wwPDB EM Map/Model Validation Report (i)

Sep 6, 2016 – 08:48 PM EDT

PDB ID	:	3JD7
EMDB ID:	:	EMD-6637
Title	:	The novel asymmetric entry intermediate of a picornavirus captured with nanodiscs
Authors	:	Lee, H.; Shingler, K.L.; Organtini, L.J.; Ashley, R.E.; Makhov, A.M.; Conway, J.F.; Hafenstein, S.
Deposited on	:	2016-04-29
Resolution	:	3.90 Å(reported)
Based on PDB ID	:	1COV

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

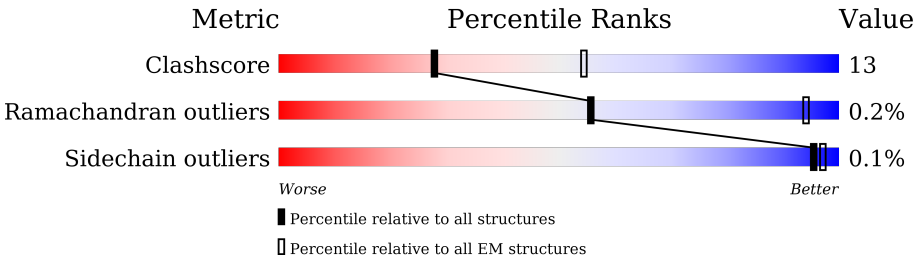
MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20027939

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 3.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)
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 <=5%

Mol	Chain	Length	Quality of chain
1	1	281	69% 27% .
2	2	263	73% 24% .
3	3	238	73% 27%
4	4	68	66% 16% 18%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	PLM	1	301	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6403 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 Capsid protein VP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	1	270	Total	C	N	O	S	0	0
			2139	1348	376	407	8		

- Molecule 2 is a protein called Capsid protein VP2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	2	256	Total	C	N	O	S	0	0
			1973	1247	335	375	16		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
2	151	THR	SER	CONFLICT	UNP Q66282
2	245	VAL	ILE	CONFLICT	UNP Q66282

- Molecule 3 is a protein called Capsid protein VP3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	3	238	Total	C	N	O	S	0	0
			1837	1174	294	352	17		

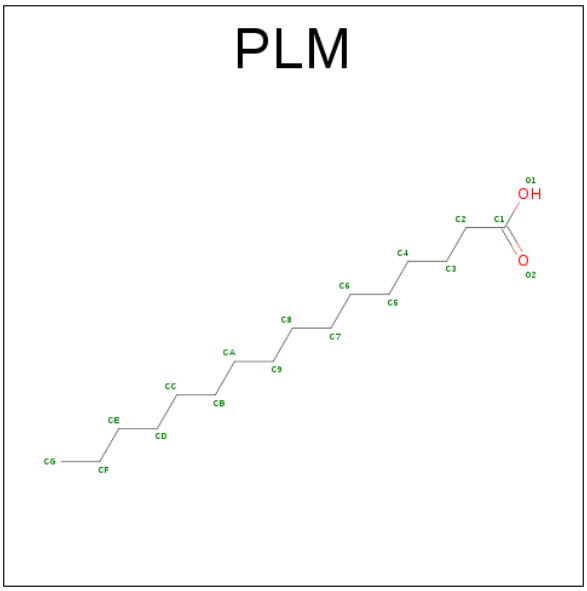
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
3	234	GLU	GLN	CONFLICT	UNP Q66282

- Molecule 4 is a protein called Capsid protein VP4.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	4	56	Total	C	N	O	S	0	0
			436	271	75	89	1		

- Molecule 5 is PALMITIC ACID (three-letter code: PLM) (formula: $C_{16}H_{32}O_2$).



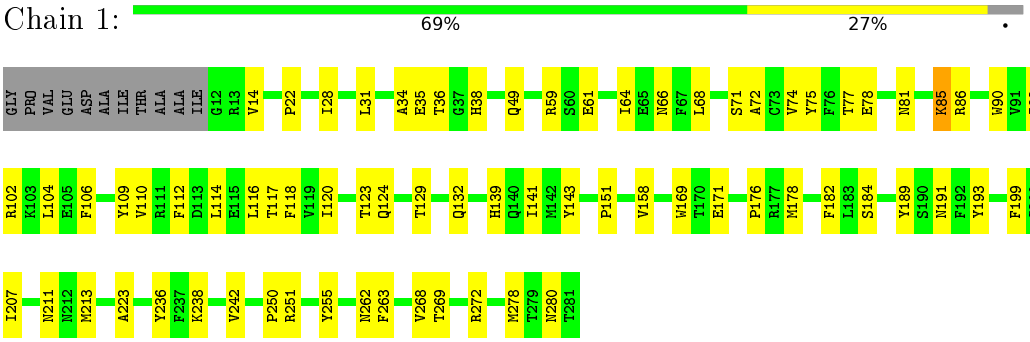
Mol	Chain	Residues	Atoms			AltConf
			Total	C	O	
5	1	1	18	16	2	0

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.

