Full wwPDB EM Validation Report ⓘ

May 4, 2024 – 10:53 am BST

PDB ID	:	5L7Q
EMDB ID	:	EMD-4009
Title	:	Structure of deformed wing virus, a honeybee pathogen
Authors	:	Skubnik, K.; Novacek, J.; Fuzik, T.; Pridal, A.; Paxton, R.; Plevka, P.
Deposited on	:	2016-06-03
Resolution	:	3.50 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:

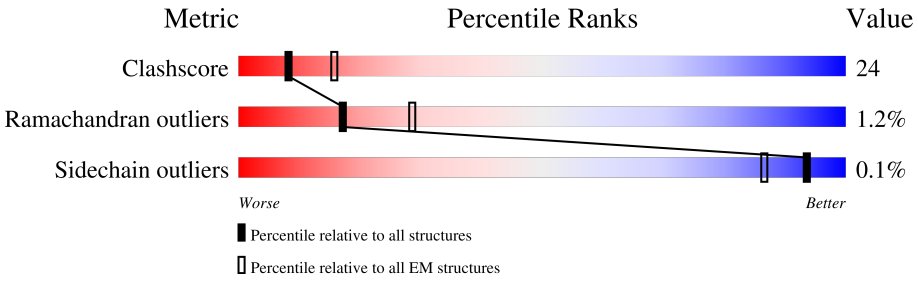
EMDB validation analysis	:	0.0.1.dev92
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	FAILED
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 3.50 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	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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 <=5%

Mol	Chain	Length	Quality of chain
1	A	258	50% 48% .
2	B	253	53% 45% ..
3	C	416	45% 50% 5%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7091 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 VP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	252	1976	1256	336	374	10	0	0

- Molecule 2 is a protein called vp2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	250	1973	1253	335	378	7	0	0

- Molecule 3 is a protein called vp3.

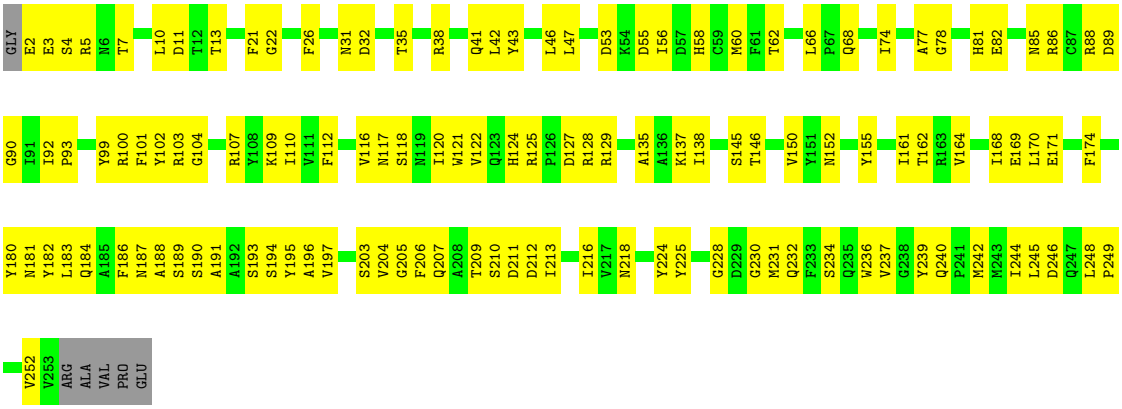
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	C	397	3142	2015	541	574	12	0	0

3 Residue-property plots (i)

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.

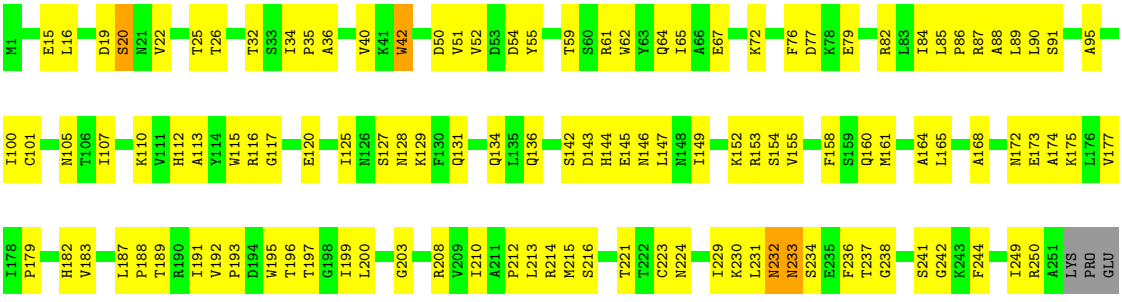
- Molecule 1: VP1

Chain A:



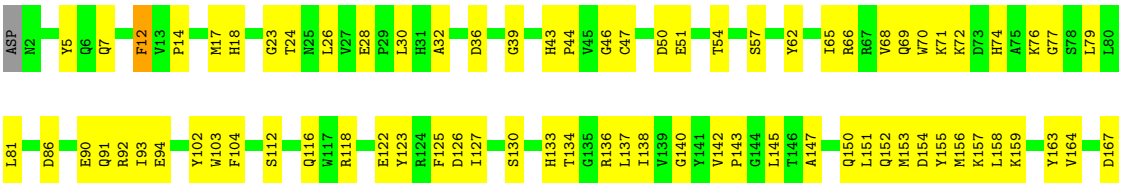
- Molecule 2: vp2

Chain B:



- Molecule 3: vp3

Chain C:




L168 S173 Y180 R184 P185 W186 W187 V188 R189 K190 Y191 G192 G193 N194 Y195 L196 P197 T200 P203 S204 T205 L206 F207 M208 Y209 V210 Q211 V212 P213 L214 I215 P216 M217 E218 A219 V220 S221 D222 T223 I224 D225 V228 Y229 V230 R231 E237 V238 C239 V240 P241 V242 Q243
L248 D253 F254 I255 L256 E261 Y262 R263 A264 K265 Y268 A269 P270 Y271 Y272 V275 W276 H277 S278 F279 N280 N281 S282 N283 S284 L285 R288 W289 A292 S293 D294 Q295 I296 A297 Q298 W299 P300 T301 I302 S303 V304 E308 L309 L312 R313 I314 K315 D316 G317 K318 Q319
V322 G323 T324 Q325 P326 W327 R328 T329 M330 V331 V332 W333 P334 Y339 N340 Y346 N347 A348 E349 R350 A351 R352 A355 Q356 H357 L358 G362 S363 L364 T365 D366 K370 F373 V374 P375 Q378 Q379 G380 P381 G382 K383 V384 S385 N386 P389 V390 W391 E392 R395 A396 P397
L398 ALA THR GLN ARG ALA HIS ILE GLN ASP PHE GLU PHE ILE GLU ALA ILE PRO GLU

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	136828	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 ($e^-/Å^2$)	15	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	4000	Depositor
Magnification	74235	Depositor
Image detector	FEI FALCON II (4k x 4k)	Depositor

5 Model quality

5.1 Standard geometry

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	0.61	0/2024	0.59	0/2754
2	B	0.61	0/2024	0.62	0/2763
3	C	0.59	0/3238	0.56	0/4418
All	All	0.60	0/7286	0.58	0/9935

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

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	1976	0	1911	113	0
2	B	1973	0	1928	96	0
3	C	3142	0	3051	172	0
All	All	7091	0	6890	342	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 24.

