



Full wwPDB EM Validation Report ⓘ

Mar 23, 2026 – 09:36 PM UTC

PDB ID : 9C1L / pdb_00009c1l
EMDB ID : EMD-45123
Title : Rhesus rotavirus (VP1 structure at 2.65 Angstrom resolution)
Authors : Jenni, S.; Herrmann, T.; De Sautu, M.; Harrison, S.C.
Deposited on : 2024-05-29
Resolution : 2.65 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 : **NOT EXECUTED**
MapQ : **FAILED**
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

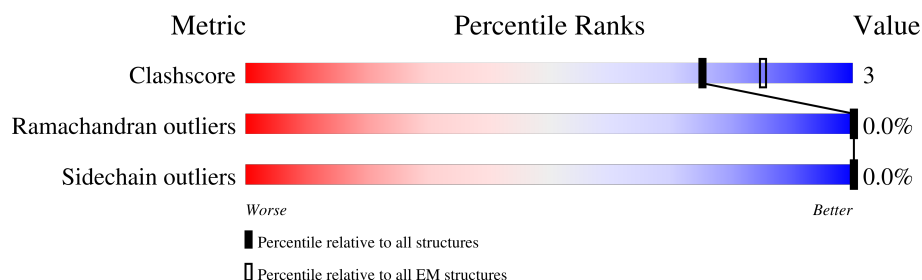
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.65 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)
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	887	80% 8% 12%
1	B	887	82% 6% 12%
1	C	887	82% 8% 9%
1	D	887	84% 10% 7%
1	E	887	82% 8% 10%
1	F	887	84% 7% 9%
1	G	887	84% 7% 9%
1	H	887	85% 6% 9%
1	I	887	83% 7% 10%

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Mol	Chain	Length	Quality of chain
1	J	887	 84% 8% 8%
2	P	1088	 90% 10%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 149101 atoms, of which 74750 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 Inner capsid protein VP2.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	782	Total	C	H	N	O	S	0	0
			12800	4057	6413	1102	1192	36		
1	B	781	Total	C	H	N	O	S	0	0
			12786	4052	6408	1101	1189	36		
1	C	803	Total	C	H	N	O	S	0	0
			13164	4174	6595	1131	1228	36		
1	D	827	Total	C	H	N	O	S	0	0
			13573	4297	6809	1163	1268	36		
1	E	795	Total	C	H	N	O	S	0	0
			13027	4132	6529	1118	1212	36		
1	F	805	Total	C	H	N	O	S	0	0
			13205	4186	6619	1134	1230	36		
1	G	804	Total	C	H	N	O	S	0	0
			13186	4180	6608	1133	1229	36		
1	H	805	Total	C	H	N	O	S	0	0
			13205	4186	6619	1134	1230	36		
1	I	799	Total	C	H	N	O	S	0	0
			13098	4154	6563	1126	1219	36		
1	J	816	Total	C	H	N	O	S	0	0
			13390	4242	6713	1149	1250	36		

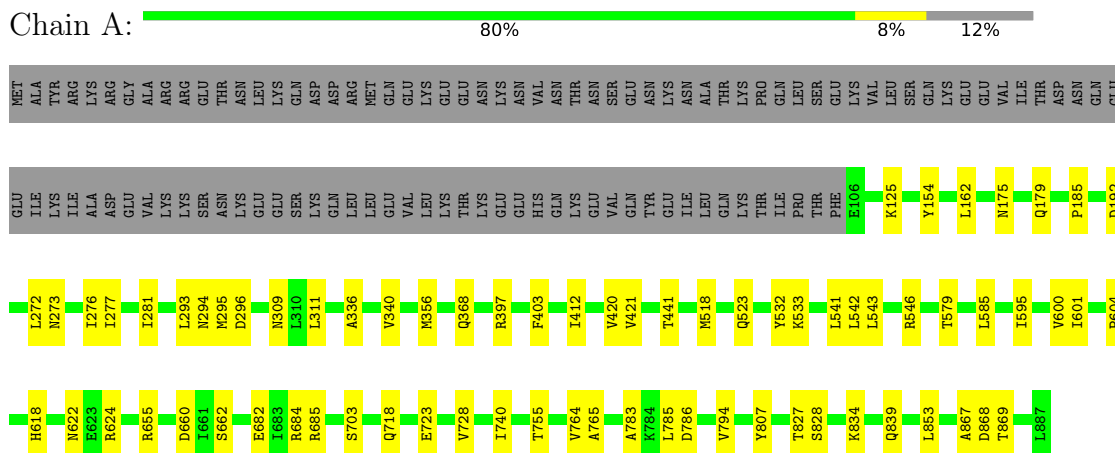
- Molecule 2 is a protein called RNA-directed RNA polymerase.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	P	1086	Total	C	H	N	O	S	0	0
			17667	5636	8874	1456	1662	39		

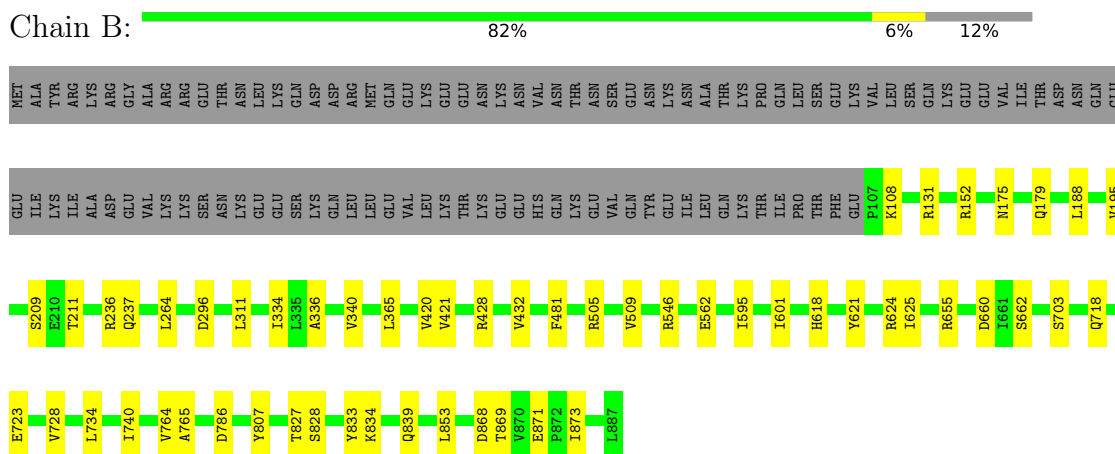
3 Residue-property plots

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. 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. 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.

