



Full wwPDB EM Validation Report ⓘ

Mar 26, 2026 – 01:57 AM UTC

PDB ID : 7M7I / pdb_00007m7i
EMDB ID : EMD-23714
Title : 6-Deoxyerythronolide B synthase (DEBS) module 1 in complex with antibody fragment 1B2 (TE-free)
Authors : Cogan, D.P.; Zhang, K.; Chiu, W.; Khosla, C.
Deposited on : 2021-03-28
Resolution : 3.40 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:

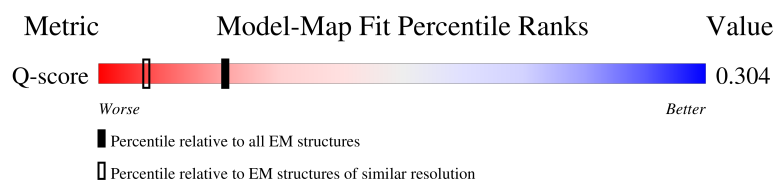
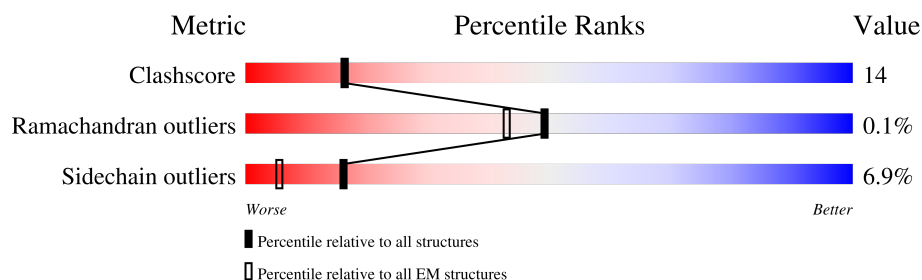
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 i

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

The reported resolution of this entry is 3.40 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	14717 (2.90 - 3.90)

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	1593	12% 58% 28% 13%
1	B	1593	7% 40% 21% 37%
2	C	249	49% 30% 18%
2	E	249	50% 31% 18%

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Mol	Chain	Length	Quality of chain
3	D	236	56%28%•13%
3	F	236	59%25%•13%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 23965 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 EryAI.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	999	Total	C	N	O	S	0	0
			7411	4610	1360	1414	27		
1	A	1390	Total	C	N	O	S	0	0
			10319	6413	1899	1972	35		

There are 112 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	expression tag	UNP Q5UNP6
B	2	ALA	-	expression tag	UNP Q5UNP6
B	3	SER	-	expression tag	UNP Q5UNP6
B	4	THR	-	expression tag	UNP Q5UNP6
B	5	ASP	-	expression tag	UNP Q5UNP6
B	6	SER	-	expression tag	UNP Q5UNP6
B	7	GLU	-	expression tag	UNP Q5UNP6
B	8	LYS	-	expression tag	UNP Q5UNP6
B	9	VAL	-	expression tag	UNP Q5UNP6
B	10	ALA	-	expression tag	UNP Q5UNP6
B	11	GLU	-	expression tag	UNP Q5UNP6
B	12	TYR	-	expression tag	UNP Q5UNP6
B	13	LEU	-	expression tag	UNP Q5UNP6
B	14	ARG	-	expression tag	UNP Q5UNP6
B	15	ARG	-	expression tag	UNP Q5UNP6
B	16	ALA	-	expression tag	UNP Q5UNP6
B	17	THR	-	expression tag	UNP Q5UNP6
B	18	LEU	-	expression tag	UNP Q5UNP6
B	19	ASP	-	expression tag	UNP Q5UNP6
B	20	LEU	-	expression tag	UNP Q5UNP6
B	21	ARG	-	expression tag	UNP Q5UNP6
B	22	ALA	-	expression tag	UNP Q5UNP6
B	23	ALA	-	expression tag	UNP Q5UNP6
B	24	ARG	-	expression tag	UNP Q5UNP6
B	25	GLN	-	expression tag	UNP Q5UNP6
B	26	ARG	-	expression tag	UNP Q5UNP6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	27	ILE	-	expression tag	UNP Q5UNP6
B	28	ARG	-	expression tag	UNP Q5UNP6
B	29	GLU	-	expression tag	UNP Q5UNP6
B	30	LEU	-	expression tag	UNP Q5UNP6
B	31	GLU	-	expression tag	UNP Q5UNP6
B	1486	THR	-	linker	UNP Q5UNP6
B	1487	SER	-	linker	UNP Q5UNP6
B	1488	GLU	-	linker	UNP Q5UNP6
B	1489	LEU	-	linker	UNP Q5UNP6
B	1490	GLY	-	linker	UNP Q5UNP6
B	1574	PRO	-	expression tag	UNP Q5UNP6
B	1575	ASN	-	expression tag	UNP Q5UNP6
B	1576	SER	-	expression tag	UNP Q5UNP6
B	1577	SER	-	expression tag	UNP Q5UNP6
B	1578	SER	-	expression tag	UNP Q5UNP6
B	1579	VAL	-	expression tag	UNP Q5UNP6
B	1580	ASP	-	expression tag	UNP Q5UNP6
B	1581	LYS	-	expression tag	UNP Q5UNP6
B	1582	LEU	-	expression tag	UNP Q5UNP6
B	1583	ALA	-	expression tag	UNP Q5UNP6
B	1584	ALA	-	expression tag	UNP Q5UNP6
B	1585	ALA	-	expression tag	UNP Q5UNP6
B	1586	LEU	-	expression tag	UNP Q5UNP6
B	1587	GLU	-	expression tag	UNP Q5UNP6
B	1588	HIS	-	expression tag	UNP Q5UNP6
B	1589	HIS	-	expression tag	UNP Q5UNP6
B	1590	HIS	-	expression tag	UNP Q5UNP6
B	1591	HIS	-	expression tag	UNP Q5UNP6
B	1592	HIS	-	expression tag	UNP Q5UNP6
B	1593	HIS	-	expression tag	UNP Q5UNP6
A	1	MET	-	expression tag	UNP Q5UNP6
A	2	ALA	-	expression tag	UNP Q5UNP6
A	3	SER	-	expression tag	UNP Q5UNP6
A	4	THR	-	expression tag	UNP Q5UNP6
A	5	ASP	-	expression tag	UNP Q5UNP6
A	6	SER	-	expression tag	UNP Q5UNP6
A	7	GLU	-	expression tag	UNP Q5UNP6
A	8	LYS	-	expression tag	UNP Q5UNP6
A	9	VAL	-	expression tag	UNP Q5UNP6
A	10	ALA	-	expression tag	UNP Q5UNP6
A	11	GLU	-	expression tag	UNP Q5UNP6
A	12	TYR	-	expression tag	UNP Q5UNP6

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Chain	Residue	Modelled	Actual	Comment	Reference
A	13	LEU	-	expression tag	UNP Q5UNP6
A	14	ARG	-	expression tag	UNP Q5UNP6
A	15	ARG	-	expression tag	UNP Q5UNP6
A	16	ALA	-	expression tag	UNP Q5UNP6
A	17	THR	-	expression tag	UNP Q5UNP6
A	18	LEU	-	expression tag	UNP Q5UNP6
A	19	ASP	-	expression tag	UNP Q5UNP6
A	20	LEU	-	expression tag	UNP Q5UNP6
A	21	ARG	-	expression tag	UNP Q5UNP6
A	22	ALA	-	expression tag	UNP Q5UNP6
A	23	ALA	-	expression tag	UNP Q5UNP6
A	24	ARG	-	expression tag	UNP Q5UNP6
A	25	GLN	-	expression tag	UNP Q5UNP6
A	26	ARG	-	expression tag	UNP Q5UNP6
A	27	ILE	-	expression tag	UNP Q5UNP6
A	28	ARG	-	expression tag	UNP Q5UNP6
A	29	GLU	-	expression tag	UNP Q5UNP6
A	30	LEU	-	expression tag	UNP Q5UNP6
A	31	GLU	-	expression tag	UNP Q5UNP6
A	1486	THR	-	linker	UNP Q5UNP6
A	1487	SER	-	linker	UNP Q5UNP6
A	1488	GLU	-	linker	UNP Q5UNP6
A	1489	LEU	-	linker	UNP Q5UNP6
A	1490	GLY	-	linker	UNP Q5UNP6
A	1574	PRO	-	expression tag	UNP Q5UNP6
A	1575	ASN	-	expression tag	UNP Q5UNP6
A	1576	SER	-	expression tag	UNP Q5UNP6
A	1577	SER	-	expression tag	UNP Q5UNP6
A	1578	SER	-	expression tag	UNP Q5UNP6
A	1579	VAL	-	expression tag	UNP Q5UNP6
A	1580	ASP	-	expression tag	UNP Q5UNP6
A	1581	LYS	-	expression tag	UNP Q5UNP6
A	1582	LEU	-	expression tag	UNP Q5UNP6
A	1583	ALA	-	expression tag	UNP Q5UNP6
A	1584	ALA	-	expression tag	UNP Q5UNP6
A	1585	ALA	-	expression tag	UNP Q5UNP6
A	1586	LEU	-	expression tag	UNP Q5UNP6
A	1587	GLU	-	expression tag	UNP Q5UNP6
A	1588	HIS	-	expression tag	UNP Q5UNP6
A	1589	HIS	-	expression tag	UNP Q5UNP6
A	1590	HIS	-	expression tag	UNP Q5UNP6
A	1591	HIS	-	expression tag	UNP Q5UNP6

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1592	HIS	-	expression tag	UNP Q5UNP6
A	1593	HIS	-	expression tag	UNP Q5UNP6

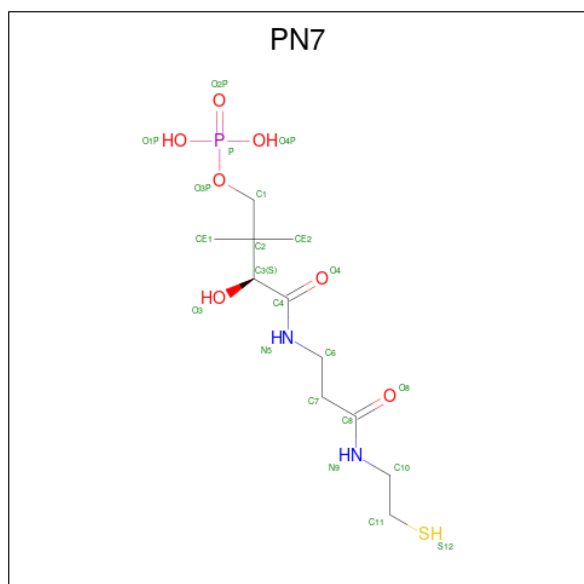
- Molecule 2 is a protein called 1B2 (light chain).

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	205	Total	C	N	O	S	0	0
			1539	978	257	298	6		
2	E	205	Total	C	N	O	S	0	0
			1539	978	257	298	6		

- Molecule 3 is a protein called 1B2 (heavy chain).

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	206	Total	C	N	O	S	0	0
			1568	983	262	317	6		
3	F	206	Total	C	N	O	S	0	0
			1568	984	262	316	6		

- Molecule 4 is N 3 -[(2S)-2-hydroxy-3,3-dimethyl-4-(phosphonooxy)butanoyl]-N-(2-sulfanylet hyl)-beta-alaninamide (CCD ID: PN7) (formula: C₁₁H₂₃N₂O₇PS).