All (342) 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:86:ARG:HD2	3:C:194:ASN:HD22	1.36	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:150:GLN:HE22	3:C:205:THR:HG21	1.37	0.88
2:B:84:ILE:HG21	2:B:87:ARG:HD3	1.55	0.86
2:B:42:TRP:HE1	3:C:46:GLY:HA3	1.44	0.82
1:A:187:ASN:HB3	2:B:147:LEU:HB3	1.62	0.81
3:C:151:LEU:HD12	3:C:152:GLN:HG3	1.61	0.80
3:C:217:MET:HG2	3:C:219:ALA:H	1.46	0.79
3:C:319:GLN:HG2	3:C:386:ASN:HD22	1.47	0.79
3:C:189:ARG:HH11	3:C:192:GLY:HA2	1.48	0.78
3:C:136:ARG:NE	3:C:167:ASP:OD1	2.17	0.77
1:A:240:GLN:HG2	2:B:152:LYS:HB2	1.66	0.77
1:A:196:ALA:HB1	2:B:183:VAL:HG11	1.65	0.77
1:A:171:GLU:HB3	3:C:32:ALA:HA	1.68	0.76
2:B:191:ILE:HG23	2:B:196:THR:HG22	1.68	0.76
1:A:101:PHE:HB2	1:A:234:SER:HB2	1.68	0.75
1:A:249:PRO:HB2	3:C:256:LEU:HD21	1.69	0.74
1:A:104:GLY:HA3	1:A:231:MET:HA	1.69	0.74
3:C:136:ARG:HB3	3:C:212:VAL:HB	1.69	0.74
3:C:122:GLU:OE1	3:C:231:ARG:NH2	2.21	0.73
3:C:112:SER:O	3:C:189:ARG:NH2	2.21	0.73
3:C:347:ASN:OD1	3:C:350:ARG:N	2.18	0.72
2:B:177:VAL:HG12	2:B:179:PRO:HD3	1.72	0.72
3:C:71:LYS:HG2	3:C:223:THR:HG22	1.71	0.72
2:B:22:VAL:HG11	2:B:62:TRP:HB2	1.72	0.71
3:C:137:LEU:HA	3:C:211:GLN:H	1.55	0.71
2:B:82:ARG:HD3	2:B:153:ARG:HD3	1.73	0.71
1:A:55:ASP:H	1:A:58:HIS:CE1	2.09	0.71
3:C:304:VAL:HA	3:C:395:ARG:HH12	1.55	0.71
1:A:5:ARG:NH1	1:A:7:THR:OG1	2.24	0.70
2:B:232:ASN:O	2:B:234:SER:N	2.25	0.69
1:A:116:VAL:HG22	1:A:118:SER:H	1.57	0.69
2:B:15:GLU:OE2	2:B:26:THR:OG1	2.11	0.69
1:A:213:ILE:HD12	1:A:216:ILE:HD12	1.75	0.68
1:A:209:THR:OG1	1:A:211:ASP:OD1	2.08	0.68
2:B:120:GLU:HB2	2:B:230:LYS:HB3	1.76	0.68
3:C:312:LEU:HB3	3:C:391:TRP:CE3	2.30	0.67
3:C:189:ARG:HG2	3:C:191:TYR:H	1.58	0.67
2:B:87:ARG:NH2	2:B:249:ILE:HB	2.10	0.67
1:A:107:ARG:NH1	3:C:28:GLU:OE1	2.28	0.67
1:A:35:THR:OG1	3:C:23:GLY:O	2.12	0.66
3:C:214:LEU:HG	3:C:216:PRO:HD3	1.77	0.66
1:A:89:ASP:OD2	3:C:248:LEU:N	2.16	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:149:ILE:HD13	2:B:153:ARG:HH21	1.61	0.66
1:A:125:ARG:HH11	1:A:152:ASN:HD22	1.43	0.66
2:B:160:GLN:NE2	3:C:104:PHE:O	2.29	0.65
3:C:66:ARG:NH2	3:C:90:GLU:O	2.29	0.65
1:A:77:ALA:HA	3:C:255:ILE:HD13	1.79	0.64
2:B:40:VAL:O	2:B:116:ARG:NH1	2.31	0.64
2:B:87:ARG:HH22	2:B:249:ILE:HB	1.61	0.64
2:B:143:ASP:OD1	2:B:144:HIS:N	2.31	0.64
3:C:116:GLN:HB2	3:C:239:CYS:SG	2.38	0.63
3:C:193:GLY:O	3:C:195:TYR:N	2.31	0.63
2:B:42:TRP:NE1	3:C:46:GLY:HA3	2.14	0.63
1:A:100:ARG:NH2	2:B:145:GLU:HB2	2.13	0.63
3:C:276:TRP:NE1	3:C:294:ASP:OD1	2.25	0.62
1:A:246:ASP:HA	3:C:194:ASN:HA	1.81	0.62
1:A:100:ARG:NH2	2:B:142:SER:O	2.28	0.62
3:C:288:ARG:NH2	3:C:292:ALA:O	2.33	0.62
2:B:77:ASP:OD2	2:B:208:ARG:NH2	2.30	0.61
3:C:309:LEU:HB2	3:C:355:ALA:HB1	1.81	0.61
