



Full wwPDB EM Validation Report ⓘ

Mar 10, 2026 – 02:07 PM UTC

PDB ID : 7WUQ / pdb_00007wuq
EMDB ID : EMD-32838
Title : Tethered peptide activation mechanism of adhesion GPCRs ADGRG2 and ADGRG4
Authors : He, Q.T.; Guo, S.C.; Xiao, P.; Sun, J.P.; Yu, X.; Gou, L.; Kong, L.L.; Zhang, L.
Deposited on : 2022-02-09
Resolution : 2.90 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

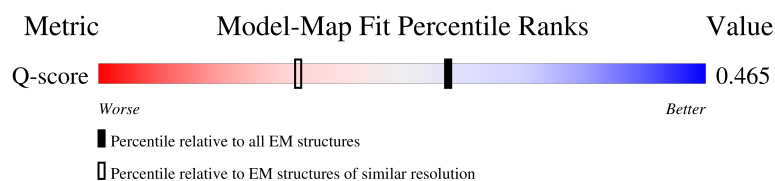
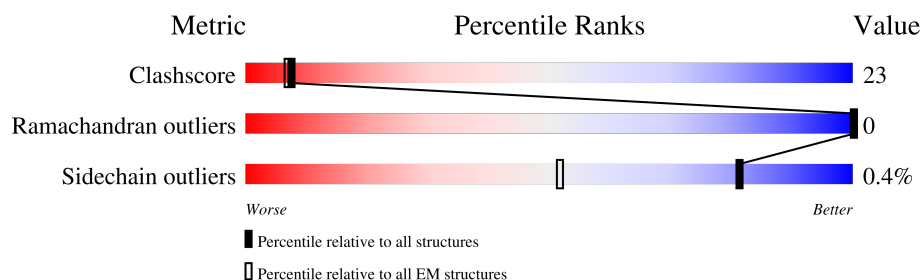
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

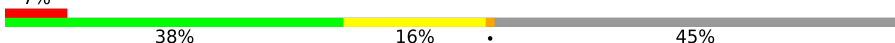
The reported resolution of this entry is 2.90 Å.

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	13054 (2.40 - 3.40)

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	394	
2	B	358	
3	G	71	
4	N	128	

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Mol	Chain	Length	Quality of chain
5	R	683	 A horizontal bar chart showing the quality of chain R. The bar is divided into three segments: a red segment at the beginning (26%), a green segment in the middle (14%), and a grey segment at the end (59%). The percentages are labeled below the bar.

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7973 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 Guanine nucleotide-binding protein G(s) subunit alpha isoforms short.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	215	Total	C	N	O	S	0	0
			1795	1135	325	330	5		

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	54	ASN	SER	engineered mutation	UNP P63092
A	226	ALA	GLY	engineered mutation	UNP P63092
A	268	ALA	GLU	engineered mutation	UNP P63092
A	271	LYS	ASN	engineered mutation	UNP P63092
A	274	ASP	LYS	engineered mutation	UNP P63092
A	280	LYS	ARG	engineered mutation	UNP P63092
A	284	ASP	THR	engineered mutation	UNP P63092
A	285	THR	ILE	engineered mutation	UNP P63092

- Molecule 2 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	334	Total	C	N	O	S	0	0
			2564	1583	462	498	21		

There are 19 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-17	MET	-	expression tag	UNP P62873
B	-16	HIS	-	expression tag	UNP P62873
B	-15	HIS	-	expression tag	UNP P62873
B	-14	HIS	-	expression tag	UNP P62873
B	-13	HIS	-	expression tag	UNP P62873
B	-12	HIS	-	expression tag	UNP P62873
B	-11	HIS	-	expression tag	UNP P62873

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	LEU	-	expression tag	UNP P62873
B	-9	GLU	-	expression tag	UNP P62873
B	-8	VAL	-	expression tag	UNP P62873
B	-7	LEU	-	expression tag	UNP P62873
B	-6	PHE	-	expression tag	UNP P62873
B	-5	GLN	-	expression tag	UNP P62873
B	-4	GLY	-	expression tag	UNP P62873
B	-3	PRO	-	expression tag	UNP P62873
B	-2	GLY	-	expression tag	UNP P62873
B	-1	SER	-	expression tag	UNP P62873
B	0	SER	-	expression tag	UNP P62873
B	1	GLN	-	expression tag	UNP P62873

- Molecule 3 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	55	Total	C	N	O	S	0	0
			421	265	75	78	3		

- Molecule 4 is a protein called Nanobody-35.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	N	127	Total	C	N	O	S	0	0
			967	602	169	190	6		

- Molecule 5 is a protein called Adhesion G-protein coupled receptor G2,mCherry.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	R	278	Total	C	N	O	S	0	0
			2226	1505	352	355	14		

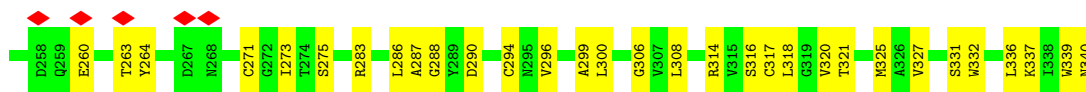
There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
R	581	MET	-	expression tag	UNP Q8CJ12
R	582	LYS	-	expression tag	UNP Q8CJ12
R	583	THR	-	expression tag	UNP Q8CJ12
R	584	ILE	-	expression tag	UNP Q8CJ12
R	585	ILE	-	expression tag	UNP Q8CJ12
R	586	ALA	-	expression tag	UNP Q8CJ12
R	587	LEU	-	expression tag	UNP Q8CJ12

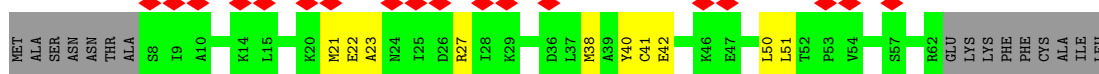
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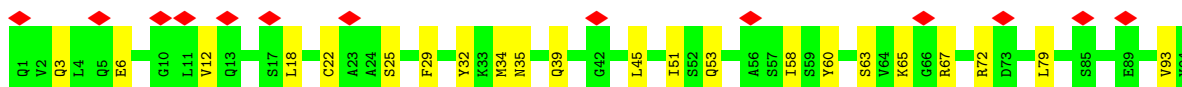
Chain	Residue	Modelled	Actual	Comment	Reference
R	588	SER	-	expression tag	UNP Q8CJ12
R	589	TYR	-	expression tag	UNP Q8CJ12
R	590	ILE	-	expression tag	UNP Q8CJ12
R	591	PHE	-	expression tag	UNP Q8CJ12
R	592	CYS	-	expression tag	UNP Q8CJ12
R	593	LEU	-	expression tag	UNP Q8CJ12
R	594	VAL	-	expression tag	UNP Q8CJ12
R	595	PHE	-	expression tag	UNP Q8CJ12
R	596	ALA	-	expression tag	UNP Q8CJ12
R	1010	HIS	-	linker	UNP Q8CJ12
R	1011	HIS	-	linker	UNP Q8CJ12
R	1012	HIS	-	linker	UNP Q8CJ12
R	1013	HIS	-	linker	UNP Q8CJ12
R	1014	HIS	-	linker	UNP Q8CJ12
R	1015	HIS	-	linker	UNP Q8CJ12
R	1016	HIS	-	linker	UNP Q8CJ12
R	1017	HIS	-	linker	UNP Q8CJ12
R	1018	GLY	-	linker	UNP Q8CJ12
R	1019	SER	-	linker	UNP Q8CJ12
R	1020	ALA	-	linker	UNP Q8CJ12
R	1021	GLU	-	linker	UNP Q8CJ12
R	1022	ASN	-	linker	UNP Q8CJ12
R	1023	LEU	-	linker	UNP Q8CJ12
R	1024	TYR	-	linker	UNP Q8CJ12
R	1025	PHE	-	linker	UNP Q8CJ12
R	1026	GLN	-	linker	UNP Q8CJ12
R	1027	GLY	-	linker	UNP Q8CJ12



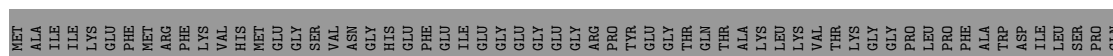
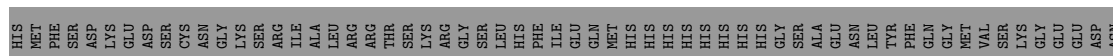
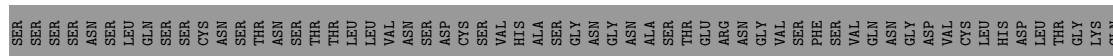
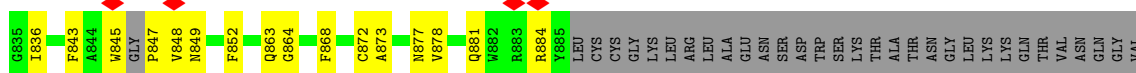
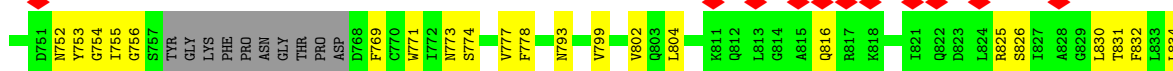
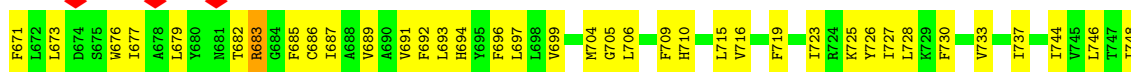
- Molecule 3: Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2



