TRP:NE1	1:D:2338:GLU:HB3	2.34	0.43
1:E:2297:SER:O	1:E:2298:PRO:C	2.61	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1710:LEU:HB2	1:C:1711:PRO:HD3	2.00	0.43
1:D:2232:LEU:HD23	1:D:2304:MET:HE3	1.99	0.43
1:E:2254:GLN:O	1:E:2258:GLU:HG2	2.19	0.43
1:E:2321:TRP:NE1	1:E:2338:GLU:HB3	2.34	0.43
1:D:2254:GLN:O	1:D:2258:GLU:HG2	2.19	0.42
1:E:2016:VAL:HG22	1:E:2069:THR:CB	2.49	0.42
1:D:1195:GLY:HA2	1:D:1198:LEU:HB3	2.02	0.42
1:D:1769:PHE:HB3	1:D:1772:ARG:HB3	2.00	0.42
1:E:1143:PHE:CD1	1:E:1144:ILE:HG23	2.54	0.42
1:E:1769:PHE:HB3	1:E:1772:ARG:HB3	2.00	0.42
1:E:2323:PHE:HB2	1:E:2336:THR:HG23	2.00	0.42
1:C:2282:VAL:CG1	1:C:2283:THR:N	2.83	0.42
1:E:1195:GLY:HA2	1:E:1198:LEU:HB3	2.02	0.42
1:E:2370:HIS:ND1	1:E:2400:GLY:CA	2.82	0.42
1:D:1759:VAL:HG13	1:D:1762:ARG:HH21	1.85	0.42
1:E:2086:LYS:O	1:E:2089:TYR:HB3	2.18	0.42
1:C:1759:VAL:HG13	1:C:1762:ARG:HH21	1.85	0.42
1:D:1143:PHE:CD1	1:D:1144:ILE:HG23	2.54	0.42
1:E:2466:ILE:O	1:E:2466:ILE:CG2	2.67	0.42
1:C:2016:VAL:HG22	1:C:2069:THR:CB	2.49	0.42
1:E:1769:PHE:CD1	1:E:1771:PRO:HD2	2.55	0.42
1:E:2238:LEU:O	1:E:2238:LEU:CD1	2.66	0.42
1:C:2321:TRP:NE1	1:C:2338:GLU:HB3	2.34	0.42
1:C:2370:HIS:ND1	1:C:2400:GLY:CA	2.82	0.42
1:E:2282:VAL:CG1	1:E:2283:THR:N	2.83	0.42
1:D:2370:HIS:ND1	1:D:2400:GLY:CA	2.82	0.42
1:E:2405:LEU:HD12	1:E:2429:TRP:O	2.20	0.42
1:C:1769:PHE:CD1	1:C:1771:PRO:HD2	2.55	0.42
1:C:2192:GLY:HA3	1:E:2141:VAL:HG21	2.01	0.42
1:C:2277:SER:N	1:C:2278:PRO:HD2	2.35	0.42
1:D:2282:VAL:CG1	1:D:2283:THR:N	2.83	0.42
1:C:1969:LYS:CG	1:C:1970:TYR:N	2.83	0.42
1:D:2016:VAL:HG22	1:D:2069:THR:CB	2.49	0.42
1:D:2277:SER:N	1:D:2278:PRO:HD2	2.35	0.41
1:E:2277:SER:N	1:E:2278:PRO:HD2	2.35	0.41
1:D:2238:LEU:O	1:D:2238:LEU:CD1	2.66	0.41
1:E:1969:LYS:CG	1:E:1970:TYR:N	2.83	0.41
1:C:1143:PHE:CD1	1:C:1144:ILE:HG23	2.54	0.41
1:C:1195:GLY:HA2	1:C:1198:LEU:HB3	2.02	0.41
1:C:2254:GLN:O	1:C:2258:GLU:HG2	2.19	0.41
1:C:2405:LEU:HD12	1:C:2429:TRP:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:1170:PHE:CD1	1:E:1787:LEU:HD13	2.55	0.41
1:E:1759:VAL:HG13	1:E:1762:ARG:HH21	1.85	0.41
1:E:2037:LEU:HD23	1:E:2046:LYS:HA	2.03	0.41
1:C:1170:PHE:CD1	1:C:1787:LEU:HD13	2.55	0.41
1:C:1670:GLY:HA2	1:C:1674:ARG:CZ	2.50	0.41
1:E:1670:GLY:HA2	1:E:1674:ARG:CZ	2.50	0.41
1:D:2174:LYS:HD2	1:D:2175:TYR:CE2	2.56	0.41
1:D:2405:LEU:HD13	1:D:2430:TRP:CE2	2.56	0.41
1:E:2174:LYS:HD2	1:E:2175:TYR:CE2	2.56	0.41
1:C:1346:LYS:HD3	1:C:1346:LYS:HA	1.96	0.41
1:C:2401:VAL:CG1	1:C:2432:ILE:HD11	2.37	0.41
1:D:1170:PHE:CD1	1:D:1787:LEU:HD13	2.55	0.41
1:D:2037:LEU:HD23	1:D:2046:LYS:HA	2.03	0.41
1:D:2530:ILE:HD13	1:D:2530:ILE:HA	1.91	0.41
1:E:2103:THR:HG23	1:E:2104:ARG:HG3	2.02	0.41
1:C:2254:GLN:CD	1:C:2254:GLN:H	2.29	0.41
1:E:1724:ARG:HH22	1:E:1805:LEU:HB3	1.86	0.41
1:C:1551:VAL:HG12	1:C:1554:GLY:H	1.86	0.41
1:C:1724:ARG:HH22	1:C:1805:LEU:HB3	1.86	0.41
1:D:1670:GLY:HA2	1:D:1674:ARG:CZ	2.50	0.41
1:E:1224:LYS:HE3	1:E:1249:PHE:HB3	2.03	0.41
1:C:2368:ILE:HG13	1:C:2369:PRO:HD2	2.03	0.41
1:D:2103:THR:HG23	1:D:2104:ARG:HG3	2.02	0.41
1:D:2106:LEU:HD23	1:D:2515:GLU:HG2	2.03	0.41
1:E:1251:LEU:HD12	1:E:1251:LEU:HA	1.88	0.41
1:E:2405:LEU:HD13	1:E:2430:TRP:CE2	2.56	0.41
1:C:2103:THR:HG23	1:C:2104:ARG:HG3	2.02	0.40
1:C:2321:TRP:CE2	1:C:2338:GLU:HB3	2.57	0.40
1:D:1749:PHE:HD1	1:D:1749:PHE:HA	1.80	0.40
1:E:1225:ASN:O	1:E:1229:LEU:HG	2.21	0.40
1:E:1553:ARG:H	1:E:1553:ARG:HG2	1.74	0.40
1:C:2020:PHE:HD1	1:C:2020:PHE:HA	1.71	0.40
1:C:2016:VAL:N	1:C:2017:PRO:HD2	2.36	0.40
1:C:2037:LEU:HD23	1:C:2046:LYS:HA	2.03	0.40
1:C:2174:LYS:HD2	1:C:2175:TYR:CE2	2.56	0.40
1:D:1654:LEU:HG	1:D:1655:HIS:N	2.36	0.40
1:D:1969:LYS:CG	1:D:1970:TYR:N	2.83	0.40
1:D:2254:GLN:H	1:D:2254:GLN:CD	2.29	0.40
1:D:2405:LEU:HD12	1:D:2429:TRP:O	2.20	0.40
1:E:2016:VAL:N	1:E:2017:PRO:HD2	2.36	0.40
1:D:1521:ARG:HB2	1:D:1528:ARG:NH1	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1551:VAL:HG12	1:D:1554:GLY:H	1.86	0.40
1:D:2321:TRP:CE2	1:D:2338:GLU:HB3	2.57	0.40
1:D:2368:ILE:HD12	1:D:2368:ILE:HA	1.96	0.40
1:E:1521:ARG:HB2	1:E:1528:ARG:NH1	2.37	0.40
1:E:2321:TRP:CE2	1:E:2338:GLU:HB3	2.57	0.40
1:D:1402:ASP:O	1:E:2179:LYS:HB3	2.21	0.40
1:D:1769:PHE:CD1	1:D:1771:PRO:HD2	2.55	0.40
1:D:2016:VAL:N	1:D:2017:PRO:HD2	2.36	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	C	1313/2561 (51%)	1197 (91%)	110 (8%)	6 (0%)	24	56
1	D	1313/2561 (51%)	1197 (91%)	110 (8%)	6 (0%)	24	56
1	E	1313/2561 (51%)	1197 (91%)	110 (8%)	6 (0%)	24	56
All	All	3939/7683 (51%)	3591 (91%)	330 (8%)	18 (0%)	26	56