2:B:128:ASN:OD1	2:B:129:LYS:N	2.32	0.61
1:A:187:ASN:ND2	2:B:144:HIS:O	2.34	0.61
2:B:50:ASP:OD1	2:B:51:VAL:N	2.33	0.61
3:C:187:TRP:CE2	3:C:203:PRO:HG3	2.35	0.61
3:C:288:ARG:HG2	3:C:296:ILE:HD13	1.81	0.61
2:B:134:GLN:HE21	2:B:165:LEU:HD22	1.65	0.61
2:B:79:GLU:HA	2:B:208:ARG:HA	1.83	0.60
2:B:19:ASP:OD2	2:B:62:TRP:NE1	2.34	0.60
1:A:10:LEU:HD13	3:C:164:VAL:HG11	1.83	0.60
3:C:370:LYS:HA	3:C:373:PHE:CD2	2.37	0.59
1:A:3:GLU:OE2	3:C:62:TYR:N	2.36	0.59
3:C:92:ARG:HH12	3:C:94:GLU:HA	1.67	0.59
2:B:32:THR:HG22	2:B:175:LYS:HB3	1.84	0.59
2:B:72:LYS:HA	2:B:213:LEU:HB2	1.83	0.59
3:C:147:ALA:HA	3:C:150:GLN:HB2	1.84	0.58
2:B:34:ILE:HG12	2:B:177:VAL:HB	1.85	0.58
1:A:2:GLU:OE1	1:A:4:SER:N	2.37	0.58
3:C:187:TRP:CD2	3:C:203:PRO:HG3	2.39	0.58
1:A:41:GLN:NE2	1:A:90:GLY:HA3	2.20	0.57
3:C:195:TYR:CZ	3:C:197:PRO:HG2	2.39	0.57
3:C:327:TRP:CH2	3:C:374:VAL:HG11	2.39	0.57
1:A:100:ARG:NE	2:B:142:SER:OG	2.34	0.57
1:A:182:TYR:CZ	2:B:145:GLU:HG3	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:95:ALA:HA	2:B:250:ARG:HH22	1.69	0.57
1:A:125:ARG:HH11	1:A:152:ASN:ND2	2.02	0.57
1:A:122:VAL:HG22	1:A:204:VAL:HG22	1.85	0.57
3:C:280:ASN:ND2	3:C:284:SER:OG	2.37	0.57
1:A:125:ARG:NH2	1:A:203:SER:OG	2.38	0.56
3:C:123:TYR:CE1	3:C:230:VAL:HG22	2.39	0.56
1:A:46:LEU:HD23	1:A:138:ILE:HD13	1.85	0.56
3:C:136:ARG:HH11	3:C:212:VAL:HG21	1.70	0.56
3:C:308:GLU:HG2	3:C:397:PRO:HA	1.87	0.56
1:A:86:ARG:HD2	3:C:194:ASN:ND2	2.14	0.56
1:A:60:MET:N	1:A:204:VAL:O	2.38	0.56
3:C:150:GLN:NE2	3:C:205:THR:HG21	2.14	0.56
2:B:79:GLU:HB3	2:B:208:ARG:HH11	1.71	0.56
2:B:182:HIS:CD2	2:B:183:VAL:H	2.24	0.56
3:C:155:TYR:O	3:C:159:LYS:HG3	2.06	0.56
2:B:64:GLN:HE21	2:B:67:GLU:HB3	1.71	0.56
1:A:68:GLN:HE21	1:A:183:LEU:HA	1.71	0.56
2:B:88:ALA:HA	2:B:91:SER:HB3	1.86	0.56
1:A:56:ILE:HD12	1:A:207:GLN:O	2.06	0.55
2:B:20:SER:HB2	2:B:61:ARG:HG2	1.89	0.55
3:C:261:GLU:HB2	3:C:390:VAL:HG13	1.88	0.55
1:A:107:ARG:NE	1:A:228:GLY:HA2	2.20	0.55
2:B:215:MET:N	3:C:221:SER:HB3	2.20	0.55
1:A:82:GLU:O	1:A:86:ARG:N	2.40	0.55
1:A:85:ASN:HD21	3:C:255:ILE:HB	1.72	0.55
3:C:272:TYR:HB3	3:C:384:VAL:HA	1.89	0.55
1:A:85:ASN:ND2	3:C:255:ILE:HB	2.22	0.55
3:C:299:TRP:O	3:C:340:ASN:HB2	2.06	0.54
3:C:239:CYS:SG	3:C:240:VAL:N	2.79	0.54
1:A:78:GLY:O	3:C:263:ARG:NE	2.36	0.54
3:C:44:PRO:HD2	3:C:47:CYS:HB2	1.89	0.54
3:C:189:ARG:NH2	3:C:242:VAL:HG23	2.23	0.54
3:C:314:ILE:HD11	3:C:389:PRO:HB2	1.90	0.54
3:C:309:LEU:HD11	3:C:396:ALA:HB3	1.89	0.54
1:A:42:LEU:HA	1:A:224:TYR:HD1	1.73	0.54
3:C:71:LYS:H	3:C:74:HIS:CE1	2.25	0.54
2:B:36:ALA:HB3	2:B:179:PRO:HG2	1.90	0.53
1:A:239:TYR:N	2:B:160:GLN:OE1	2.41	0.53
3:C:271:TYR:O	3:C:385:SER:N	2.28	0.53
3:C:190:LYS:HA	3:C:200:THR:O	2.08	0.53