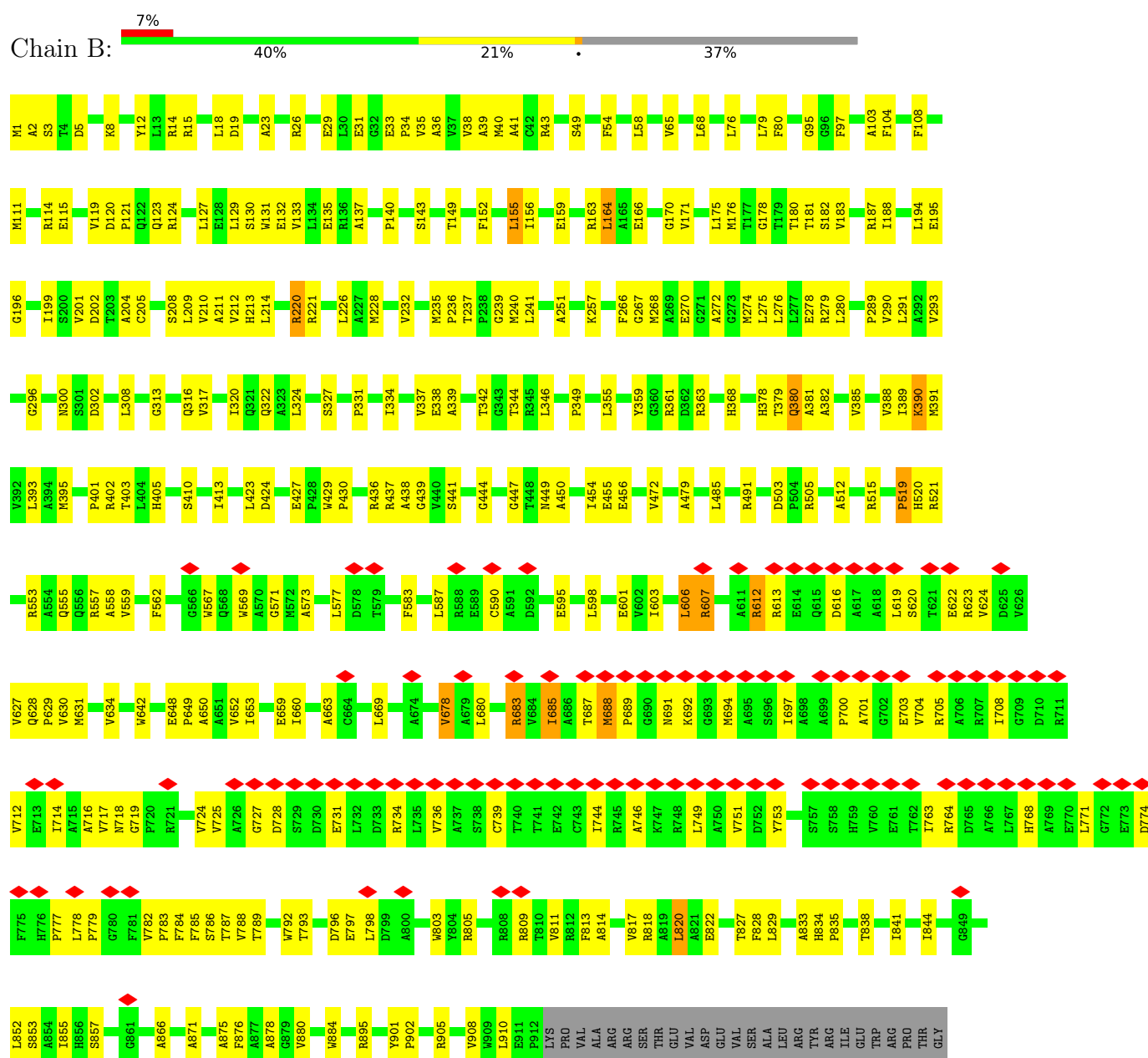


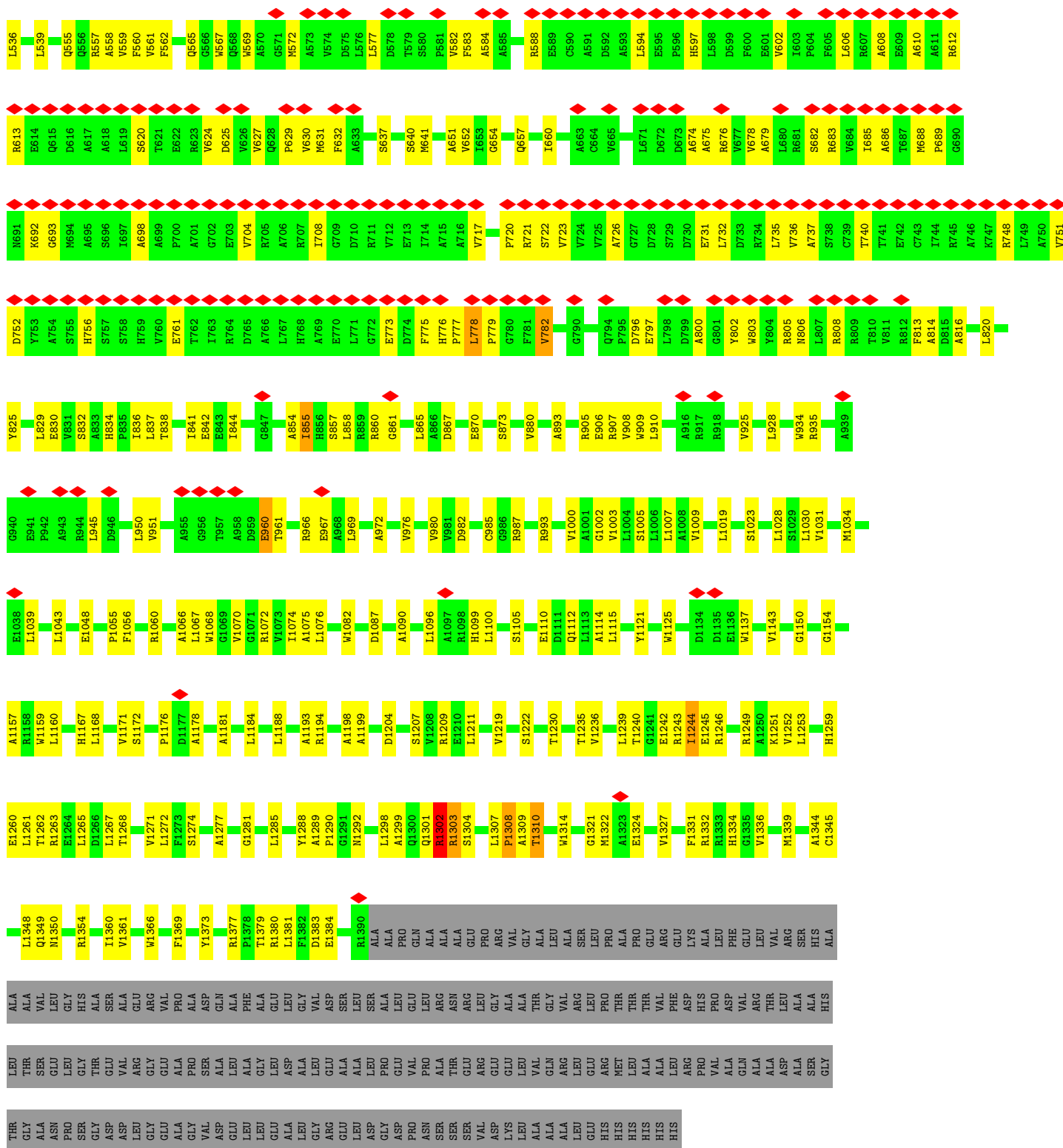
Mol	Chain	Residues	Atoms					AltConf
4	B	1	Total	C	N	O	P S	0
			21	11	2	6	1 1	

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

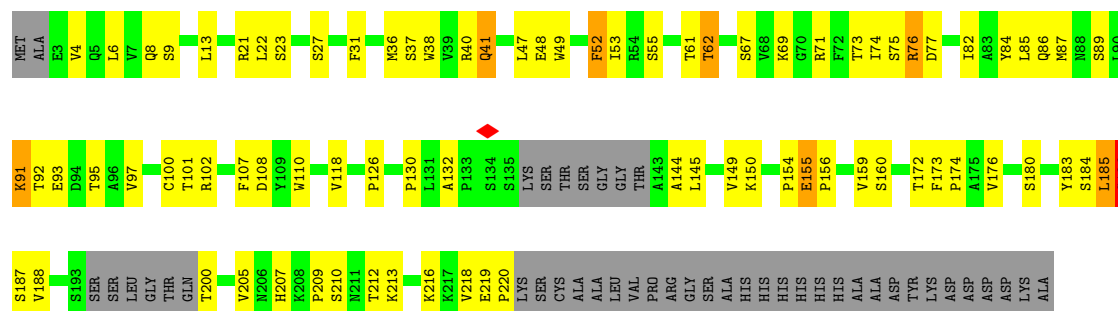
• Molecule 1: EryAI





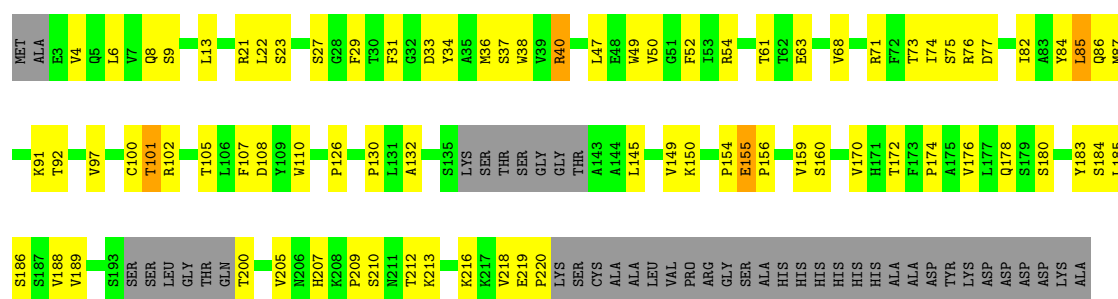
• Molecule 2: 1B2 (light chain)

Chain C: 49% 30% 18%



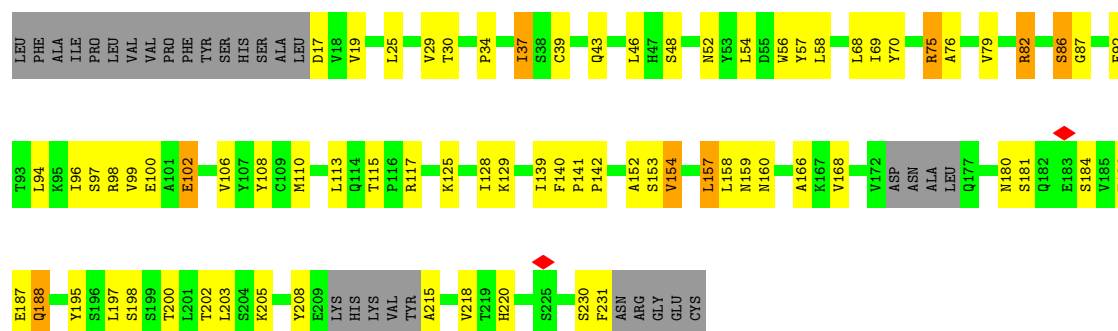
• Molecule 2: 1B2 (light chain)

Chain E: 50% 31% 18%



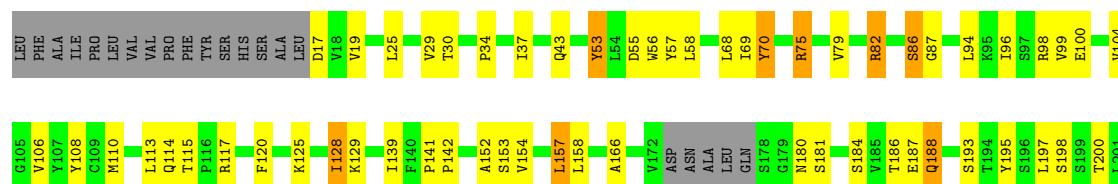
• Molecule 3: 1B2 (heavy chain)

Chain D: 56% 28% 13%



• Molecule 3: 1B2 (heavy chain)

Chain F: 59% 25% 13%





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	69566	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	4.722	Depositor
Minimum map value	-2.257	Depositor
Average map value	0.010	Depositor
Map value standard deviation	0.123	Depositor
Recommended contour level	0.5	Depositor
Map size (Å)	336.0, 336.0, 336.0	wwPDB
Map dimensions	336, 336, 336	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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PN7

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	0.34	0/10522	0.62	16/14323 (0.1%)
1	B	0.45	2/7560 (0.0%)	0.70	14/10286 (0.1%)
2	C	0.57	1/1575 (0.1%)	0.86	7/2141 (0.3%)
2	E	0.49	0/1575	0.82	6/2141 (0.3%)
3	D	0.42	0/1601	0.64	0/2175
3	F	0.35	0/1601	0.59	1/2174 (0.0%)
All	All	0.41	3/24434 (0.0%)	0.68	44/33240 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
2	C	0	1
2	E	0	1
3	D	0	1
3	F	0	1
All	All	0	5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	519	PRO	CG-CD	-13.61	1.04	1.50
1	B	1449	SER	CA-CB	-6.76	1.42	1.53
2	C	187	SER	CA-CB	-5.93	1.44	1.53

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	519	PRO	N-CD-CG	-18.24	75.84	103.20
1	A	233	THR	CB-CA-C	-18.17	83.19	110.62
1	B	519	PRO	CA-CB-CG	-11.37	82.91	104.50
1	A	221	ARG	CB-CA-C	-10.86	92.94	110.86
1	B	519	PRO	CA-N-CD	-8.75	99.75	112.00
1	B	685	ILE	N-CA-C	-8.48	105.22	113.53
1	A	1310	THR	CB-CA-C	-8.27	96.41	109.72
2	E	40	ARG	CA-C-N	7.72	134.57	123.14
2	E	40	ARG	C-N-CA	7.72	134.57	123.14
2	C	48	GLU	CB-CA-C	-7.51	99.27	110.24
1	B	519	PRO	CB-CG-CD	7.36	129.66	106.10
1	B	390	LYS	N-CA-C	-7.01	103.33	110.97
3	F	53	TYR	CB-CA-C	-7.00	102.15	111.74
2	E	101	THR	CB-CA-C	6.47	121.20	109.37
1	A	1301	GLN	CB-CA-C	-6.46	98.73	110.56
2	E	85	LEU	CA-C-N	-6.43	113.61	122.30
2	E	85	LEU	C-N-CA	-6.43	113.61	122.30
1	A	782	VAL	CA-C-O	-6.42	115.36	119.15
2	C	185	LEU	O-C-N	6.42	131.10	123.33
1	A	1308	PRO	CB-CA-C	-6.39	100.10	110.25
1	B	1411	GLU	N-CA-C	-6.25	104.57	111.82
2	E	40	ARG	CA-C-O	-6.24	114.06	121.11
2	C	41	GLN	CB-CA-C	-6.11	101.04	111.31
1	A	1302	ARG	N-CA-C	-5.96	104.78	111.28
1	B	221	ARG	N-CA-C	-5.95	106.64	114.31
1	A	1308	PRO	N-CD-CG	-5.94	94.29	103.20
1	A	8	LYS	N-CA-C	-5.75	102.78	111.34
2	C	52	PHE	CA-CB-CG	5.73	119.53	113.80
1	B	774	ASP	N-CA-C	-5.69	108.13	114.62
2	C	41	GLN	CB-CG-CD	-5.68	102.95	112.60
1	A	960	GLU	N-CA-C	-5.67	106.20	113.01
1	B	380	GLN	CB-CA-C	-5.66	109.53	117.23
1	A	1304	SER	N-CA-C	-5.61	105.70	112.54
1	A	15	ARG	N-CA-C	-5.59	105.10	111.14
1	A	8	LYS	CB-CA-C	-5.41	104.04	112.31
2	C	40	ARG	N-CA-CB	-5.34	101.72	110.68
1	B	774	ASP	CB-CA-C	5.24	116.25	109.28
1	B	221	ARG	CB-CA-C	-5.10	100.45	109.02
1	A	44	LEU	CA-C-O	-5.09	115.63	120.97
1	A	44	LEU	CA-C-N	5.06	126.16	119.84
1	A	44	LEU	C-N-CA	5.06	126.16	119.84
1	B	220	ARG	N-CA-C	-5.05	106.70	112.92
2	C	186	SER	N-CA-C	-5.05	100.05	108.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1427	VAL	N-CA-C	-5.02	107.04	113.22

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	797	GLU	Peptide
2	C	155	GLU	Peptide
3	D	115	THR	Peptide
2	E	155	GLU	Peptide
3	F	115	THR	Peptide

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	10319	0	10164	307	0
1	B	7411	0	7295	237	0
2	C	1539	0	1511	48	0
2	E	1539	0	1511	44	0
3	D	1568	0	1528	38	0
3	F	1568	0	1533	37	0
4	B	21	0	22	3	0
All	All	23965	0	23564	675	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 14.