- Molecule 4: Nanobody-35



- Molecule 5: Adhesion G-protein coupled receptor G2,mCherry



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

LYS
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ARG
ALA
GLU
GLY
ARG
HIS
SER
THR
GLY
GLY
MET
ASP
GLU
LEU
TYR
LYS

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	3743997	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	64	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.226	Depositor
Minimum map value	-0.155	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.008	Depositor
Recommended contour level	0.026	Depositor
Map size (Å)	192.0, 192.0, 192.0	wwPDB
Map dimensions	192, 192, 192	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.0, 1.0, 1.0	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	0.29	1/1828 (0.1%)	0.59	1/2456 (0.0%)
2	B	0.31	0/2611	0.64	0/3539
3	G	0.21	0/427	0.49	0/576
4	N	0.24	0/987	0.55	0/1337
5	R	0.29	0/2284	0.69	1/3101 (0.0%)
All	All	0.29	1/8137 (0.0%)	0.63	2/11009 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	356	ARG	CA-CB	5.09	1.62	1.53

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	R	683	ARG	N-CA-C	6.18	118.02	111.28
1	A	40	THR	N-CA-C	5.49	117.34	111.36

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	A	1795	0	1766	61	0
2	B	2564	0	2474	102	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	421	0	434	13	0
4	N	967	0	937	49	0
5	R	2226	0	2302	168	0
All	All	7973	0	7913	363	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 23.