3:C:301:THR:OG1	3:C:340:ASN:ND2	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:314:ILE:HG12	3:C:316:ASP:H	1.73	0.53
2:B:223:CYS:SG	2:B:224:ASN:N	2.82	0.53
1:A:102:TYR:CE2	1:A:181:ASN:HB2	2.44	0.53
1:A:211:ASP:OD1	1:A:212:ASP:N	2.41	0.53
2:B:117:GLY:HA3	2:B:234:SER:HA	1.90	0.53
2:B:152:LYS:O	2:B:154:SER:N	2.42	0.53
3:C:50:ASP:OD1	3:C:51:GLU:N	2.42	0.53
2:B:77:ASP:CG	2:B:208:ARG:HH21	2.12	0.52
3:C:91:GLN:HB3	3:C:103:TRP:HB2	1.91	0.52
3:C:133:HIS:CG	3:C:224:ILE:HD11	2.44	0.52
1:A:129:ARG:NH2	1:A:195:TYR:OH	2.43	0.52
3:C:275:VAL:HB	3:C:282:SER:HB3	1.91	0.52
3:C:127:ILE:HG21	3:C:168:LEU:HD22	1.90	0.52
3:C:322:VAL:HB	3:C:383:LYS:HD3	1.89	0.52
2:B:188:PRO:HB3	2:B:192:VAL:HG21	1.92	0.52
3:C:133:HIS:CD2	3:C:224:ILE:HD11	2.44	0.52
3:C:333:TRP:CD2	3:C:334:PRO:HD2	2.44	0.52
1:A:190:SER:HB2	1:A:193:SER:HB2	1.91	0.52
3:C:138:ILE:HG13	3:C:211:GLN:HB2	1.91	0.52
3:C:261:GLU:HB2	3:C:390:VAL:CG1	2.39	0.51
1:A:32:ASP:O	1:A:35:THR:HG22	2.10	0.51
3:C:276:TRP:CG	3:C:296:ILE:HD11	2.45	0.51
3:C:36:ASP:OD2	3:C:39:GLY:N	2.43	0.51
3:C:268:TYR:HD1	3:C:289:TRP:CD1	2.29	0.51
3:C:352:ARG:O	3:C:356:GLN:HG3	2.11	0.51
3:C:363:SER:N	3:C:366:ASP:OD2	2.41	0.51
1:A:11:ASP:OD1	1:A:13:THR:OG1	2.22	0.51
2:B:61:ARG:HB2	2:B:105:ASN:OD1	2.09	0.51
1:A:21:PHE:CD2	1:A:31:ASN:HB2	2.46	0.51
1:A:150:VAL:O	1:A:152:ASN:ND2	2.44	0.51
1:A:103:ARG:O	1:A:232:GLN:N	2.42	0.51
2:B:192:VAL:HG13	2:B:193:PRO:HD2	1.93	0.51
1:A:35:THR:HA	1:A:38:ARG:HE	1.77	0.50
1:A:43:TYR:HB2	1:A:225:TYR:CD2	2.45	0.50
3:C:325:GLN:HE22	3:C:346:TYR:HA	1.75	0.50
1:A:125:ARG:HH22	1:A:203:SER:CB	2.24	0.50
3:C:302:ILE:HD11	3:C:312:LEU:HD21	1.93	0.50
3:C:349:GLU:OE1	3:C:352:ARG:NH1	2.43	0.50
1:A:104:GLY:O	1:A:174:PHE:HB2	2.10	0.50
3:C:12:PHE:HE1	3:C:14:PRO:HG3	1.76	0.50
1:A:100:ARG:NH1	1:A:184:GLN:HB2	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:312:LEU:HD12	3:C:391:TRP:HB3	1.92	0.50
3:C:118:ARG:NH1	3:C:237:GLU:OE2	2.44	0.50
3:C:308:GLU:HA	3:C:398:LEU:H	1.77	0.50
1:A:145:SER:OG	1:A:146:THR:N	2.45	0.50
1:A:174:PHE:HE1	1:A:181:ASN:ND2	2.10	0.49
2:B:212:PRO:HD2	3:C:225:ASP:OD2	2.12	0.49
3:C:72:LYS:HA	3:C:214:LEU:HB3	1.94	0.49
3:C:270:PRO:HB3	3:C:385:SER:O	2.11	0.49
3:C:180:TYR:OH	3:C:186:TRP:O	2.30	0.49
2:B:16:LEU:HA	2:B:25:THR:HG22	1.94	0.49
3:C:5:TYR:CZ	3:C:7:GLN:HB2	2.48	0.49
3:C:265:LYS:HE3	3:C:299:TRP:C	2.32	0.49
3:C:325:GLN:NE2	3:C:346:TYR:HA	2.28	0.49
2:B:65:ILE:HG22	2:B:89:LEU:HA	1.93	0.49
3:C:69:GLN:NE2	3:C:223:THR:HB	2.26	0.49
3:C:312:LEU:N	3:C:330:MET:O	2.46	0.49
1:A:171:GLU:HB2	3:C:30:LEU:HD13	1.94	0.49
3:C:292:ALA:N	3:C:295:GLN:OE1	2.45	0.49
1:A:60:MET:H	1:A:205:GLY:HA2	1.76	0.48
2:B:55:TYR:HB2	2:B:59:THR:HG23	1.95	0.48
2:B:86:PRO:HB2	2:B:200:LEU:HG	1.94	0.48