All (675) 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:A:1302:ARG:HG2	1:A:1309:ALA:HB2	1.66	0.77
1:B:558:ALA:HA	1:B:827:THR:HG23	1.68	0.76
1:A:565:GLN:HE21	1:A:751:VAL:HG11	1.53	0.74
1:A:777:PRO:HB3	1:A:796:ASP:HA	1.71	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:565:GLN:OE1	1:A:657:GLN:NE2	2.24	0.71
3:F:158:LEU:HB3	3:F:197:LEU:HB3	1.73	0.70
1:B:577:LEU:HD12	1:B:587:LEU:HD23	1.74	0.70
1:B:235:MET:HE1	1:B:241:LEU:HG	1.72	0.69
1:A:148:PRO:HB2	1:A:224:SER:HA	1.75	0.69
1:A:602:VAL:HG13	1:A:630:VAL:HG22	1.73	0.69
1:B:123:GLN:NE2	1:B:180:THR:O	2.25	0.69
1:B:557:ARG:HG3	1:B:650:ALA:HB2	1.76	0.68
1:A:367:LEU:HB3	1:A:420:ILE:HG12	1.76	0.68
1:A:302:ASP:OD1	1:A:449:ASN:ND2	2.26	0.67
1:B:559:VAL:HG11	1:B:820:LEU:HD11	1.76	0.67
1:B:205:CYS:HB2	1:B:444:GLY:HA2	1.77	0.67
1:B:214:LEU:HD11	1:A:214:LEU:HD22	1.76	0.67
1:B:232:VAL:HG23	1:B:272:ALA:HB2	1.75	0.67
1:B:491:ARG:HD3	1:B:902:PRO:HG3	1.77	0.67
1:A:368:HIS:HB3	1:A:423:LEU:HD21	1.75	0.67
3:D:184:SER:HB3	3:D:198:SER:HB3	1.77	0.66
3:F:17:ASP:OD1	3:F:43:GLN:NE2	2.28	0.66
1:A:572:MET:HE3	1:A:631:MET:HE1	1.77	0.66
3:F:184:SER:HB3	3:F:198:SER:HB3	1.77	0.66
1:A:121:PRO:HA	1:A:124:ARG:HG2	1.76	0.66
3:D:17:ASP:OD1	3:D:43:GLN:NE2	2.28	0.66
2:E:149:VAL:HG13	2:E:205:VAL:HG11	1.78	0.65
1:A:625:ASP:HB3	1:A:686:ALA:HB2	1.77	0.65
1:A:1260:GLU:O	1:A:1263:ARG:NH1	2.29	0.65
1:B:388:VAL:HG22	1:B:454:ILE:HD11	1.78	0.65
1:A:980:VAL:O	1:A:993:ARG:NH1	2.30	0.65
2:C:149:VAL:HG13	2:C:205:VAL:HG11	1.78	0.65
3:D:158:LEU:HB3	3:D:197:LEU:HB3	1.78	0.64
1:A:39:ALA:HB2	1:A:137:ALA:HB2	1.80	0.64
1:B:58:LEU:HD21	1:B:390:LYS:HG3	1.78	0.64
1:A:69:PRO:HB3	1:A:72:ARG:HH11	1.62	0.64
1:A:1184:LEU:O	1:A:1188:LEU:HG	1.97	0.64
1:B:555:GLN:NE2	1:B:878:ALA:O	2.30	0.64
1:B:1416:LEU:HD21	1:B:1485:LEU:HB3	1.79	0.64
1:A:505:ARG:HH11	1:A:893:ALA:HA	1.62	0.64
1:B:43:ARG:HG2	1:B:129:LEU:HD21	1.80	0.64
1:A:474:PRO:HB3	1:A:873:SER:HB2	1.80	0.64
1:A:526:PRO:HG3	1:A:532:ALA:HB2	1.80	0.64
1:A:72:ARG:NH2	1:A:98:LEU:O	2.28	0.63
1:A:209:LEU:HD11	1:A:440:VAL:HG12	1.78	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:SER:HB3	1:A:188:ILE:HD13	1.78	0.63
1:B:181:THR:OG1	1:A:202:ASP:OD2	2.14	0.63
1:A:1143:VAL:HB	1:A:1168:LEU:HD12	1.80	0.63
1:B:289:PRO:HG3	1:B:895:ARG:HE	1.64	0.63
1:B:553:ARG:NH2	1:B:1444:GLU:HG3	2.15	0.62
1:A:92:GLN:HB2	1:A:250:LEU:HD23	1.82	0.62
1:B:182:SER:OG	1:A:202:ASP:OD1	2.14	0.62
2:E:9:SER:OG	2:E:23:SER:OG	2.18	0.62
2:E:132:ALA:HB1	2:E:220:PRO:HA	1.81	0.62
1:B:130:SER:HB2	1:B:228:MET:HE1	1.82	0.62
2:E:8:GLN:NE2	2:E:100:CYS:H	1.98	0.62
1:B:553:ARG:NH1	1:B:1440:GLN:HE22	1.98	0.62
1:A:57:LEU:HG	1:A:62:ARG:HB3	1.82	0.62
1:A:300:ASN:HD21	1:A:318:ARG:HB2	1.65	0.62
1:A:329:LEU:O	3:F:98:ARG:NH1	2.33	0.62
1:A:1176:PRO:HD2	1:A:1199:ALA:HB2	1.81	0.62
2:E:174:PRO:HD3	3:F:186:THR:HG22	1.82	0.62
1:A:175:LEU:O	1:A:179:THR:OG1	2.17	0.61
1:B:379:THR:HB	1:B:382:ALA:HB3	1.82	0.61
1:A:1000:VAL:HG12	1:A:1002:GLY:H	1.66	0.61
1:B:627:VAL:HG22	1:B:631:MET:SD	2.41	0.61
2:C:9:SER:OG	2:C:23:SER:OG	2.18	0.61
1:A:1048:GLU:HG3	1:A:1090:ALA:HA	1.82	0.61
2:C:22:LEU:HD12	2:C:85:LEU:HD23	1.81	0.61
2:C:132:ALA:HB1	2:C:220:PRO:HA	1.82	0.61
1:B:817:VAL:HG21	1:B:841:ILE:HD12	1.83	0.60
1:B:1427:VAL:HG22	1:B:1451:SER:HA	1.84	0.60
2:C:8:GLN:NE2	2:C:100:CYS:H	2.00	0.60
1:B:829:LEU:HD11	1:B:880:VAL:HG21	1.83	0.60
1:A:478:SER:O	1:A:488:GLN:NE2	2.34	0.60
1:B:736:VAL:HG22	1:B:746:ALA:HB3	1.83	0.60
2:C:126:PRO:HD3	2:C:207:HIS:HB2	1.84	0.60
1:B:124:ARG:NH1	1:B:908:VAL:O	2.35	0.60
1:B:1428:LEU:HD21	1:B:1447:VAL:HG22	1.82	0.60
1:A:1072:ARG:HG2	1:A:1112:GLN:HE21	1.67	0.60
1:B:308:LEU:HD21	1:A:183:VAL:HB	1.83	0.60
1:A:338:GLU:N	1:A:439:GLY:O	2.32	0.60
1:A:1000:VAL:HB	1:A:1039:LEU:HD21	1.83	0.60
1:A:268:MET:HA	1:A:378:HIS:HB3	1.83	0.59
1:A:637:SER:O	1:A:641:MET:HG2	2.02	0.59
1:A:704:VAL:HG21	1:A:723:VAL:HG21	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:VAL:N	1:B:276:LEU:O	2.30	0.59
1:B:155:LEU:HD13	1:B:181:THR:HB	1.84	0.59
1:A:560:PHE:HB2	1:A:652:VAL:HG12	1.85	0.59
1:B:569:TRP:HD1	1:B:571:GLY:H	1.50	0.59
1:A:141:PRO:HG2	1:A:516:ALA:HB2	1.84	0.59
1:A:522:ALA:HB1	1:A:539:LEU:HD11	1.85	0.59
2:E:22:LEU:HD12	2:E:85:LEU:HD23	1.85	0.59
1:A:1168:LEU:HD23	1:A:1188:LEU:HD13	1.84	0.58
1:A:1262:THR:HA	1:A:1265:LEU:HD13	1.85	0.58
1:A:390:LYS:HE3	1:A:401:PRO:HG2	1.84	0.58
2:E:126:PRO:HD3	2:E:207:HIS:HB2	1.84	0.58
1:B:121:PRO:HA	1:B:124:ARG:HG2	1.85	0.58
1:A:1259:HIS:CE1	1:A:1302:ARG:HB2	2.37	0.58
1:A:1171:VAL:HG21	1:A:1211:LEU:HD21	1.85	0.58
2:E:188:VAL:HG11	3:F:157:LEU:HD11	1.86	0.58
1:B:140:PRO:HB2	1:B:143:SER:HB3	1.85	0.58
1:A:340:HIS:CD2	1:A:443:PHE:H	2.22	0.58
1:B:813:PHE:CE2	1:B:841:ILE:HD11	2.39	0.58
1:A:1075:ALA:HB2	1:A:1082:TRP:HD1	1.68	0.58
2:C:41:GLN:HG2	2:C:47:LEU:HD23	1.85	0.58
1:A:717:VAL:O	1:A:814:ALA:N	2.36	0.58
1:A:126:MET:HE2	1:A:188:ILE:HD12	1.85	0.57
1:A:124:ARG:NH1	1:A:908:VAL:O	2.37	0.57
2:C:174:PRO:HD3	3:D:186:THR:HG22	1.87	0.57
1:B:555:GLN:HE21	1:B:827:THR:HB	1.69	0.57
1:B:835:PRO:O	1:B:838:THR:OG1	2.21	0.57
1:B:855:ILE:HG12	1:B:875:ALA:HA	1.86	0.57
1:A:43:ARG:HG2	1:A:125:LEU:HD22	1.84	0.57
1:A:103:ALA:HB1	1:A:905:ARG:HD3	1.85	0.57
2:E:176:VAL:HB	2:E:184:SER:HB2	1.85	0.57
1:B:649:PRO:HB2	1:B:782:VAL:HG11	1.87	0.57
1:A:1:MET:HE1	2:C:52:PHE:CZ	2.39	0.57
1:B:15:ARG:NH1	2:E:102:ARG:HH22	2.03	0.57
1:A:1075:ALA:HB2	1:A:1082:TRP:CD1	2.40	0.57
1:B:1:MET:SD	1:B:2:ALA:N	2.77	0.57
2:C:176:VAL:HB	2:C:184:SER:HB2	1.85	0.57
2:E:38:TRP:CD1	2:E:85:LEU:HD13	2.40	0.56
1:A:1087:ASP:HB3	1:A:1114:ALA:HA	1.87	0.56
1:B:26:ARG:HA	1:B:29:GLU:HG2	1.87	0.56
1:B:120:ASP:HB3	1:B:123:GLN:HG3	1.86	0.56
1:B:627:VAL:O	1:B:630:VAL:HG22	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:830:GLU:HB2	1:A:838:THR:HG23	1.86	0.56
3:D:180:ASN:ND2	3:D:202:THR:O	2.39	0.56
1:B:788:VAL:HG13	1:B:789:THR:HG23	1.86	0.56
1:A:244:PHE:HB3	1:A:250:LEU:HD13	1.87	0.56
1:A:208:SER:HB2	1:A:385:VAL:HB	1.87	0.56
1:A:582:VAL:HG23	1:A:641:MET:HE1	1.87	0.56
1:A:1274:SER:HB2	1:A:1292:ASN:HB3	1.88	0.56
3:F:180:ASN:ND2	3:F:202:THR:O	2.39	0.56
1:B:1:MET:SD	1:B:3:SER:N	2.78	0.56
1:A:610:ALA:HB2	1:A:613:ARG:HH21	1.70	0.56
2:C:95:THR:HG23	2:C:118:VAL:H	1.71	0.56
3:F:58:LEU:HB2	3:F:68:LEU:HD11	1.88	0.56
1:B:36:ALA:N	1:B:278:GLU:O	2.37	0.55
1:B:694:MET:HG3	1:B:751:VAL:HG11	1.87	0.55
1:A:1159:TRP:CZ2	1:A:1349:GLN:HB2	2.42	0.55
1:A:1272:LEU:HD13	1:A:1299:ALA:HB2	1.87	0.55
2:E:22:LEU:HB2	2:E:85:LEU:HB3	1.87	0.55
1:B:120:ASP:HB2	1:B:178:GLY:O	2.06	0.55
1:B:163:ARG:HB3	1:B:166:GLU:HG3	1.87	0.55
1:A:92:GLN:HE22	1:A:254:GLY:HA2	1.71	0.55
1:A:1350:ASN:OD1	1:A:1354:ARG:NH1	2.29	0.55
1:A:1019:LEU:HG	1:A:1252:VAL:HG11	1.89	0.55
1:A:1150:GLY:O	1:A:1154:GLY:N	2.36	0.55
1:A:608:ALA:HB1	1:A:612:ARG:HG3	1.87	0.55
2:C:156:PRO:HG2	2:C:209:PRO:HD3	1.89	0.55
3:F:56:TRP:HB2	3:F:69:ILE:HB	1.88	0.55
1:B:339:ALA:HA	1:B:441:SER:HB2	1.89	0.55
1:B:512:ALA:HB1	1:B:884:TRP:CG	2.42	0.55
3:D:58:LEU:HB2	3:D:68:LEU:HD11	1.88	0.55
1:B:853:SER:HB2	1:B:878:ALA:HB1	1.89	0.55
2:C:155:GLU:HG2	2:C:183:TYR:CD2	2.42	0.55
2:E:155:GLU:HG2	2:E:183:TYR:CD2	2.42	0.55
1:B:133:VAL:HG22	1:B:276:LEU:HB2	1.89	0.54
1:B:201:VAL:HG22	1:A:199:ILE:HD11	1.88	0.54
1:A:340:HIS:HD2	1:A:443:PHE:H	1.56	0.54
1:A:945:LEU:HD21	1:A:1105:SER:HB3	1.89	0.54
1:B:503:ASP:OD2	1:B:505:ARG:HG2	2.06	0.54
2:C:37:SER:HB2	2:C:101:THR:OG1	2.06	0.54
2:E:49:TRP:HB3	3:F:117:ARG:O	2.06	0.54
1:A:205:CYS:SG	1:A:378:HIS:NE2	2.74	0.54
1:A:756:HIS:ND1	1:A:806:ASN:O	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1271:VAL:HA	1:A:1310:THR:HB	1.88	0.54
1:A:1230:THR:HA	1:A:1322:MET:HE2	1.88	0.54
1:B:97:PHE:HA	1:B:270:GLU:HG2	1.90	0.54
1:A:15:ARG:HH21	3:D:76:ALA:HB1	1.71	0.54
1:A:68:LEU:HD22	1:A:74:TRP:HZ3	1.73	0.54
1:A:562:PHE:HB2	1:A:654:GLY:HA2	1.90	0.54
1:A:177:THR:O	1:A:187:ARG:NH2	2.41	0.54
1:B:739:CYS:HB2	1:B:744:ILE:HG23	1.90	0.54
1:A:300:ASN:ND2	1:A:318:ARG:HB2	2.23	0.53
1:A:567:TRP:CD2	1:A:832:SER:HB2	2.43	0.53
1:B:624:VAL:HA	1:B:627:VAL:HG12	1.90	0.53
1:A:492:LEU:HD23	1:A:536:LEU:HD21	1.89	0.53
1:B:567:TRP:HB3	1:B:834:HIS:HB3	1.91	0.53
1:A:149:THR:HB	1:A:194:LEU:HD23	1.89	0.53
1:A:736:VAL:HG11	1:A:748:ARG:HG2	1.90	0.53
2:C:144:ALA:HB3	3:D:140:PHE:HZ	1.73	0.53
1:A:493:ALA:HB2	1:A:536:LEU:HB3	1.91	0.53
1:A:1193:ALA:HB1	1:A:1194:ARG:HH11	1.73	0.53
1:B:181:THR:H	1:A:157:PRO:HG2	1.74	0.53
1:B:268:MET:HE1	1:B:380:GLN:HG3	1.90	0.53
1:A:64:ALA:HB3	1:A:376:LEU:HA	1.89	0.53
1:A:613:ARG:NH2	1:A:620:SER:OG	2.42	0.53
1:B:29:GLU:HA	1:B:33:GLU:HB2	1.91	0.53
1:B:237:THR:HG23	1:B:239:GLY:H	1.74	0.53
1:A:1321:GLY:HA2	1:A:1324:GLU:HB3	1.91	0.53
1:A:174:TYR:HA	1:A:177:THR:HG22	1.90	0.53
1:B:472:VAL:HG21	1:B:866:ALA:HA	1.91	0.53
1:A:122:GLN:HE22	1:A:184:ALA:HB1	1.73	0.53
1:A:126:MET:HE3	1:A:153:VAL:HG21	1.91	0.53
2:E:156:PRO:HG2	2:E:209:PRO:HD3	1.89	0.52
1:B:557:ARG:NH2	1:B:648:GLU:O	2.43	0.52
1:B:719:GLY:HA2	1:B:844:ILE:HD11	1.92	0.52
1:A:629:PRO:HG3	1:A:679:ALA:HA	1.91	0.52
1:A:1068:TRP:HZ2	1:A:1087:ASP:HB2	1.74	0.52
1:A:624:VAL:HA	1:A:627:VAL:HG12	1.90	0.52
1:A:300:ASN:HB3	1:A:319:VAL:HB	1.91	0.52
3:D:142:PRO:HD3	3:D:154:VAL:HB	1.91	0.52
2:E:38:TRP:HE1	2:E:74:ILE:HG23	1.75	0.52
1:A:1236:VAL:HA	1:A:1239:LEU:HB2	1.92	0.52
1:B:35:VAL:HA	1:B:279:ARG:HA	1.91	0.52
1:A:141:PRO:O	1:A:192:LEU:HD11	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:266:PHE:HD2	1:B:378:HIS:HD1	1.57	0.52
1:B:324:LEU:HD21	1:B:334:ILE:HG12	1.92	0.52
1:A:100:GLU:HB3	1:A:125:LEU:HD21	1.91	0.52
1:A:558:ALA:HB2	1:A:880:VAL:HG13	1.92	0.52
1:A:1031:VAL:HG21	1:A:1074:ILE:HD13	1.92	0.52
2:C:38:TRP:HE1	2:C:74:ILE:HG23	1.75	0.52
2:E:4:VAL:HA	2:E:27:SER:O	2.10	0.52
1:B:209:LEU:O	1:B:212:VAL:HG22	2.10	0.52
1:B:783:PRO:HB3	1:B:792:TRP:HE1	1.74	0.52
1:A:1299:ALA:O	1:A:1303:ARG:HB3	2.10	0.52
2:C:4:VAL:HA	2:C:27:SER:O	2.10	0.52
1:B:199:ILE:HG23	1:A:201:VAL:HG23	1.91	0.51
1:B:334:ILE:HG23	1:B:438:ALA:HA	1.92	0.51
1:A:682:SER:HA	1:A:685:ILE:HG22	1.91	0.51
2:E:145:LEU:HD23	2:E:218:VAL:HG13	1.92	0.51
1:B:183:VAL:HG12	1:A:446:SER:HA	1.91	0.51
1:B:403:THR:N	1:B:423:LEU:O	2.43	0.51