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	919	PRO
1	D	919	PRO
1	E	919	PRO
1	C	1177	ILE
1	D	1177	ILE
1	E	1177	ILE
1	C	2144	ASP
1	D	2144	ASP
1	E	2144	ASP

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Mol	Chain	Res	Type
1	C	1081	TYR
1	D	1081	TYR
1	E	1081	TYR
1	C	918	TYR
1	D	918	TYR
1	E	918	TYR
1	C	842	TYR
1	D	842	TYR
1	E	842	TYR

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	C	940/2260 (42%)	932 (99%)	8 (1%)	70	75
1	D	940/2260 (42%)	932 (99%)	8 (1%)	70	75
1	E	940/2260 (42%)	932 (99%)	8 (1%)	70	75
All	All	2820/6780 (42%)	2796 (99%)	24 (1%)	68	75

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	1184	TYR
1	C	1258	TYR
1	C	1300	HIS
1	C	1699	HIS
1	C	1768	TYR
1	C	1971	ARG
1	C	2020	PHE
1	C	2339	LYS
1	D	1184	TYR
1	D	1258	TYR
1	D	1300	HIS
1	D	1699	HIS
1	D	1768	TYR

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Mol	Chain	Res	Type
1	D	1971	ARG
1	D	2020	PHE
1	D	2339	LYS
1	E	1184	TYR
1	E	1258	TYR
1	E	1300	HIS
1	E	1699	HIS
1	E	1768	TYR
1	E	1971	ARG
1	E	2020	PHE
1	E	2339	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	1225	ASN
1	C	1241	ASN
1	C	1348	GLN
1	C	2076	ASN
1	C	2116	HIS
1	C	2310	ASN
1	C	2347	ASN
1	C	2385	ASN
1	C	2490	HIS
1	D	1225	ASN
1	D	1241	ASN
1	D	1348	GLN
1	D	2076	ASN
1	D	2080	GLN
1	D	2116	HIS
1	D	2161	ASN
1	D	2310	ASN
1	D	2347	ASN
1	D	2385	ASN
1	E	1225	ASN
1	E	1241	ASN
1	E	1344	GLN
1	E	1799	GLN
1	E	2076	ASN
1	E	2080	GLN
1	E	2116	HIS
1	E	2310	ASN

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Mol	Chain	Res	Type
1	E	2347	ASN
1	E	2385	ASN

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.

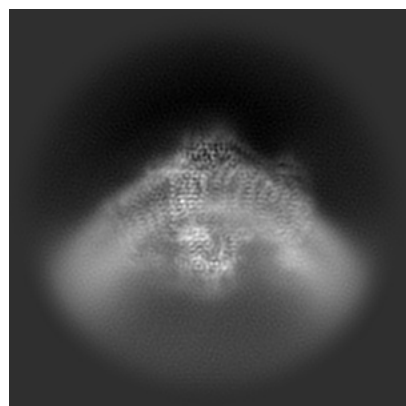
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-76184. These allow visual inspection of the internal detail of the map and identification of artifacts.

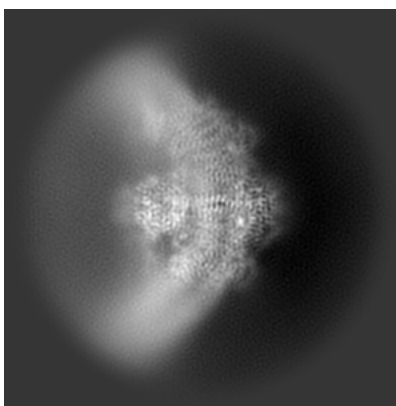
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

