



wwPDB EM Map/Model Validation Report ⓘ

Jul 6, 2016 – 08:32 AM EDT

PDB ID : 3JB8
Title : Insight into Three-dimensional structure of Maize Chlorotic Mottle Virus Revealed by Single Particle Analysis
Authors : Wang, C.Y.; Zhang, Q.F.; Gao, Y.Z.; Zhou, X.P.; Ji, G.; Huang, X.J.; Hong, J.; Zhang, C.X.
Deposited on : 2015-08-04
Resolution : 3.60 Å(reported)

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.
For rigid body fitted models, validation errors reported here could stem from errors in the original structure(s) used in the fitting.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/EMValidationReportHelp>

MolProbity : 4.02b-467
Mogul : unknown
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027790

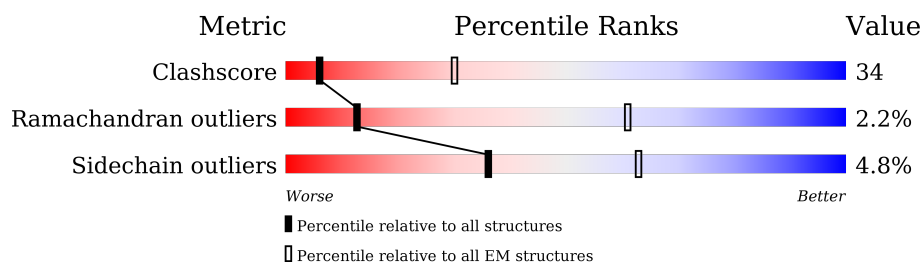
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.60 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	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	185	 64% 30% 6%
1	B	185	 61% 32% 6% •
1	C	185	 62% 31% 6% •

2 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 4098 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.

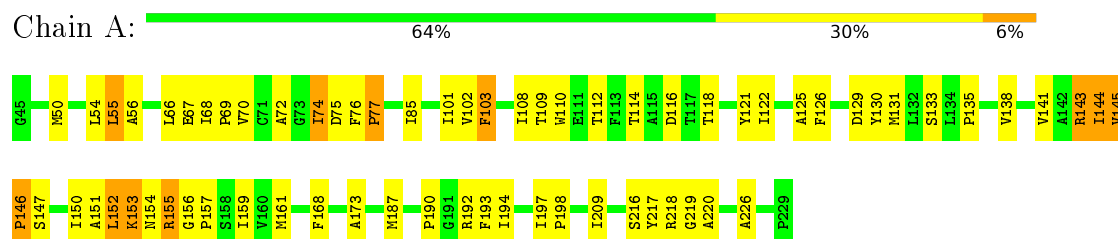
- Molecule 1 is a protein called Coat protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	185	Total	C	N	O	S	0	0
			1378	881	231	257	9		
1	B	182	Total	C	N	O	S	0	0
			1356	865	228	254	9		
1	C	183	Total	C	N	O	S	0	0
			1364	871	229	255	9		

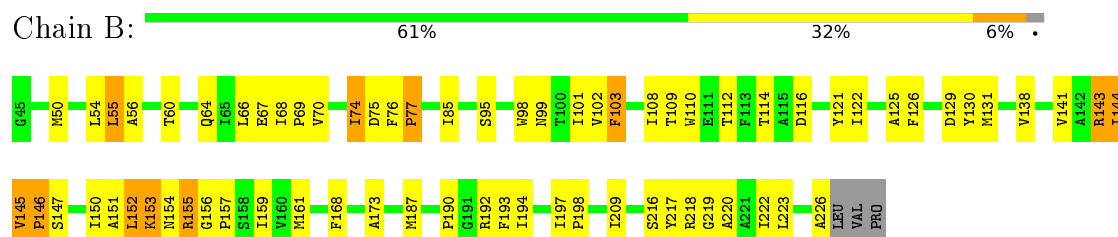
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Coat protein



• Molecule 1: Coat protein



• Molecule 1: Coat protein



4 Experimental information

Property	Value	Source
Reconstruction method	Not provided	Depositor
Imposed symmetry	POINT, I	Depositor
Number of images used	53600	Depositor
Resolution determination method	FSC 0.143	Depositor
CTF correction method	Not provided	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	Not provided	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	Depositor
Image detector	GATAN ULTRASCAN 4000 (4k 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	$\# Z > 2$	RMSZ	$\# Z > 2$
1	A	0.51	0/1408	0.59	1/1922 (0.1%)
1	B	0.51	0/1385	0.60	1/1889 (0.1%)
1	C	0.51	0/1393	0.60	1/1900 (0.1%)
All	All	0.51	0/4186	0.60	3/5711 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	145	VAL	C-N-CD	-5.95	107.51	120.60
1	A	145	VAL	C-N-CD	-5.94	107.54	120.60
1	C	145	VAL	C-N-CD	-5.94	107.54	120.60

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	1378	0	1397	103	0
1	B	1356	0	1370	106	0
1	C	1364	0	1381	109	0
All	All	4098	0	4148	279	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 34.