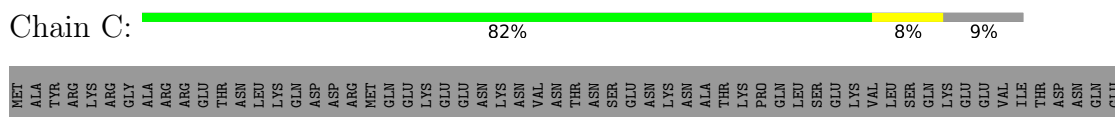
• Molecule 1: Inner capsid protein VP2

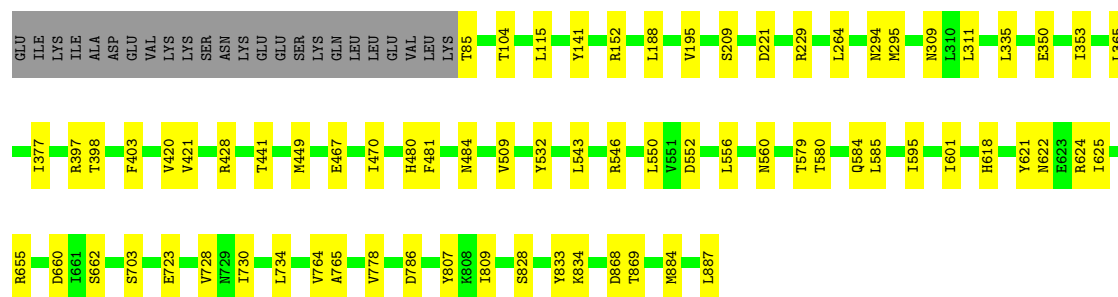


• Molecule 1: Inner capsid protein VP2



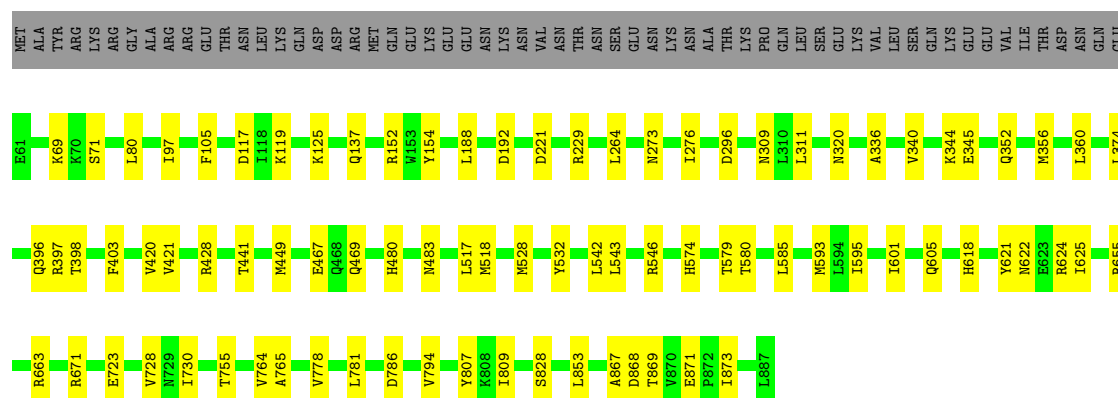
• Molecule 1: Inner capsid protein VP2





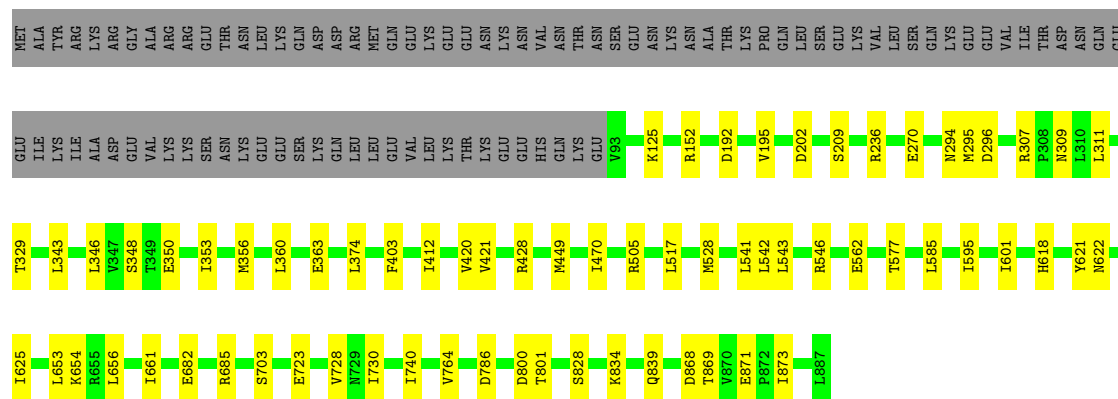
• Molecule 1: Inner capsid protein VP2

Chain D: 84% 10% 7%



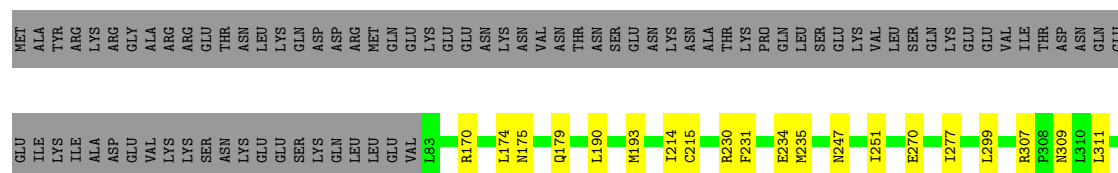
• Molecule 1: Inner capsid protein VP2

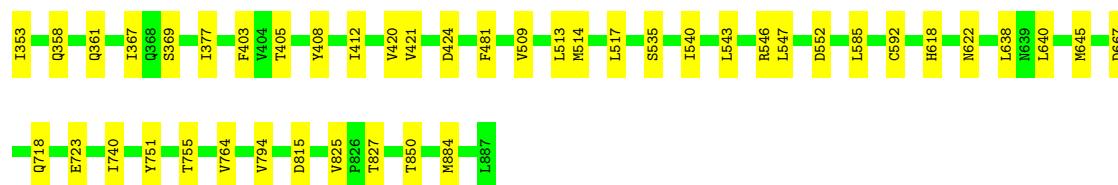
Chain E: 82% 8% 10%



• Molecule 1: Inner capsid protein VP2

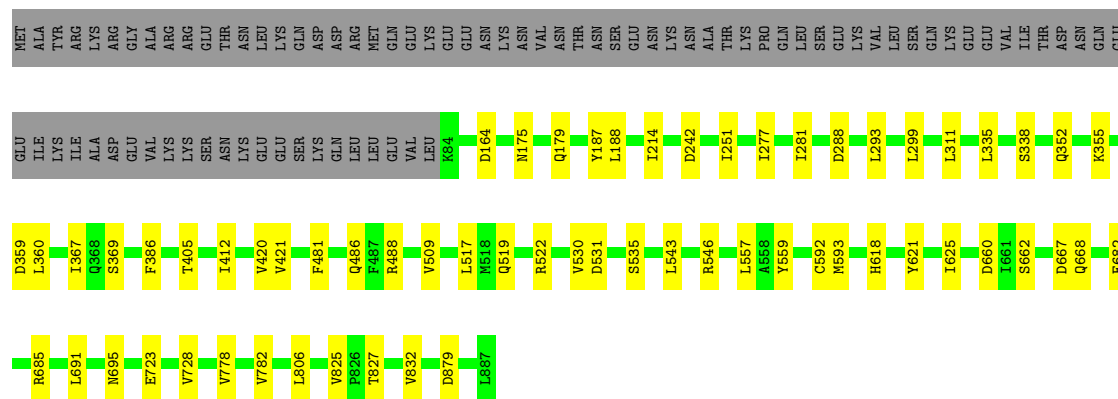
Chain F: 84% 7% 9%





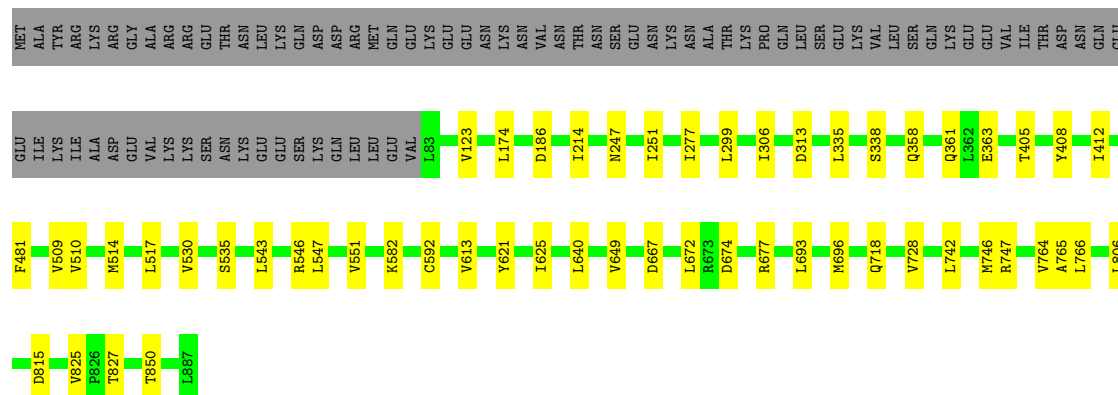
• Molecule 1: Inner capsid protein VP2

Chain G: 84% 7% 9%



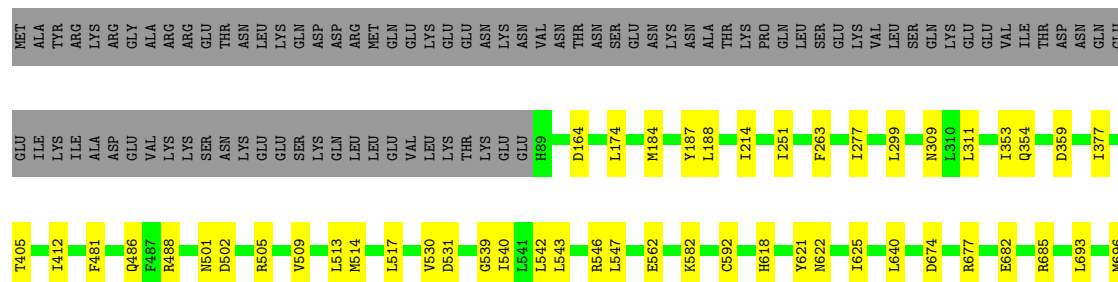
• Molecule 1: Inner capsid protein VP2

Chain H: 85% 6% 9%



• Molecule 1: Inner capsid protein VP2

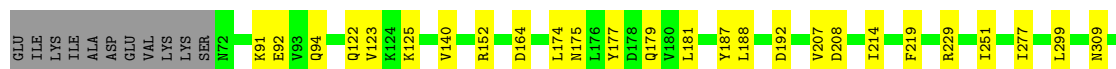
Chain I: 83% 7% 10%





• Molecule 1: Inner capsid protein VP2

Chain J: 84% 8% 8%



• Molecule 2: RNA-directed RNA polymerase

Chain P: 90% 10%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	469892	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)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	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	0.14	0/6503	0.27	0/8823
1	B	0.14	0/6494	0.26	0/8810
1	C	0.14	0/6689	0.28	0/9074
1	D	0.14	0/6884	0.26	0/9332
1	E	0.14	0/6617	0.27	0/8979
1	F	0.13	0/6706	0.27	0/9096
1	G	0.13	0/6698	0.26	0/9085
1	H	0.13	0/6706	0.27	0/9096
1	I	0.13	0/6655	0.26	0/9029
1	J	0.14	0/6797	0.27	0/9217
2	P	0.13	0/8967	0.28	0/12124
All	All	0.14	0/75716	0.27	0/102665

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6387	6413	6413	44	0
1	B	6378	6408	6408	36	0
1	C	6569	6595	6595	43	0
1	D	6764	6809	6809	52	0
1	E	6498	6529	6529	45	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	6586	6619	6619	39	0
1	G	6578	6608	6608	38	0
1	H	6586	6619	6619	40	0
1	I	6535	6563	6563	36	0
1	J	6677	6713	6713	44	0
2	P	8793	8874	8872	66	0
All	All	74351	74750	74748	457	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 3.