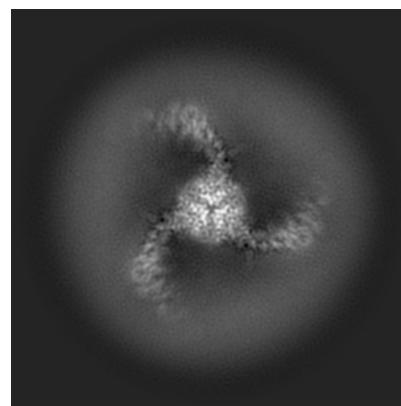
6.1.1 Primary map



X

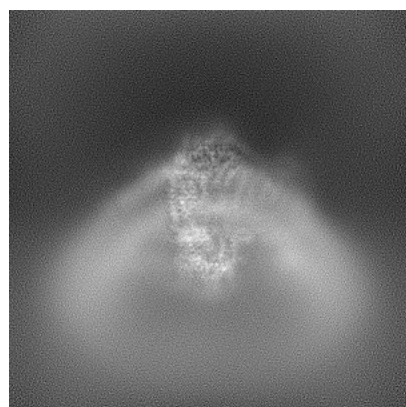


Y

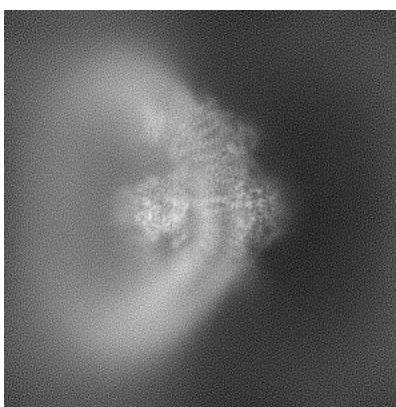


Z

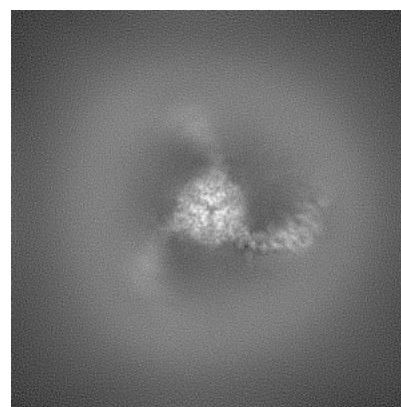
6.1.2 Raw map



X



Y

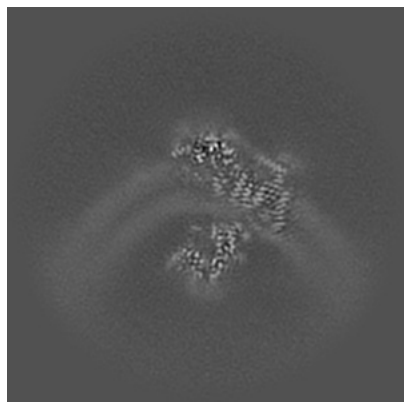


Z

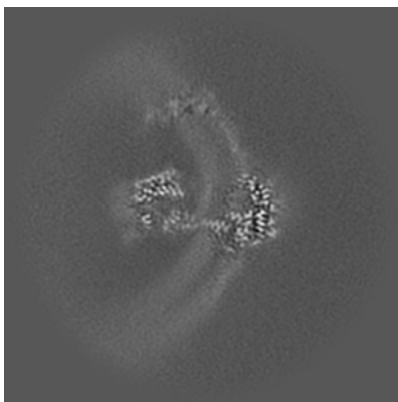
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

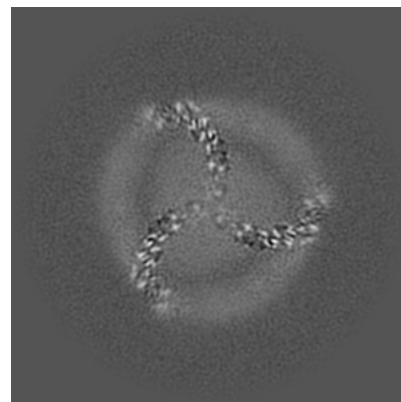
6.2.1 Primary map



X Index: 200

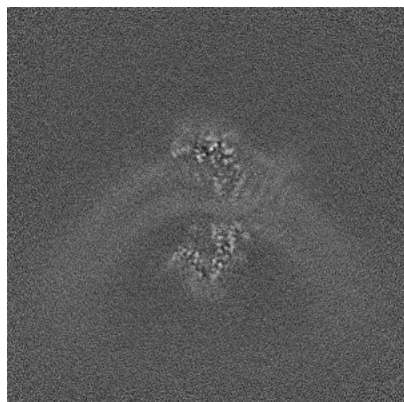


Y Index: 200

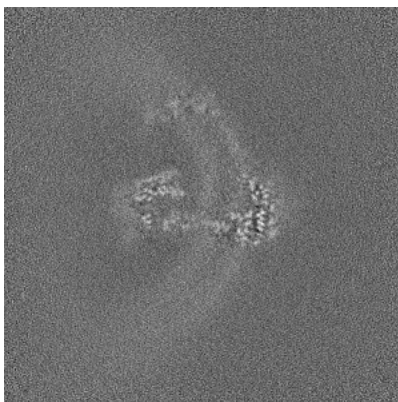


Z Index: 200

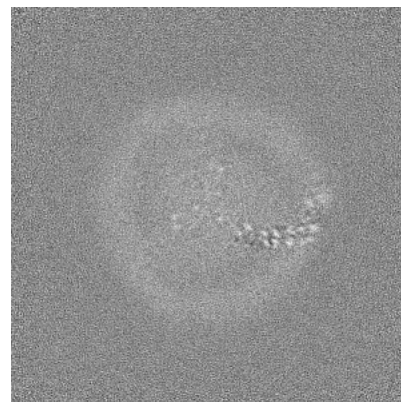
6.2.2 Raw map



X Index: 200



Y Index: 200

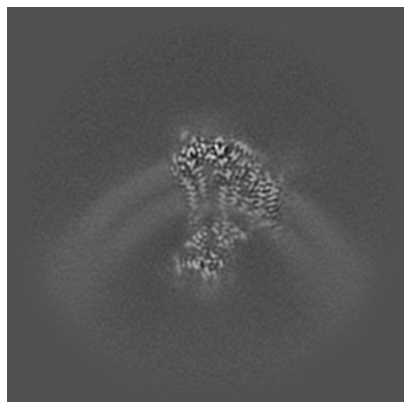


Z Index: 200

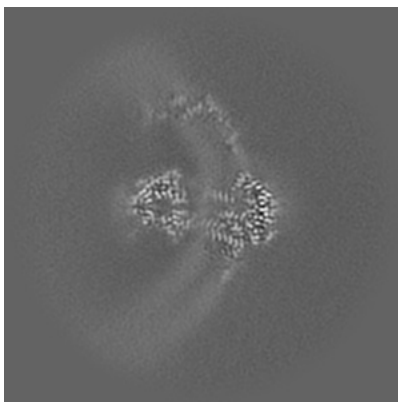
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