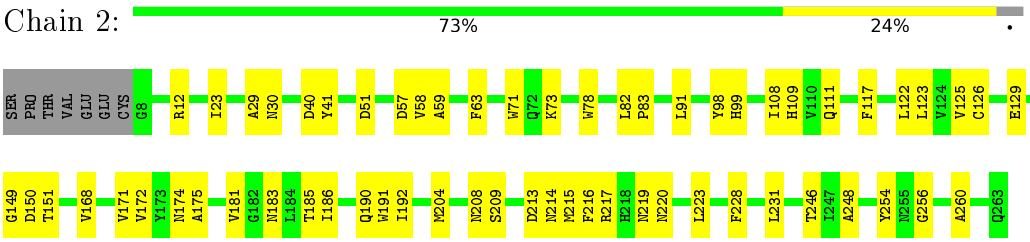
- Molecule 1: Capsid protein VP1

Chain 1:



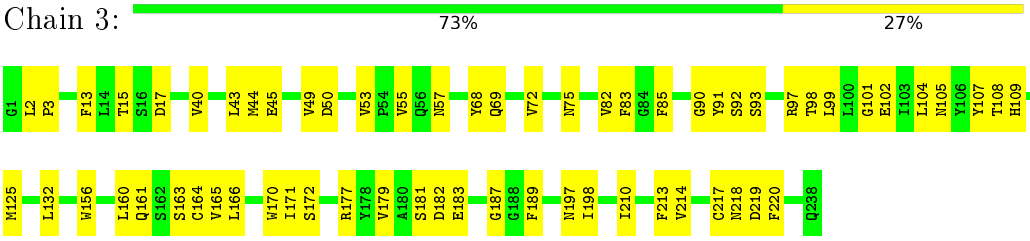
- Molecule 2: Capsid protein VP2

Chain 2:



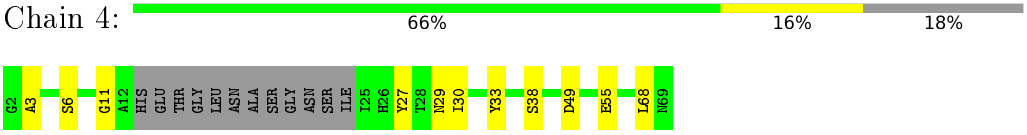
- Molecule 3: Capsid protein VP3

Chain 3:



- Molecule 4: Capsid protein VP4

Chain 4:



4 Experimental information (i)

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, I	Depositor
Number of particles used	57203	Depositor
Resolution determination method	FSC 0.143	Depositor
CTF correction method	Each particle	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	Not provided	Depositor
Minimum defocus (nm)	1362	Depositor
Maximum defocus (nm)	5875	Depositor
Magnification	Not provided	Depositor
Image detector	FEI FALCON II (4k x 4k)	Depositor

5 Model quality

5.1 Standard geometry

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

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	1	0.45	0/2198	0.51	0/2999
2	2	0.45	0/2025	0.53	0/2770
3	3	0.47	0/1888	0.50	0/2573
4	4	0.40	0/443	0.49	0/597
All	All	0.45	0/6554	0.51	0/8939

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	1	2139	0	2072	81	0
2	2	1973	0	1906	48	0
3	3	1837	0	1766	59	0
4	4	436	0	422	8	0
5	1	18	0	31	26	0
All	All	6403	0	6197	160	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 13.