All (363) 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)
5:R:845:TRP:C	5:R:847:PRO:HD2	1.32	1.52
4:N:95:TYR:CE1	4:N:121:GLY:HA3	1.55	1.41
5:R:845:TRP:CA	5:R:849:ASN:HB2	1.57	1.34
5:R:679:LEU:CD1	5:R:689:VAL:HG11	1.56	1.31
5:R:648:ILE:HD12	5:R:877:ASN:ND2	1.46	1.31
1:A:41:HIS:CE1	1:A:383:ILE:HD13	1.71	1.24
5:R:648:ILE:CD1	5:R:877:ASN:HD22	1.49	1.24
4:N:35:ASN:HB2	4:N:97:ALA:O	1.38	1.21
5:R:598:LEU:HD21	5:R:617:MET:CG	1.74	1.17
5:R:598:LEU:HD21	5:R:617:MET:HG3	1.19	1.14
5:R:630:LEU:HD12	5:R:633:ILE:HD11	1.32	1.11
5:R:634:PHE:CE1	5:R:868:PHE:CE2	2.40	1.10
5:R:845:TRP:C	5:R:847:PRO:CD	2.25	1.10
5:R:845:TRP:O	5:R:847:PRO:HD2	1.52	1.10
5:R:679:LEU:HD11	5:R:689:VAL:HG11	1.34	1.07
5:R:679:LEU:HD12	5:R:689:VAL:HG11	1.32	1.07
1:A:41:HIS:CE1	1:A:383:ILE:CD1	2.37	1.06
5:R:679:LEU:CD1	5:R:689:VAL:CG1	2.33	1.05
5:R:845:TRP:C	5:R:849:ASN:HB2	1.81	1.05
2:B:81:ILE:CG2	2:B:105:TYR:HE2	1.68	1.05
5:R:670:ILE:CD1	5:R:692:PHE:HB3	1.87	1.04
5:R:845:TRP:HA	5:R:849:ASN:HB2	1.32	1.03
5:R:670:ILE:CD1	5:R:692:PHE:CB	2.37	1.02
4:N:95:TYR:CE1	4:N:121:GLY:CA	2.44	1.01
5:R:670:ILE:HD11	5:R:692:PHE:CB	1.92	0.99
5:R:670:ILE:HD12	5:R:692:PHE:HB3	1.45	0.98
2:B:81:ILE:HG23	2:B:105:TYR:HE2	1.26	0.97
5:R:648:ILE:HD12	5:R:877:ASN:HD22	0.83	0.97
5:R:670:ILE:HD11	5:R:692:PHE:HB2	1.42	0.97
5:R:744:ILE:O	5:R:748:ILE:HG12	1.65	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:N:95:TYR:HE1	4:N:121:GLY:HA3	1.26	0.95
2:B:318:LEU:HD12	2:B:318:LEU:O	1.68	0.94
1:A:43:LEU:HD23	1:A:243:ALA:HB3	1.48	0.93
1:A:42:ARG:HG2	1:A:222:PHE:HE2	1.34	0.93
2:B:37:ILE:HD13	3:G:38:MET:CE	1.99	0.92
2:B:156:GLN:HE21	2:B:168:LEU:HD11	1.31	0.92
5:R:845:TRP:HA	5:R:849:ASN:CB	1.99	0.92
2:B:81:ILE:HG23	2:B:105:TYR:CE2	2.04	0.92
2:B:81:ILE:CG2	2:B:105:TYR:CE2	2.55	0.90
2:B:37:ILE:HD11	3:G:38:MET:HE3	1.53	0.90
4:N:95:TYR:CD1	4:N:121:GLY:HA3	2.05	0.90
4:N:93:VAL:HG11	4:N:95:TYR:OH	1.72	0.89
2:B:37:ILE:CD1	3:G:38:MET:HE3	2.02	0.89
2:B:320:VAL:HG22	2:B:327:VAL:HG22	1.53	0.89
1:A:41:HIS:HE1	1:A:383:ILE:HD13	1.36	0.88
5:R:634:PHE:HE1	5:R:868:PHE:CE2	1.90	0.88
5:R:845:TRP:CA	5:R:849:ASN:CB	2.49	0.88
5:R:634:PHE:CD1	5:R:868:PHE:CD2	2.63	0.87
5:R:845:TRP:N	5:R:849:ASN:HB2	1.88	0.86
5:R:670:ILE:HD13	5:R:692:PHE:C	2.01	0.86
2:B:157:ILE:O	2:B:157:ILE:HG23	1.77	0.85
5:R:673:LEU:O	5:R:677:ILE:HG12	1.76	0.85
4:N:95:TYR:HE1	4:N:120:GLN:O	1.59	0.84
5:R:694:HIS:HD1	5:R:753:TYR:HH	1.26	0.84
5:R:634:PHE:CD1	5:R:868:PHE:CE2	2.67	0.82
5:R:755:ILE:HD11	5:R:769:PHE:O	1.81	0.81
4:N:95:TYR:CD1	4:N:121:GLY:CA	2.64	0.80
5:R:843:PHE:O	5:R:848:VAL:HG23	1.82	0.79
5:R:744:ILE:O	5:R:748:ILE:CG1	2.31	0.79
2:B:288:GLY:N	2:B:318:LEU:HD21	1.98	0.79
2:B:37:ILE:CD1	3:G:38:MET:CE	2.58	0.78
5:R:845:TRP:N	5:R:849:ASN:HD22	1.82	0.78
5:R:673:LEU:O	5:R:677:ILE:CG1	2.32	0.78
5:R:648:ILE:HG22	5:R:648:ILE:O	1.83	0.77
5:R:679:LEU:HD11	5:R:689:VAL:CG1	2.06	0.77
5:R:634:PHE:HE1	5:R:868:PHE:CZ	2.03	0.76
4:N:93:VAL:HG12	4:N:95:TYR:CZ	2.22	0.75
4:N:93:VAL:CG1	4:N:95:TYR:CZ	2.71	0.74
2:B:183:HIS:CE1	2:B:209:LYS:HG2	2.22	0.73
4:N:97:ALA:HB1	4:N:110:VAL:HB	1.71	0.73
5:R:744:ILE:O	5:R:748:ILE:CD1	2.36	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:LEU:HD21	1:A:379:CYS:SG	2.28	0.72
1:A:43:LEU:CD2	1:A:243:ALA:HB3	2.18	0.72
5:R:682:THR:HG23	5:R:682:THR:O	1.87	0.71
5:R:845:TRP:O	5:R:849:ASN:N	2.20	0.71
5:R:598:LEU:HD21	5:R:617:MET:HG2	1.68	0.71
1:A:41:HIS:CE1	1:A:383:ILE:HD11	2.25	0.71
4:N:35:ASN:CB	4:N:97:ALA:O	2.29	0.71
5:R:670:ILE:CD1	5:R:692:PHE:C	2.64	0.70
5:R:754:GLY:O	5:R:773:ASN:ND2	2.24	0.70
5:R:845:TRP:C	5:R:849:ASN:CB	2.63	0.69
4:N:95:TYR:CE1	4:N:120:GLN:O	2.44	0.68
5:R:845:TRP:H	5:R:849:ASN:HD22	1.40	0.68
2:B:71:VAL:HG12	2:B:105:TYR:HB2	1.75	0.68
5:R:634:PHE:CE1	5:R:868:PHE:CZ	2.80	0.68
5:R:755:ILE:CD1	5:R:769:PHE:O	2.42	0.67
5:R:693:LEU:HD21	5:R:769:PHE:CZ	2.30	0.67
5:R:774:SER:OG	5:R:777:VAL:HB	1.95	0.67
2:B:318:LEU:HD12	2:B:318:LEU:C	2.20	0.66
5:R:679:LEU:HD13	5:R:686:CYS:HA	1.76	0.66
5:R:673:LEU:O	5:R:677:ILE:CG2	2.43	0.66
5:R:845:TRP:O	5:R:849:ASN:HB2	1.94	0.66
1:A:392:GLU:OE1	5:R:825:ARG:NH2	2.27	0.65
4:N:95:TYR:HE1	4:N:121:GLY:CA	2.00	0.65
5:R:793:ASN:ND2	5:R:834:LEU:O	2.30	0.65
5:R:634:PHE:HD1	5:R:868:PHE:CD2	2.15	0.64
5:R:845:TRP:HA	5:R:849:ASN:ND2	2.13	0.64
5:R:848:VAL:HB	5:R:852:PHE:CE2	2.33	0.64
4:N:97:ALA:CB	4:N:110:VAL:HB	2.28	0.63
5:R:752:ASN:O	5:R:773:ASN:HB2	1.99	0.63
2:B:75:GLN:NE2	2:B:100:VAL:O	2.31	0.63
5:R:635:LEU:HD13	5:R:864:GLY:CA	2.28	0.63
5:R:848:VAL:HB	5:R:852:PHE:CD2	2.33	0.63
4:N:95:TYR:HE1	4:N:120:GLN:C	2.07	0.62
1:A:362:HIS:NE2	1:A:378:ASP:OD2	2.33	0.62
5:R:679:LEU:HD12	5:R:689:VAL:CG1	2.14	0.62
5:R:673:LEU:O	5:R:677:ILE:HG23	2.00	0.62
5:R:612:LEU:HB2	5:R:617:MET:HE2	1.81	0.61
2:B:158:VAL:HG13	2:B:190:LEU:HD21	1.82	0.61
2:B:283:ARG:N	3:G:41:CYS:SG	2.74	0.61
5:R:845:TRP:H	5:R:849:ASN:ND2	1.97	0.61
2:B:156:GLN:NE2	2:B:168:LEU:HD11	2.08	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:R:748:ILE:HD13	5:R:748:ILE:N	2.15	0.61
5:R:845:TRP:N	5:R:849:ASN:CB	2.63	0.61
5:R:753:TYR:C	5:R:773:ASN:ND2	2.59	0.60
4:N:35:ASN:ND2	4:N:107:CYS:SG	2.73	0.60
1:A:42:ARG:HG2	1:A:222:PHE:CE2	2.26	0.60
2:B:156:GLN:HE21	2:B:168:LEU:CD1	2.09	0.60
4:N:35:ASN:ND2	4:N:99:CYS:SG	2.72	0.60
1:A:41:HIS:ND1	1:A:383:ILE:HD11	2.16	0.59
2:B:157:ILE:O	2:B:157:ILE:CG2	2.49	0.59
5:R:845:TRP:HA	5:R:849:ASN:CG	2.27	0.59
5:R:598:LEU:CD2	5:R:617:MET:HG3	2.13	0.59
4:N:39:GLN:HB2	4:N:45:LEU:HD13	1.85	0.59