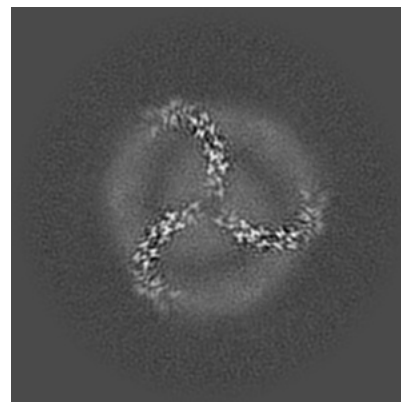
6.3.1 Primary map



X Index: 208

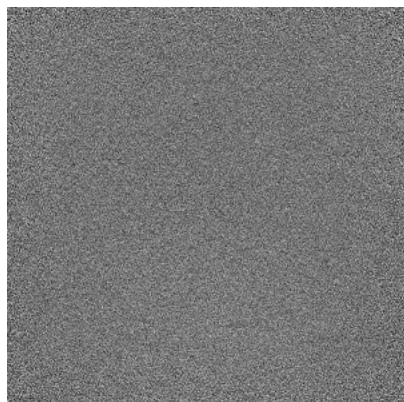


Y Index: 197

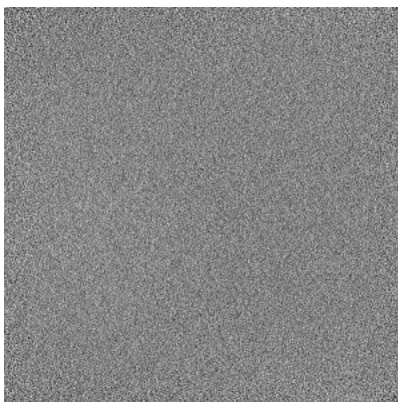


Z Index: 207

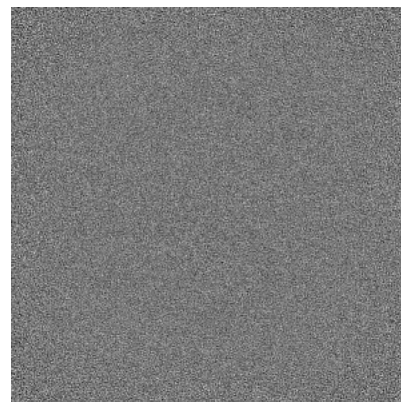
6.3.2 Raw map



X Index: 0



Y Index: 0

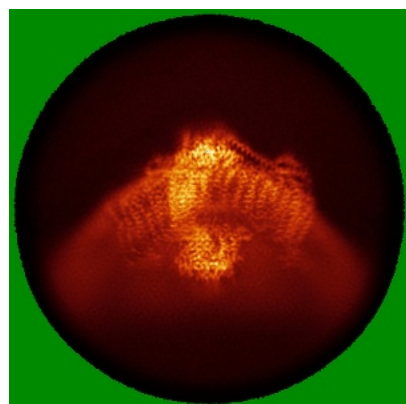


Z Index: 399

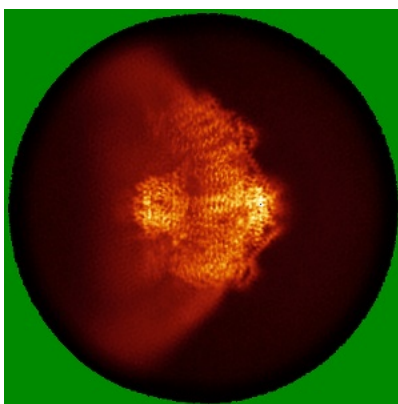
The images above show the largest variance slices of the map in three orthogonal directions.