All (457) 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:541:LEU:HD13	1:E:528:MET:HE1	1.56	0.87
2:P:937:TYR:OH	2:P:971:ASP:OD2	1.97	0.83
1:A:546:ARG:NH2	1:A:595:ILE:O	2.12	0.82
1:E:868:ASP:OD2	1:I:405:THR:OG1	1.99	0.80
1:J:91:LYS:NZ	2:P:642:GLU:OE2	2.15	0.80
1:A:368:GLN:NE2	2:P:990:GLU:OE2	2.18	0.77
1:E:270:GLU:OE1	1:E:307:ARG:NH1	2.18	0.76
2:P:768:SER:OG	2:P:774:GLU:OE2	2.00	0.76
1:H:247:ASN:OD1	1:H:850:THR:OG1	2.01	0.76
1:C:546:ARG:NH2	1:C:595:ILE:O	2.20	0.75
1:I:674:ASP:OD1	1:I:677:ARG:NH2	2.20	0.75
2:P:880:ASP:OD2	2:P:1069:LYS:NZ	2.19	0.75
1:B:868:ASP:OD2	1:F:405:THR:OG1	2.02	0.75
1:A:523:GLN:O	1:A:533:LYS:NZ	2.18	0.74
1:B:546:ARG:NH2	1:B:595:ILE:O	2.21	0.74
1:E:546:ARG:NH2	1:E:595:ILE:O	2.20	0.74
1:I:486:GLN:O	1:I:488:ARG:NH1	2.21	0.73
2:P:729:ARG:NH2	2:P:767:ASP:O	2.23	0.72
1:E:348:SER:OG	1:E:350:GLU:OE1	2.06	0.71
2:P:528:GLN:O	2:P:531:THR:OG1	2.08	0.71
1:H:514:MET:HE2	1:H:547:LEU:HD22	1.71	0.71
2:P:44:LYS:NZ	2:P:61:GLU:OE2	2.23	0.71
1:I:786:ASP:OD2	1:I:830:THR:OG1	2.08	0.70
1:J:175:ASN:OD1	1:J:179:GLN:NE2	2.25	0.70
1:G:682:GLU:OE1	1:G:685:ARG:NH1	2.25	0.70
1:E:152:ARG:NH1	1:E:723:GLU:OE1	2.25	0.70
1:F:247:ASN:OD1	1:F:850:THR:OG1	2.05	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:486:GLN:O	1:G:488:ARG:NH1	2.25	0.69
1:H:412:ILE:HG22	1:H:543:LEU:HD22	1.74	0.69
1:D:296:ASP:OD2	1:H:582:LYS:NZ	2.26	0.68
1:H:408:TYR:OH	1:H:535:SER:OG	2.11	0.68
1:B:152:ARG:NH1	1:B:723:GLU:OE1	2.26	0.68
1:H:674:ASP:OD1	1:H:677:ARG:NH2	2.27	0.67
1:I:682:GLU:OE1	1:I:685:ARG:NH1	2.27	0.67
1:C:152:ARG:NH1	1:C:723:GLU:OE1	2.26	0.67
1:D:221:ASP:O	1:D:229:ARG:NH2	2.28	0.67
1:C:868:ASP:OD2	1:G:405:THR:OG1	2.09	0.67
1:E:723:GLU:OE2	1:E:839:GLN:NE2	2.28	0.67
1:C:350:GLU:OE2	1:D:71:SER:OG	2.08	0.66
2:P:571:SER:OG	2:P:586:TYR:O	2.14	0.66
1:A:175:ASN:OD1	1:A:179:GLN:NE2	2.29	0.66
1:D:546:ARG:NH2	1:D:595:ILE:O	2.29	0.65
1:F:353:ILE:HD11	1:F:377:ILE:HB	1.78	0.65
1:J:481:PHE:CE1	1:J:509:VAL:HG21	2.31	0.65
1:F:546:ARG:NH1	1:F:592:CYS:O	2.30	0.65
2:P:495:SER:OG	2:P:499:GLN:O	2.12	0.65
1:A:412:ILE:CG2	1:A:543:LEU:HD22	2.27	0.65
1:F:412:ILE:HG22	1:F:543:LEU:HD22	1.77	0.65
1:G:412:ILE:HG22	1:G:543:LEU:HD22	1.77	0.65
1:B:188:LEU:HD11	1:B:264:LEU:HD21	1.79	0.65
1:C:309:ASN:O	1:C:622:ASN:ND2	2.30	0.64
1:G:355:LYS:NZ	1:G:359:ASP:OD2	2.31	0.64
1:J:376:GLY:O	1:J:590:SER:OG	2.16	0.64
1:B:188:LEU:HD11	1:B:264:LEU:CD2	2.28	0.64
1:D:69:LYS:HA	1:D:80:LEU:HD11	1.79	0.64
1:C:556:LEU:O	1:C:560:ASN:ND2	2.31	0.63
1:E:296:ASP:OD2	1:I:582:LYS:NZ	2.31	0.63
1:C:104:THR:OG1	1:C:584:GLN:OE1	2.17	0.62
1:H:408:TYR:HH	1:H:535:SER:HG	1.41	0.61
1:C:703:SER:O	1:C:834:LYS:NZ	2.31	0.61
1:E:412:ILE:CG2	1:E:543:LEU:HD22	2.31	0.61
1:A:311:LEU:O	1:A:618:HIS:NE2	2.33	0.61
1:I:481:PHE:CE1	1:I:509:VAL:HG21	2.35	0.61
1:J:486:GLN:O	1:J:488:ARG:NH1	2.33	0.61
1:J:723:GLU:N	1:J:723:GLU:OE1	2.32	0.61
2:P:791:LEU:HD22	2:P:856:ALA:HB1	1.82	0.61
1:E:311:LEU:O	1:E:618:HIS:NE2	2.33	0.61
1:B:723:GLU:OE2	1:B:839:GLN:NE2	2.33	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:LYS:NZ	1:A:192:ASP:OD2	2.33	0.60
1:G:660:ASP:OD1	1:G:662:SER:OG	2.19	0.60
1:D:528:MET:HE1	1:E:541:LEU:HD13	1.81	0.60
1:E:601:ILE:HD11	1:E:869:THR:CG2	2.31	0.60
1:J:349:THR:OG1	2:P:358:GLU:OE2	2.11	0.60
2:P:77:TYR:CZ	2:P:748:LEU:HD13	2.37	0.60
1:H:693:LEU:HD12	1:H:696:MET:SD	2.42	0.60
1:D:311:LEU:O	1:D:618:HIS:NE2	2.33	0.60
1:J:125:LYS:NZ	1:J:192:ASP:OD2	2.33	0.59
1:C:601:ILE:HD11	1:C:869:THR:HG21	1.84	0.59
1:G:175:ASN:ND2	1:G:179:GLN:OE1	2.35	0.59
1:H:174:LEU:HD22	1:H:640:LEU:HD12	1.84	0.59
2:P:261:PHE:O	2:P:507:THR:OG1	2.11	0.59
1:B:601:ILE:HD11	1:B:869:THR:CG2	2.33	0.59
2:P:680:ARG:HD3	2:P:687:ILE:HD11	1.84	0.59
1:A:154:TYR:OH	1:A:723:GLU:OE2	2.20	0.59
2:P:717:TYR:OH	2:P:770:ASP:OD2	2.21	0.59
2:P:24:ILE:HD11	2:P:49:SER:OG	2.02	0.58
1:B:195:VAL:HG21	1:B:209:SER:HA	1.86	0.58
1:C:428:ARG:HG2	1:H:530:VAL:HG11	1.85	0.58
1:D:97:ILE:HD11	1:D:374:LEU:HB2	1.86	0.58
1:G:481:PHE:CD1	1:G:509:VAL:HG21	2.39	0.58
1:F:175:ASN:OD1	1:F:179:GLN:NE2	2.36	0.58
1:A:412:ILE:HG22	1:A:543:LEU:HD22	1.86	0.58
1:G:481:PHE:CE1	1:G:509:VAL:HG21	2.38	0.58
1:A:660:ASP:OD2	1:A:662:SER:OG	2.22	0.57
1:C:601:ILE:HD11	1:C:869:THR:CG2	2.35	0.57
1:D:137:GLN:NE2	1:D:221:ASP:OD2	2.35	0.57
1:D:663:ARG:NE	1:I:354:GLN:OE1	2.37	0.57
1:F:309:ASN:O	1:F:622:ASN:ND2	2.36	0.57
1:F:481:PHE:CE1	1:F:509:VAL:HG21	2.39	0.57
1:H:481:PHE:CE1	1:H:509:VAL:HG21	2.38	0.57
1:H:546:ARG:NH1	1:H:592:CYS:O	2.36	0.57
1:A:703:SER:O	1:A:834:LYS:NZ	2.34	0.57
1:H:517:LEU:HD11	1:H:543:LEU:HD23	1.87	0.57
1:J:309:ASN:O	1:J:622:ASN:ND2	2.38	0.56
1:C:480:HIS:O	1:C:484:ASN:ND2	2.37	0.56
1:G:311:LEU:O	1:G:618:HIS:NE2	2.36	0.56
1:J:355:LYS:NZ	1:J:359:ASP:OD2	2.37	0.56
1:F:270:GLU:OE1	1:F:307:ARG:NH1	2.38	0.56
2:P:363:TYR:CZ	2:P:367:ILE:HD11	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:729:ARG:NE	2:P:770:ASP:OD1	2.38	0.56
1:H:825:VAL:O	1:H:827:THR:HG23	2.06	0.56
1:H:363:GLU:OE1	1:H:363:GLU:N	2.37	0.56
1:D:398:THR:HB	1:D:580:THR:HG22	1.87	0.56
1:E:786:ASP:OD2	1:E:828:SER:OG	2.24	0.56
1:G:517:LEU:HD11	1:G:543:LEU:HD23	1.88	0.56