1:A:244:ILE:HD11	3:C:90:GLU:OE2	2.14	0.48
3:C:155:TYR:HB3	3:C:209:TYR:CG	2.48	0.48
1:A:102:TYR:OH	1:A:197:VAL:O	2.27	0.48
1:A:56:ILE:HD11	1:A:209:THR:HG22	1.96	0.48
2:B:214:ARG:NH2	3:C:69:GLN:OE1	2.47	0.48
3:C:70:TRP:HD1	3:C:74:HIS:HD1	1.61	0.48
1:A:100:ARG:HH22	2:B:145:GLU:HB2	1.78	0.48
2:B:113:ALA:N	2:B:238:GLY:O	2.42	0.48
3:C:155:TYR:O	3:C:158:LEU:HB3	2.13	0.48
1:A:42:LEU:HA	1:A:224:TYR:CD1	2.48	0.47
2:B:172:ASN:OD1	2:B:173:GLU:N	2.47	0.47
3:C:284:SER:HB3	3:C:373:PHE:HB3	1.96	0.47
2:B:232:ASN:CG	2:B:233:ASN:H	2.17	0.47
3:C:375:PRO:HG2	3:C:378:GLN:HG2	1.95	0.47
3:C:298:GLN:HG2	3:C:339:TYR:CD1	2.49	0.47
2:B:127:SER:OG	2:B:168:ALA:O	2.14	0.47
3:C:154:ASP:OD1	3:C:157:LYS:N	2.24	0.47
3:C:195:TYR:CE2	3:C:197:PRO:HG2	2.50	0.47
1:A:135:ALA:HA	1:A:137:LYS:HE3	1.97	0.47
1:A:228:GLY:O	1:A:230:GLY:N	2.40	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:85:LEU:HB2	2:B:86:PRO:HD3	1.97	0.47
3:C:357:HIS:CE1	3:C:362:GLY:HA3	2.50	0.47
2:B:113:ALA:O	2:B:189:THR:N	2.36	0.46
2:B:136:GLN:NE2	2:B:158:PHE:HB2	2.30	0.46
3:C:92:ARG:HA	3:C:102:TYR:CD1	2.49	0.46
3:C:333:TRP:CZ3	3:C:358:LEU:HB3	2.50	0.46
2:B:115:TRP:HZ2	2:B:231:LEU:HD11	1.79	0.46
1:A:88:ARG:NH2	3:C:253:ASP:O	2.48	0.46
3:C:217:MET:CE	3:C:219:ALA:HB3	2.45	0.46
2:B:101:CYS:HA	2:B:107:ILE:HG22	1.97	0.46
1:A:2:GLU:O	1:A:5:ARG:HG2	2.15	0.46
1:A:47:LEU:HB3	1:A:218:ASN:HA	1.98	0.46
2:B:196:THR:OG1	2:B:197:THR:N	2.49	0.46
2:B:52:VAL:HG22	2:B:237:THR:HB	1.97	0.46
1:A:32:ASP:OD1	1:A:32:ASP:N	2.49	0.46
1:A:210:SER:O	1:A:213:ILE:HG22	2.15	0.46
1:A:99:TYR:CD1	1:A:236:TRP:HA	2.51	0.46
3:C:125:PHE:O	3:C:173:SER:OG	2.29	0.46
1:A:66:LEU:HD13	1:A:68:GLN:HB2	1.98	0.46
1:A:103:ARG:HB2	1:A:180:TYR:HB3	1.98	0.46
1:A:124:HIS:O	1:A:155:TYR:HB3	2.16	0.46
1:A:240:GLN:O	1:A:242:MET:HE3	2.16	0.45
1:A:109:LYS:HA	1:A:168:ILE:O	2.16	0.45
3:C:86:ASP:OD2	3:C:190:LYS:HD2	2.16	0.45
1:A:125:ARG:HE	1:A:152:ASN:ND2	2.15	0.45
3:C:357:HIS:HE2	3:C:366:ASP:HB2	1.82	0.45
2:B:241:SER:HB3	2:B:244:PHE:CE2	2.51	0.45
3:C:65:ILE:HG22	3:C:228:VAL:O	2.17	0.45
2:B:34:ILE:HA	2:B:35:PRO:HD2	1.84	0.45
3:C:92:ARG:NH1	3:C:93:ILE:O	2.50	0.45
1:A:62:THR:HG22	1:A:203:SER:CB	2.46	0.45
1:A:55:ASP:HB3	1:A:58:HIS:NE2	2.31	0.45
3:C:156:MET:SD	3:C:159:LYS:HD3	2.57	0.45
2:B:112:HIS:CG	2:B:236:PHE:HB3	2.51	0.45
3:C:346:TYR:OH	3:C:375:PRO:HD2	2.17	0.44
1:A:121:TRP:CZ2	1:A:205:GLY:HA3	2.53	0.44
2:B:143:ASP:HB2	2:B:149:ILE:CD1	2.47	0.44
3:C:357:HIS:NE2	3:C:366:ASP:HB2	2.32	0.44
2:B:110:LYS:O	2:B:242:GLY:HA2	2.17	0.44
2:B:143:ASP:H	2:B:203:GLY:HA2	1.82	0.44
2:B:215:MET:HG3	2:B:216:SER:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:36:ALA:O	2:B:179:PRO:HG3	2.17	0.44
3:C:184:ARG:HD2	3:C:187:TRP:HA	1.99	0.44