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

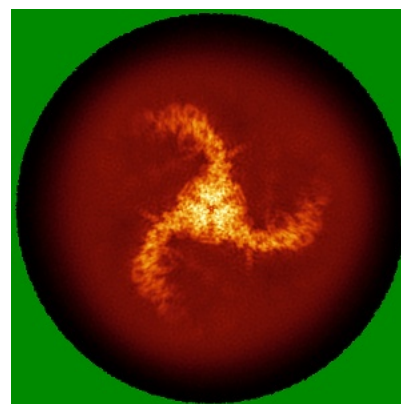
6.4.1 Primary map



X

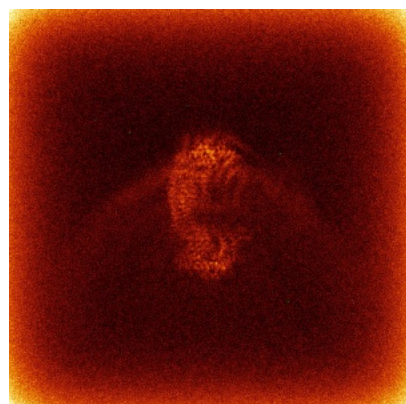


Y

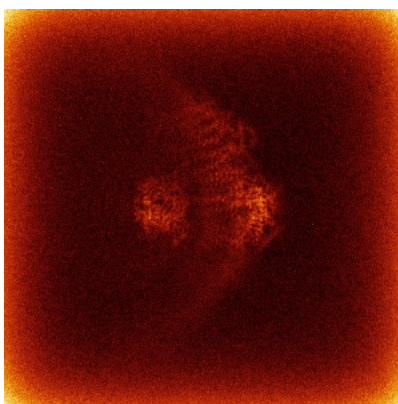


Z

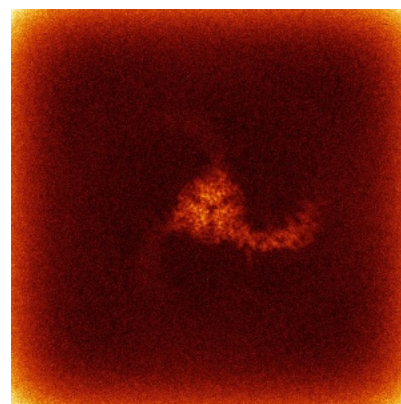
6.4.2 Raw map



X



Y



Z

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

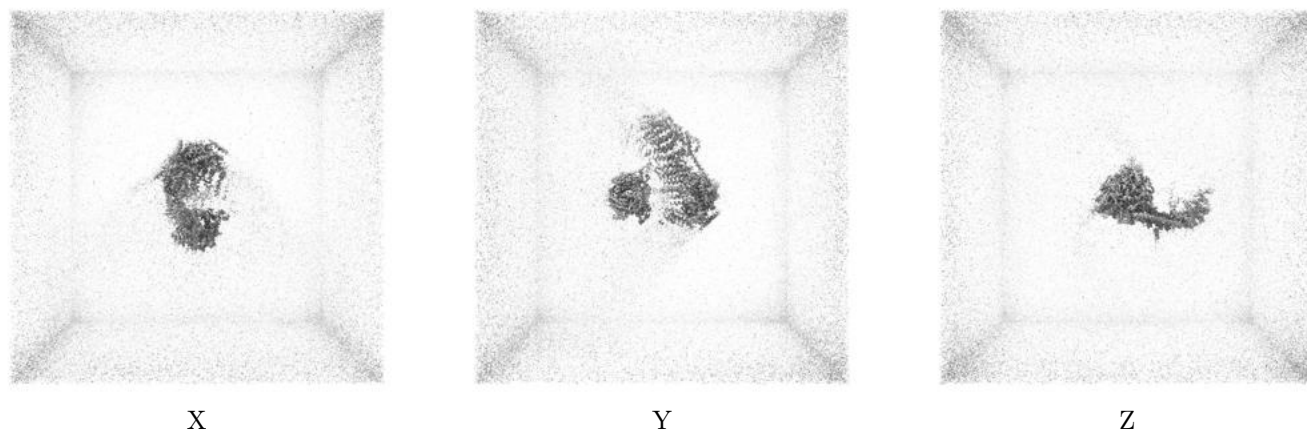
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 4.5. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

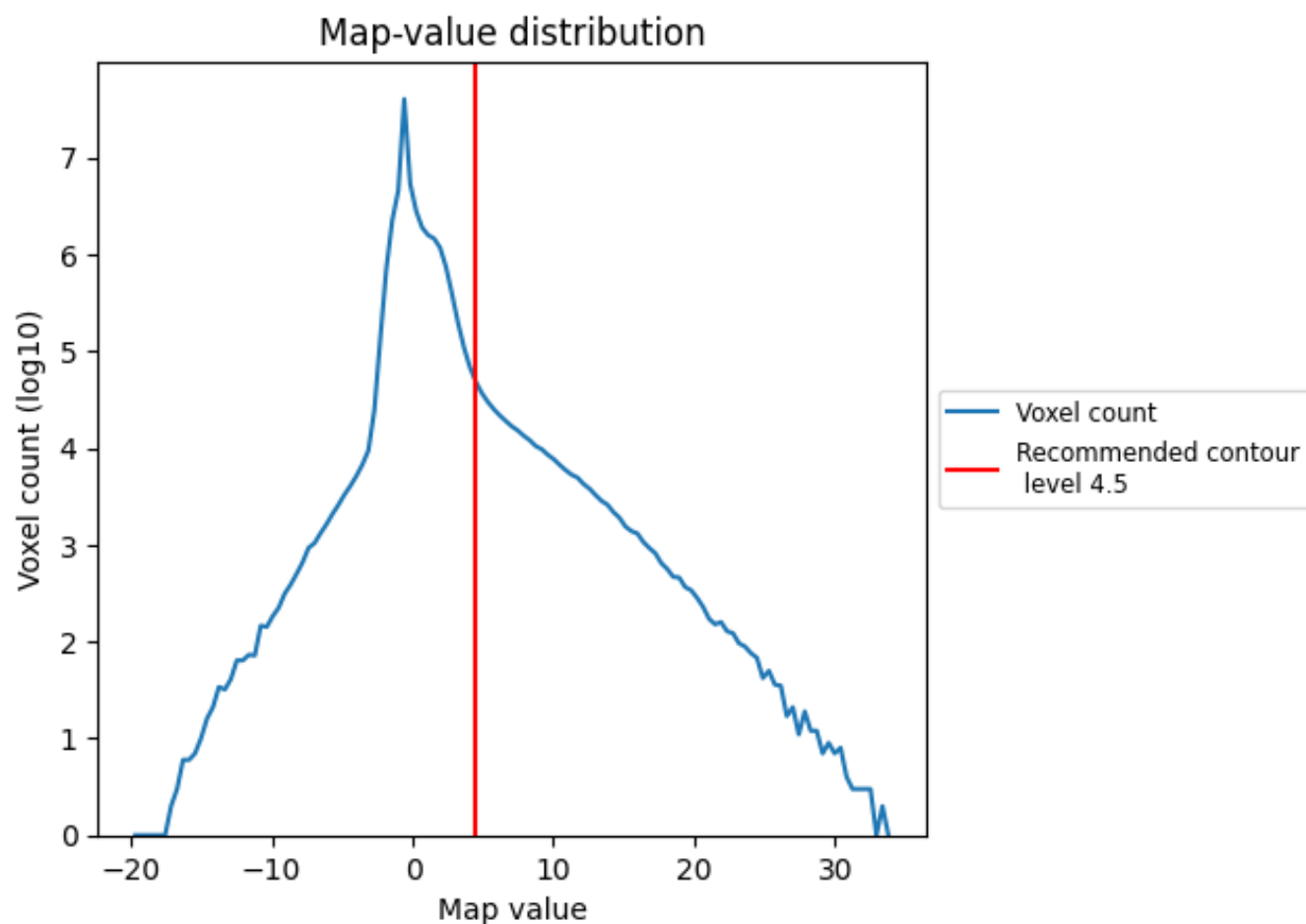
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