1:A:1019:LEU:HD13	1:A:1244:ILE:HG22	1.91	0.51
1:B:104:PHE:HB2	1:B:124:ARG:HD3	1.92	0.51
1:B:798:LEU:HD11	1:B:803:TRP:CH2	2.46	0.51
3:D:39:CYS:HB2	3:D:56:TRP:CH2	2.45	0.51
1:A:65:VAL:HG12	1:A:96:GLY:N	2.26	0.51
1:A:133:VAL:HG22	1:A:276:LEU:HB2	1.93	0.51
1:B:569:TRP:CD1	1:B:571:GLY:H	2.29	0.51
1:B:805:ARG:HG2	1:B:809:ARG:HH12	1.76	0.51
1:A:800:ALA:HA	1:A:803:TRP:CD1	2.46	0.51
1:A:820:LEU:HB3	1:A:825:TYR:HB2	1.93	0.51
2:E:101:THR:HB	2:E:107:PHE:HB3	1.92	0.51
1:A:1377:ARG:NH1	1:A:1379:THR:HG21	2.26	0.50
1:B:688:MET:HE3	1:B:753:TYR:HB2	1.93	0.50
1:B:704:VAL:HG13	1:B:714:ILE:HD11	1.93	0.50
1:A:24:ARG:CZ	3:F:75:ARG:HB2	2.41	0.50
1:A:945:LEU:N	1:A:972:ALA:O	2.43	0.50
1:B:828:PHE:HE1	1:B:852:LEU:HB3	1.77	0.50
1:A:982:ASP:HB2	1:A:985:CYS:HB3	1.93	0.50
2:C:145:LEU:HD23	2:C:218:VAL:HG13	1.92	0.50
1:B:716:ALA:HB3	1:B:724:VAL:HB	1.93	0.50
1:B:135:GLU:O	1:B:515:ARG:NH1	2.45	0.50
1:A:208:SER:OG	1:A:383:ALA:O	2.28	0.50
1:A:559:VAL:HG22	1:A:651:ALA:HB3	1.94	0.50
1:B:19:ASP:O	1:B:23:ALA:N	2.43	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:867:ASP:HA	1:A:870:GLU:HG2	1.94	0.50
1:B:519:PRO:HB3	1:B:553:ARG:HD2	1.94	0.50
1:A:51:PRO:HD2	1:A:136:ARG:CZ	2.42	0.50
1:A:340:HIS:CE1	1:A:373:LYS:HE2	2.47	0.50
3:D:152:ALA:O	3:D:203:LEU:N	2.45	0.50
1:A:265:GLY:HA2	1:A:344:THR:HG22	1.94	0.49
1:A:369:LEU:HB3	1:A:422:LEU:HD23	1.94	0.49
1:A:584:ALA:O	1:A:588:ARG:HG2	2.12	0.49
2:C:6:LEU:HD11	2:C:102:ARG:HB2	1.94	0.49
3:D:69:ILE:HD12	3:D:94:LEU:HD12	1.94	0.49
1:B:176:MET:HE1	1:A:445:ILE:HG21	1.93	0.49
1:A:577:LEU:HB3	1:A:583:PHE:HD2	1.77	0.49
3:F:69:ILE:HD12	3:F:94:LEU:HD12	1.94	0.49
1:B:171:VAL:HG21	1:B:910:LEU:HD21	1.94	0.49
1:B:777:PRO:HB3	1:B:798:LEU:H	1.77	0.49
1:A:108:PHE:HZ	1:A:131:TRP:CE2	2.31	0.49
1:A:1076:LEU:HD13	1:A:1281:GLY:HA3	1.93	0.49
1:A:1350:ASN:CG	1:A:1354:ARG:HH22	2.19	0.49
3:F:57:TYR:HB2	3:F:108:TYR:HB2	1.93	0.49
1:A:567:TRP:HB3	1:A:836:ILE:HG12	1.93	0.49
1:A:842:GLU:HG3	1:A:854:ALA:HB3	1.94	0.49
1:A:950:LEU:HB2	1:A:1003:VAL:HG22	1.94	0.49
1:A:660:ILE:HG22	1:A:674:ALA:HB1	1.94	0.49
3:D:57:TYR:HB2	3:D:108:TYR:HB2	1.93	0.49
3:D:188:GLN:HB3	3:D:195:TYR:CE1	2.48	0.49
2:E:6:LEU:HD11	2:E:102:ARG:HB2	1.94	0.49
1:A:1048:GLU:O	1:A:1060:ARG:NE	2.43	0.49
2:C:22:LEU:HB2	2:C:85:LEU:HB3	1.94	0.49
2:E:150:LYS:HE2	2:E:184:SER:HB3	1.94	0.49
3:F:152:ALA:O	3:F:203:LEU:N	2.45	0.49
1:B:731:GLU:HG2	1:B:734:ARG:HH11	1.78	0.49
2:C:150:LYS:HE2	2:C:184:SER:HB3	1.94	0.49
1:B:36:ALA:HB1	1:B:290:VAL:HG13	1.95	0.49
1:B:410:SER:HB2	1:B:413:ILE:HD12	1.94	0.49
1:B:712:VAL:HG22	1:B:727:GLY:HA3	1.94	0.49
1:B:1427:VAL:HG11	1:B:1455:LEU:HB2	1.93	0.49
3:D:52:ASN:HB3	3:D:92:PHE:HE2	1.78	0.49
1:B:209:LEU:HD22	1:B:450:ALA:HB1	1.94	0.49
1:B:603:ILE:HG23	1:B:607:ARG:HH21	1.76	0.49
1:A:802:TYR:HA	1:A:805:ARG:HE	1.78	0.49
1:A:152:PHE:O	1:A:229:ALA:HA	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1066:ALA:HB1	1:A:1290:PRO:HB3	1.94	0.49
1:A:1172:SER:O	1:A:1199:ALA:HA	2.13	0.49
1:B:553:ARG:HH11	1:B:1440:GLN:HE22	1.59	0.48
1:A:337:VAL:HG23	1:A:367:LEU:HD11	1.94	0.48
1:A:1068:TRP:O	1:A:1072:ARG:HG3	2.14	0.48
1:B:115:GLU:O	1:B:119:VAL:HG23	2.13	0.48
1:A:489:ALA:HB3	1:A:539:LEU:HD22	1.93	0.48
1:A:720:PRO:HB2	1:A:721:ARG:HD2	1.95	0.48
2:C:155:GLU:HG2	2:C:183:TYR:CE2	2.48	0.48
1:A:54:PHE:HE1	1:A:58:LEU:HD11	1.78	0.48
3:D:29:VAL:O	3:D:129:LYS:N	2.47	0.48
2:E:54:ARG:HH22	2:E:63:GLU:HB2	1.78	0.48
1:A:1222:SER:HB3	1:A:1268:THR:H	1.78	0.48
3:F:188:GLN:HB3	3:F:195:TYR:CE1	2.48	0.48
1:B:15:ARG:HB3	3:F:70:TYR:HE2	1.77	0.48
1:A:103:ALA:HB1	1:A:905:ARG:HB3	1.95	0.48
1:B:778:LEU:HD12	1:B:779:PRO:HD2	1.94	0.48
4:B:1601:PN7:H15	4:B:1601:PN7:H19	1.56	0.48
1:A:156:ILE:HG13	1:A:381:ALA:HB2	1.96	0.48
2:C:49:TRP:HB3	3:D:117:ARG:O	2.14	0.48
1:B:114:ARG:HB2	1:B:1429:GLY:HA3	1.95	0.48
3:D:37:ILE:HD11	3:D:56:TRP:CZ3	2.49	0.48
2:E:155:GLU:HG2	2:E:183:TYR:CE2	2.48	0.48
1:A:1159:TRP:CG	1:A:1345:CYS:HB2	2.49	0.48
2:E:37:SER:HB2	2:E:101:THR:OG1	2.13	0.48
1:B:33:GLU:HB3	1:B:220:ARG:HG2	1.95	0.48
1:A:466:VAL:HG23	1:A:505:ARG:HH21	1.78	0.48
1:B:127:LEU:HD23	1:B:188:ILE:HG13	1.96	0.48
1:A:1:MET:HE1	2:C:52:PHE:HZ	1.77	0.48
1:A:40:MET:HG2	1:A:275:LEU:HD22	1.96	0.48
1:A:688:MET:N	1:A:689:PRO:HD2	2.29	0.47
1:A:761:GLU:HG3	1:A:808:ARG:HD2	1.96	0.47
1:A:1009:VAL:HA	1:A:1048:GLU:HB2	1.96	0.47
1:B:430:PRO:O	1:B:436:ARG:NE	2.46	0.47
1:B:764:ARG:HE	1:B:768:HIS:CD2	2.32	0.47
1:A:331:PRO:O	1:A:363:ARG:NH1	2.47	0.47
1:A:319:VAL:HG11	1:A:449:ASN:HB3	1.96	0.47
1:A:625:ASP:HB2	1:A:683:ARG:HA	1.96	0.47
2:E:149:VAL:HG21	2:E:185:LEU:HD22	1.96	0.47
3:F:29:VAL:O	3:F:129:LYS:N	2.47	0.47
1:B:337:VAL:HG11	1:B:355:LEU:HD22	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:573:ALA:HB1	1:B:583:PHE:CE2	2.49	0.47
1:A:778:LEU:HG	1:A:779:PRO:HD2	1.96	0.47
1:A:1070:VAL:HB	1:A:1290:PRO:HG3	1.97	0.47
1:B:14:ARG:NH1	2:E:33:ASP:OD1	2.48	0.47
1:B:212:VAL:HG11	1:B:275:LEU:HD12	1.96	0.47
1:B:555:GLN:NE2	1:B:827:THR:HB	2.28	0.47
1:B:663:ALA:HB2	1:B:798:LEU:HD21	1.97	0.47
1:B:712:VAL:HG21	1:B:731:GLU:HB2	1.96	0.47
1:A:22:ALA:O	1:A:26:ARG:HG3	2.14	0.47
1:A:752:ASP:OD1	1:A:752:ASP:N	2.48	0.47
2:E:205:VAL:O	2:E:213:LYS:HA	2.15	0.47
1:B:385:VAL:O	1:B:389:ILE:HD12	2.13	0.47
1:A:124:ARG:HH12	1:A:908:VAL:HB	1.80	0.47
1:A:316:GLN:O	1:A:320:ILE:HG12	2.15	0.47
1:A:704:VAL:O	1:A:708:ILE:HG23	2.15	0.47
1:A:1125:TRP:CH2	1:A:1360:ILE:HG21	2.49	0.47
1:A:1272:LEU:HD21	1:A:1298:LEU:HD23	1.97	0.47
2:C:149:VAL:HG21	2:C:185:LEU:HD22	1.96	0.47
3:F:203:LEU:HD21	3:F:208:TYR:HA	1.96	0.47
1:B:857:SER:O	1:B:871:ALA:HB2	2.14	0.47
1:A:123:GLN:NE2	1:A:127:LEU:HG	2.29	0.47
1:A:1209:ARG:HA	1:A:1261:LEU:HD21	1.96	0.47
1:B:624:VAL:HG11	1:B:685:ILE:HG22	1.97	0.47
1:A:797:GLU:HB2	1:A:805:ARG:HH22	1.80	0.47
1:B:39:ALA:HB2	1:B:137:ALA:HB2	1.97	0.47
1:B:114:ARG:HD3	1:B:170:GLY:O	2.14	0.47
2:C:107:PHE:HB2	2:C:110:TRP:CZ2	2.50	0.47
2:C:205:VAL:O	2:C:213:LYS:HA	2.15	0.47
1:B:103:ALA:HB1	1:B:905:ARG:HB3	1.96	0.46
1:A:162:PRO:HD2	1:A:910:LEU:HG	1.97	0.46
1:A:471:VAL:HG11	1:A:865:LEU:HD23	1.98	0.46
1:B:15:ARG:CB	3:F:70:TYR:HE2	2.28	0.46
1:A:26:ARG:O	1:A:30:LEU:HB2	2.14	0.46
1:B:208:SER:HB3	1:B:381:ALA:O	2.14	0.46
1:A:57:LEU:HD11	1:A:64:ALA:HB2	1.96	0.46
1:A:359:TYR:OH	1:A:451:HIS:NE2	2.49	0.46
1:A:1034:MET:HE1	1:A:1043:LEU:HB2	1.97	0.46
1:B:195:GLU:HG2	1:A:300:ASN:OD1	2.15	0.46
1:A:286:ASN:HB2	1:A:288:HIS:CE1	2.50	0.46
1:A:693:GLY:HA3	1:A:752:ASP:HB2	1.96	0.46
1:A:1171:VAL:HG22	1:A:1198:ALA:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:102:ARG:NH2	2:E:108:ASP:OD2	2.49	0.46
2:E:107:PHE:HB2	2:E:110:TRP:CZ2	2.50	0.46
1:B:210:VAL:O	1:B:213:HIS:HB3	2.16	0.46
1:A:969:LEU:HB2	1:A:976:VAL:HG22	1.98	0.46
1:A:72:ARG:HE	1:A:102:THR:HG23	1.81	0.46
3:F:30:THR:HB	3:F:129:LYS:HE3	1.98	0.46
1:A:1:MET:HB2	1:A:7:GLU:HB2	1.98	0.46
1:A:50:THR:HG23	1:A:52:GLU:HG3	1.97	0.46
1:A:597:HIS:CE1	1:A:676:ARG:HB2	2.51	0.46
1:A:816:ALA:O	1:A:820:LEU:HG	2.16	0.46
1:A:1178:ALA:HB3	1:A:1181:ALA:HB2	1.97	0.46
1:A:1271:VAL:HG11	1:A:1348:LEU:HD11	1.97	0.46
1:B:15:ARG:NH1	1:B:18:LEU:HD22	2.31	0.46
1:B:363:ARG:HD2	1:B:363:ARG:HA	1.83	0.46
1:A:987:ARG:NH1	1:A:1239:LEU:O	2.48	0.46
2:C:102:ARG:NH2	2:C:108:ASP:OD2	2.49	0.46
1:B:40:MET:SD	1:B:389:ILE:HG13	2.56	0.45
2:C:53:ILE:HA	2:C:62:THR:HA	1.98	0.45
2:C:55:SER:HA	2:C:76:ARG:HH12	1.80	0.45
2:E:130:PRO:HG3	2:E:216:LYS:HE2	1.99	0.45
1:B:8:LYS:HG2	2:E:105:THR:HG21	1.97	0.45
1:B:519:PRO:O	1:B:520:HIS:ND1	2.49	0.45
1:B:660:ILE:HG21	1:B:678:VAL:HB	1.97	0.45
1:A:123:GLN:HE21	1:A:127:LEU:HG	1.81	0.45
1:A:126:MET:CE	1:A:153:VAL:HG21	2.46	0.45
1:A:235:MET:HE2	1:A:235:MET:HB2	1.53	0.45
1:A:627:VAL:HG22	1:A:631:MET:HE3	1.97	0.45
1:B:58:LEU:HG	1:B:401:PRO:HB3	1.97	0.45
1:A:1143:VAL:HG11	1:A:1160:LEU:HD13	1.98	0.45
1:A:1285:LEU:HG	1:A:1327:VAL:HG21	1.98	0.45
2:C:38:TRP:CE2	2:C:85:LEU:HB2	2.51	0.45
2:C:130:PRO:HG3	2:C:216:LYS:HE2	1.99	0.45
1:B:391:MET:HE1	1:B:439:GLY:C	2.41	0.45
1:A:1048:GLU:HB3	1:A:1060:ARG:HE	1.80	0.45
1:A:1339:MET:HE1	1:A:1361:VAL:HG12	1.98	0.45
2:E:50:VAL:HG13	2:E:68:VAL:HG21	1.98	0.45
1:B:194:LEU:HD11	1:B:228:MET:HE2	1.97	0.45
1:A:339:ALA:HB2	1:A:369:LEU:HD11	1.99	0.45
1:A:505:ARG:NH1	1:A:893:ALA:HA	2.28	0.45
3:D:203:LEU:HD21	3:D:208:TYR:HA	1.97	0.45
1:B:601:GLU:HG2	1:B:603:ILE:H	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:642:TRP:NE1	1:B:857:SER:OG	2.46	0.45
1:B:731:GLU:HA	1:B:734:ARG:HG2	1.98	0.45
1:B:1421:ARG:NH2	1:B:1436:VAL:H	2.14	0.45
1:A:395:MET:O	1:A:436:ARG:NH2	2.49	0.45
1:A:610:ALA:HA	1:A:613:ARG:HE	1.81	0.45
2:C:55:SER:HA	2:C:76:ARG:NH1	2.32	0.45
3:D:102:GLU:H	3:D:102:GLU:HG3	1.59	0.45
3:D:106:VAL:HG22	3:D:125:LYS:HD2	1.99	0.45
3:F:25:LEU:O	3:F:125:LYS:N	2.44	0.45
1:B:108:PHE:HZ	1:B:131:TRP:CE2	2.34	0.45
1:B:5:ASP:O	1:A:6:SER:OG	2.35	0.45
1:B:208:SER:O	1:B:211:ALA:HB3	2.16	0.45
1:B:402:ARG:NE	1:B:424:ASP:O	2.42	0.45
1:B:777:PRO:HB2	1:B:796:ASP:OD1	2.16	0.45
1:B:1466:ARG:HD2	1:B:1466:ARG:HA	1.65	0.45
1:A:1007:LEU:HD23	1:A:1007:LEU:H	1.82	0.45
2:E:170:VAL:HA	2:E:189:VAL:HG12	1.98	0.45
3:F:106:VAL:HG22	3:F:125:LYS:HD2	1.99	0.45
1:B:68:LEU:HB2	1:A:1056:PHE:CD2	2.51	0.45
1:B:1448:ASP:O	1:B:1449:SER:C	2.59	0.45
1:A:1068:TRP:CZ2	1:A:1087:ASP:HB2	2.51	0.45
2:C:31:PHE:HE2	2:C:77:ASP:H	1.65	0.45
1:B:378:HIS:CD2	1:B:380:GLN:H	2.35	0.45
1:A:123:GLN:HE22	1:A:188:ILE:HG13	1.81	0.45
2:E:154:PRO:HD2	2:E:209:PRO:HG2	1.99	0.45
1:B:300:ASN:O	1:B:449:ASN:N	2.49	0.44
1:A:1204:ASP:HB3	1:A:1207:SER:HB3	1.98	0.44
1:B:12:TYR:HE1	3:F:70:TYR:HB2	1.82	0.44
1:B:787:THR:HB	1:B:811:VAL:HG13	2.00	0.44
1:A:1222:SER:HA	1:A:1267:LEU:HA	1.99	0.44
1:B:338:GLU:HB2	1:B:391:MET:CE	2.47	0.44
1:A:103:ALA:HA	1:A:906:GLU:O	2.17	0.44
1:A:837:LEU:O	1:A:841:ILE:HG12	2.18	0.44
1:B:680:LEU:HD23	1:B:771:LEU:HA	1.99	0.44
1:A:1099:HIS:HB2	1:A:1115:LEU:HD13	1.99	0.44
3:D:30:THR:HB	3:D:129:LYS:HE3	1.98	0.44
1:A:312:ASN:HB3	1:A:315:ALA:HB3	2.00	0.44
1:A:1096:LEU:HA	1:A:1100:LEU:HG	2.00	0.44
2:C:75:SER:OG	2:C:84:TYR:HB2	2.18	0.44
2:C:154:PRO:HD2	2:C:209:PRO:HG2	1.99	0.44