All (279) 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:145:VAL:HG13	1:A:146:PRO:CD	1.08	1.54
1:B:145:VAL:HG13	1:B:146:PRO:CD	1.08	1.53
1:C:145:VAL:HG13	1:C:146:PRO:CD	1.08	1.50
1:B:145:VAL:CG1	1:B:146:PRO:CD	1.93	1.45
1:C:145:VAL:CG1	1:C:146:PRO:CD	1.93	1.44
1:A:145:VAL:CG1	1:A:146:PRO:CD	1.93	1.42
1:B:187:MET:SD	1:C:187:MET:HE3	1.61	1.39
1:A:187:MET:HE3	1:C:187:MET:SD	1.64	1.37
1:B:145:VAL:CG1	1:B:146:PRO:HD3	1.55	1.29
1:C:145:VAL:CG1	1:C:146:PRO:HD2	1.57	1.26
1:B:145:VAL:CG1	1:B:146:PRO:HD2	1.58	1.25
1:A:145:VAL:CG1	1:A:146:PRO:HD2	1.57	1.25
1:C:145:VAL:CG1	1:C:146:PRO:HD3	1.55	1.25
1:A:187:MET:SD	1:B:187:MET:HE3	1.78	1.22
1:A:145:VAL:CG1	1:A:146:PRO:HD3	1.55	1.20
1:B:187:MET:SD	1:C:187:MET:CE	2.34	1.15
1:A:187:MET:CE	1:C:187:MET:SD	2.37	1.12
1:A:187:MET:SD	1:B:187:MET:CE	2.41	1.08
1:A:72:ALA:CB	1:A:101:ILE:HD11	1.83	1.08
1:C:72:ALA:CB	1:C:101:ILE:HD11	1.83	1.07
1:C:145:VAL:HG12	1:C:146:PRO:HD2	1.44	0.99
1:C:145:VAL:HG13	1:C:146:PRO:HD2	1.14	0.98
1:A:145:VAL:HG13	1:A:146:PRO:HD2	1.14	0.98
1:B:143:ARG:O	1:B:144:ILE:HG12	1.62	0.98
1:A:143:ARG:O	1:A:144:ILE:HG12	1.62	0.98
1:C:143:ARG:O	1:C:144:ILE:HG12	1.62	0.97
1:A:145:VAL:HG12	1:A:146:PRO:HD2	1.44	0.96
1:B:145:VAL:HG12	1:B:146:PRO:HD2	1.44	0.96
1:B:187:MET:CE	1:C:187:MET:HE3	1.96	0.94
1:B:101:ILE:HG12	1:B:217:TYR:CD1	2.03	0.94
1:B:145:VAL:HG13	1:B:146:PRO:HD2	1.14	0.91
1:C:72:ALA:HB2	1:C:101:ILE:HD11	1.52	0.89
1:A:72:ALA:HB2	1:A:101:ILE:HD11	1.52	0.89
1:C:72:ALA:CB	1:C:101:ILE:CD1	2.52	0.87
1:A:101:ILE:HG22	1:A:217:TYR:CD1	2.10	0.87
1:A:72:ALA:CB	1:A:101:ILE:CD1	2.52	0.87
1:B:220:ALA:O	1:C:145:VAL:HB	1.75	0.86
1:C:101:ILE:HG22	1:C:217:TYR:CD1	2.10	0.86
1:B:125:ALA:HB1	1:B:144:ILE:HD12	1.58	0.85
1:B:101:ILE:CD1	1:B:217:TYR:HE1	1.88	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:ALA:HB1	1:A:144:ILE:HD12	1.58	0.85
1:B:101:ILE:HD11	1:B:217:TYR:HE1	1.41	0.84
1:A:116:ASP:HA	1:A:153:LYS:HD2	1.60	0.84
1:B:116:ASP:HA	1:B:153:LYS:HD2	1.60	0.84
1:C:72:ALA:HB1	1:C:101:ILE:CD1	2.08	0.84
1:C:125:ALA:HB1	1:C:144:ILE:HD12	1.58	0.83
1:A:72:ALA:HB1	1:A:101:ILE:CD1	2.08	0.82
1:C:116:ASP:HA	1:C:153:LYS:HD2	1.60	0.82
1:A:187:MET:CE	1:B:187:MET:HE3	2.09	0.82