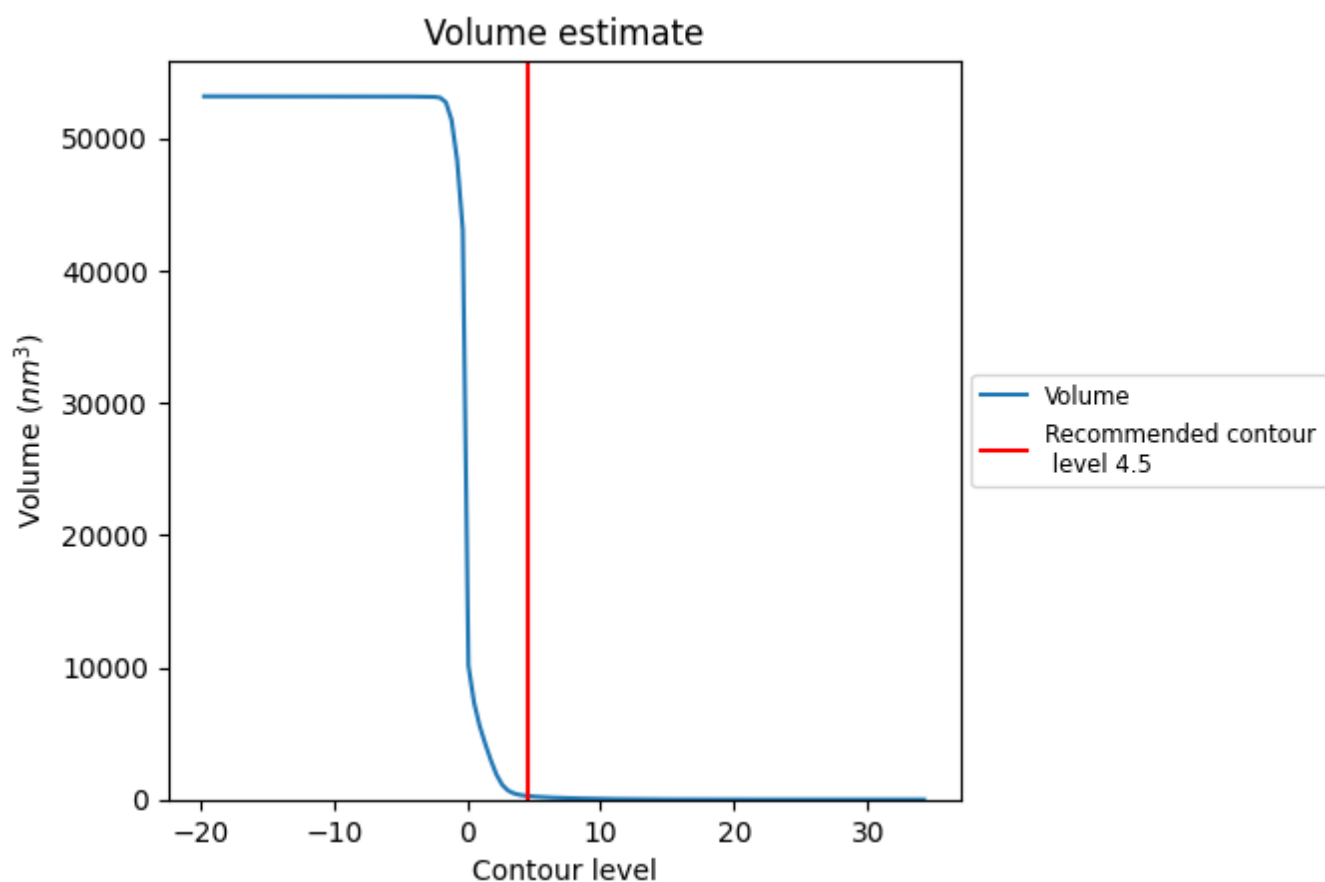
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

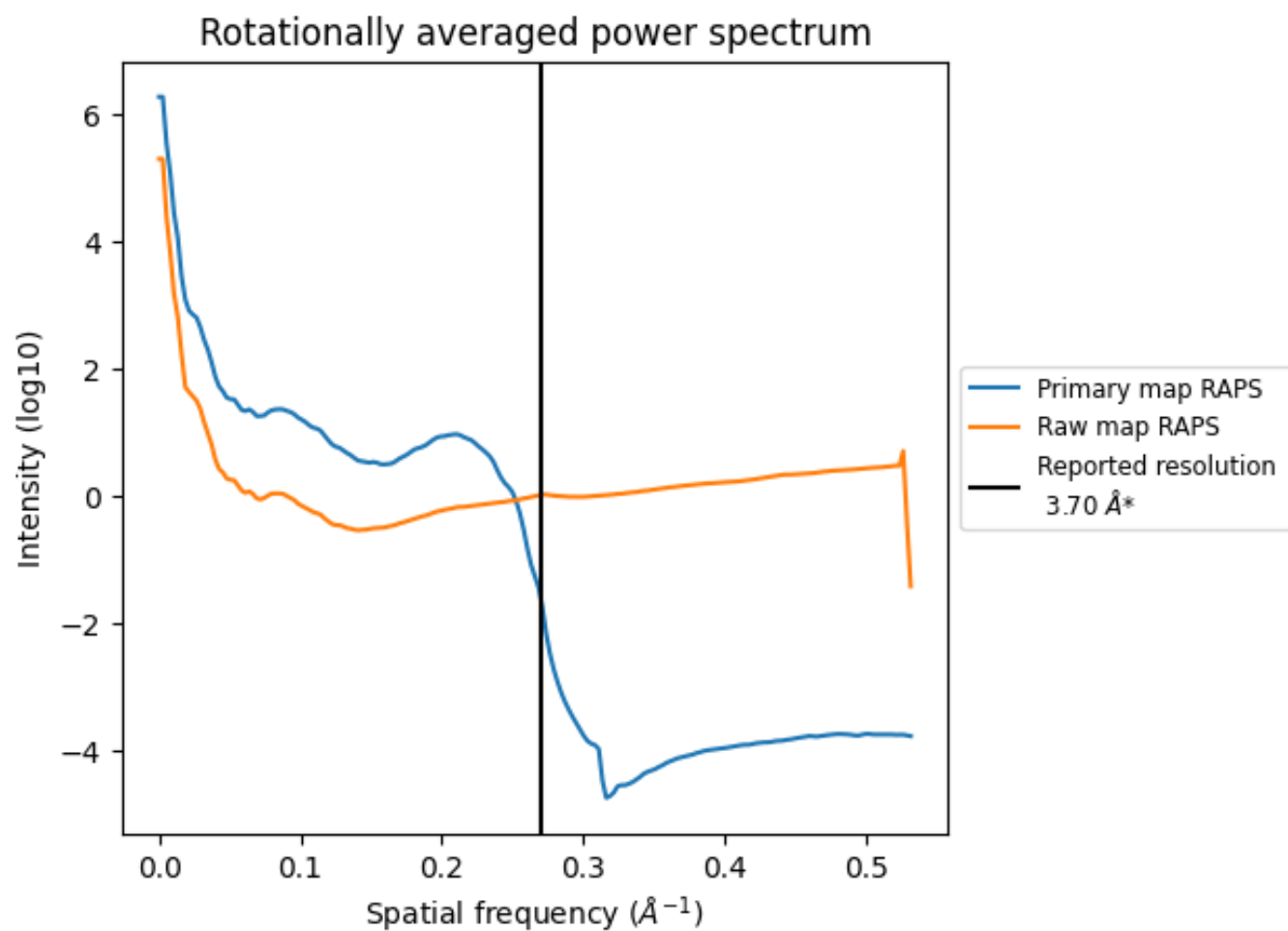
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 276 nm³; this corresponds to an approximate mass of 249 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