1:D:601:ILE:HD11	1:D:869:THR:OG1	2.05	0.55
2:P:887:VAL:HG22	2:P:1054:LEU:CD2	2.35	0.55
1:D:396:GLN:OE1	1:D:574:HIS:ND1	2.38	0.55
1:F:481:PHE:CD1	1:F:509:VAL:HG21	2.41	0.55
1:E:309:ASN:O	1:E:622:ASN:ND2	2.37	0.55
1:C:397:ARG:NH2	1:C:579:THR:OG1	2.40	0.55
1:D:403:PHE:HB3	1:D:585:LEU:HD12	1.89	0.55
1:H:742:LEU:HD21	1:H:766:LEU:HB3	1.89	0.55
2:P:1084:ASN:O	2:P:1087:GLN:NE2	2.36	0.55
1:B:311:LEU:O	1:B:618:HIS:NE2	2.37	0.55
1:D:868:ASP:OD2	1:H:405:THR:OG1	2.14	0.55
1:H:481:PHE:CD1	1:H:509:VAL:HG21	2.42	0.55
1:E:412:ILE:HG22	1:E:543:LEU:HD22	1.88	0.55
1:D:336:ALA:O	1:D:340:VAL:HG23	2.06	0.55
2:P:103:THR:HG22	2:P:103:THR:O	2.06	0.55
1:A:868:ASP:OD1	1:A:869:THR:N	2.40	0.55
1:D:152:ARG:NH1	1:D:723:GLU:OE1	2.40	0.55
1:I:546:ARG:NH1	1:I:592:CYS:O	2.40	0.54
1:I:621:TYR:CZ	1:I:625:ILE:HD11	2.41	0.54
1:J:377:ILE:O	1:J:377:ILE:HG22	2.07	0.54
1:A:718:GLN:HA	1:A:827:THR:HG22	1.90	0.54
1:D:449:MET:HE1	1:D:469:GLN:HB3	1.90	0.54
1:E:125:LYS:NZ	1:E:192:ASP:OD2	2.41	0.54
1:I:481:PHE:CD1	1:I:509:VAL:HG21	2.42	0.54
1:D:352:GLN:NE2	1:D:593:MET:O	2.41	0.54
1:G:667:ASP:OD1	1:G:668:GLN:N	2.41	0.54
2:P:855:SER:O	2:P:859:THR:HG23	2.07	0.54
2:P:787:ALA:O	2:P:790:SER:OG	2.18	0.54
1:I:164:ASP:N	1:I:164:ASP:OD1	2.40	0.54
1:J:546:ARG:NH1	1:J:592:CYS:O	2.40	0.54
2:P:505:ASP:OD1	2:P:505:ASP:N	2.41	0.54
2:P:561:TYR:O	2:P:564:THR:OG1	2.22	0.53
1:D:344:LYS:NZ	1:D:345:GLU:OE2	2.41	0.53
1:E:601:ILE:HD11	1:E:869:THR:HG21	1.89	0.53
1:F:277:ILE:HD11	1:F:299:LEU:HD11	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:815:ASP:OD1	1:H:815:ASP:N	2.39	0.53
1:B:601:ILE:HD11	1:B:869:THR:HG21	1.90	0.53
1:C:311:LEU:O	1:C:618:HIS:NE2	2.37	0.53
1:J:363:GLU:N	1:J:363:GLU:OE1	2.41	0.53
2:P:186:ASN:ND2	2:P:192:GLU:OE2	2.40	0.53
1:A:336:ALA:O	1:A:340:VAL:HG23	2.09	0.53
1:B:420:VAL:HG23	1:B:421:VAL:HG23	1.91	0.53
1:D:320:ASN:O	1:D:671:ARG:NH2	2.39	0.53
1:C:398:THR:HB	1:C:580:THR:HG22	1.91	0.53
1:D:397:ARG:NH2	1:D:579:THR:OG1	2.40	0.53
2:P:24:ILE:HD12	2:P:77:TYR:HE1	1.73	0.53
2:P:141:ASN:ND2	2:P:208:ASN:OD1	2.41	0.53
1:D:441:THR:O	1:D:532:TYR:OH	2.27	0.52
1:E:195:VAL:HG21	1:E:209:SER:HA	1.91	0.52
1:B:621:TYR:CZ	1:B:625:ILE:HD11	2.43	0.52
1:D:117:ASP:O	1:D:119:LYS:NZ	2.39	0.52
1:J:728:VAL:CG2	1:J:806:LEU:HD11	2.40	0.52
2:P:24:ILE:HD12	2:P:77:TYR:CE1	2.44	0.52
1:E:420:VAL:HG23	1:E:421:VAL:HG23	1.90	0.52
2:P:1023:LYS:NZ	2:P:1058:TYR:O	2.37	0.52
1:C:221:ASP:O	1:C:229:ARG:NH2	2.42	0.52
1:B:718:GLN:HA	1:B:827:THR:HG22	1.92	0.52
1:C:543:LEU:HG	1:C:550:LEU:HD11	1.91	0.52
2:P:887:VAL:HG22	2:P:1054:LEU:HD21	1.91	0.51
2:P:51:ASN:OD1	2:P:53:LEU:HD13	2.10	0.51
1:A:441:THR:O	1:A:532:TYR:OH	2.26	0.51
1:E:800:ASP:OD1	1:E:801:THR:N	2.44	0.51
1:J:481:PHE:CD1	1:J:509:VAL:HG21	2.45	0.51
1:D:125:LYS:NZ	1:D:192:ASP:OD2	2.41	0.51
1:F:815:ASP:OD1	1:F:815:ASP:N	2.44	0.51
1:H:514:MET:HE2	1:H:547:LEU:CD2	2.39	0.51
1:G:621:TYR:CZ	1:G:625:ILE:HD11	2.46	0.51
1:I:309:ASN:O	1:I:622:ASN:ND2	2.39	0.51
2:P:357:ASP:OD2	2:P:361:ARG:NE	2.43	0.51
1:G:367:ILE:HG22	1:G:369:SER:H	1.76	0.51
1:E:682:GLU:OE1	1:E:685:ARG:NH1	2.44	0.51
1:B:432:VAL:HG12	1:C:887:LEU:HD22	1.91	0.51
1:C:728:VAL:HG23	1:C:807:TYR:O	2.10	0.51
1:D:517:LEU:CD1	1:D:543:LEU:HD23	2.41	0.51
1:D:528:MET:HE1	1:E:541:LEU:CD1	2.41	0.50
1:J:92:GLU:OE2	1:J:94:GLN:NE2	2.40	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:408:TYR:OH	1:F:535:SER:OG	2.12	0.50
1:I:184:MET:HE3	1:I:263:PHE:HD1	1.76	0.50
1:H:277:ILE:HD11	1:H:299:LEU:HD11	1.94	0.50
1:E:202:ASP:OD2	1:E:236:ARG:NH1	2.44	0.50
1:G:288:ASP:OD2	1:G:559:TYR:OH	2.29	0.50
1:A:601:ILE:HD11	1:A:869:THR:CG2	2.41	0.50
1:F:190:LEU:HD23	1:F:193:MET:HE2	1.94	0.50
1:B:336:ALA:O	1:B:340:VAL:HG23	2.11	0.50
1:F:358:GLN:O	1:F:361:GLN:NE2	2.38	0.50
1:G:352:GLN:NE2	1:G:593:MET:O	2.40	0.50
1:H:621:TYR:CZ	1:H:625:ILE:HD11	2.47	0.50
1:J:621:TYR:CZ	1:J:625:ILE:HD11	2.46	0.50
1:I:184:MET:HE3	1:I:263:PHE:CD1	2.47	0.49
1:J:815:ASP:OD1	1:J:815:ASP:N	2.44	0.49
1:A:397:ARG:NH2	1:A:579:THR:OG1	2.44	0.49
1:B:660:ASP:OD2	1:B:662:SER:OG	2.30	0.49
1:B:871:GLU:HG3	1:B:873:ILE:HG22	1.94	0.49
1:I:187:TYR:C	1:I:188:LEU:HD12	2.37	0.49
1:A:356:MET:HG2	1:A:542:LEU:HD22	1.94	0.49
1:J:360:LEU:HD12	1:J:374:LEU:HD21	1.95	0.49
2:P:194:GLU:N	2:P:194:GLU:OE1	2.46	0.49
1:E:740:ILE:HD12	1:E:764:VAL:HG21	1.93	0.49
1:F:514:MET:HE2	1:F:547:LEU:HD22	1.94	0.49
1:I:723:GLU:OE1	1:I:723:GLU:N	2.42	0.49
1:F:718:GLN:HA	1:F:827:THR:HG22	1.95	0.49
1:A:420:VAL:HG23	1:A:421:VAL:HG23	1.94	0.49
1:J:140:VAL:HG11	1:J:152:ARG:HD2	1.94	0.49
1:A:786:ASP:OD2	1:A:828:SER:OG	2.31	0.49
1:I:514:MET:HE2	1:I:547:LEU:HD22	1.93	0.49
1:G:187:TYR:C	1:G:188:LEU:HD12	2.38	0.49
1:G:214:ILE:HG21	1:G:251:ILE:HG21	1.95	0.49
1:C:188:LEU:HD23	1:C:264:LEU:CD2	2.42	0.48
1:E:505:ARG:NH2	1:E:562:GLU:OE1	2.45	0.48
2:P:401:SER:O	2:P:420:LYS:NZ	2.39	0.48
1:G:546:ARG:NH1	1:G:592:CYS:O	2.46	0.48
2:P:453:ASP:OD1	2:P:459:THR:OG1	2.27	0.48
1:B:728:VAL:HG23	1:B:807:TYR:O	2.13	0.48
1:D:309:ASN:O	1:D:622:ASN:ND2	2.43	0.48
1:F:230:ARG:NH1	1:F:234:GLU:OE2	2.47	0.48
1:I:174:LEU:HD22	1:I:640:LEU:CD1	2.43	0.48