5:R:633:ILE:HG13	5:R:634:PHE:N	2.16	0.59
4:N:99:CYS:HG	4:N:107:CYS:CB	2.10	0.58
5:R:673:LEU:O	5:R:677:ILE:CD1	2.51	0.58
2:B:325:MET:O	2:B:340:ASN:ND2	2.36	0.58
4:N:35:ASN:HD21	4:N:107:CYS:HG	1.51	0.58
1:A:46:LEU:HD23	1:A:231:ARG:HH11	1.67	0.58
2:B:114:CYS:SG	2:B:122:SER:OG	2.61	0.58
5:R:631:SER:O	5:R:635:LEU:N	2.28	0.58
5:R:845:TRP:O	5:R:849:ASN:CB	2.52	0.58
5:R:845:TRP:O	5:R:847:PRO:CD	2.42	0.58
4:N:3:GLN:NE2	4:N:25:SER:OG	2.36	0.58
1:A:313:PRO:HG2	4:N:65:LYS:HB3	1.86	0.57
5:R:687:ILE:HG23	5:R:753:TYR:HB2	1.86	0.57
1:A:289:LEU:HB3	1:A:361:PRO:HA	1.85	0.57
5:R:845:TRP:H	5:R:849:ASN:CB	2.17	0.57
5:R:648:ILE:O	5:R:648:ILE:CG2	2.51	0.56
2:B:106:ALA:HB2	2:B:111:TYR:O	2.05	0.56
2:B:288:GLY:HA3	2:B:318:LEU:HD23	1.87	0.56
5:R:609:ARG:HD3	5:R:778:PHE:CE2	2.40	0.56
5:R:693:LEU:HD21	5:R:769:PHE:CE2	2.39	0.56
2:B:107:PRO:HD3	2:B:151:PHE:HB2	1.86	0.56
5:R:635:LEU:HD13	5:R:864:GLY:HA3	1.87	0.56
1:A:40:THR:HG23	1:A:220:HIS:HB2	1.87	0.56
5:R:606:ASP:OD1	5:R:609:ARG:HD2	2.05	0.56
2:B:283:ARG:HD3	2:B:300:LEU:HD12	1.87	0.56
5:R:881:GLN:OE1	5:R:884:ARG:NH2	2.38	0.56
4:N:93:VAL:HG11	4:N:95:TYR:CZ	2.39	0.55
5:R:845:TRP:CA	5:R:849:ASN:HD22	2.19	0.55
5:R:648:ILE:HG21	5:R:878:VAL:HG22	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:ARG:NH2	5:R:716:VAL:O	2.39	0.55
1:A:384:GLN:NE2	5:R:715:LEU:O	2.40	0.55
2:B:170:ASP:HB3	2:B:173:THR:HG22	1.89	0.55
5:R:647:LYS:HD3	5:R:881:GLN:HE22	1.72	0.55
1:A:232:ARG:HG3	1:A:233:LYS:HG2	1.89	0.55
1:A:274:ASP:OD1	1:A:278:ASN:ND2	2.40	0.55
2:B:106:ALA:CB	2:B:111:TYR:O	2.55	0.55
4:N:63:SER:O	4:N:67:ARG:NH1	2.40	0.55
5:R:618:MET:HE3	5:R:618:MET:O	2.07	0.55
2:B:47:THR:HG23	2:B:337:LYS:HB3	1.88	0.54
5:R:673:LEU:O	5:R:677:ILE:HD13	2.07	0.54
2:B:59:TYR:CD2	2:B:101:MET:HG2	2.42	0.54
2:B:107:PRO:HD3	2:B:151:PHE:CB	2.38	0.54
2:B:149:CYS:O	2:B:150:ARG:NH1	2.39	0.54
5:R:604:LEU:HD13	5:R:697:LEU:HD13	1.88	0.54
1:A:232:ARG:HH21	2:B:204:CYS:HB3	1.72	0.54
1:A:215:ASP:OD2	1:A:373:ARG:NH1	2.40	0.54
1:A:227:GLN:NE2	2:B:119:ASN:OD1	2.41	0.54
5:R:754:GLY:O	5:R:773:ASN:CG	2.51	0.54
1:A:283:ARG:O	1:A:357:HIS:ND1	2.41	0.54
2:B:49:ARG:NH1	2:B:85:TYR:O	2.41	0.53
2:B:156:GLN:NE2	2:B:168:LEU:HD21	2.22	0.53
2:B:145:TYR:O	2:B:162:GLY:N	2.41	0.53
5:R:843:PHE:O	5:R:848:VAL:CG2	2.53	0.53
1:A:386:MET:SD	1:A:390:GLN:NE2	2.82	0.53
5:R:609:ARG:NH2	5:R:778:PHE:CG	2.76	0.53
5:R:843:PHE:CD1	5:R:843:PHE:N	2.73	0.53
2:B:251:ARG:NH1	2:B:263:THR:OG1	2.41	0.53
5:R:598:LEU:CD2	5:R:617:MET:CG	2.68	0.53
5:R:754:GLY:O	5:R:773:ASN:OD1	2.27	0.53
1:A:234:TRP:HB2	2:B:117:LEU:HD21	1.91	0.53
2:B:271:CYS:HB2	2:B:290:ASP:HB2	1.90	0.52
4:N:95:TYR:CE1	4:N:120:GLN:C	2.88	0.52
5:R:606:ASP:OD2	5:R:609:ARG:HD2	2.09	0.52
2:B:54:HIS:NE2	2:B:72:SER:OG	2.37	0.52
5:R:598:LEU:HA	5:R:621:THR:HG21	1.91	0.52
5:R:704:MET:HE2	5:R:863:GLN:HE22	1.74	0.52
2:B:58:ILE:O	2:B:316:SER:OG	2.28	0.52
5:R:670:ILE:CD1	5:R:692:PHE:HB2	2.17	0.52
5:R:771:TRP:NE1	5:R:778:PHE:CE1	2.71	0.52
5:R:676:TRP:HA	5:R:676:TRP:CE3	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:288:GLY:CA	2:B:318:LEU:HD23	2.40	0.52
1:A:30:LEU:O	1:A:34:LYS:HB2	2.10	0.52
5:R:606:ASP:CG	5:R:609:ARG:HD2	2.34	0.52
5:R:683:ARG:CZ	5:R:687:ILE:HD11	2.39	0.52
5:R:639:LEU:HD12	5:R:665:LEU:HD13	1.92	0.52
5:R:755:ILE:HG22	5:R:756:GLY:N	2.24	0.52
2:B:250:CYS:HB3	2:B:273:ILE:HD13	1.91	0.51
5:R:752:ASN:O	5:R:773:ASN:ND2	2.41	0.51
5:R:753:TYR:O	5:R:773:ASN:ND2	2.43	0.51
1:A:393:LEU:HD23	5:R:826:SER:HB2	1.93	0.51
4:N:99:CYS:SG	4:N:107:CYS:CB	2.99	0.51
5:R:706:LEU:HD11	5:R:733:VAL:HG13	1.92	0.51
2:B:331:SER:OG	2:B:332:TRP:N	2.44	0.51
5:R:799:VAL:HA	5:R:802:VAL:HG12	1.93	0.51
5:R:683:ARG:NE	5:R:687:ILE:HD11	2.26	0.51
1:A:393:LEU:HD11	5:R:830:LEU:HD13	1.92	0.51
4:N:51:ILE:HG13	4:N:58:ILE:HG12	1.92	0.51
5:R:733:VAL:HG23	5:R:737:ILE:HD12	1.93	0.51
5:R:845:TRP:HA	5:R:849:ASN:HD22	1.76	0.51
2:B:187:VAL:HA	2:B:203:ALA:HA	1.94	0.50
5:R:725:LYS:HB3	5:R:728:LEU:HB3	1.92	0.50
2:B:73:ALA:HB2	2:B:103:CYS:HB3	1.92	0.50
5:R:670:ILE:HD13	5:R:693:LEU:N	2.26	0.50
5:R:696:PHE:HA	5:R:699:VAL:HG22	1.94	0.50
5:R:831:THR:HG23	5:R:836:ILE:HB	1.92	0.50
2:B:100:VAL:HG23	2:B:115:GLY:H	1.77	0.50
5:R:635:LEU:HD13	5:R:864:GLY:C	2.37	0.49
5:R:655:LYS:HD3	5:R:727:ILE:HG13	1.93	0.49
2:B:220:GLN:HA	3:G:22:GLU:HG3	1.94	0.49
4:N:115:TYR:HD2	4:N:118:ARG:HB2	1.77	0.49
4:N:93:VAL:HG12	4:N:95:TYR:CE2	2.47	0.49
5:R:671:PHE:CE2	5:R:769:PHE:CZ	3.01	0.49
1:A:263:THR:OG1	1:A:264:ASN:N	2.45	0.48
5:R:744:ILE:O	5:R:748:ILE:HD13	2.10	0.48
1:A:233:LYS:NZ	2:B:246:ASP:OD2	2.37	0.48
2:B:288:GLY:CA	2:B:318:LEU:CD2	2.91	0.48
5:R:687:ILE:O	5:R:691:VAL:HG23	2.12	0.48
1:A:317:ARG:O	1:A:317:ARG:HG2	2.14	0.48
1:A:356:ARG:HB2	1:A:389:ARG:HD2	1.96	0.48
1:A:390:GLN:HB3	5:R:653:PRO:HG2	1.96	0.48
2:B:212:ASP:HB3	2:B:215:GLU:HB2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:288:GLY:N	2:B:318:LEU:CD2	2.74	0.48
5:R:647:LYS:CD	5:R:881:GLN:HE22	2.26	0.48
1:A:346:LEU:HD11	1:A:361:PRO:HG3	1.96	0.48
2:B:168:LEU:O	2:B:177:THR:N	2.47	0.48
5:R:832:PHE:HB3	5:R:863:GLN:HG3	1.95	0.48
2:B:219:ARG:HA	3:G:21:MET:HE1	1.94	0.48
5:R:705:GLY:O	5:R:709:PHE:N	2.42	0.48
5:R:609:ARG:NH2	5:R:778:PHE:CB	2.77	0.48
2:B:318:LEU:O	2:B:318:LEU:CD1	2.53	0.47
5:R:604:LEU:HD22	5:R:697:LEU:HD22	1.94	0.47
2:B:42:ARG:HH11	2:B:306:GLY:HA2	1.80	0.47
2:B:321:THR:OG1	2:B:325:MET:N	2.46	0.47
2:B:83:ASP:OD2	2:B:86:THR:OG1	2.33	0.47
2:B:275:SER:OG	2:B:317:CYS:HA	2.15	0.47
1:A:44:LEU:HD22	1:A:238:PHE:CG	2.50	0.47
2:B:34:THR:HG21	2:B:300:LEU:HB3	1.96	0.47