2:C:188:VAL:HG21	3:D:157:LEU:HD21	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:717:VAL:O	1:B:814:ALA:N	2.39	0.44
4:B:1601:PN7:H22	1:A:205:CYS:HG	1.64	0.44
1:B:251:ALA:H	1:B:267:GLY:H	1.65	0.44
1:A:210:VAL:O	1:A:214:LEU:HG	2.18	0.44
1:A:502:GLN:HB3	1:A:507:ILE:HD11	1.99	0.44
1:A:627:VAL:O	1:A:631:MET:HG2	2.17	0.44
2:E:75:SER:OG	2:E:84:TYR:HB2	2.18	0.44
1:B:41:ALA:HB1	1:B:129:LEU:HD22	1.99	0.44
1:B:209:LEU:HD22	1:B:450:ALA:CB	2.48	0.44
1:A:732:LEU:HD12	1:A:735:LEU:HD12	1.99	0.44
1:B:111:MET:HE2	1:B:187:ARG:HH11	1.82	0.44
1:B:683:ARG:HE	1:B:683:ARG:HB3	1.62	0.44
1:B:788:VAL:HG21	1:B:805:ARG:HB3	2.00	0.44
1:A:232:VAL:HA	1:A:272:ALA:HA	1.99	0.44
1:A:282:ASP:O	1:A:283:ALA:C	2.61	0.44
3:D:158:LEU:HD22	3:D:218:VAL:HG11	2.00	0.44
1:B:479:ALA:HA	1:B:901:TYR:HE2	1.83	0.43
1:A:72:ARG:HH21	1:A:101:ALA:H	1.65	0.43
1:A:951:VAL:HG13	1:A:951:VAL:O	2.18	0.43
1:A:1167:HIS:CE1	1:A:1219:VAL:HG11	2.52	0.43
3:F:86:SER:OG	3:F:87:GLY:N	2.49	0.43
1:A:154:GLY:HA3	1:A:207:SER:HB3	2.00	0.43
1:A:316:GLN:O	1:A:319:VAL:HG12	2.18	0.43
1:A:462:GLU:HB3	1:A:465:ARG:HE	1.83	0.43
1:A:1331:PHE:HB3	1:A:1336:VAL:HG12	2.00	0.43
2:C:101:THR:HB	2:C:107:PHE:HB3	2.00	0.43
1:B:149:THR:H	1:B:196:GLY:H	1.67	0.43
1:B:562:PHE:CE1	1:B:652:VAL:HG13	2.53	0.43
1:B:660:ILE:HD13	1:B:803:TRP:CZ3	2.53	0.43
1:A:121:PRO:HD3	1:A:160:TYR:CD1	2.53	0.43
1:A:146:ALA:HA	1:A:195:GLU:HG3	2.00	0.43
2:C:173:PHE:CE2	3:D:198:SER:HB2	2.54	0.43
1:B:612:ARG:HE	1:B:612:ARG:HB3	1.60	0.43
1:A:149:THR:HG23	1:A:226:LEU:C	2.44	0.43
1:A:312:ASN:HD22	1:A:315:ALA:HB2	1.84	0.43
1:A:698:ALA:HA	1:A:722:SER:HA	2.00	0.43
1:A:1159:TRP:HZ2	1:A:1349:GLN:HB2	1.84	0.43
3:D:141:PRO:HB3	3:D:231:PHE:HB2	2.00	0.43
1:B:34:PRO:HB2	1:B:293:VAL:HG11	2.01	0.43
1:B:291:LEU:O	1:B:395:MET:HE1	2.19	0.43
1:A:476:VAL:O	1:A:511:LEU:HG	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:594:LEU:HG	1:A:675:ALA:HB1	2.01	0.43
1:A:1125:TRP:HD1	1:A:1381:LEU:HG	1.83	0.43
1:A:1302:ARG:HE	1:A:1307:LEU:HB3	1.83	0.43
3:F:215:ALA:HA	3:F:230:SER:HA	2.00	0.43
1:B:313:GLY:O	1:B:317:VAL:HG23	2.18	0.43
1:B:607:ARG:H	1:B:607:ARG:HG2	1.54	0.43
1:A:279:ARG:O	1:A:280:LEU:C	2.61	0.43
1:A:935:ARG:NH1	1:A:1384:GLU:OE2	2.52	0.43
1:A:1314:TRP:CZ3	1:A:1344:ALA:HB1	2.52	0.43
1:B:121:PRO:HG2	1:B:236:PRO:HB3	2.00	0.43
1:B:235:MET:HG2	1:B:240:MET:HE2	2.00	0.43
1:B:663:ALA:HB1	1:B:669:LEU:HB2	1.99	0.43
1:B:692:LYS:HE2	1:B:728:ASP:HB3	2.00	0.43
1:A:33:GLU:HB3	1:A:220:ARG:HG2	1.99	0.43
1:A:54:PHE:CE1	1:A:58:LEU:HD11	2.54	0.43
1:A:69:PRO:HG3	1:A:97:PHE:HB3	2.01	0.43
1:A:720:PRO:HB2	1:A:721:ARG:HH11	1.84	0.43
3:D:34:PRO:HA	3:D:96:ILE:O	2.19	0.43
1:B:34:PRO:O	1:B:280:LEU:HB2	2.19	0.43
1:B:188:ILE:HD12	1:B:188:ILE:H	1.84	0.43
1:B:302:ASP:N	1:B:447:GLY:O	2.50	0.43
3:D:215:ALA:HA	3:D:230:SER:HA	2.00	0.43
2:E:47:LEU:HB2	3:F:120:PHE:CG	2.53	0.43
1:B:43:ARG:HH12	1:B:132:GLU:CD	2.27	0.43
1:B:300:ASN:OD1	1:B:322:GLN:NE2	2.51	0.43
1:A:235:MET:HA	1:A:236:PRO:HD3	1.84	0.43
1:A:685:ILE:HD12	1:A:688:MET:HE3	2.00	0.43
2:C:21:ARG:HD3	2:C:84:TYR:CE1	2.54	0.43
3:D:86:SER:OG	3:D:87:GLY:N	2.49	0.43
1:B:54:PHE:HD2	1:B:393:LEU:HD12	1.83	0.42
1:B:368:HIS:CE1	1:B:427:GLU:HG3	2.54	0.42
1:A:569:TRP:CZ3	1:A:861:GLY:HA2	2.53	0.42
1:A:906:GLU:HG2	1:A:907:ARG:H	1.82	0.42
1:A:934:TRP:HB3	1:A:1121:TYR:HB3	2.01	0.42
4:B:1601:PN7:O3	4:B:1601:PN7:O3P	2.37	0.42
2:C:91:LYS:HB3	2:C:91:LYS:HE3	1.38	0.42
3:F:34:PRO:HA	3:F:96:ILE:O	2.19	0.42
3:F:141:PRO:HB3	3:F:231:PHE:HB2	2.00	0.42
1:B:577:LEU:HD13	1:B:606:LEU:HD22	2.01	0.42
1:A:637:SER:O	1:A:640:SER:OG	2.27	0.42
2:C:185:LEU:HD23	2:C:186:SER:C	2.44	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:159:VAL:HG22	2:E:205:VAL:HG13	2.02	0.42
3:F:19:VAL:HG22	3:F:43:GLN:HB2	2.01	0.42
1:B:164:LEU:HG	1:B:175:LEU:HD22	2.01	0.42
1:B:1437:PRO:HB2	1:B:1440:GLN:HB2	2.02	0.42
1:A:125:LEU:O	1:A:129:LEU:HD13	2.19	0.42
1:B:18:LEU:HD11	2:E:29:PHE:CZ	2.55	0.42
1:B:41:ALA:HB3	1:B:274:MET:HB2	2.02	0.42
1:A:384:GLY:O	1:A:388:VAL:HG23	2.18	0.42
1:A:632:PHE:CG	1:A:678:VAL:HG11	2.55	0.42
1:B:156:ILE:HD11	1:B:204:ALA:HA	2.01	0.42
1:B:257:LYS:HG3	1:B:405:HIS:HB3	2.01	0.42
1:B:712:VAL:HA	1:B:727:GLY:HA3	2.02	0.42
1:A:855:ILE:HD13	1:A:855:ILE:HA	1.67	0.42
2:C:159:VAL:HG22	2:C:205:VAL:HG13	2.02	0.42
3:D:205:LYS:HE2	3:D:205:LYS:HB2	1.83	0.42
2:E:149:VAL:HB	2:E:185:LEU:HB3	2.02	0.42
1:B:65:VAL:HG13	1:B:95:GLY:H	1.84	0.42
1:B:455:GLU:HG2	1:B:456:GLU:N	2.34	0.42
1:B:583:PHE:HZ	1:B:634:VAL:HG13	1.85	0.42
1:B:763:ILE:O	1:B:763:ILE:HG22	2.20	0.42
1:B:1409:ALA:C	1:B:1411:GLU:N	2.78	0.42
1:A:470:ASP:CG	1:A:471:VAL:H	2.28	0.42
1:A:1005:SER:OG	1:A:1043:LEU:HD11	2.20	0.42
1:A:1289:ALA:HB3	1:A:1290:PRO:HD3	2.02	0.42
2:E:21:ARG:HD3	2:E:84:TYR:CE1	2.54	0.42
1:B:202:ASP:O	1:A:200:SER:HB3	2.19	0.42
1:B:485:LEU:HD13	1:B:520:HIS:O	2.20	0.42
1:B:697:ILE:HD12	1:B:746:ALA:HB2	2.01	0.42
1:A:102:THR:HG22	1:A:909:TRP:CD2	2.55	0.42
1:A:560:PHE:CE1	1:A:829:LEU:HD23	2.55	0.42
1:A:841:ILE:HA	1:A:844:ILE:HG12	2.01	0.42
1:A:1331:PHE:HA	1:A:1334:HIS:CE1	2.55	0.42
1:B:331:PRO:HB3	1:B:359:TYR:HA	2.02	0.42
1:B:429:TRP:CD1	1:B:436:ARG:HH11	2.38	0.42
1:B:437:ARG:HG3	1:B:455:GLU:HG3	2.01	0.42
1:A:97:PHE:HA	1:A:270:GLU:HG2	2.02	0.42
1:A:925:VAL:HA	1:A:928:LEU:HD12	2.01	0.42
1:A:1332:ARG:HA	1:A:1336:VAL:O	2.20	0.42
1:B:573:ALA:HB1	1:B:583:PHE:HE2	1.85	0.42
1:B:1471:THR:HG23	1:B:1475:HIS:HD2	1.84	0.42
1:A:692:LYS:NZ	1:A:726:ALA:HB1	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:834:HIS:HA	1:A:858:LEU:O	2.20	0.42
1:A:830:GLU:HB3	1:A:857:SER:HB2	2.02	0.41
1:A:1157:ALA:HB1	1:A:1168:LEU:HD21	2.01	0.41
2:C:8:GLN:HE21	2:C:100:CYS:HB3	1.85	0.41
2:C:87:MET:HE2	2:C:87:MET:HB3	1.89	0.41
2:C:149:VAL:HB	2:C:185:LEU:HB3	2.02	0.41
1:B:316:GLN:O	1:B:320:ILE:HG13	2.19	0.41
1:B:595:GLU:HA	1:B:598:LEU:O	2.20	0.41
1:B:712:VAL:HG12	1:B:725:VAL:HG22	2.01	0.41
1:B:1465:VAL:HG21	1:B:1485:LEU:HD11	2.02	0.41
1:A:197:PRO:HD2	1:A:224:SER:HB3	2.03	0.41
1:A:302:ASP:O	1:A:312:ASN:ND2	2.54	0.41
3:D:30:THR:HB	3:D:129:LYS:HB3	2.03	0.41
3:D:82:ARG:O	3:D:96:ILE:HA	2.20	0.41
2:E:31:PHE:HE2	2:E:77:ASP:H	1.68	0.41
1:B:80:PHE:HD2	1:A:1055:PRO:HB2	1.84	0.41
1:A:555:GLN:O	1:A:557:ARG:HG2	2.21	0.41
1:A:577:LEU:HD23	1:A:577:LEU:H	1.86	0.41
1:A:1028:LEU:HD21	1:A:1070:VAL:HG21	2.02	0.41
1:A:1267:LEU:O	1:A:1308:PRO:HG2	2.19	0.41
2:C:53:ILE:HD13	2:C:76:ARG:HG2	2.02	0.41
1:B:76:LEU:H	1:B:79:LEU:HD13	1.85	0.41
1:B:296:GLY:HA3	1:B:327:SER:HB3	2.01	0.41
1:A:68:LEU:HD22	1:A:74:TRP:CZ3	2.54	0.41
1:A:149:THR:HA	1:A:224:SER:HB2	2.01	0.41
1:A:1377:ARG:HH11	1:A:1379:THR:HG21	1.85	0.41
3:D:75:ARG:CZ	3:D:79:VAL:O	2.68	0.41
1:A:38:VAL:N	1:A:276:LEU:O	2.53	0.41
1:A:625:ASP:HA	1:A:682:SER:HB2	2.01	0.41
1:A:737:ALA:O	1:A:740:THR:HG22	2.20	0.41
1:A:1137:TRP:HE1	1:A:1268:THR:HG23	1.84	0.41
1:A:1307:LEU:HA	1:A:1308:PRO:HD3	1.87	0.41
3:D:19:VAL:HG22	3:D:43:GLN:HB2	2.02	0.41
3:F:75:ARG:CZ	3:F:79:VAL:O	2.68	0.41
1:B:15:ARG:HH11	2:E:102:ARG:HH22	1.67	0.41
1:B:562:PHE:N	1:B:653:ILE:O	2.50	0.41
1:B:652:VAL:HB	1:B:784:PHE:HD1	1.84	0.41
1:A:731:GLU:O	1:A:735:LEU:N	2.53	0.41
1:A:1366:TRP:HA	1:A:1369:PHE:HB3	2.02	0.41
3:D:168:VAL:HG22	3:D:218:VAL:HG12	2.03	0.41
1:B:14:ARG:NH1	2:E:34:TYR:CE1	2.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:378:HIS:HD2	1:B:380:GLN:H	1.67	0.41
1:B:659:GLU:OE2	1:B:787:THR:OG1	2.33	0.41
1:B:700:PRO:HG2	1:B:703:GLU:HB2	2.02	0.41
1:A:98:LEU:HD23	1:A:98:LEU:HA	1.86	0.41
1:A:1235:THR:O	1:A:1239:LEU:HG	2.20	0.41
3:D:166:ALA:HB2	3:D:220:HIS:HB2	2.03	0.41
3:F:82:ARG:O	3:F:96:ILE:HA	2.20	0.41
3:F:166:ALA:HB2	3:F:220:HIS:HB2	2.03	0.41
1:B:187:ARG:HA	1:A:308:LEU:HD11	2.02	0.41
1:A:285:ARG:HE	1:A:285:ARG:HB3	1.66	0.41
1:B:361:ARG:HD2	1:B:361:ARG:HA	1.80	0.41
1:B:701:ALA:HA	1:B:704:VAL:HG12	2.02	0.41
1:B:718:ASN:HA	1:B:813:PHE:HD2	1.85	0.41
1:B:1409:ALA:C	1:B:1411:GLU:H	2.28	0.41
1:A:602:VAL:HG12	1:A:606:LEU:HD23	2.03	0.41
1:A:1110:GLU:OE2	1:A:1380:ARG:HG3	2.21	0.41
1:A:1373:TYR:OH	1:A:1377:ARG:NH2	2.54	0.41
2:E:22:LEU:HD23	2:E:22:LEU:HA	1.93	0.41
1:B:201:VAL:HG21	1:B:211:ALA:HA	2.02	0.41
1:B:276:LEU:HD12	1:B:276:LEU:HA	1.82	0.41
1:B:521:ARG:HD3	1:B:876:PHE:O	2.20	0.41
1:B:786:SER:OG	1:B:793:THR:OG1	2.33	0.41
1:A:561:VAL:HG23	1:A:813:PHE:HZ	1.86	0.41
3:F:128:ILE:HG22	3:F:193:SER:HB2	2.03	0.41
3:F:142:PRO:HD3	3:F:154:VAL:HB	2.03	0.41
3:F:217:GLU:OE1	3:F:217:GLU:N	2.54	0.41
1:B:214:LEU:HD12	1:B:214:LEU:HA	1.92	0.40
1:B:349:PRO:HA	1:B:413:ILE:HG12	2.02	0.40
1:B:1409:ALA:HB3	1:B:1411:GLU:CG	2.51	0.40
1:B:1419:LEU:HG	1:B:1423:HIS:CE1	2.56	0.40
1:A:1030:LEU:CD2	1:A:1034:MET:HE3	2.50	0.40
1:A:1159:TRP:CD1	1:A:1345:CYS:HB2	2.57	0.40
1:B:587:LEU:HA	1:B:590:CYS:SG	2.61	0.40
1:B:705:ARG:O	1:B:708:ILE:HG12	2.20	0.40
1:A:375:ASN:OD1	1:A:404:LEU:HB2	2.20	0.40
1:A:834:HIS:ND1	1:A:860:ARG:HB2	2.35	0.40
1:A:1277:ALA:HB2	1:A:1285:LEU:HD13	2.02	0.40
2:C:38:TRP:HZ3	2:C:100:CYS:HB2	1.85	0.40
3:D:25:LEU:O	3:D:125:LYS:N	2.44	0.40
1:B:40:MET:HG2	1:B:291:LEU:HD12	2.03	0.40
1:B:152:PHE:HA	1:B:199:ILE:O	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:338:GLU:HB2	1:B:391:MET:HE2	2.04	0.40
1:B:567:TRP:CD2	1:B:833:ALA:HB3	2.56	0.40
1:B:694:MET:SD	1:B:749:LEU:HB3	2.61	0.40
1:A:241:LEU:H	1:A:241:LEU:HG	1.49	0.40
3:D:56:TRP:HB2	3:D:69:ILE:HB	2.03	0.40
3:F:30:THR:HB	3:F:129:LYS:HB3	2.03	0.40
3:F:82:ARG:HE	3:F:82:ARG:HB3	1.57	0.40
1:B:43:ARG:NH1	1:B:49:SER:O	2.54	0.40
1:B:149:THR:HG23	1:B:226:LEU:HG	2.03	0.40
1:A:154:GLY:CA	1:A:207:SER:HB3	2.50	0.40
1:B:429:TRP:CD1	1:B:436:ARG:NH1	2.90	0.40
1:B:628:GLN:N	1:B:629:PRO:HD2	2.36	0.40
1:A:29:GLU:HA	1:A:33:GLU:HB2	2.04	0.40
1:A:86:ARG:HB3	1:A:89:THR:OG1	2.22	0.40
1:A:1023:SER:HB2	1:A:1067:LEU:HD21	2.04	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	1388/1593 (87%)	1311 (94%)	76 (6%)	1 (0%)	48	78
1	B	995/1593 (62%)	921 (93%)	72 (7%)	2 (0%)	43	71
2	C	199/249 (80%)	182 (92%)	17 (8%)	0	100	100
2	E	199/249 (80%)	182 (92%)	17 (8%)	0	100	100
3	D	200/236 (85%)	191 (96%)	9 (4%)	0	100	100
3	F	200/236 (85%)	191 (96%)	9 (4%)	0	100	100
All	All	3181/4156 (76%)	2978 (94%)	200 (6%)	3 (0%)	49	78