1:A:220:ALA:H	1:B:145:VAL:HB	1.46	0.80
1:B:101:ILE:CD1	1:B:217:TYR:CE1	2.63	0.80
1:B:187:MET:CE	1:C:187:MET:CE	2.60	0.78
1:B:95:SER:OG	1:B:101:ILE:CD1	2.32	0.78
1:B:95:SER:OG	1:B:101:ILE:HD11	1.84	0.78
1:C:121:TYR:HA	1:C:151:ALA:HB3	1.66	0.78
1:A:187:MET:HE3	1:C:187:MET:CE	2.15	0.77
1:A:121:TYR:HA	1:A:151:ALA:HB3	1.66	0.76
1:B:121:TYR:HA	1:B:151:ALA:HB3	1.66	0.75
1:C:101:ILE:HD12	1:C:103:PHE:HE1	1.52	0.74
1:A:101:ILE:HD12	1:A:103:PHE:HE1	1.52	0.74
1:A:72:ALA:HB1	1:A:101:ILE:HD13	1.68	0.74
1:A:220:ALA:N	1:B:145:VAL:HB	2.02	0.73
1:C:72:ALA:HB1	1:C:101:ILE:HD13	1.68	0.73
1:A:219:GLY:HA2	1:B:145:VAL:HG21	1.72	0.72
1:A:145:VAL:CB	1:A:146:PRO:HD3	2.21	0.70
1:B:145:VAL:CB	1:B:146:PRO:HD3	2.21	0.70
1:A:145:VAL:HG13	1:A:146:PRO:HD3	0.71	0.70
1:B:145:VAL:HG13	1:B:146:PRO:HD3	0.71	0.70
1:A:108:ILE:HG22	1:A:159:ILE:HB	1.74	0.70
1:C:145:VAL:CB	1:C:146:PRO:HD3	2.21	0.69
1:C:108:ILE:HG22	1:C:159:ILE:HB	1.74	0.69
1:C:145:VAL:HG13	1:C:146:PRO:HD3	0.70	0.69
1:B:108:ILE:HG22	1:B:159:ILE:HB	1.74	0.69
1:B:101:ILE:HD11	1:B:217:TYR:CE1	2.26	0.68
1:C:55:LEU:C	1:C:55:LEU:HD23	2.14	0.68
1:B:55:LEU:C	1:B:55:LEU:HD23	2.14	0.68
1:A:55:LEU:HD23	1:A:55:LEU:C	2.14	0.67
1:A:187:MET:CE	1:B:187:MET:CE	2.73	0.66
1:C:72:ALA:CA	1:C:101:ILE:HD11	2.27	0.65
1:A:72:ALA:CA	1:A:101:ILE:HD11	2.27	0.64
1:B:101:ILE:HG12	1:B:217:TYR:CE1	2.32	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:55:LEU:HD23	1:C:56:ALA:N	2.13	0.64
1:A:55:LEU:HD23	1:A:56:ALA:N	2.13	0.63
1:B:55:LEU:HD23	1:B:56:ALA:N	2.13	0.62
1:A:187:MET:CE	1:C:187:MET:CE	2.74	0.62
1:A:145:VAL:HB	1:C:220:ALA:H	1.64	0.62
1:A:143:ARG:C	1:A:144:ILE:HG12	2.21	0.61
1:B:143:ARG:C	1:B:144:ILE:HG12	2.21	0.61
1:B:125:ALA:CB	1:B:144:ILE:HD12	2.31	0.60
1:B:226:ALA:HB3	1:C:131:MET:O	2.01	0.60
1:A:220:ALA:H	1:B:145:VAL:CB	2.15	0.58
1:A:125:ALA:CB	1:A:144:ILE:HD12	2.31	0.58
1:C:125:ALA:CB	1:C:144:ILE:HD12	2.31	0.57
1:A:133:SER:OG	1:C:224:GLN:O	2.23	0.56
1:A:226:ALA:HB3	1:B:131:MET:O	2.04	0.56
1:C:108:ILE:HG21	1:C:193:PHE:CE1	2.41	0.56
1:A:108:ILE:HG21	1:A:193:PHE:CE1	2.41	0.55
1:B:108:ILE:HG21	1:B:193:PHE:CE1	2.41	0.55
1:B:112:THR:HG21	1:B:153:LYS:CB	2.36	0.55
1:B:220:ALA:O	1:C:145:VAL:CB	2.52	0.55
1:A:112:THR:HG21	1:A:153:LYS:CB	2.36	0.55