2:B:81:ILE:HG21	2:B:105:TYR:CE2	2.43	0.47
5:R:723:ILE:HB	5:R:726:TYR:HD1	1.79	0.47
1:A:385:ARG:HA	1:A:388:LEU:HD12	1.98	0.46
2:B:57:LYS:HD2	2:B:332:TRP:CE2	2.50	0.46
2:B:207:SER:HA	2:B:223:THR:HA	1.95	0.46
5:R:646:GLU:HA	5:R:649:ARG:HE	1.80	0.46
2:B:287:ALA:C	2:B:318:LEU:HD21	2.40	0.46
5:R:670:ILE:CD1	5:R:693:LEU:N	2.79	0.46
2:B:79:LEU:HB3	2:B:93:ILE:HB	1.98	0.46
4:N:98:ARG:O	4:N:115:TYR:OH	2.26	0.46
5:R:634:PHE:HD1	5:R:868:PHE:CG	2.33	0.46
5:R:753:TYR:C	5:R:773:ASN:HD21	2.24	0.46
1:A:371:ASN:OD1	1:A:374:ARG:NH1	2.48	0.46
4:N:6:GLU:HA	4:N:22:CYS:HA	1.97	0.46
4:N:98:ARG:HH21	4:N:100:PRO:HA	1.81	0.46
2:B:45:MET:HE2	3:G:50:LEU:HD22	1.98	0.45
5:R:673:LEU:HA	5:R:677:ILE:HG23	1.98	0.45
5:R:868:PHE:HA	5:R:872:CYS:HB3	1.98	0.45
1:A:43:LEU:HD21	1:A:379:CYS:HG	1.81	0.45
2:B:63:TRP:O	2:B:105:TYR:HB3	2.15	0.45
1:A:252:SER:HB3	1:A:266:LEU:HB3	1.98	0.45
1:A:384:GLN:HG2	5:R:715:LEU:HA	1.99	0.45
2:B:320:VAL:HG12	2:B:321:THR:N	2.32	0.45
5:R:693:LEU:CD2	5:R:769:PHE:CZ	2.98	0.45
4:N:22:CYS:SG	4:N:79:LEU:HB3	2.57	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:R:635:LEU:HD23	5:R:665:LEU:HB2	1.99	0.45
5:R:683:ARG:CZ	5:R:687:ILE:CD1	2.94	0.45
1:A:364:THR:HG22	1:A:371:ASN:HD21	1.80	0.45
1:A:276:ILE:HG23	1:A:282:LEU:HD13	1.97	0.45
5:R:754:GLY:C	5:R:773:ASN:HD21	2.25	0.45
5:R:653:PRO:HD3	5:R:710:HIS:CE1	2.52	0.45
1:A:271:LYS:HD2	1:A:271:LYS:HA	1.82	0.45
2:B:52:ARG:O	2:B:89:LYS:NZ	2.50	0.44
4:N:35:ASN:ND2	4:N:110:VAL:HG11	2.33	0.44
2:B:51:LEU:HB2	2:B:336:LEU:HB2	1.99	0.44
1:A:41:HIS:CD2	5:R:719:PHE:HD1	2.34	0.44
1:A:370:GLU:O	1:A:370:GLU:HG3	2.17	0.44
1:A:278:ASN:OD1	4:N:105:ARG:NE	2.47	0.44
2:B:23:LYS:HA	2:B:26:ALA:HB2	2.00	0.44
2:B:327:VAL:O	2:B:339:TRP:N	2.49	0.44
5:R:679:LEU:CD1	5:R:689:VAL:HG12	2.37	0.44
1:A:389:ARG:HG2	1:A:394:LEU:HB2	2.00	0.44
2:B:294:CYS:HB2	2:B:308:LEU:HB2	1.99	0.44
4:N:99:CYS:SG	4:N:107:CYS:HB2	2.58	0.43
5:R:635:LEU:HB3	5:R:665:LEU:HD12	2.00	0.43
1:A:44:LEU:O	1:A:245:ILE:N	2.52	0.43
1:A:364:THR:CG2	1:A:371:ASN:HD21	2.31	0.43
2:B:260:GLU:OE2	2:B:263:THR:OG1	2.35	0.43
4:N:93:VAL:CG1	4:N:95:TYR:CE2	3.01	0.43
4:N:32:TYR:O	4:N:72:ARG:NH2	2.45	0.43
2:B:34:THR:HG23	3:G:38:MET:HE3	1.99	0.43
2:B:264:TYR:OH	2:B:299:ALA:O	2.34	0.43
2:B:57:LYS:HB2	2:B:332:TRP:CD1	2.53	0.43
2:B:275:SER:O	2:B:318:LEU:HD21	2.18	0.43
4:N:60:TYR:HB2	4:N:65:LYS:HG3	2.00	0.43
1:A:275:SER:OG	4:N:107:CYS:O	2.36	0.43
2:B:245:SER:OG	2:B:247:ASP:OD1	2.36	0.43
4:N:34:MET:HB3	4:N:79:LEU:HD22	2.01	0.43
1:A:212:PHE:HE1	1:A:221:MET:HB2	1.83	0.43
4:N:53:GLN:NE2	4:N:102:PRO:O	2.52	0.43
1:A:343:ASP:HA	1:A:346:LEU:HB2	2.01	0.42
3:G:23:ALA:HA	3:G:27:ARG:HH22	1.84	0.42
5:R:693:LEU:HG	5:R:771:TRP:CZ3	2.54	0.42
1:A:388:LEU:HD22	5:R:804:LEU:HD13	2.01	0.42
2:B:318:LEU:C	2:B:318:LEU:CD1	2.91	0.42
5:R:682:THR:O	5:R:682:THR:CG2	2.59	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:48:ARG:HE	2:B:340:ASN:HB2	1.83	0.42
5:R:679:LEU:HD23	5:R:679:LEU:HA	1.71	0.42
5:R:679:LEU:HD21	5:R:685:PHE:CD1	2.54	0.42
1:A:274:ASP:HA	1:A:348:ILE:HG21	2.02	0.42
4:N:12:VAL:HG21	4:N:18:LEU:HD12	2.01	0.42
1:A:45:LEU:CD2	1:A:247:VAL:HG23	2.50	0.42
5:R:599:THR:H	5:R:621:THR:HG23	1.84	0.42
5:R:730:PHE:HA	5:R:733:VAL:HG12	2.02	0.42
5:R:868:PHE:O	5:R:873:ALA:N	2.51	0.42
2:B:290:ASP:HA	2:B:314:ARG:HG2	2.02	0.41
5:R:848:VAL:HB	5:R:852:PHE:HE2	1.83	0.41
5:R:613:PRO:HA	5:R:614:PRO:HD3	1.91	0.41
2:B:42:ARG:HG3	2:B:44:GLN:HG2	2.02	0.41
2:B:152:LEU:HB2	2:B:156:GLN:HB3	2.03	0.41
2:B:286:LEU:HD13	2:B:296:VAL:HB	2.02	0.41
2:B:156:GLN:HE21	2:B:168:LEU:HD21	1.83	0.41
5:R:618:MET:O	5:R:618:MET:SD	2.79	0.41
1:A:393:LEU:HA	5:R:826:SER:HB2	2.02	0.41
2:B:48:ARG:HA	2:B:48:ARG:HD3	1.87	0.41
2:B:275:SER:CB	2:B:318:LEU:H	2.33	0.41
4:N:6:GLU:HG3	4:N:22:CYS:HB3	2.02	0.41
5:R:746:LEU:HD23	5:R:746:LEU:HA	1.91	0.41
2:B:70:LEU:O	2:B:82:TRP:N	2.51	0.41
4:N:29:PHE:HE1	4:N:72:ARG:HB2	1.85	0.41
2:B:40:VAL:HG21	3:G:51:LEU:HD22	2.02	0.41
4:N:96:CYS:O	4:N:96:CYS:SG	2.78	0.41
1:A:231:ARG:H	4:N:109:ASP:CG	2.29	0.41
2:B:57:LYS:HE2	2:B:75:GLN:HG3	2.01	0.41
4:N:35:ASN:HD22	4:N:110:VAL:HG11	1.86	0.41
1:A:216:LYS:HE2	1:A:216:LYS:HB2	1.86	0.41
2:B:103:CYS:O	2:B:103:CYS:SG	2.79	0.41
5:R:673:LEU:CA	5:R:677:ILE:HG23	2.51	0.41
5:R:752:ASN:O	5:R:773:ASN:CB	2.67	0.41
2:B:101:MET:HE1	2:B:145:TYR:HB2	2.03	0.41
2:B:217:MET:HE1	3:G:21:MET:HE3	2.03	0.41
1:A:247:VAL:HG13	1:A:292:ASN:HB2	2.02	0.40
2:B:209:LYS:HB3	2:B:221:THR:HG23	2.04	0.40
5:R:816:GLN:H	5:R:816:GLN:HG3	1.72	0.40
2:B:104:ALA:HB2	2:B:149:CYS:O	2.21	0.40
2:B:151:PHE:CZ	2:B:157:ILE:HD12	2.57	0.40
2:B:317:CYS:SG	2:B:317:CYS:O	2.79	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:370:GLU:O	1:A:370:GLU:CG	2.70	0.40
5:R:679:LEU:HD13	5:R:689:VAL:CG1	2.44	0.40
5:R:609:ARG:NH2	5:R:778:PHE:CD2	2.89	0.40
5:R:848:VAL:HB	5:R:852:PHE:HD2	1.83	0.40
2:B:236:PRO:HB2	3:G:40:TYR:CE1	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	205/394 (52%)	199 (97%)	6 (3%)	0	100	100
2	B	332/358 (93%)	307 (92%)	25 (8%)	0	100	100
3	G	53/71 (75%)	52 (98%)	1 (2%)	0	100	100
4	N	125/128 (98%)	116 (93%)	9 (7%)	0	100	100
5	R	272/683 (40%)	264 (97%)	8 (3%)	0	100	100
All	All	987/1634 (60%)	938 (95%)	49 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	193/349 (55%)	193 (100%)	0	100	100
2	B	276/299 (92%)	276 (100%)	0	100	100
3	G	44/58 (76%)	43 (98%)	1 (2%)	44	76
4	N	105/106 (99%)	105 (100%)	0	100	100
5	R	147/592 (25%)	145 (99%)	2 (1%)	59	85
All	All	765/1404 (54%)	762 (100%)	3 (0%)	81	94