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	689	PRO
1	B	1449	SER
1	A	236	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	1028/1175 (88%)	992 (96%)	36 (4%)	32	56
1	B	745/1175 (63%)	704 (94%)	41 (6%)	19	46
2	C	170/203 (84%)	146 (86%)	24 (14%)	3	13
2	E	170/203 (84%)	147 (86%)	23 (14%)	4	15
3	D	182/208 (88%)	156 (86%)	26 (14%)	3	13
3	F	182/208 (88%)	161 (88%)	21 (12%)	5	20
All	All	2477/3172 (78%)	2306 (93%)	171 (7%)	16	40

All (171) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	31	GLU
1	B	155	LEU
1	B	159	GLU
1	B	164	LEU
1	B	342	THR
1	B	344	THR
1	B	346	LEU
1	B	606	LEU
1	B	607	ARG
1	B	612	ARG
1	B	613	ARG
1	B	616	ASP
1	B	619	LEU
1	B	620	SER
1	B	622	GLU
1	B	623	ARG

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Mol	Chain	Res	Type
1	B	678	VAL
1	B	683	ARG
1	B	687	THR
1	B	688	MET
1	B	691	ASN
1	B	785	PHE
1	B	818	ARG
1	B	820	LEU
1	B	822	GLU
1	B	1407	LEU
1	B	1411	GLU
1	B	1412	ARG
1	B	1414	LYS
1	B	1435	ARG
1	B	1436	VAL
1	B	1447	VAL
1	B	1448	ASP
1	B	1453	LEU
1	B	1455	LEU
1	B	1456	ARG
1	B	1458	ARG
1	B	1459	LEU
1	B	1463	THR
1	B	1466	ARG
1	B	1485	LEU
1	A	11	GLU
1	A	43	ARG
1	A	44	LEU
1	A	232	VAL
1	A	234	VAL
1	A	235	MET
1	A	237	THR
1	A	240	MET
1	A	241	LEU
1	A	243	ASP
1	A	247	MET
1	A	279	ARG
1	A	286	ASN
1	A	773	GLU
1	A	775	PHE
1	A	776	HIS
1	A	778	LEU

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Mol	Chain	Res	Type
1	A	782	VAL
1	A	855	ILE
1	A	960	GLU
1	A	961	THR
1	A	966	ARG
1	A	967	GLU
1	A	1240	THR
1	A	1242	GLU
1	A	1243	ARG
1	A	1244	ILE
1	A	1245	GLU
1	A	1246	ARG
1	A	1249	ARG
1	A	1251	LYS
1	A	1253	LEU
1	A	1288	TYR
1	A	1302	ARG
1	A	1303	ARG
1	A	1383	ASP
2	C	13	LEU
2	C	36	MET
2	C	61	THR
2	C	62	THR
2	C	67	SER
2	C	69	LYS
2	C	71	ARG
2	C	73	THR
2	C	76	ARG
2	C	82	ILE
2	C	86	GLN
2	C	89	SER
2	C	91	LYS
2	C	92	THR
2	C	93	GLU
2	C	97	VAL
2	C	160	SER
2	C	172	THR
2	C	180	SER
2	C	186	SER
2	C	200	THR
2	C	210	SER
2	C	212	THR

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Mol	Chain	Res	Type
2	C	219	GLU
3	D	37	ILE
3	D	46	LEU
3	D	48	SER
3	D	54	LEU
3	D	70	TYR
3	D	75	ARG
3	D	82	ARG
3	D	86	SER
3	D	97	SER
3	D	98	ARG
3	D	99	VAL
3	D	100	GLU
3	D	102	GLU
3	D	110	MET
3	D	113	LEU
3	D	128	ILE
3	D	139	ILE
3	D	153	SER
3	D	154	VAL
3	D	157	LEU
3	D	159	ASN
3	D	160	ASN
3	D	181	SER
3	D	187	GLU
3	D	188	GLN
3	D	200	THR
2	E	13	LEU
2	E	36	MET
2	E	40	ARG
2	E	52	PHE
2	E	61	THR
2	E	71	ARG
2	E	73	THR
2	E	76	ARG
2	E	82	ILE
2	E	86	GLN
2	E	87	MET
2	E	91	LYS
2	E	92	THR
2	E	97	VAL
2	E	160	SER

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Mol	Chain	Res	Type
2	E	172	THR
2	E	178	GLN
2	E	180	SER
2	E	186	SER
2	E	200	THR
2	E	210	SER
2	E	212	THR
2	E	219	GLU
3	F	37	ILE
3	F	53	TYR
3	F	55	ASP
3	F	70	TYR
3	F	75	ARG
3	F	82	ARG
3	F	86	SER
3	F	99	VAL
3	F	100	GLU
3	F	104	VAL
3	F	110	MET
3	F	113	LEU
3	F	114	GLN
3	F	128	ILE
3	F	139	ILE
3	F	153	SER
3	F	157	LEU
3	F	181	SER
3	F	187	GLU
3	F	188	GLN
3	F	200	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	322	GLN
1	B	368	HIS
1	B	405	HIS
1	B	568	GLN
1	B	691	ASN
1	A	123	GLN
1	A	288	HIS
1	A	312	ASN
1	A	488	GLN

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Mol	Chain	Res	Type
1	A	502	GLN
1	A	597	HIS
1	A	657	GLN
1	A	776	HIS
1	A	806	ASN
1	A	1064	HIS
1	A	1112	GLN
1	A	1292	ASN
1	A	1349	GLN
2	C	88	ASN
2	C	178	GLN
3	D	43	GLN
3	D	114	GLN
3	D	182	GLN
2	E	41	GLN
3	F	43	GLN

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](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	PN7	B	1601	1	14,20,21	2.07	4 (28%)	18,26,29	2.42	9 (50%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	PN7	B	1601	1	-	8/24/26/27	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1601	PN7	C8-N9	4.46	1.44	1.33
4	B	1601	PN7	C4-N5	4.32	1.43	1.33
4	B	1601	PN7	O4-C4	-2.80	1.18	1.23
4	B	1601	PN7	O8-C8	-2.62	1.18	1.23

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1601	PN7	C6-N5-C4	-4.25	114.91	122.55
4	B	1601	PN7	CE2-C2-C1	-3.77	101.99	108.22
4	B	1601	PN7	O3-C3-C2	-3.39	102.33	110.18
4	B	1601	PN7	CE1-C2-C1	3.20	113.51	108.22
4	B	1601	PN7	O4-C4-N5	-3.00	116.63	122.98
4	B	1601	PN7	C6-C7-C8	2.90	117.22	112.39
4	B	1601	PN7	CE1-C2-C3	2.55	113.11	108.77
4	B	1601	PN7	C10-N9-C8	-2.53	118.11	122.82
4	B	1601	PN7	C3-C4-N5	2.35	120.94	116.48

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1601	PN7	N5-C6-C7-C8
4	B	1601	PN7	C7-C8-N9-C10
4	B	1601	PN7	O8-C8-N9-C10
4	B	1601	PN7	CE2-C2-C3-O3
4	B	1601	PN7	O3P-C1-C2-CE2
4	B	1601	PN7	N9-C10-C11-S12

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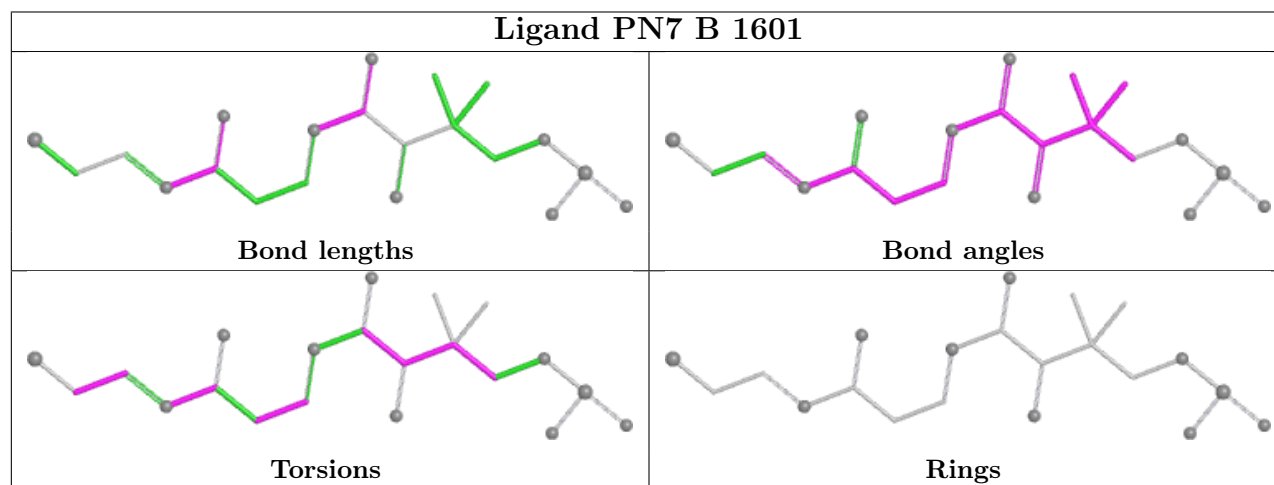
Mol	Chain	Res	Type	Atoms
4	B	1601	PN7	CE2-C2-C3-C4
4	B	1601	PN7	O3-C3-C4-N5

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1601	PN7	3	0

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

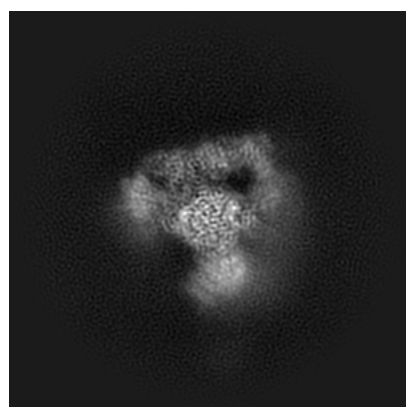
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23714. These allow visual inspection of the internal detail of the map and identification of artifacts.