1:I:728:VAL:CG2	1:I:806:LEU:HD11	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:621:TYR:CZ	1:E:625:ILE:HD11	2.49	0.48
1:F:517:LEU:HB2	1:F:540:ILE:HG23	1.94	0.48
1:G:335:LEU:O	1:G:338:SER:OG	2.30	0.48
1:A:518:MET:HE2	1:A:518:MET:HA	1.95	0.48
1:D:403:PHE:CB	1:D:585:LEU:HD12	2.43	0.48
1:A:294:ASN:OD1	1:A:295:MET:N	2.46	0.48
1:A:723:GLU:OE2	1:A:839:GLN:NE2	2.46	0.48
1:B:786:ASP:OD2	1:B:828:SER:OG	2.31	0.48
1:C:195:VAL:HG21	1:C:209:SER:HA	1.96	0.48
1:C:552:ASP:HB3	1:C:884:MET:HE2	1.96	0.48
1:C:188:LEU:HD23	1:C:264:LEU:HG	1.96	0.48
1:G:281:ILE:HD13	1:G:293:LEU:HD13	1.96	0.47
2:P:930:ILE:N	2:P:984:ASP:OD2	2.45	0.47
1:G:360:LEU:O	1:G:535:SER:OG	2.32	0.47
1:C:778:VAL:HG22	1:C:809:ILE:HG23	1.96	0.47
2:P:591:SER:HA	2:P:596:THR:HG21	1.96	0.47
1:H:718:GLN:HA	1:H:827:THR:HG22	1.95	0.47
1:C:621:TYR:CZ	1:C:625:ILE:HD11	2.50	0.47
1:C:786:ASP:OD2	1:C:828:SER:OG	2.32	0.47
1:J:122:GLN:OE1	1:J:123:VAL:N	2.47	0.47
1:C:728:VAL:HG22	1:C:730:ILE:HG12	1.95	0.47
1:D:621:TYR:CZ	1:D:625:ILE:HD11	2.50	0.47
1:I:311:LEU:O	1:I:618:HIS:NE2	2.41	0.47
1:I:353:ILE:HD11	1:I:377:ILE:HB	1.96	0.47
2:P:145:VAL:HG11	2:P:170:PHE:CD1	2.49	0.47
1:E:403:PHE:HB3	1:E:585:LEU:HD12	1.97	0.47
2:P:789:MET:HE2	2:P:1074:THR:HG22	1.96	0.47
1:H:746:MET:HE3	1:H:747:ARG:NH1	2.30	0.47
1:J:412:ILE:HG22	1:J:543:LEU:HD22	1.96	0.47
2:P:646:GLN:NE2	2:P:650:ASP:OD1	2.47	0.47
1:E:363:GLU:O	2:P:983:GLN:NE2	2.48	0.46
1:C:420:VAL:HG23	1:C:421:VAL:HG23	1.97	0.46
1:C:660:ASP:OD2	1:C:662:SER:OG	2.34	0.46
1:F:552:ASP:HB3	1:F:884:MET:HE2	1.97	0.46
1:C:403:PHE:CB	1:C:585:LEU:HD12	2.46	0.46
1:D:786:ASP:OD2	1:D:828:SER:OG	2.33	0.46
1:C:467:GLU:OE2	1:C:480:HIS:ND1	2.40	0.46
1:F:367:ILE:HG22	1:F:369:SER:H	1.81	0.46
1:D:428:ARG:HG2	1:I:530:VAL:HG11	1.97	0.46
1:I:746:MET:SD	1:I:747:ARG:NH1	2.89	0.46
2:P:745:THR:HG22	2:P:746:GLY:H	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:853:LEU:HD23	1:J:667:ASP:HA	1.98	0.45
1:F:751:TYR:O	1:F:755:THR:OG1	2.22	0.45
1:G:519:GLN:OE1	1:G:522:ARG:NH2	2.47	0.45
1:D:356:MET:HE2	1:D:374:LEU:HD23	1.98	0.45
1:H:174:LEU:HD22	1:H:640:LEU:CD1	2.47	0.45
1:D:778:VAL:HG22	1:D:809:ILE:HG23	1.97	0.45
1:A:296:ASP:OD2	1:J:582:LYS:NZ	2.49	0.45
1:A:403:PHE:HB3	1:A:585:LEU:HD12	1.99	0.45
2:P:387:GLN:NE2	2:P:551:ASP:OD2	2.46	0.45
2:P:575:ILE:HD12	2:P:584:ILE:HD12	1.97	0.45
1:A:728:VAL:HG23	1:A:807:TYR:O	2.17	0.45
1:F:311:LEU:O	1:F:618:HIS:NE2	2.47	0.45
1:F:755:THR:HG23	1:F:794:VAL:HG13	1.97	0.45
2:P:542:LEU:HD11	2:P:561:TYR:HD1	1.82	0.45
1:B:703:SER:O	1:B:834:LYS:NZ	2.45	0.45
1:D:764:VAL:HG22	1:D:765:ALA:N	2.31	0.45
1:J:177:TYR:CZ	1:J:181:LEU:HD11	2.51	0.45
2:P:75:LEU:HD22	2:P:748:LEU:HD11	1.97	0.45
1:D:467:GLU:OE1	1:D:483:ASN:ND2	2.45	0.45
1:J:335:LEU:O	1:J:338:SER:OG	2.33	0.45
1:C:481:PHE:CZ	1:C:509:VAL:HG11	2.52	0.44
1:A:185:PRO:O	1:A:684:ARG:NH2	2.50	0.44
1:I:214:ILE:HG21	1:I:251:ILE:HG21	1.99	0.44
1:D:360:LEU:HD21	1:D:542:LEU:HD11	1.98	0.44
1:E:703:SER:O	1:E:834:LYS:NZ	2.51	0.44
1:H:728:VAL:CG2	1:H:806:LEU:HD11	2.48	0.44
2:P:86:GLU:O	2:P:90:VAL:HG13	2.17	0.44
1:E:517:LEU:CD1	1:E:543:LEU:HD23	2.47	0.44
1:F:718:GLN:HG2	1:F:825:VAL:HG11	1.99	0.44
1:H:277:ILE:CD1	1:H:299:LEU:HD11	2.47	0.44
1:A:403:PHE:CB	1:A:585:LEU:HD12	2.47	0.44
1:C:353:ILE:HD11	1:C:377:ILE:HB	1.99	0.44
1:E:428:ARG:HG2	1:J:530:VAL:HG21	2.00	0.44
1:G:825:VAL:O	1:G:827:THR:HG23	2.17	0.44
1:I:531:ASP:N	1:I:531:ASP:OD1	2.51	0.44
1:J:800:ASP:N	1:J:800:ASP:OD1	2.50	0.44
1:B:740:ILE:HD12	1:B:764:VAL:HG21	2.00	0.44
1:J:703:SER:O	1:J:834:LYS:NZ	2.50	0.44
1:F:174:LEU:HD22	1:F:640:LEU:HD12	2.00	0.44
1:D:853:LEU:HD23	1:H:667:ASP:HA	2.00	0.43
1:J:214:ILE:HG21	1:J:251:ILE:HG21	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:309:ASN:O	1:A:622:ASN:ND2	2.46	0.43
1:D:154:TYR:OH	1:D:723:GLU:OE2	2.33	0.43
1:D:624:ARG:HD2	1:D:655:ARG:HB3	2.00	0.43
1:E:871:GLU:HG3	1:E:873:ILE:HG22	1.99	0.43
1:G:386:PHE:CE1	1:G:557:LEU:HD12	2.54	0.43
1:H:649:VAL:HG13	1:H:672:LEU:HD23	2.00	0.43
1:I:517:LEU:HB2	1:I:540:ILE:HG23	1.99	0.43
1:D:755:THR:HG23	1:D:794:VAL:HG13	2.01	0.43
1:F:214:ILE:HG21	1:F:251:ILE:HG21	1.99	0.43
1:G:242:ASP:OD1	1:G:242:ASP:N	2.51	0.43
1:G:778:VAL:O	1:G:782:VAL:HG23	2.18	0.43
1:J:349:THR:HG23	1:J:377:ILE:O	2.18	0.43
2:P:528:GLN:C	2:P:531:THR:HG1	2.25	0.43
1:A:755:THR:HG23	1:A:794:VAL:HG13	2.01	0.43
1:B:505:ARG:NH2	1:B:562:GLU:OE1	2.47	0.43
2:P:613:SER:O	2:P:616:SER:OG	2.35	0.43
1:E:353:ILE:HG13	2:P:974:THR:HG23	2.00	0.43
1:E:517:LEU:HD13	1:E:543:LEU:HD23	2.01	0.43
1:I:412:ILE:HG22	1:I:543:LEU:HD22	1.99	0.43
1:I:513:LEU:O	1:I:517:LEU:HG	2.19	0.43
1:J:174:LEU:HD22	1:J:640:LEU:CD1	2.48	0.43
2:P:374:MET:SD	2:P:602:ILE:HD11	2.58	0.43
1:A:272:LEU:HD11	1:A:277:ILE:HG13	2.01	0.43
1:B:734:LEU:HD21	1:B:833:TYR:CG	2.53	0.43
1:I:764:VAL:HG22	1:I:765:ALA:N	2.33	0.43
2:P:145:VAL:CG2	2:P:211:ILE:HG23	2.49	0.43
1:F:420:VAL:HG23	1:F:421:VAL:HG23	2.00	0.43
1:G:691:LEU:O	1:G:695:ASN:ND2	2.52	0.43
2:P:184:ARG:H	2:P:195:VAL:HG21	1.83	0.43
2:P:319:PHE:N	2:P:320:PRO:HD2	2.33	0.43
1:F:277:ILE:CD1	1:F:299:LEU:HD11	2.49	0.43
1:G:420:VAL:HG23	1:G:421:VAL:HG23	2.00	0.43
1:H:510:VAL:HB	1:H:551:VAL:HG22	2.01	0.43