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	G	42	GLU
5	R	616	GLN
5	R	618	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	GLN
1	A	23	ASN
1	A	29	GLN
1	A	41	HIS
1	A	267	GLN
1	A	387	HIS
1	A	390	GLN
2	B	16	ASN
2	B	75	GLN
2	B	156	GLN
2	B	175	GLN
2	B	293	ASN
4	N	5	GLN
4	N	35	ASN
4	N	74	ASN
4	N	84	ASN
4	N	123	GLN
5	R	616	GLN
5	R	849	ASN
5	R	877	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.

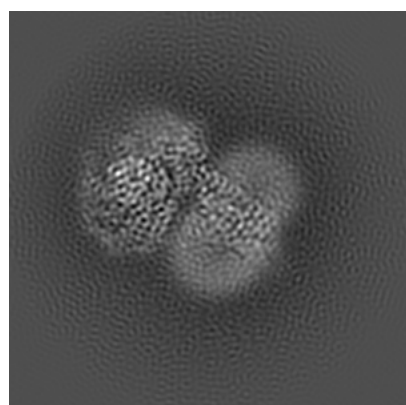
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-32838. These allow visual inspection of the internal detail of the map and identification of artifacts.

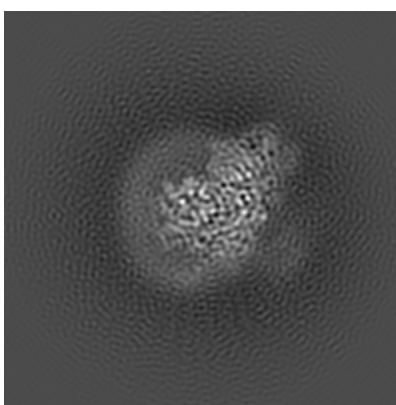
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

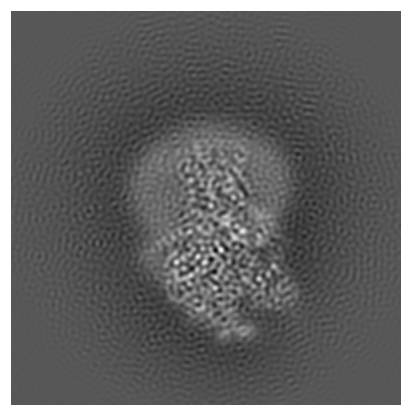
6.1.1 Primary map



X



Y

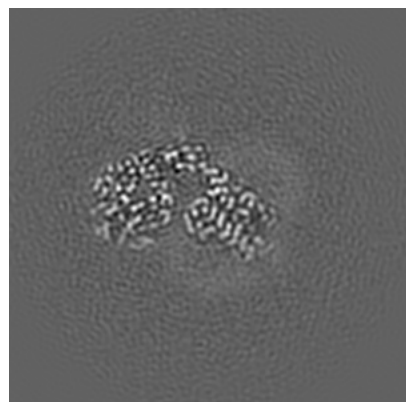


Z

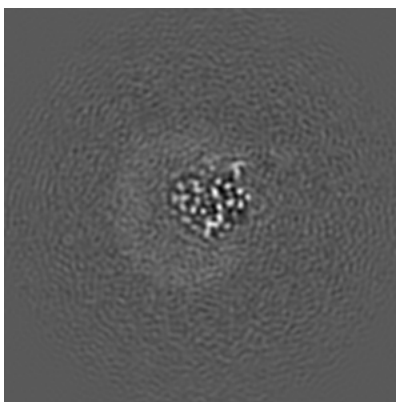
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

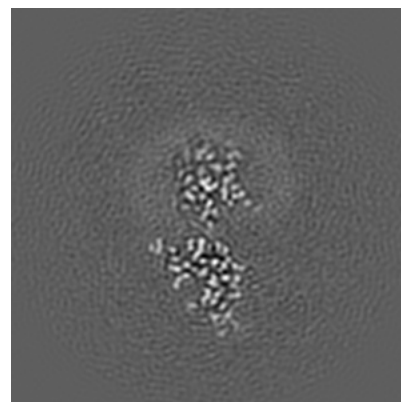
6.2.1 Primary map



X Index: 96



Y Index: 96

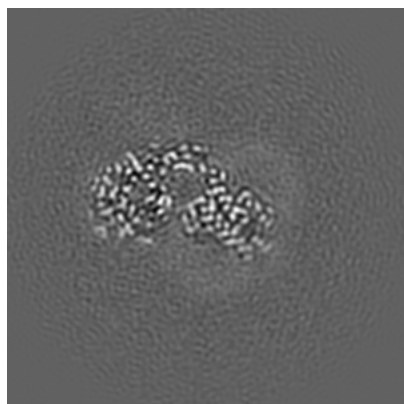


Z Index: 96

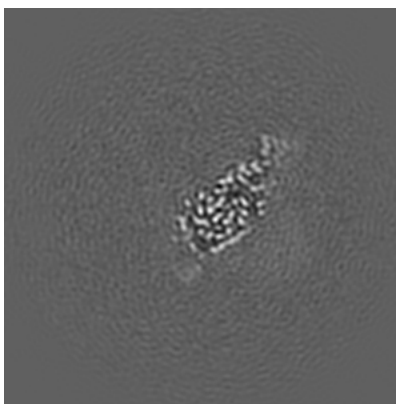
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

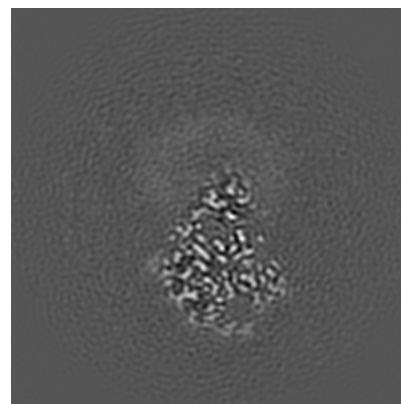
6.3.1 Primary map



X Index: 95



Y Index: 73

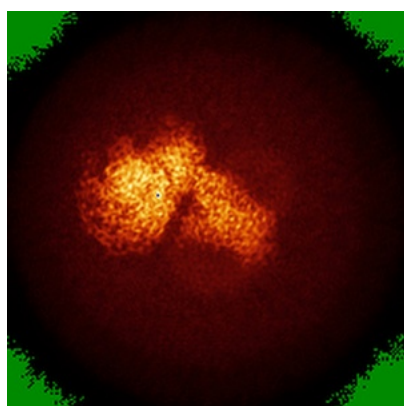


Z Index: 109

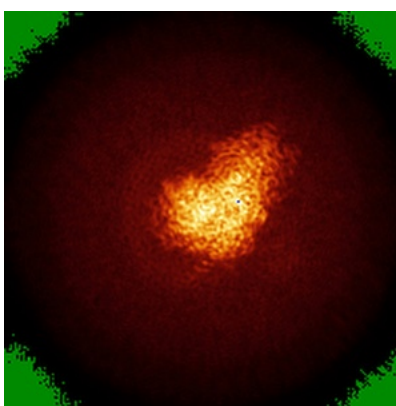
The images above show the largest variance slices of the map in three orthogonal directions.

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

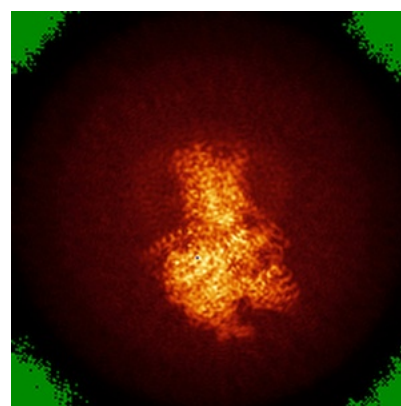
6.4.1 Primary map



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



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.026. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

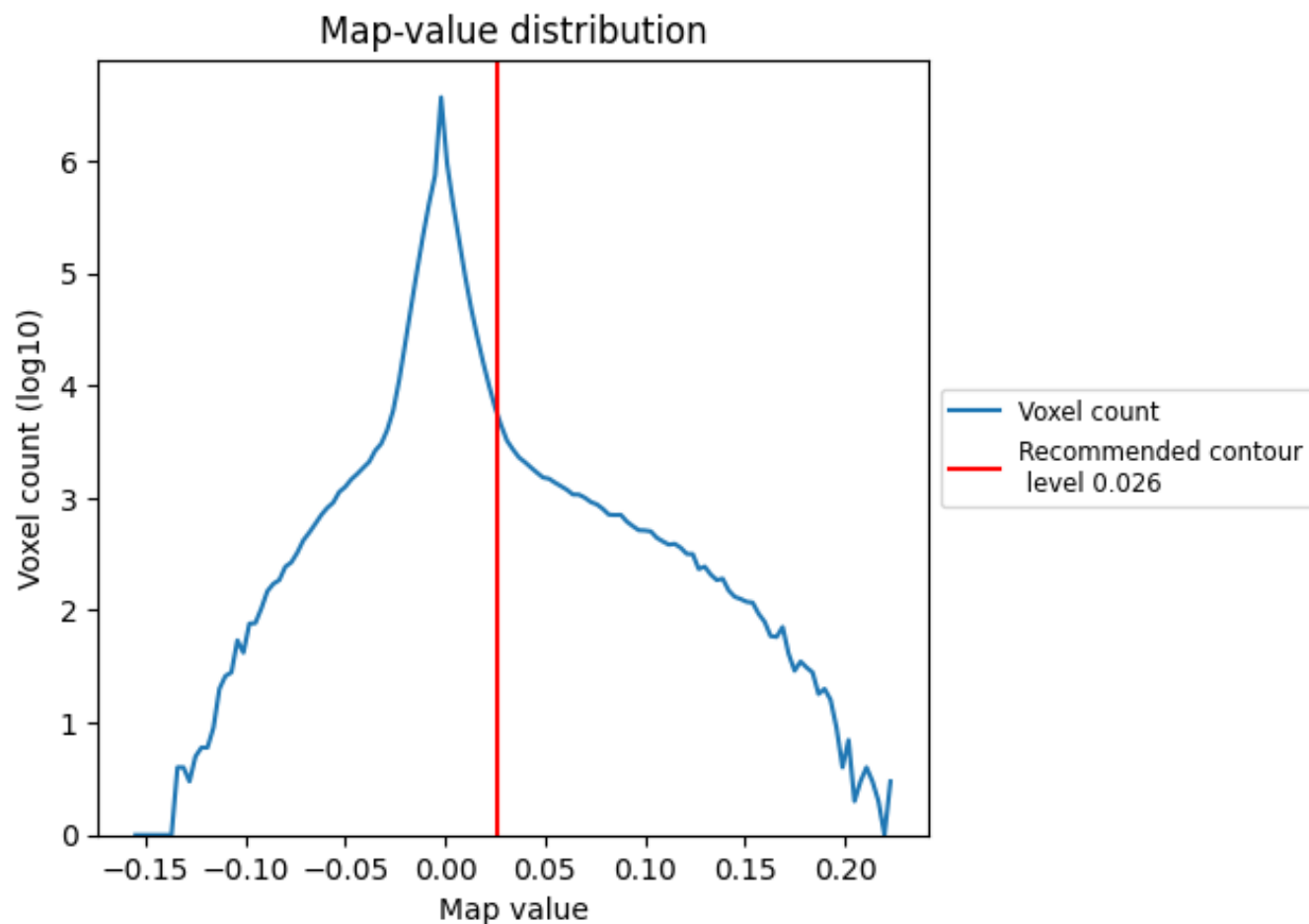
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