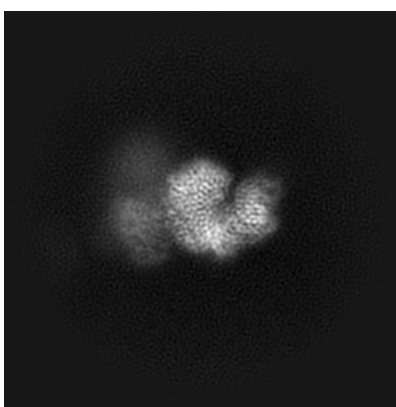
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

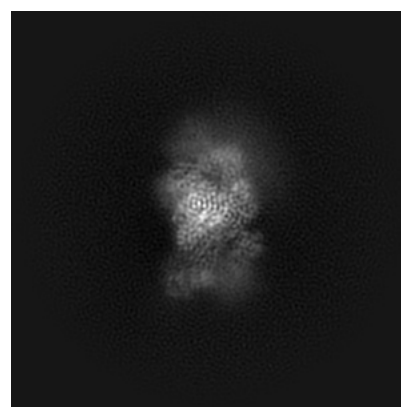
6.1.1 Primary map



X



Y

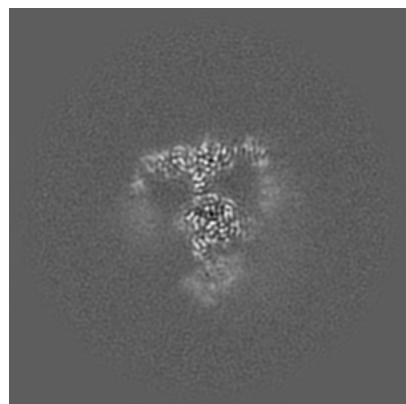


Z

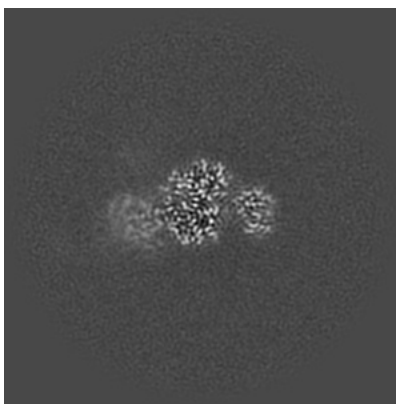
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

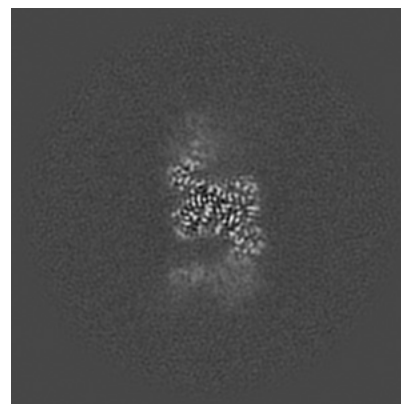
6.2.1 Primary map



X Index: 168



Y Index: 168

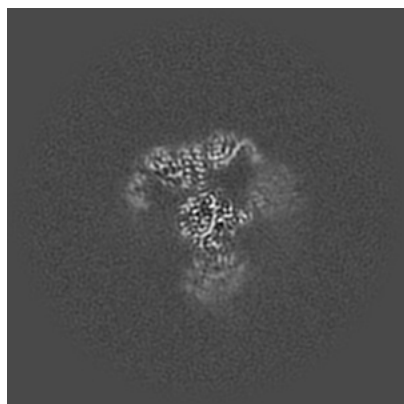


Z Index: 168

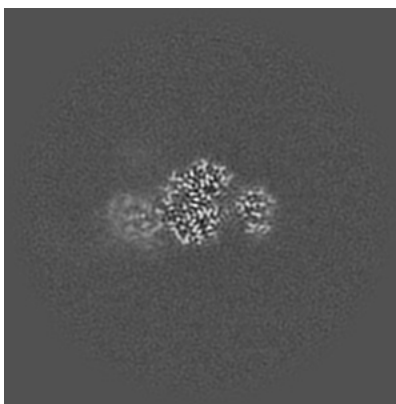
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

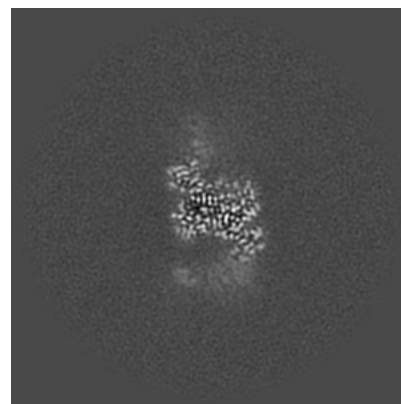
6.3.1 Primary map



X Index: 158



Y Index: 167

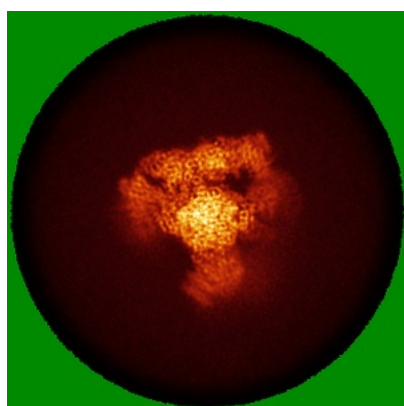


Z Index: 166

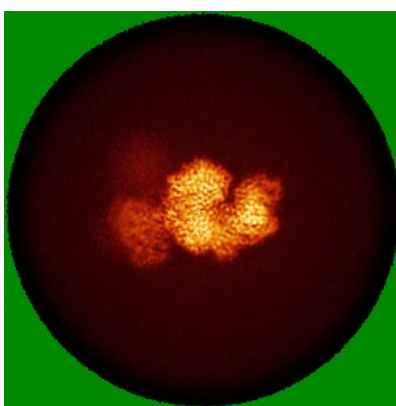
The images above show the largest variance slices of the map in three orthogonal directions.

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

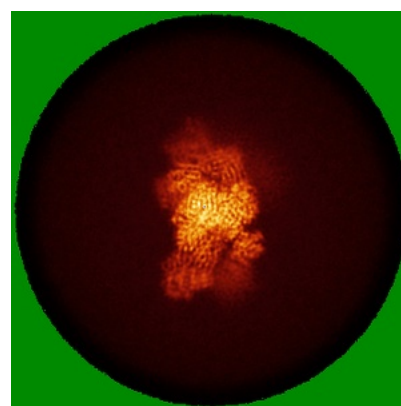
6.4.1 Primary map



X



Y

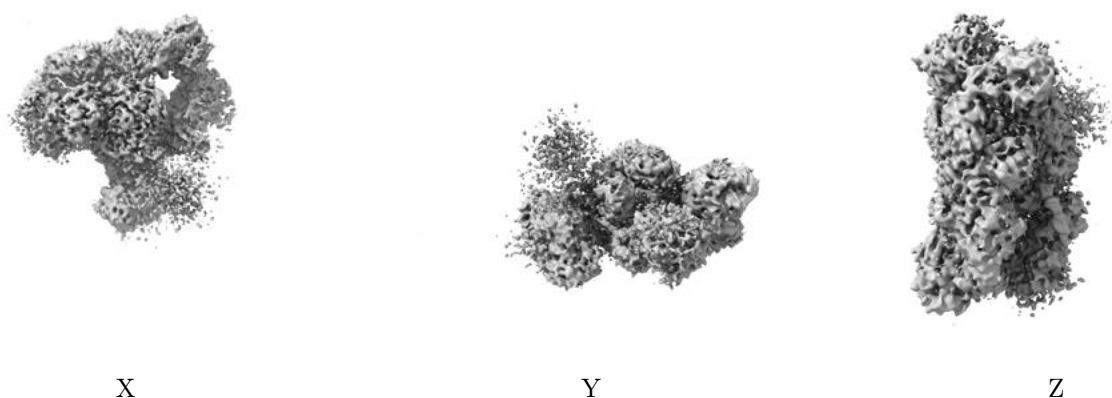


Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.5. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

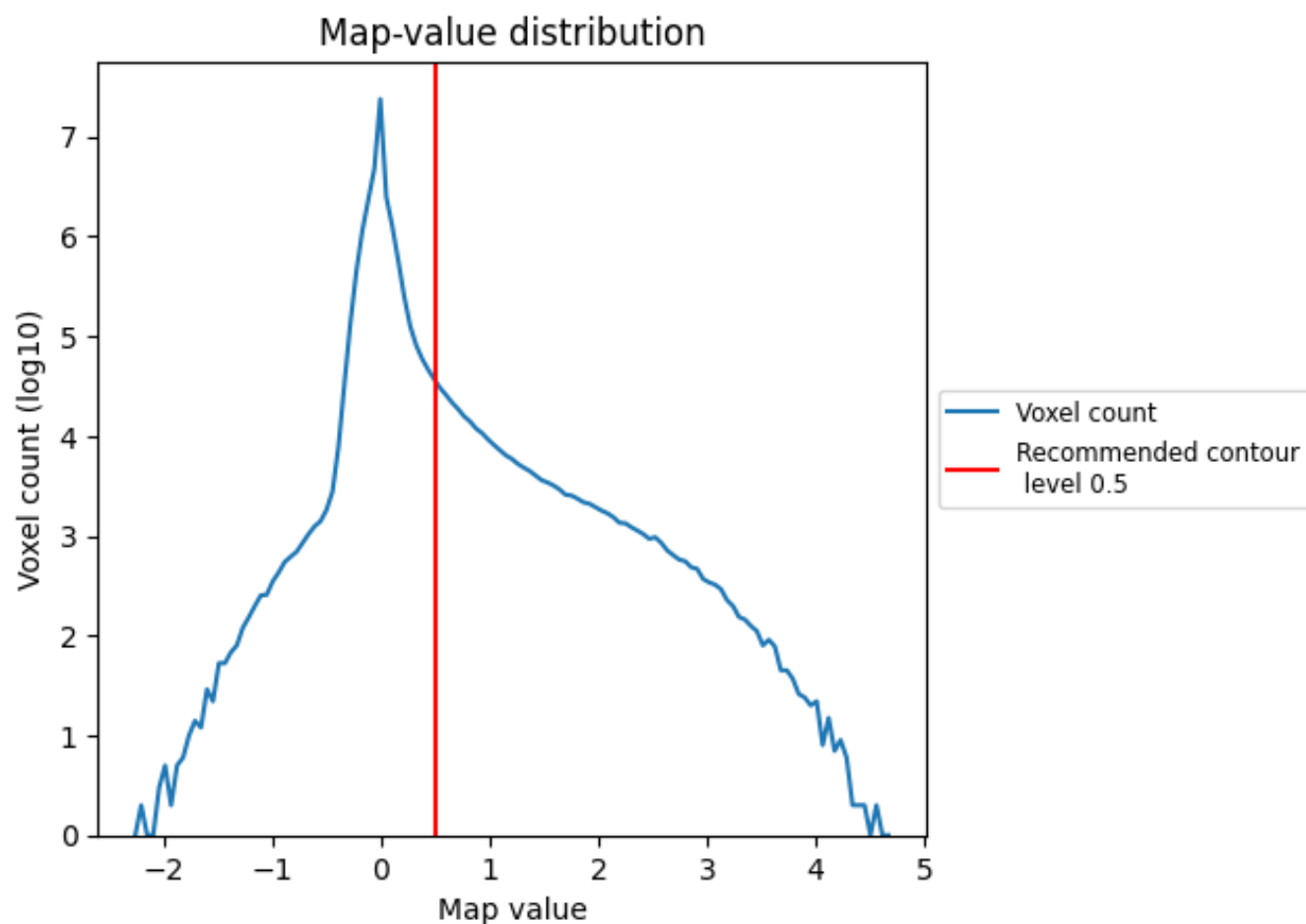
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

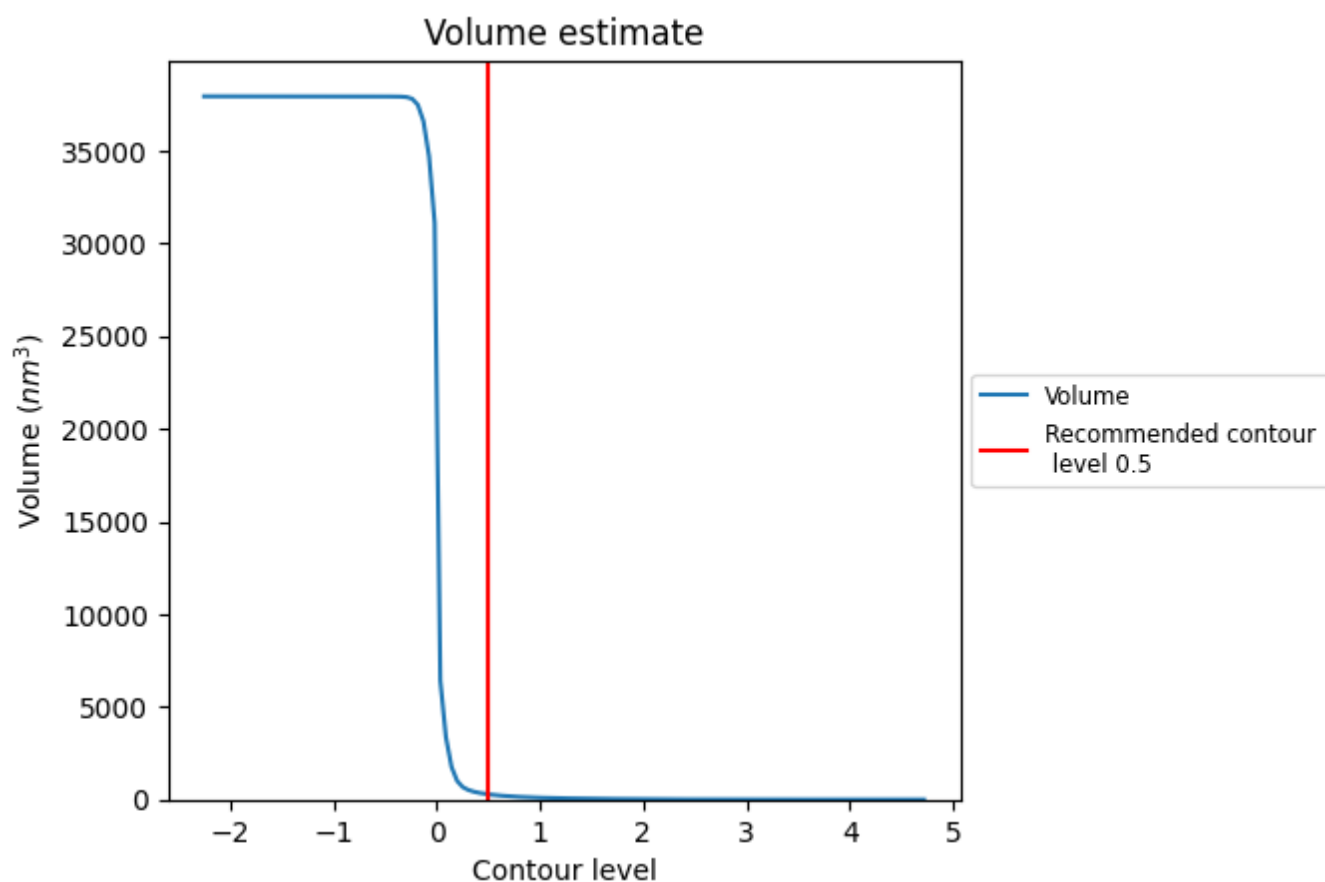
This section contains the results of statistical analysis of the map.