1:I:825:VAL:O	1:I:827:THR:HG23	2.19	0.43
1:J:621:TYR:CE2	1:J:625:ILE:HD11	2.54	0.43
1:A:764:VAL:HG22	1:A:765:ALA:H	1.84	0.42
1:G:723:GLU:OE1	1:G:723:GLU:N	2.50	0.42
2:P:337:GLN:O	2:P:451:ARG:NH1	2.46	0.42
1:A:604:PRO:HB2	1:A:867:ALA:HB3	2.01	0.42
1:A:624:ARG:HD2	1:A:655:ARG:HB3	2.01	0.42
1:A:682:GLU:OE1	1:A:685:ARG:NH1	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:LYS:HG3	1:B:334:ILE:HD12	2.01	0.42
1:D:728:VAL:HG23	1:D:807:TYR:O	2.19	0.42
1:F:740:ILE:HD12	1:F:764:VAL:HG21	2.01	0.42
1:H:186:ASP:OD2	1:H:186:ASP:N	2.51	0.42
1:E:621:TYR:CE2	1:E:625:ILE:HD11	2.54	0.42
1:F:215:CYS:SG	1:F:235:MET:HE1	2.59	0.42
1:F:424:ASP:OD1	1:F:424:ASP:N	2.52	0.42
1:H:335:LEU:O	1:H:338:SER:OG	2.36	0.42
1:E:329:THR:HG21	1:E:577:THR:HG21	2.01	0.42
1:E:449:MET:SD	1:E:470:ILE:HD13	2.59	0.42
1:E:353:ILE:HG23	1:E:374:LEU:HD23	2.02	0.42
1:H:214:ILE:HG21	1:H:251:ILE:HG21	2.01	0.42
1:H:517:LEU:CD1	1:H:543:LEU:HD23	2.49	0.42
2:P:426:ASP:OD2	2:P:432:ARG:NH2	2.51	0.42
1:B:764:VAL:HG22	1:B:765:ALA:H	1.84	0.42
1:J:531:ASP:OD1	1:J:531:ASP:N	2.52	0.42
1:D:605:GLN:NE2	1:D:867:ALA:O	2.50	0.42
1:I:539:GLY:HA2	1:I:542:LEU:HD12	2.02	0.42
1:B:624:ARG:HD2	1:B:655:ARG:HB3	2.02	0.42
1:F:170:ARG:NE	1:F:638:LEU:O	2.52	0.42
1:H:123:VAL:HG22	1:H:306:ILE:HD12	2.01	0.42
1:I:277:ILE:CD1	1:I:299:LEU:HD11	2.50	0.42
1:I:501:ASN:OD1	1:I:502:ASP:N	2.53	0.42
1:A:740:ILE:HD12	1:A:764:VAL:HG21	2.01	0.42
1:C:294:ASN:OD1	1:C:295:MET:N	2.52	0.42
1:D:781:LEU:HD13	1:D:807:TYR:CD2	2.54	0.42
1:D:871:GLU:HG3	1:D:873:ILE:HG22	2.01	0.42
1:G:879:ASP:OD1	1:G:879:ASP:N	2.51	0.42
1:J:164:ASP:OD1	1:J:164:ASP:N	2.42	0.42
1:B:175:ASN:OD1	1:B:179:GLN:NE2	2.49	0.42
1:D:728:VAL:HG22	1:D:730:ILE:HG13	2.02	0.42
1:F:645:MET:SD	1:F:645:MET:N	2.93	0.42
1:J:187:TYR:C	1:J:188:LEU:HD12	2.44	0.42
2:P:35:GLU:OE2	2:P:866:THR:HG22	2.19	0.42
1:A:601:ILE:HD11	1:A:869:THR:HG21	2.01	0.41
1:B:365:LEU:HD13	1:C:365:LEU:HG	2.00	0.41
1:J:600:VAL:O	1:J:600:VAL:HG23	2.20	0.41
2:P:878:LEU:HD22	2:P:1036:TYR:CD1	2.55	0.41
1:A:281:ILE:HD13	1:A:293:LEU:HD13	2.02	0.41
1:C:624:ARG:HD2	1:C:655:ARG:HB3	2.02	0.41
1:D:420:VAL:HG23	1:D:421:VAL:HG23	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:518:MET:HE2	1:D:518:MET:HA	2.02	0.41
1:E:654:LYS:HG3	1:E:661:ILE:HD13	2.03	0.41
1:G:531:ASP:OD1	1:G:531:ASP:N	2.50	0.41
2:P:38:CYS:SG	2:P:65:VAL:HG13	2.60	0.41
1:A:783:ALA:HB3	1:A:785:LEU:HD13	2.02	0.41
1:B:131:ARG:NH2	1:B:211:THR:OG1	2.52	0.41
1:B:188:LEU:HD11	1:B:264:LEU:HD23	2.02	0.41
1:D:467:GLU:OE2	1:D:480:HIS:ND1	2.42	0.41
1:E:343:LEU:HB3	1:E:346:LEU:HD21	2.00	0.41
1:G:164:ASP:OD1	1:G:164:ASP:N	2.53	0.41
1:J:207:VAL:HA	1:J:208:ASP:HA	1.88	0.41
1:D:188:LEU:HD22	1:D:264:LEU:HD21	2.02	0.41
1:D:273:ASN:ND2	1:D:276:ILE:HD12	2.35	0.41
1:H:313:ASP:OD1	1:H:621:TYR:OH	2.35	0.41
1:H:335:LEU:CD1	1:H:613:VAL:HG11	2.51	0.41
1:I:693:LEU:HD12	1:I:696:MET:SD	2.61	0.41
1:C:141:TYR:N	1:C:723:GLU:O	2.51	0.41
1:C:441:THR:O	1:C:532:TYR:OH	2.34	0.41
1:J:219:PHE:O	1:J:229:ARG:NH1	2.51	0.41
1:B:296:ASP:OD1	1:B:296:ASP:N	2.48	0.41
1:H:764:VAL:HG22	1:H:765:ALA:N	2.35	0.41
1:J:429:GLU:OE1	1:J:429:GLU:N	2.53	0.41
1:A:600:VAL:HG23	1:A:600:VAL:O	2.21	0.41
1:A:764:VAL:HG22	1:A:765:ALA:N	2.36	0.41
1:E:356:MET:HG2	1:E:542:LEU:HD22	2.03	0.41
1:G:806:LEU:HD13	1:G:832:VAL:HB	2.02	0.41
1:I:505:ARG:NH2	1:I:562:GLU:OE1	2.46	0.41
1:F:403:PHE:HB3	1:F:585:LEU:HD12	2.01	0.41
1:J:532:TYR:CE2	1:J:536:ILE:HD11	2.56	0.41
1:C:449:MET:SD	1:C:470:ILE:HD13	2.61	0.41
1:D:356:MET:HE2	1:D:374:LEU:CD2	2.51	0.41
1:G:188:LEU:HD12	1:G:188:LEU:N	2.36	0.41
1:J:517:LEU:HD11	1:J:543:LEU:HD23	2.03	0.41
1:C:764:VAL:HG22	1:C:765:ALA:N	2.35	0.41
1:B:428:ARG:HG2	1:G:530:VAL:HG11	2.02	0.40
1:B:853:LEU:HD23	1:F:667:ASP:HA	2.01	0.40
1:E:653:LEU:HD23	1:E:656:LEU:HD12	2.03	0.40
1:F:513:LEU:O	1:F:517:LEU:HG	2.20	0.40
1:F:723:GLU:OE1	1:F:723:GLU:N	2.49	0.40
1:G:728:VAL:CG2	1:G:806:LEU:HD11	2.50	0.40
2:P:77:TYR:CE1	2:P:748:LEU:HD13	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:123:GLU:OE1	2:P:123:GLU:N	2.48	0.40
1:A:273:ASN:OD1	1:A:276:ILE:HD12	2.21	0.40
1:B:621:TYR:CE2	1:B:625:ILE:HD11	2.57	0.40
1:C:115:LEU:HD23	1:C:335:LEU:HB2	2.01	0.40
1:C:403:PHE:HB3	1:C:585:LEU:HD12	2.04	0.40
1:E:728:VAL:HG22	1:E:730:ILE:HG13	2.02	0.40
1:J:277:ILE:CD1	1:J:299:LEU:HD11	2.50	0.40
2:P:900:MET:HE3	2:P:903:LEU:HD11	2.03	0.40
2:P:1084:ASN:ND2	2:P:1086:PHE:O	2.54	0.40
1:B:236:ARG:NH2	1:B:237:GLN:O	2.53	0.40
2:P:542:LEU:HD11	2:P:561:TYR:CD1	2.56	0.40
1:B:481:PHE:CZ	1:B:509:VAL:HG11	2.56	0.40
1:C:734:LEU:HD21	1:C:833:TYR:CG	2.56	0.40
1:E:294:ASN:OD1	1:E:295:MET:N	2.55	0.40
1:E:360:LEU:HD21	1:E:542:LEU:HD11	2.02	0.40
1:J:624:ARG:HD2	1:J:655:ARG:HB3	2.04	0.40
1:F:231:PHE:O	1:F:235:MET:N	2.51	0.40
1:G:277:ILE:HD11	1:G:299:LEU:HD11	2.03	0.40
1:H:358:GLN:O	1:H:361:GLN:NE2	2.54	0.40
1:H:412:ILE:CG2	1:H:543:LEU:HD22	2.47	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	780/887 (88%)	767 (98%)	13 (2%)	0	100	100
1	B	779/887 (88%)	767 (98%)	12 (2%)	0	100	100
1	C	801/887 (90%)	787 (98%)	14 (2%)	0	100	100
1	D	825/887 (93%)	811 (98%)	13 (2%)	1 (0%)	48	66