1:A:74:ILE:HG23	1:A:75:ASP:H	1.72	0.55
1:C:112:THR:HG21	1:C:153:LYS:CB	2.36	0.55
1:A:226:ALA:CB	1:B:131:MET:O	2.55	0.55
1:B:138:VAL:HA	1:B:141:VAL:HG12	1.88	0.55
1:B:101:ILE:CG1	1:B:217:TYR:CE1	2.90	0.55
1:A:138:VAL:HA	1:A:141:VAL:HG12	1.89	0.54
1:B:74:ILE:HG23	1:B:75:ASP:H	1.72	0.54
1:C:74:ILE:HG23	1:C:75:ASP:H	1.72	0.54
1:B:144:ILE:HG22	1:B:146:PRO:HD2	1.90	0.54
1:C:144:ILE:HG22	1:C:146:PRO:HD2	1.90	0.54
1:B:101:ILE:CG1	1:B:217:TYR:CD1	2.85	0.54
1:C:129:ASP:OD1	1:C:130:TYR:N	2.41	0.54
1:C:143:ARG:C	1:C:144:ILE:HG12	2.21	0.54
1:B:129:ASP:OD1	1:B:130:TYR:N	2.41	0.54
1:A:217:TYR:O	1:A:218:ARG:HG2	2.08	0.54
1:A:129:ASP:OD1	1:A:130:TYR:N	2.41	0.53
1:C:138:VAL:HA	1:C:141:VAL:HG12	1.88	0.53
1:A:69:PRO:HG2	1:A:76:PHE:CE2	2.43	0.53
1:A:133:SER:CB	1:C:224:GLN:O	2.57	0.53
1:A:74:ILE:O	1:A:173:ALA:HB2	2.09	0.53
1:C:217:TYR:O	1:C:218:ARG:HG2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:ALA:HA	1:A:101:ILE:HD11	1.90	0.53
1:A:144:ILE:HG22	1:A:146:PRO:HD2	1.90	0.53
1:C:72:ALA:HA	1:C:101:ILE:HD11	1.91	0.53
1:C:121:TYR:HA	1:C:151:ALA:CB	2.38	0.53
1:A:114:THR:HG21	1:A:152:LEU:CD2	2.39	0.53
1:C:144:ILE:HG22	1:C:145:VAL:N	2.24	0.53
1:B:69:PRO:HG2	1:B:76:PHE:CE2	2.43	0.53
1:C:74:ILE:O	1:C:173:ALA:HB2	2.09	0.53
1:C:112:THR:HG21	1:C:153:LYS:HB2	1.91	0.52
1:B:217:TYR:O	1:B:218:ARG:HG2	2.08	0.52
1:C:114:THR:HG21	1:C:152:LEU:CD2	2.39	0.52
1:B:114:THR:HG21	1:B:152:LEU:CD2	2.39	0.52
1:A:112:THR:HG21	1:A:153:LYS:HB2	1.91	0.52
1:B:220:ALA:N	1:C:145:VAL:HB	2.25	0.52
1:C:69:PRO:HG2	1:C:76:PHE:CE2	2.43	0.52
1:B:112:THR:HG21	1:B:153:LYS:HB2	1.91	0.52
1:A:144:ILE:HG22	1:A:145:VAL:N	2.24	0.52
1:B:144:ILE:HG22	1:B:145:VAL:N	2.24	0.52
1:B:74:ILE:O	1:B:173:ALA:HB2	2.09	0.52
1:A:109:THR:HG22	1:A:110:TRP:N	2.25	0.51
1:B:109:THR:HG22	1:B:110:TRP:N	2.25	0.51
1:A:121:TYR:HA	1:A:151:ALA:CB	2.38	0.51
1:C:109:THR:HG22	1:C:110:TRP:N	2.25	0.51
1:B:220:ALA:H	1:C:145:VAL:CG2	2.23	0.50
1:C:101:ILE:HG22	1:C:217:TYR:HD1	1.73	0.50
1:A:101:ILE:HG22	1:A:217:TYR:CE1	2.46	0.50
1:C:150:ILE:HG23	1:C:156:GLY:HA3	1.94	0.50
1:C:101:ILE:HG22	1:C:217:TYR:CE1	2.46	0.49
1:A:116:ASP:O	1:A:153:LYS:NZ	2.43	0.49
1:A:150:ILE:HG23	1:A:156:GLY:HA3	1.94	0.49