2:B:54:ASP:HB3	2:B:234:SER:O	2.17	0.44
2:B:164:ALA:HB1	2:B:174:ALA:HB1	1.99	0.44
2:B:241:SER:HB3	2:B:244:PHE:CD2	2.53	0.44
1:A:2:GLU:CD	1:A:4:SER:H	2.20	0.44
1:A:88:ARG:HB3	1:A:245:LEU:HD21	2.00	0.44
1:A:186:PHE:CZ	1:A:188:ALA:HA	2.53	0.44
1:A:191:ALA:O	1:A:194:SER:OG	2.27	0.44
2:B:187:LEU:HA	2:B:188:PRO:HD2	1.81	0.44
3:C:77:GLY:HA2	3:C:155:TYR:CE2	2.53	0.44
3:C:140:GLY:HA3	3:C:158:LEU:HD11	2.00	0.44
1:A:127:ASP:HB2	1:A:128:ARG:NH1	2.32	0.44
2:B:115:TRP:HE1	2:B:234:SER:HB2	1.83	0.44
2:B:214:ARG:NH1	3:C:223:THR:OG1	2.51	0.44
3:C:126:ASP:OD1	3:C:173:SER:OG	2.36	0.44
3:C:145:LEU:HD12	3:C:205:THR:OG1	2.17	0.43
3:C:285:LEU:HD21	3:C:379:GLN:HB2	1.99	0.43
1:A:74:ILE:O	1:A:81:HIS:N	2.34	0.43
1:A:125:ARG:HG2	1:A:155:TYR:CD2	2.52	0.43
2:B:50:ASP:O	2:B:52:VAL:HG23	2.19	0.43
3:C:12:PHE:CD1	3:C:12:PHE:C	2.92	0.43
3:C:65:ILE:HG21	3:C:230:VAL:HG23	2.00	0.43
3:C:322:VAL:HG22	3:C:326:PRO:HA	2.01	0.43
3:C:347:ASN:OD1	3:C:350:ARG:HG2	2.18	0.43
3:C:130:SER:OG	3:C:133:HIS:ND1	2.40	0.43
3:C:277:HIS:O	3:C:281:ASN:HA	2.18	0.43
1:A:82:GLU:HA	1:A:85:ASN:HB3	2.00	0.43
2:B:215:MET:SD	2:B:221:THR:HA	2.59	0.43
3:C:134:THR:O	3:C:214:LEU:HD12	2.19	0.43
1:A:129:ARG:HD3	1:A:194:SER:OG	2.19	0.43
1:A:237:VAL:HG13	2:B:161:MET:HE1	1.99	0.43
2:B:72:LYS:HE2	2:B:221:THR:HG23	2.01	0.43
2:B:131:GLN:HB3	2:B:213:LEU:HD22	2.01	0.43
1:A:117:ASN:O	1:A:162:THR:HB	2.19	0.43
1:A:125:ARG:HD3	1:A:128:ARG:NH2	2.33	0.43
1:A:152:ASN:HB2	1:A:155:TYR:HD2	1.82	0.43
3:C:319:GLN:HB3	3:C:389:PRO:HB3	2.01	0.42
3:C:333:TRP:HD1	3:C:364:LEU:HD11	1.83	0.42
2:B:125:ILE:HG12	2:B:127:SER:H	1.83	0.42
3:C:280:ASN:ND2	3:C:283:ASN:HB2	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:317:GLY:HA2	3:C:328:ARG:HG2	2.01	0.42
1:A:187:ASN:OD1	2:B:145:GLU:HA	2.19	0.42
3:C:159:LYS:HA	3:C:163:TYR:CD2	2.54	0.42
1:A:135:ALA:HA	1:A:137:LYS:HG2	2.01	0.42
1:A:248:LEU:HA	1:A:249:PRO:HD3	1.82	0.42
3:C:217:MET:HE2	3:C:219:ALA:HB3	2.02	0.42
3:C:265:LYS:HG3	3:C:300:PRO:HD3	2.01	0.42
3:C:380:GLY:HA3	3:C:381:PRO:HD2	1.83	0.42
1:A:122:VAL:HG11	1:A:170:LEU:HD11	2.02	0.42
2:B:77:ASP:OD1	2:B:155:VAL:HG13	2.19	0.42
2:B:115:TRP:CG	2:B:116:ARG:N	2.88	0.42
3:C:17:MET:HG2	3:C:18:HIS:O	2.18	0.42
3:C:276:TRP:CH2	3:C:278:SER:HB2	2.54	0.42
3:C:332:VAL:CG1	3:C:340:ASN:HB3	2.49	0.42
1:A:240:GLN:NE2	2:B:146:ASN:HD21	2.18	0.42
2:B:85:LEU:HA	2:B:85:LEU:HD23	1.79	0.42
3:C:289:TRP:HD1	3:C:297:ALA:HA	1.85	0.42
1:A:99:TYR:HD1	1:A:236:TRP:HA	1.85	0.42
3:C:240:VAL:HA	3:C:241:PRO:HD3	1.69	0.42
3:C:358:LEU:HD23	3:C:362:GLY:O	2.20	0.42
1:A:120:ILE:HG12	1:A:206:PHE:CD1	2.55	0.41
2:B:90:LEU:HD22	2:B:100:ILE:HD12	2.01	0.41
3:C:288:ARG:HH22	3:C:293:SER:HA	1.85	0.41