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	793/887 (89%)	782 (99%)	11 (1%)	0	100	100
1	F	803/887 (90%)	789 (98%)	14 (2%)	0	100	100
1	G	802/887 (90%)	789 (98%)	13 (2%)	0	100	100
1	H	803/887 (90%)	790 (98%)	13 (2%)	0	100	100
1	I	797/887 (90%)	787 (99%)	10 (1%)	0	100	100
1	J	814/887 (92%)	800 (98%)	14 (2%)	0	100	100
2	P	1084/1088 (100%)	1054 (97%)	30 (3%)	0	100	100
All	All	9081/9958 (91%)	8923 (98%)	157 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	105	PHE

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	718/818 (88%)	717 (100%)	1 (0%)	88	95
1	B	717/818 (88%)	717 (100%)	0	100	100
1	C	739/818 (90%)	738 (100%)	1 (0%)	88	95
1	D	762/818 (93%)	762 (100%)	0	100	100
1	E	731/818 (89%)	731 (100%)	0	100	100
1	F	741/818 (91%)	741 (100%)	0	100	100
1	G	740/818 (90%)	740 (100%)	0	100	100
1	H	741/818 (91%)	741 (100%)	0	100	100
1	I	735/818 (90%)	734 (100%)	1 (0%)	88	95
1	J	752/818 (92%)	752 (100%)	0	100	100
2	P	987/989 (100%)	986 (100%)	1 (0%)	88	95