1:B:150:ILE:HG23	1:B:156:GLY:HA3	1.94	0.49
1:A:187:MET:HE3	1:B:187:MET:HE3	1.94	0.49
1:B:121:TYR:HA	1:B:151:ALA:CB	2.38	0.49
1:B:226:ALA:CB	1:C:131:MET:O	2.59	0.49
1:A:144:ILE:HB	1:A:147:SER:HB3	1.95	0.48
1:C:103:PHE:CE2	1:C:168:PHE:HD2	2.31	0.48
1:C:67:GLU:OE1	1:C:192:ARG:HD2	2.13	0.48
1:B:187:MET:HE3	1:C:187:MET:HE3	1.89	0.48
1:A:145:VAL:HB	1:C:220:ALA:N	2.28	0.48
1:B:220:ALA:H	1:C:145:VAL:HB	1.78	0.48
1:A:101:ILE:CD1	1:A:103:PHE:HE1	2.24	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:GLU:OE1	1:A:192:ARG:HD2	2.13	0.48
1:B:103:PHE:CE2	1:B:168:PHE:HD2	2.31	0.48
1:B:67:GLU:OE1	1:B:192:ARG:HD2	2.13	0.48
1:C:144:ILE:HB	1:C:147:SER:HB3	1.95	0.48
1:B:122:ILE:HB	1:B:151:ALA:HB2	1.95	0.48
1:B:144:ILE:HB	1:B:147:SER:HB3	1.95	0.48
1:C:122:ILE:HB	1:C:151:ALA:HB2	1.95	0.48
1:A:122:ILE:HB	1:A:151:ALA:HB2	1.95	0.48
1:B:154:ASN:O	1:B:155:ARG:C	2.53	0.47
1:A:220:ALA:H	1:B:145:VAL:CG2	2.27	0.47
1:A:154:ASN:O	1:A:155:ARG:C	2.53	0.47
1:A:103:PHE:CE2	1:A:168:PHE:HD2	2.31	0.47
1:C:151:ALA:CB	1:C:152:LEU:HA	2.42	0.47
1:B:150:ILE:HD12	1:B:157:PRO:HD2	1.97	0.47
1:A:151:ALA:HB1	1:A:152:LEU:HA	1.97	0.46
1:A:150:ILE:HD12	1:A:157:PRO:HD2	1.97	0.46
1:A:101:ILE:CG2	1:A:217:TYR:CE1	2.98	0.46
1:C:154:ASN:O	1:C:155:ARG:C	2.53	0.46
1:C:101:ILE:CG2	1:C:217:TYR:CE1	2.98	0.46
1:B:151:ALA:HB1	1:B:152:LEU:HA	1.97	0.46
1:A:68:ILE:HG13	1:A:193:PHE:HB3	1.98	0.46
1:C:150:ILE:HD12	1:C:157:PRO:HD2	1.97	0.46
1:C:101:ILE:CD1	1:C:103:PHE:HE1	2.24	0.46
1:C:68:ILE:HG13	1:C:193:PHE:HB3	1.98	0.46
1:B:126:PHE:CE2	1:B:161:MET:HG3	2.51	0.46
1:A:126:PHE:CE2	1:A:161:MET:HG3	2.51	0.46
1:C:126:PHE:CE2	1:C:161:MET:HG3	2.51	0.46
1:A:76:PHE:HB3	1:A:77:PRO:HA	1.99	0.45
1:C:151:ALA:HB1	1:C:152:LEU:HA	1.97	0.45
1:A:151:ALA:CB	1:A:152:LEU:HA	2.42	0.45
1:A:220:ALA:O	1:B:145:VAL:HB	2.15	0.45
1:B:76:PHE:HB3	1:B:77:PRO:HA	1.99	0.45
1:B:68:ILE:HG13	1:B:193:PHE:HB3	1.98	0.45
1:A:131:MET:HG2	1:C:171:ILE:HG22	1.98	0.45
1:C:101:ILE:CG2	1:C:217:TYR:CD1	2.93	0.45
1:C:76:PHE:HB3	1:C:77:PRO:HA	1.99	0.44
1:C:101:ILE:HD12	1:C:103:PHE:CE1	2.42	0.44
1:A:72:ALA:HA	1:A:101:ILE:CD1	2.48	0.44
1:B:116:ASP:O	1:B:153:LYS:NZ	2.43	0.44
1:B:187:MET:CE	1:C:187:MET:HE1	2.45	0.44