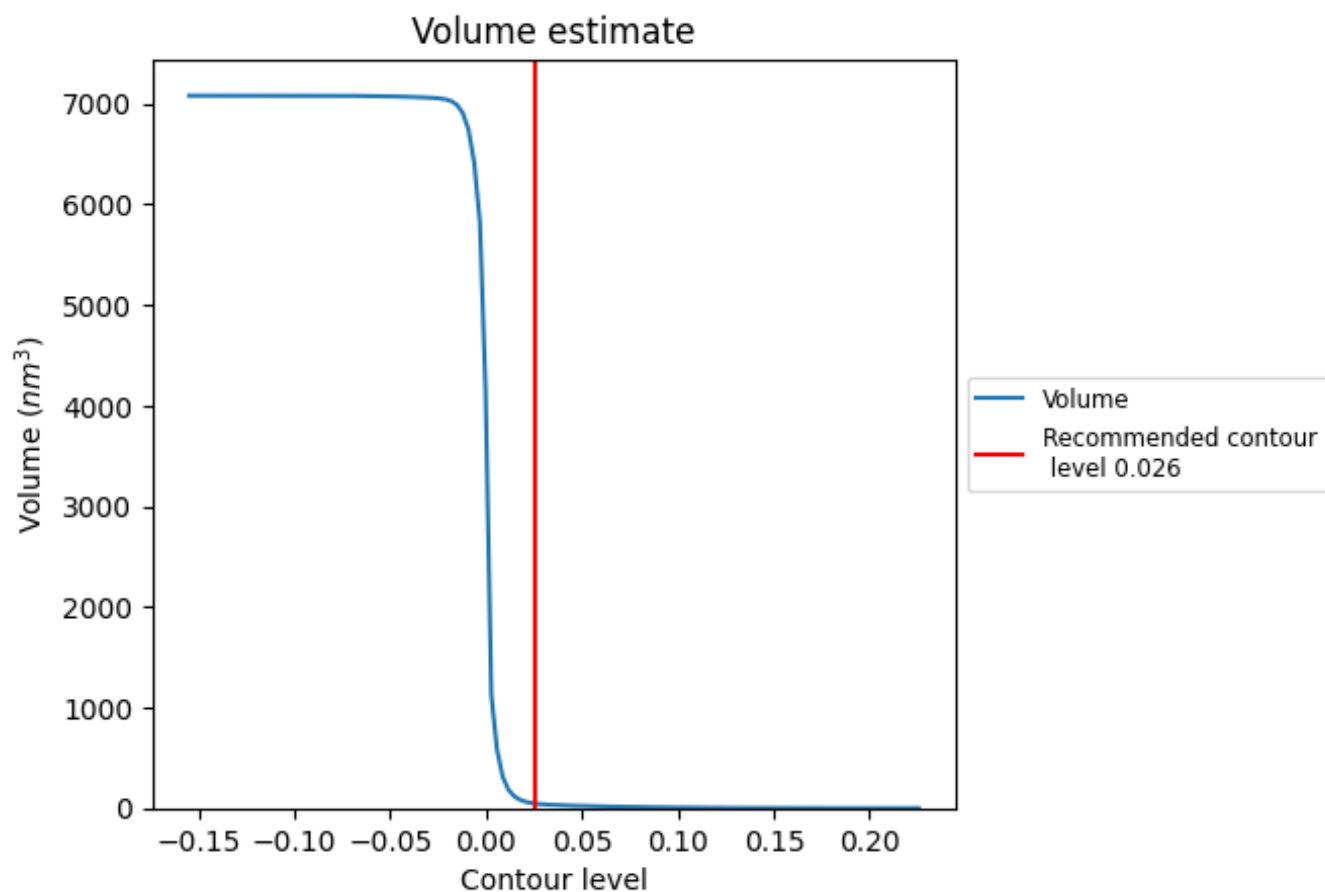
This section contains the results of statistical analysis of the map.

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.

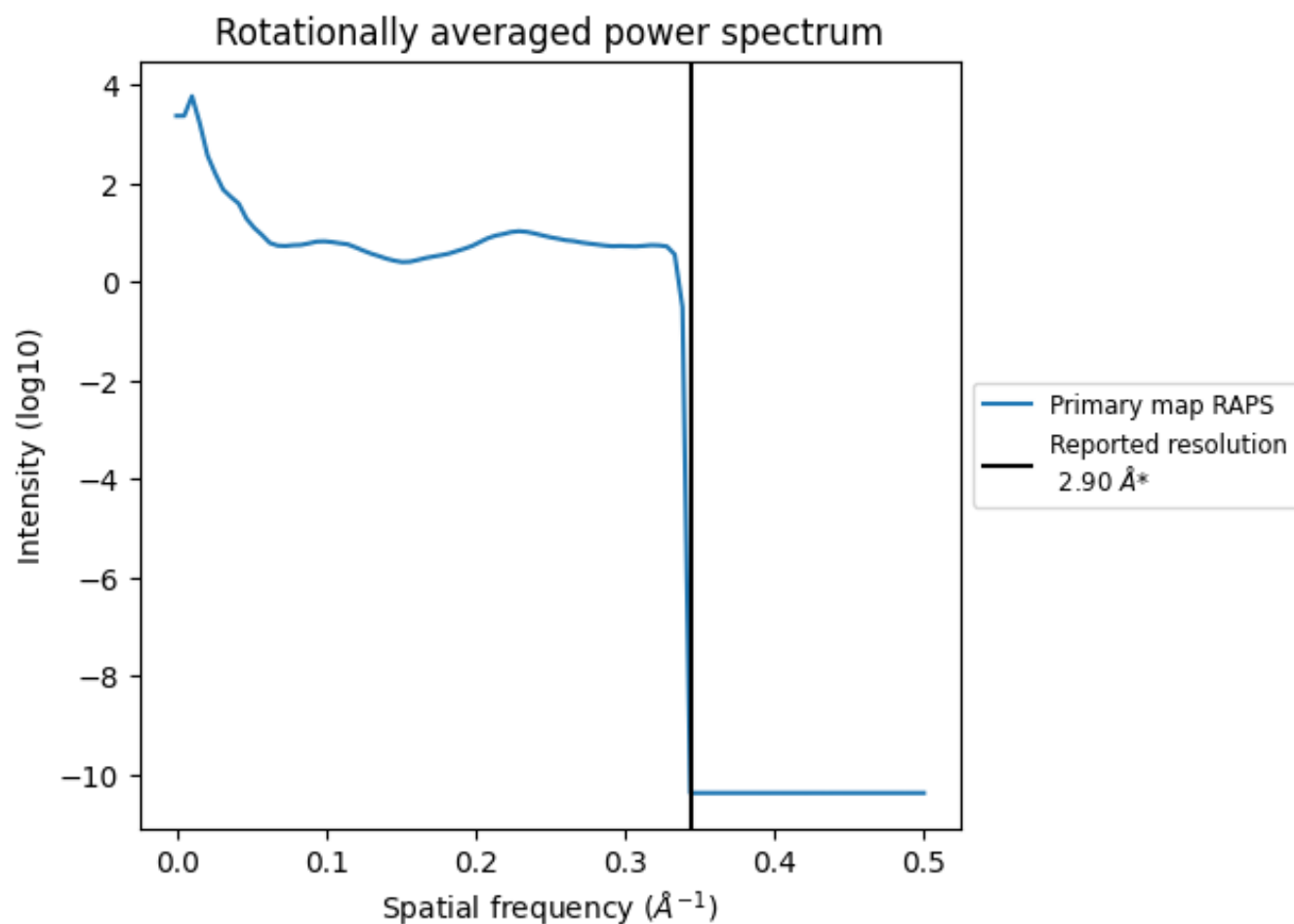
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 45 nm^3 ; this corresponds to an approximate mass of 40 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 0.345 Å⁻¹