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	8363/9169 (91%)	8359 (100%)	4 (0%)	100	100

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162	LEU
1	C	85	THR
1	I	359	ASP
2	P	139	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (101) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	199	ASN
1	A	257	HIS
1	A	266	HIS
1	A	287	ASN
1	A	354	GLN
1	A	361	GLN
1	A	460	GLN
1	A	468	GLN
1	A	668	GLN
1	A	695	ASN
1	A	733	ASN
1	A	760	ASN
1	A	835	GLN
1	A	847	HIS
1	B	257	HIS
1	B	266	HIS
1	B	372	GLN
1	B	497	ASN
1	B	503	ASN
1	B	605	GLN
1	B	657	GLN
1	B	708	GLN
1	B	756	ASN
1	B	760	ASN
1	B	885	ASN
1	C	94	GLN
1	C	450	GLN
1	C	460	GLN

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Mol	Chain	Res	Type
1	C	468	GLN
1	C	695	ASN
1	C	708	GLN
1	C	835	GLN
1	D	90	GLN
1	D	257	HIS
1	D	318	HIS
1	D	368	GLN
1	D	423	ASN
1	D	460	GLN
1	D	503	ASN
1	D	537	GLN
1	D	584	GLN
1	D	695	ASN
1	D	733	ASN
1	D	792	GLN
1	D	847	HIS
1	E	257	HIS
1	E	266	HIS
1	E	273	ASN
1	E	287	ASN
1	E	423	ASN
1	E	450	GLN
1	E	503	ASN
1	E	511	ASN
1	E	537	GLN
1	E	642	GLN
1	E	847	HIS
1	F	94	GLN
1	F	294	ASN
1	F	460	GLN
1	F	657	GLN
1	F	668	GLN
1	F	733	ASN
1	G	294	ASN
1	G	460	GLN
1	G	668	GLN
1	G	756	ASN
1	G	847	HIS
1	G	885	ASN
1	H	90	GLN
1	H	294	ASN

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Mol	Chain	Res	Type
1	H	512	GLN
1	H	792	GLN
1	H	852	ASN
1	H	874	ASN
1	I	126	GLN
1	I	273	ASN
1	I	294	ASN
1	I	372	GLN
1	I	460	GLN
1	I	512	GLN
1	I	668	GLN
1	I	733	ASN
1	I	753	GLN
1	I	756	ASN
1	J	240	GLN
1	J	287	ASN
1	J	372	GLN
1	J	460	GLN
1	J	477	ASN
1	J	524	GLN
1	J	792	GLN
1	J	847	HIS
1	J	874	ASN
2	P	290	GLN
2	P	430	ASN
2	P	498	ASN
2	P	529	HIS
2	P	638	GLN
2	P	825	ASN
2	P	862	GLN
2	P	1084	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

This section contains visualisations of the EMDB entry EMD-45123. These allow visual inspection of the internal detail of the map and identification of artifacts.

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

6.1 Orthogonal projections

This section was not generated.

6.2 Central slices

This section was not generated.

6.3 Largest variance slices

This section was not generated.

6.4 Orthogonal standard-deviation projections (False-color)

This section was not generated.

6.5 Orthogonal surface views

This section was not generated.

6.6 Mask visualisation

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

7 Map analysis ⓘ

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

7.1 Map-value distribution ⓘ

This section was not generated.

7.2 Volume estimate versus contour level ⓘ

This section was not generated.

7.3 Rotationally averaged power spectrum ⓘ

This section was not generated. The rotationally averaged power spectrum had issues being displayed.

8 Fourier-Shell correlation ⓘ

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

9 Map-model fit

This section was not generated.