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.

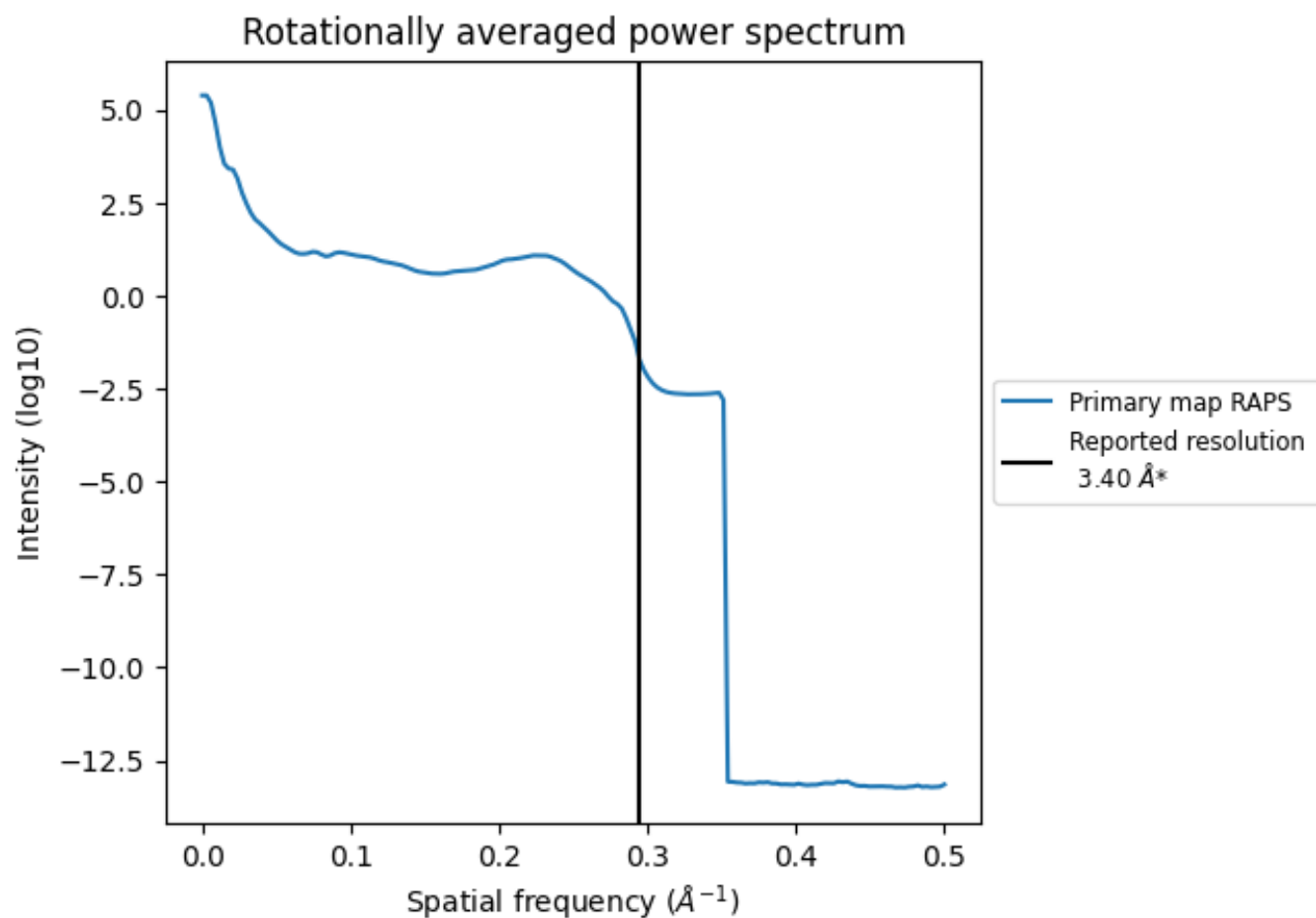
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 283 nm³; this corresponds to an approximate mass of 255 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 0.294 $\text{\AA}^{-1}$

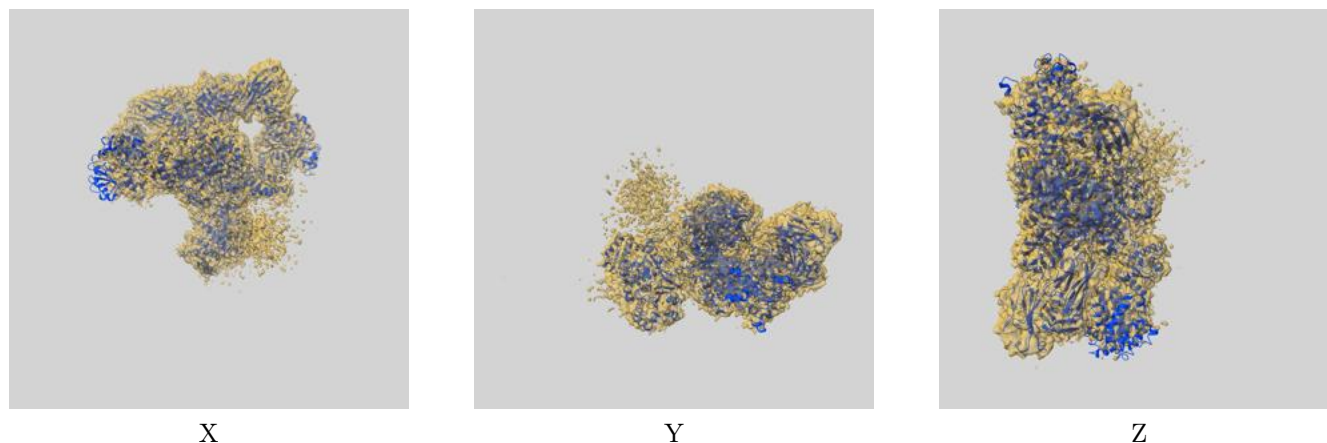
8 Fourier-Shell correlation

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

9 Map-model fit [i](#)

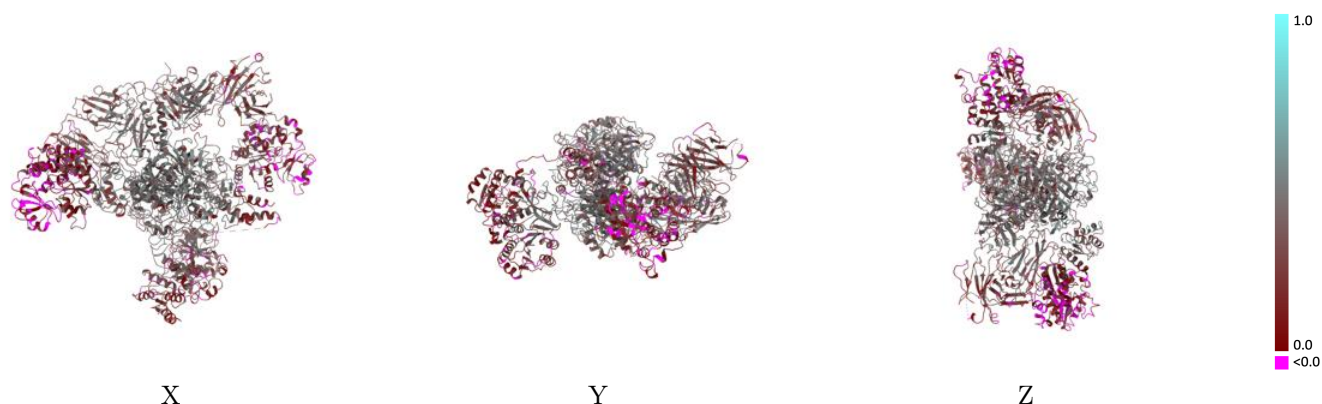
This section contains information regarding the fit between EMDB map EMD-23714 and PDB model 7M7I. Per-residue inclusion information can be found in [section 3](#) on [page 8](#).

9.1 Map-model overlay [i](#)



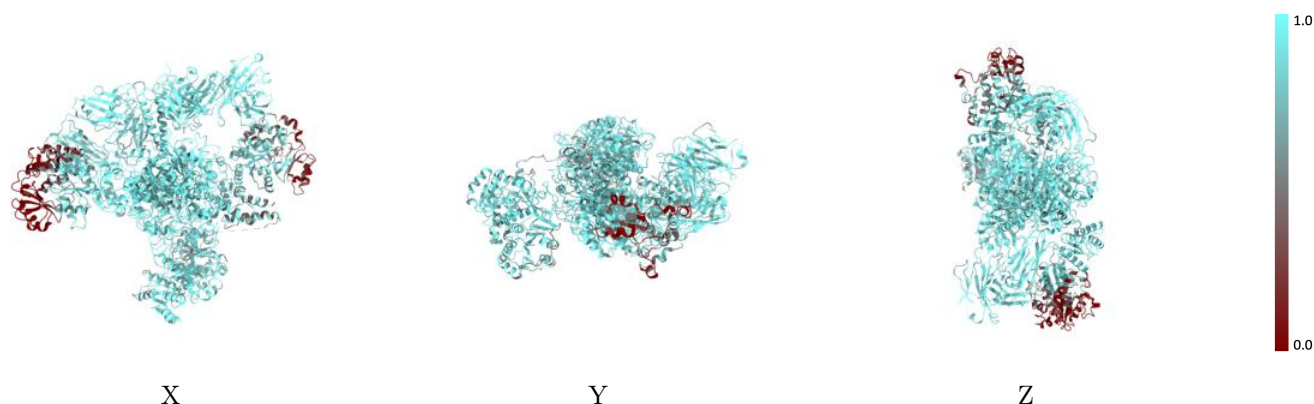
The images above show the 3D surface view of the map at the recommended contour level 0.5 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



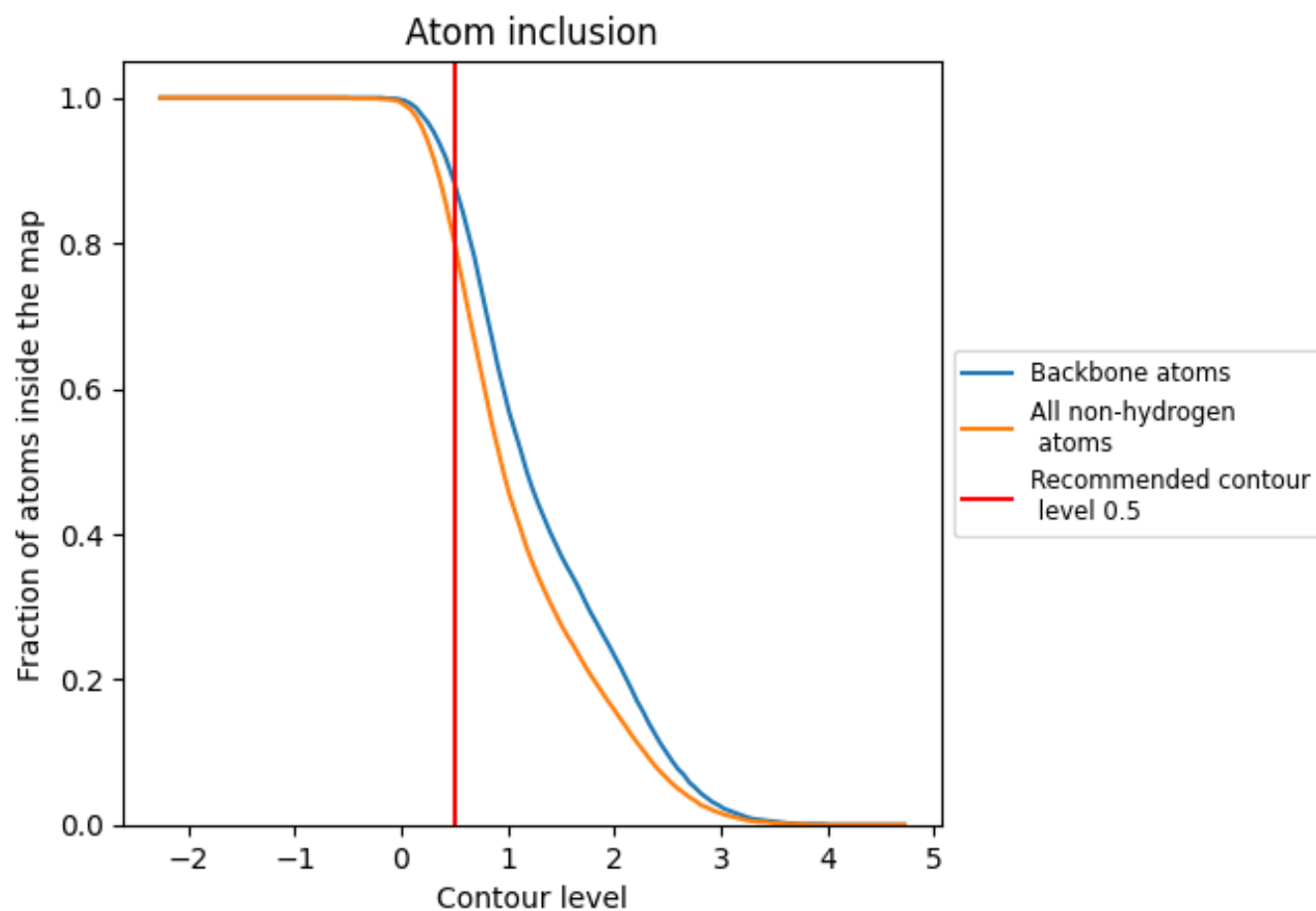
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.5).

9.4 Atom inclusion [i](#)



At the recommended contour level, 88% of all backbone atoms, 80% 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 (0.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8010	0.3040
A	0.7670	0.2770
B	0.7800	0.3180
C	0.8780	0.3240
D	0.8700	0.3190
E	0.8820	0.3340
F	0.8940	0.3480

1.0

0.0

<0.0