1:C:70:VAL:HG13	1:C:190:PRO:HD2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:101:ILE:CG2	1:A:217:TYR:CD1	2.93	0.44
1:A:55:LEU:CD2	1:A:56:ALA:N	2.81	0.44
1:B:108:ILE:HG12	1:B:209:ILE:HD11	2.00	0.44
1:C:151:ALA:HA	1:C:152:LEU:HA	1.74	0.44
1:C:116:ASP:O	1:C:153:LYS:NZ	2.43	0.43
1:C:108:ILE:HG12	1:C:209:ILE:HD11	2.00	0.43
1:A:101:ILE:HG22	1:A:217:TYR:HD1	1.73	0.43
1:A:108:ILE:HG12	1:A:209:ILE:HD11	2.00	0.43
1:C:55:LEU:CD2	1:C:56:ALA:N	2.81	0.43
1:A:101:ILE:HD12	1:A:103:PHE:CE1	2.42	0.43
1:A:54:LEU:HD23	1:A:54:LEU:HA	1.87	0.43
1:A:133:SER:O	1:C:223:LEU:HD23	2.18	0.43
1:B:54:LEU:HA	1:B:54:LEU:HD23	1.87	0.43
1:B:98:TRP:HB3	1:B:99:ASN:H	1.54	0.43
1:C:108:ILE:HD13	1:C:193:PHE:CE1	2.54	0.43
1:A:138:VAL:O	1:A:141:VAL:HG12	2.19	0.43
1:B:108:ILE:HD13	1:B:193:PHE:CE1	2.54	0.43
1:B:70:VAL:HG13	1:B:190:PRO:HD2	2.00	0.43
1:B:55:LEU:CD2	1:B:56:ALA:N	2.81	0.43
1:C:66:LEU:HG	1:C:85:ILE:HD11	2.01	0.43
1:A:114:THR:HG21	1:A:152:LEU:HD23	2.01	0.42
1:B:138:VAL:O	1:B:141:VAL:HG12	2.19	0.42
1:A:108:ILE:HD13	1:A:193:PHE:CE1	2.54	0.42
1:A:144:ILE:HG22	1:A:146:PRO:CD	2.49	0.42
1:A:70:VAL:HG13	1:A:190:PRO:HD2	2.00	0.42
1:B:114:THR:HG21	1:B:152:LEU:HD23	2.01	0.42
1:B:74:ILE:HG23	1:B:75:ASP:N	2.34	0.42
1:B:66:LEU:HG	1:B:85:ILE:HD11	2.01	0.42
1:A:135:PRO:HD3	1:C:223:LEU:CD2	2.49	0.42
1:C:114:THR:HG21	1:C:152:LEU:HD23	2.01	0.42
1:A:66:LEU:HG	1:A:85:ILE:HD11	2.01	0.42
1:C:138:VAL:O	1:C:141:VAL:HG12	2.19	0.42
1:B:95:SER:HG	1:B:101:ILE:HD11	1.84	0.42
1:B:122:ILE:HG12	1:B:197:ILE:HG22	2.02	0.42
1:A:150:ILE:CG2	1:A:156:GLY:HA3	2.50	0.41
1:B:219:GLY:HA2	1:C:145:VAL:HG21	2.01	0.41
1:B:187:MET:HE1	1:C:187:MET:HE1	2.02	0.41
1:C:54:LEU:HA	1:C:54:LEU:HD23	1.87	0.41
1:C:122:ILE:HG12	1:C:197:ILE:HG22	2.02	0.41
1:C:72:ALA:CA	1:C:101:ILE:CD1	2.96	0.41
1:B:144:ILE:HG22	1:B:146:PRO:CD	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:101:ILE:HD13	1:B:217:TYR:CE1	2.50	0.41
1:C:74:ILE:HG23	1:C:75:ASP:N	2.35	0.41
1:A:102:VAL:HG12	1:A:216:SER:O	2.21	0.41
1:A:122:ILE:HG12	1:A:197:ILE:HG22	2.02	0.41
1:B:151:ALA:CB	1:B:152:LEU:HA	2.42	0.41
1:B:222:ILE:C	1:B:223:LEU:HD12	2.41	0.41
1:B:108:ILE:HG21	1:B:193:PHE:HE1	1.85	0.41
1:A:118:THR:O	1:A:153:LYS:NZ	2.46	0.41
1:C:102:VAL:HG12	1:C:216:SER:O	2.21	0.41