1:A:62:THR:HG22	1:A:203:SER:HB3	2.02	0.41
3:C:62:TYR:CD1	3:C:231:ARG:HB2	2.55	0.41
3:C:79:LEU:HD13	3:C:209:TYR:CZ	2.56	0.41
3:C:151:LEU:O	3:C:152:GLN:HB2	2.21	0.41
1:A:92:ILE:N	1:A:93:PRO:HD2	2.35	0.41
2:B:120:GLU:O	2:B:229:ILE:HG13	2.20	0.41
3:C:184:ARG:HH11	3:C:188:VAL:H	1.69	0.41
3:C:313:ARG:O	3:C:392:GLU:HB3	2.20	0.41
1:A:169:GLU:OE1	3:C:18:HIS:HD2	2.04	0.41
3:C:76:LYS:HA	3:C:210:VAL:O	2.21	0.41
3:C:142:VAL:HA	3:C:143:PRO:HD2	1.85	0.41
3:C:329:THR:HG21	3:C:348:ALA:N	2.35	0.41
2:B:115:TRP:CZ2	2:B:231:LEU:HD11	2.55	0.41
3:C:140:GLY:O	3:C:206:LEU:HD12	2.21	0.41
3:C:189:ARG:NH1	3:C:243:GLN:O	2.50	0.41
1:A:56:ILE:HG23	1:A:207:GLN:HA	2.03	0.41
1:A:112:PHE:CD1	1:A:120:ILE:HD13	2.56	0.41
3:C:12:PHE:CE1	3:C:14:PRO:HG3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:71:LYS:HA	3:C:223:THR:HA	2.03	0.41
3:C:153:MET:HG3	3:C:207:PHE:CZ	2.56	0.41
3:C:262:TYR:CE1	3:C:300:PRO:HB2	2.55	0.41
3:C:24:THR:O	3:C:26:LEU:N	2.53	0.40
1:A:22:GLY:O	1:A:26:PHE:N	2.54	0.40
1:A:110:ILE:HB	1:A:168:ILE:HG22	2.04	0.40
1:A:183:LEU:HA	1:A:183:LEU:HD12	1.88	0.40
3:C:54:THR:HB	3:C:57:SER:HB3	2.02	0.40
3:C:68:VAL:CG2	3:C:81:LEU:HD23	2.52	0.40
1:A:161:ILE:HG22	1:A:164:VAL:H	1.85	0.40
2:B:188:PRO:HG2	2:B:199:ILE:HD12	2.03	0.40
1:A:103:ARG:HD3	3:C:43:HIS:CD2	2.57	0.40
3:C:323:GLY:N	3:C:327:TRP:HE1	2.20	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	250/258 (97%)	227 (91%)	20 (8%)	3 (1%)	13 50
2	B	248/253 (98%)	221 (89%)	21 (8%)	6 (2%)	6 35
3	C	395/416 (95%)	362 (92%)	31 (8%)	2 (0%)	29 68
All	All	893/927 (96%)	810 (91%)	72 (8%)	11 (1%)	17 50

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	76	PHE
2	B	233	ASN
3	C	12	PHE
3	C	194	ASN

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Mol	Chain	Res	Type
1	A	252	VAL
2	B	232	ASN
1	A	189	SER
1	A	53	ASP
2	B	20	SER
2	B	42	TRP
2	B	210	ILE

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	215/219 (98%)	215 (100%)	0	100 100
2	B	221/224 (99%)	220 (100%)	1 (0%)	88 94
3	C	338/354 (96%)	338 (100%)	0	100 100
All	All	774/797 (97%)	773 (100%)	1 (0%)	93 98

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	195	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	GLN
1	A	124	HIS
1	A	140	ASN
1	A	148	GLN
1	A	152	ASN
2	B	64	GLN
2	B	134	GLN
2	B	136	GLN
2	B	146	ASN

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Mol	Chain	Res	Type
3	C	18	HIS
3	C	116	GLN
3	C	150	GLN
3	C	227	ASN
3	C	243	GLN
3	C	280	ASN
3	C	325	GLN
3	C	386	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no monosaccharides 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-4009. 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

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.