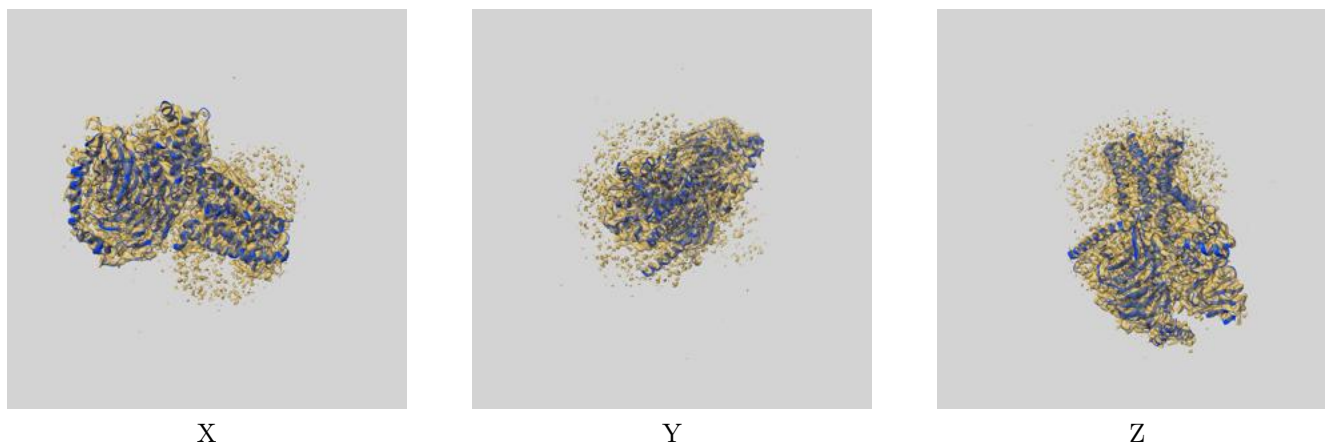
8 Fourier-Shell correlation

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

9 Map-model fit [i](#)

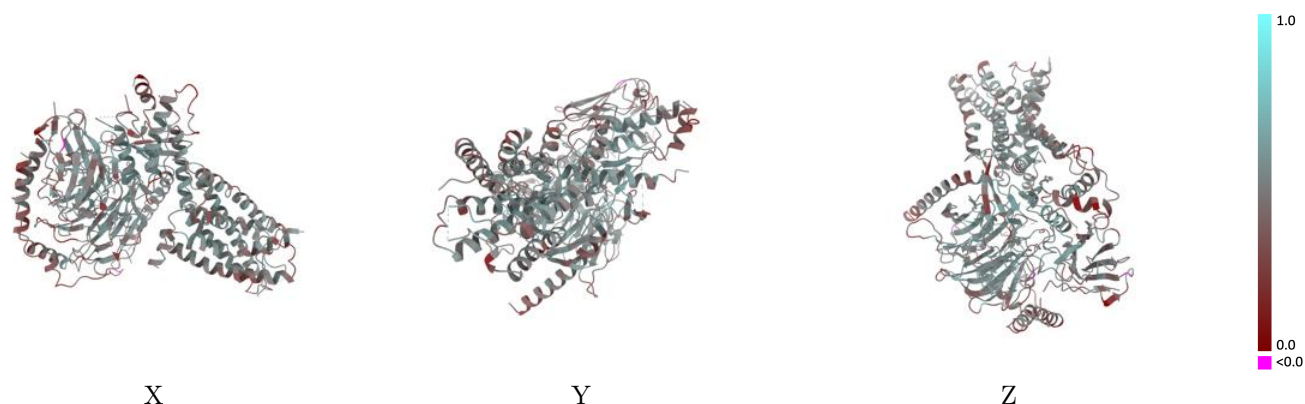
This section contains information regarding the fit between EMDB map EMD-32838 and PDB model 7WUQ. Per-residue inclusion information can be found in section [3](#) on page [7](#).

9.1 Map-model overlay [i](#)



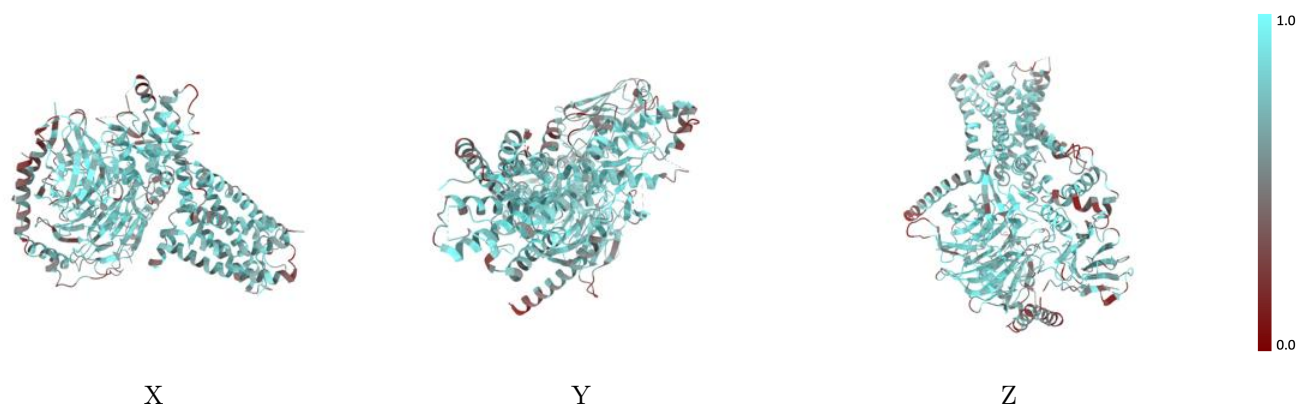
The images above show the 3D surface view of the map at the recommended contour level 0.026 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



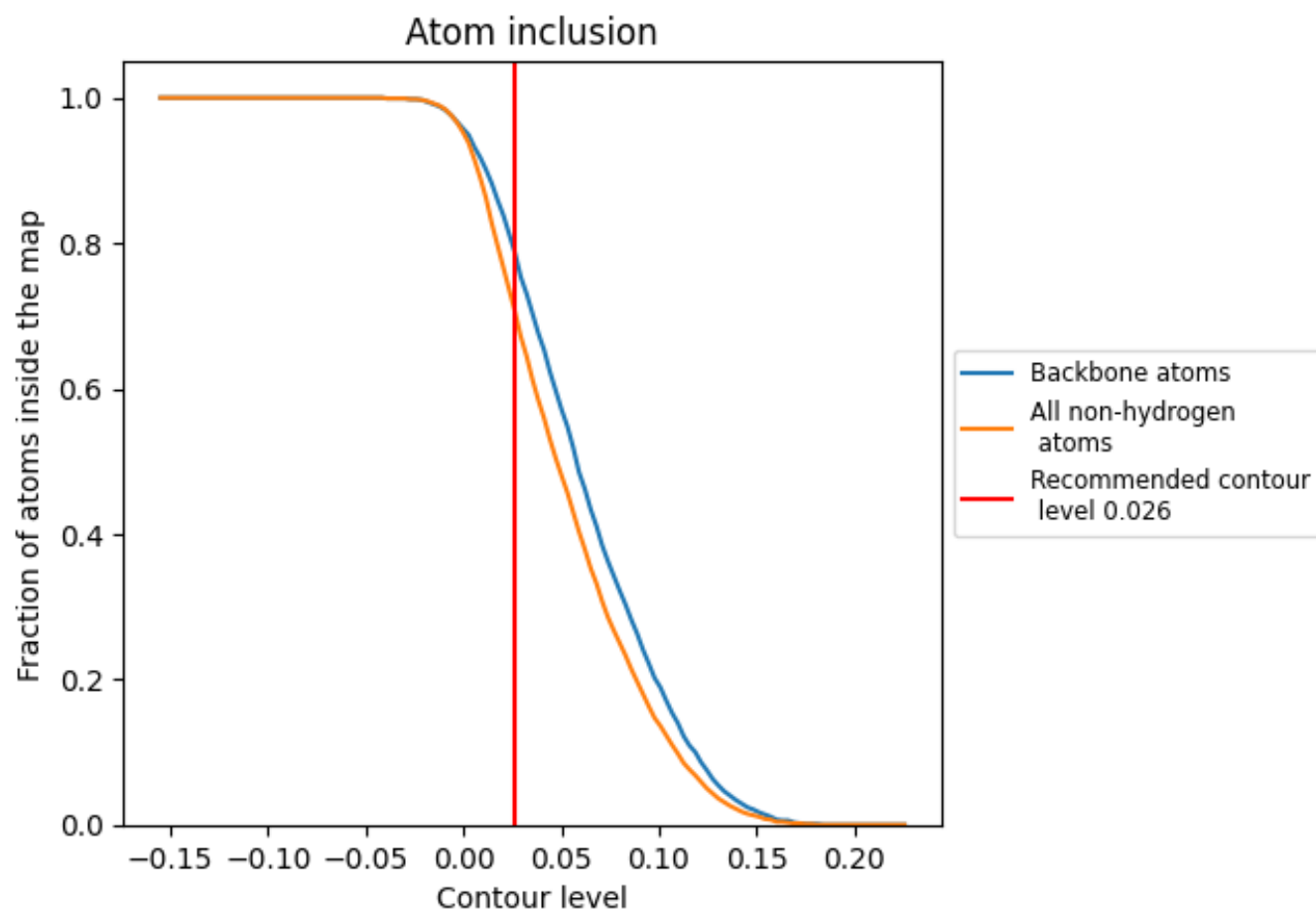
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.026).

9.4 Atom inclusion [i](#)



At the recommended contour level, 79% of all backbone atoms, 71% 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 (0.026) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7100	0.4650
A	0.7100	0.4660
B	0.7310	0.4690
G	0.5630	0.3790
N	0.7070	0.4620
R	0.7150	0.4780