1:C:68:ILE:HG13	1:C:68:ILE:O	2.21	0.41
1:A:102:VAL:HG13	1:A:102:VAL:O	2.21	0.41
1:A:121:TYR:CE1	1:A:198:PRO:HG3	2.56	0.41
1:A:68:ILE:HG13	1:A:68:ILE:O	2.21	0.41
1:A:74:ILE:HG23	1:A:75:ASP:N	2.34	0.41
1:C:222:ILE:C	1:C:223:LEU:HD12	2.41	0.41
1:C:72:ALA:HA	1:C:101:ILE:CD1	2.48	0.41
1:B:102:VAL:HG13	1:B:102:VAL:O	2.21	0.41
1:C:144:ILE:HG22	1:C:146:PRO:CD	2.50	0.41
1:B:102:VAL:HG12	1:B:216:SER:O	2.20	0.41
1:B:150:ILE:CG2	1:B:156:GLY:HA3	2.50	0.41
1:B:64:GLN:HB2	1:B:197:ILE:HG12	2.03	0.41
1:A:72:ALA:CA	1:A:101:ILE:CD1	2.96	0.40
1:A:66:LEU:O	1:A:194:ILE:HG13	2.21	0.40
1:B:66:LEU:O	1:B:194:ILE:HG13	2.21	0.40
1:B:68:ILE:HG13	1:B:68:ILE:O	2.21	0.40
1:C:102:VAL:O	1:C:102:VAL:HG13	2.21	0.40
1:B:121:TYR:CE1	1:B:198:PRO:HG3	2.56	0.40
1:B:60:THR:HB	1:B:64:GLN:HE21	1.86	0.40
1:C:55:LEU:CG	1:C:56:ALA:N	2.84	0.40
1:C:98:TRP:HB3	1:C:99:ASN:H	1.54	0.40
1:C:118:THR:O	1:C:153:LYS:NZ	2.45	0.40
1:C:150:ILE:CG2	1:C:156:GLY:HA3	2.50	0.40
1:C:168:PHE:CE2	1:C:190:PRO:HG3	2.56	0.40
1:C:121:TYR:CE1	1:C:198:PRO:HG3	2.56	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	183/185 (99%)	161 (88%)	18 (10%)	4 (2%)	8	51
1	B	180/185 (97%)	158 (88%)	18 (10%)	4 (2%)	8	51
1	C	181/185 (98%)	159 (88%)	18 (10%)	4 (2%)	8	51
All	All	544/555 (98%)	478 (88%)	54 (10%)	12 (2%)	13	51

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	155	ARG
1	B	155	ARG
1	C	155	ARG
1	A	74	ILE
1	A	146	PRO
1	B	74	ILE
1	B	146	PRO
1	C	74	ILE
1	C	146	PRO
1	A	77	PRO
1	B	77	PRO
1	C	77	PRO

5.3.2 Protein sidechains [i](#)

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

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

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	146/146 (100%)	139 (95%)	7 (5%)	31 72
1	B	143/146 (98%)	136 (95%)	7 (5%)	31 71
1	C	144/146 (99%)	137 (95%)	7 (5%)	31 71
All	All	433/438 (99%)	412 (95%)	21 (5%)	36 72

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	MET
1	A	55	LEU
1	A	103	PHE
1	A	143	ARG
1	A	144	ILE
1	A	152	LEU
1	A	153	LYS
1	B	50	MET
1	B	55	LEU
1	B	103	PHE
1	B	143	ARG
1	B	144	ILE
1	B	152	LEU
1	B	153	LYS
1	C	50	MET
1	C	55	LEU
1	C	103	PHE
1	C	143	ARG
1	C	144	ILE
1	C	152	LEU
1	C	153	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.