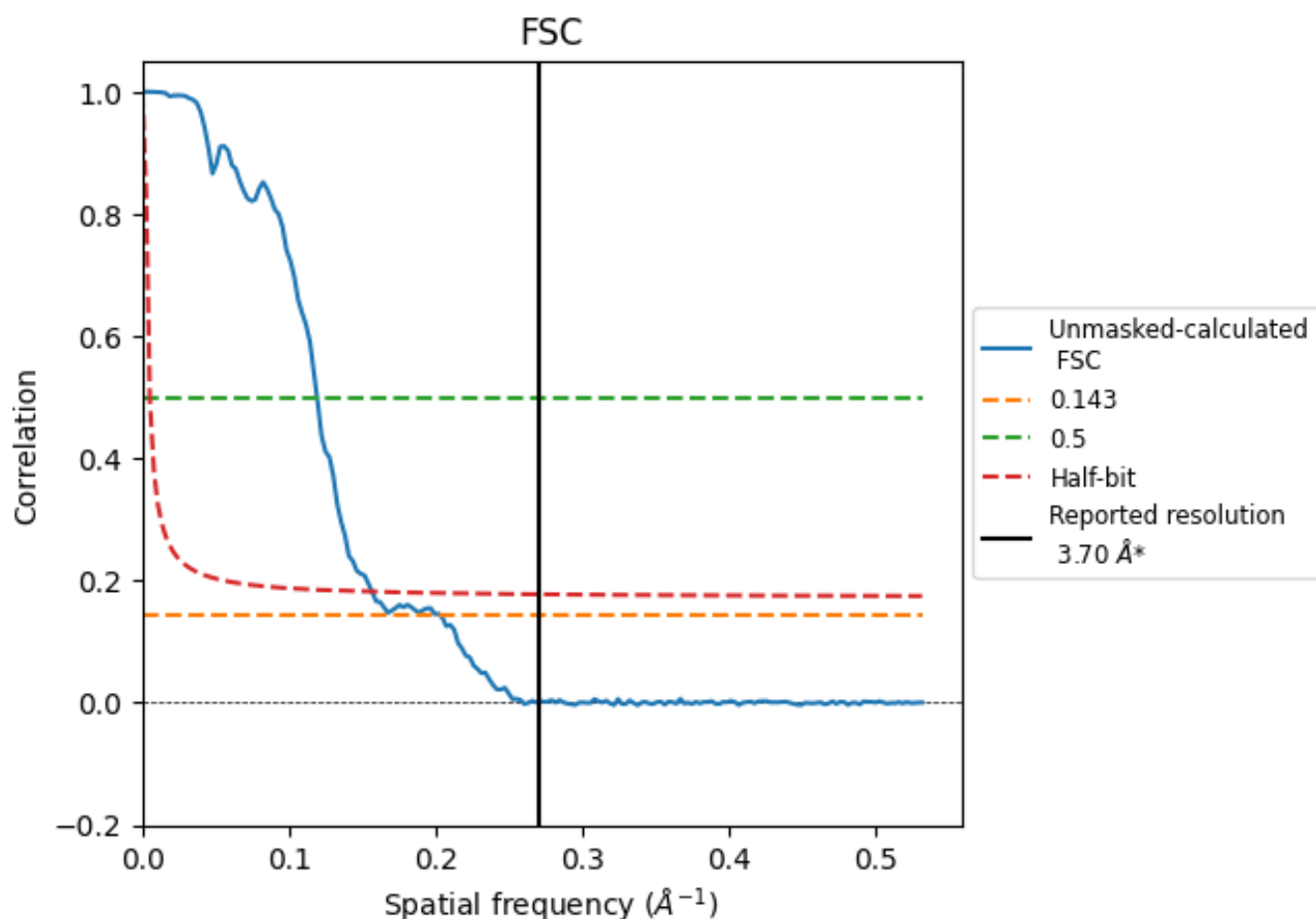


*Reported resolution corresponds to spatial frequency of 0.270 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



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

8.2 Resolution estimates [i](#)

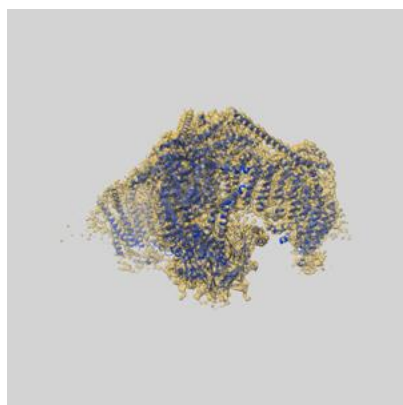
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.70	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.92	8.39	6.38

*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 4.92 differs from the reported value 3.7 by more than 10 %

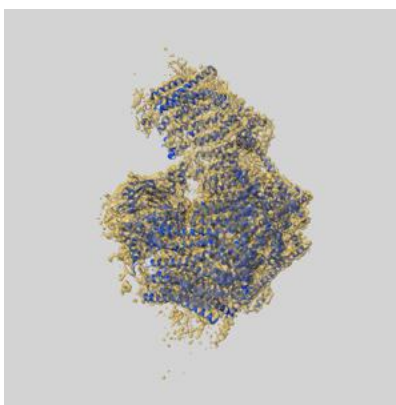
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-76184 and PDB model 11YE. Per-residue inclusion information can be found in section [3](#) on page [5](#).

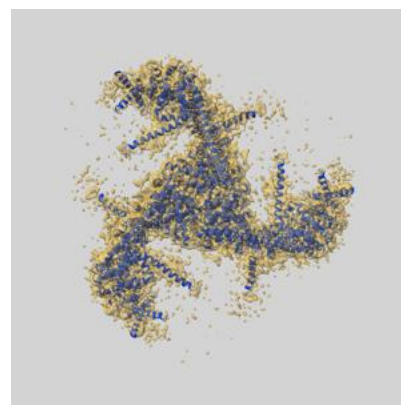
9.1 Map-model overlay [i](#)



X



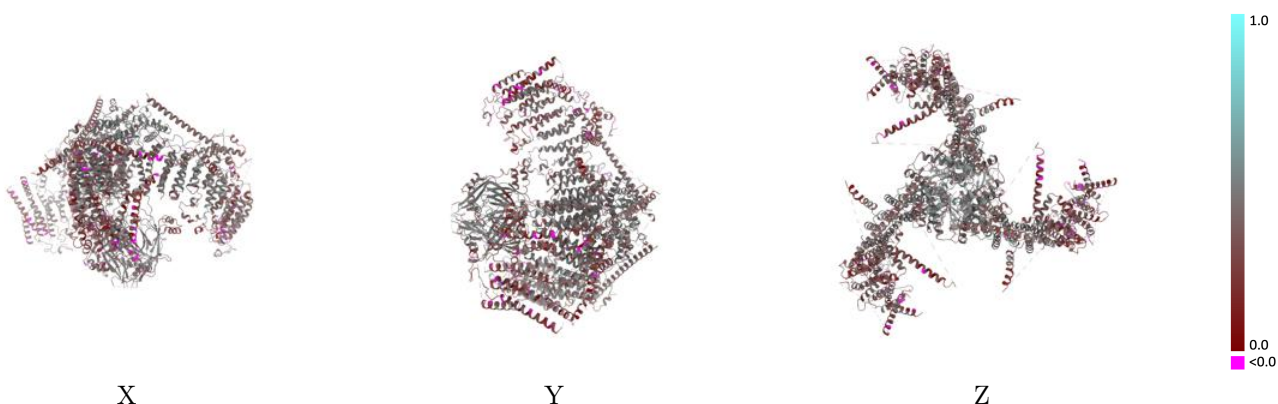
Y



Z

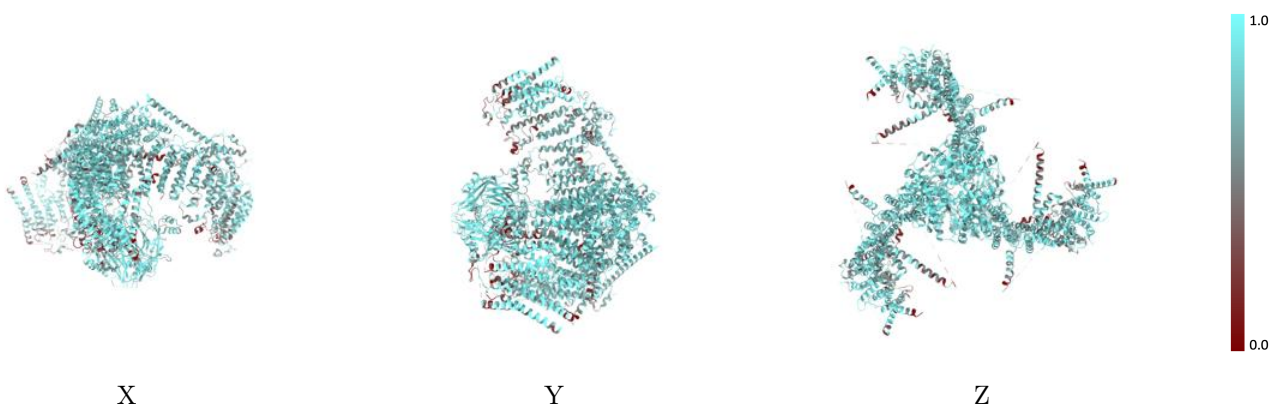
The images above show the 3D surface view of the map at the recommended contour level 4.5 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



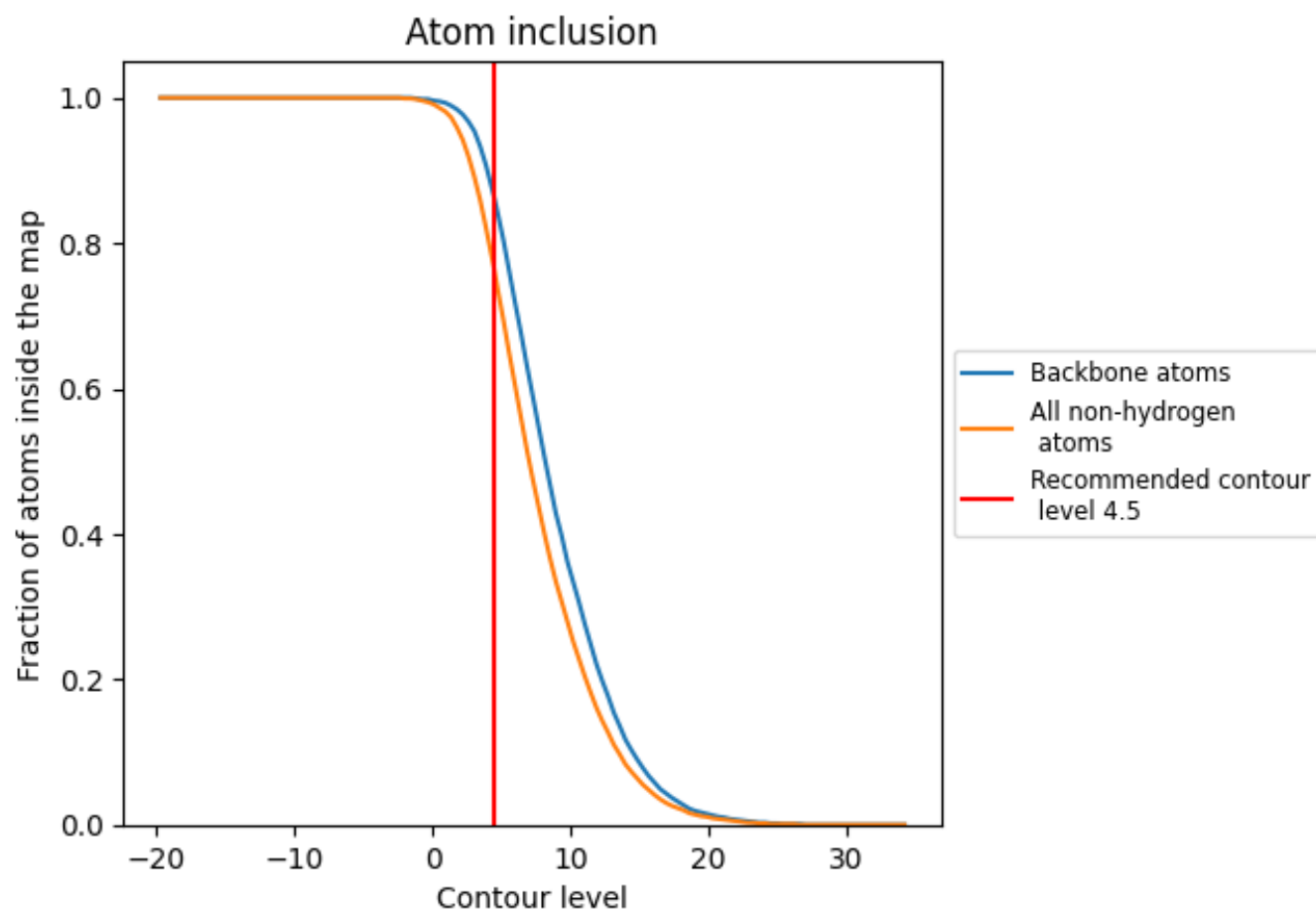
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (4.5).

9.4 Atom inclusion [i](#)



At the recommended contour level, 86% of all backbone atoms, 76% 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 (4.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7630	0.3750
C	0.7690	0.3790
D	0.7510	0.3680
E	0.7690	0.3780