All (160) 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)
1:1:116:LEU:HD11	5:1:301:PLM:HC2	1.16	1.13
1:1:213:MET:HE1	5:1:301:PLM:H71	1.14	1.13
1:1:114:LEU:CD1	5:1:301:PLM:H82	1.85	1.06
1:1:92:ILE:HD13	5:1:301:PLM:HA1	1.07	1.05
1:1:213:MET:CE	5:1:301:PLM:H71	1.86	1.05
1:1:213:MET:HE1	5:1:301:PLM:C7	1.89	1.00
1:1:213:MET:CE	5:1:301:PLM:C7	2.39	1.00
1:1:92:ILE:CD1	5:1:301:PLM:HA1	1.98	0.94
1:1:116:LEU:CD1	5:1:301:PLM:HC2	1.99	0.91
1:1:116:LEU:HD11	5:1:301:PLM:CC	2.04	0.88
1:1:213:MET:HE2	5:1:301:PLM:C7	2.09	0.80
1:1:92:ILE:HD13	5:1:301:PLM:CA	2.03	0.80
1:1:213:MET:CE	5:1:301:PLM:H72	2.13	0.77
1:1:213:MET:HE2	5:1:301:PLM:H72	1.66	0.75
1:1:114:LEU:HD11	5:1:301:PLM:H61	1.70	0.74
3:3:172:SER:HB2	3:3:177:ARG:HH21	1.53	0.73
2:2:12:ARG:HE	4:4:68:LEU:HB3	1.53	0.73
1:1:102:ARG:NH2	3:3:102:GLU:OE1	2.22	0.73
3:3:117:THR:HB	3:3:213:PHE:HB2	1.72	0.72
1:1:68:LEU:HD12	1:1:242:VAL:HG11	1.72	0.71
1:1:114:LEU:HD12	5:1:301:PLM:H82	1.70	0.70
1:1:238:LYS:NZ	3:3:17:ASP:O	2.22	0.70
3:3:75:ASN:O	3:3:197:ASN:ND2	2.26	0.68
1:1:114:LEU:HD11	5:1:301:PLM:H82	1.72	0.67
1:1:143:TYR:CE1	5:1:301:PLM:HB2	2.29	0.67
2:2:117:PHE:HE1	3:3:125:MET:HG3	1.61	0.66
2:2:213:ASP:OD1	2:2:214:ASN:N	2.28	0.66
1:1:71:SER:HB2	3:3:15:THR:HG23	1.78	0.66
1:1:36:THR:HG22	2:2:29:ALA:HB1	1.78	0.65
1:1:118:PHE:HE2	1:1:178:MET:HE3	1.62	0.65
1:1:268:VAL:HG12	1:1:269:THR:HG23	1.80	0.64
1:1:199:PHE:HE1	2:2:217:ARG:HD2	1.62	0.63
1:1:59:ARG:NH2	3:3:218:ASN:O	2.30	0.63
2:2:228:PHE:O	3:3:69:GLN:NE2	2.31	0.63
2:2:82:LEU:HD11	2:2:108:ILE:HD11	1.81	0.62
1:1:81:ASN:O	1:1:86:ARG:NH1	2.33	0.62
2:2:83:PRO:HG2	2:2:219:ASN:HA	1.81	0.62
1:1:114:LEU:HD11	5:1:301:PLM:C6	2.29	0.61
3:3:2:LEU:HD12	3:3:3:PRO:HD2	1.82	0.60
3:3:101:GLY:O	3:3:105:ASN:ND2	2.34	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:3:90:GLY:HA3	3:3:179:VAL:HG22	1.83	0.60
3:3:171:ILE:O	3:3:177:ARG:NH2	2.36	0.58
3:3:44:MET:HE1	3:3:220:PHE:HB3	1.85	0.58
2:2:40:ASP:OD1	2:2:41:TYR:N	2.31	0.58
2:2:208:ASN:OD1	2:2:209:SER:N	2.37	0.57
2:2:12:ARG:HH22	3:3:160:LEU:HD11	1.68	0.57
3:3:164:CYS:SG	3:3:165:VAL:N	2.76	0.57
3:3:156:TRP:NE1	3:3:164:CYS:HB2	2.19	0.57
3:3:197:ASN:OD1	3:3:198:ILE:N	2.38	0.57
1:1:250:PRO:HG2	3:3:102:GLU:HG3	1.88	0.56
1:1:14:VAL:HG13	3:3:219:ASP:HA	1.87	0.56
2:2:23:ILE:HD11	2:2:246:THR:HG21	1.87	0.56
1:1:151:PRO:HG3	1:1:158:VAL:HB	1.87	0.55
1:1:269:THR:HG22	3:3:97:ARG:HB3	1.86	0.55
2:2:108:ILE:HD13	2:2:223:LEU:HD13	1.87	0.55
1:1:64:ILE:HD11	3:3:40:VAL:HG12	1.89	0.54
1:1:117:THR:HG21	3:3:13:PHE:CE1	2.43	0.53
3:3:45:GLU:OE2	4:4:49:ASP:N	2.41	0.53
2:2:141:PRO:HB3	2:2:168:VAL:HG22	1.91	0.53
1:1:199:PHE:CE1	2:2:217:ARG:HD2	2.43	0.52
2:2:51:ASP:OD2	2:2:256:GLY:N	2.39	0.52
1:1:129:THR:HB	1:1:132:GLN:HE21	1.75	0.52
1:1:104:LEU:O	1:1:191:ASN:ND2	2.44	0.51
1:1:120:ILE:HG21	1:1:139:HIS:CD2	2.46	0.51
3:3:132:LEU:HB3	3:3:166:LEU:HD22	1.93	0.51
3:3:107:TYR:O	3:3:179:VAL:HG11	2.11	0.51
1:1:184:SER:HB2	1:1:189:TYR:CD1	2.45	0.51
1:1:280:ASN:HD21	3:3:92:SER:HA	1.76	0.50
1:1:110:VAL:HG12	1:1:112:PHE:HD1	1.76	0.50
3:3:98:THR:HG22	3:3:99:LEU:H	1.76	0.50
1:1:262:ASN:OD1	2:2:134:CYS:N	2.44	0.50
1:1:38:HIS:NE2	4:4:55:GLU:HG2	2.26	0.50
1:1:251:ARG:NH2	1:1:263:PHE:HB3	2.26	0.49
2:2:149:GLY:O	2:2:151:THR:N	2.46	0.49
3:3:115:LYS:HG3	3:3:217:CYS:SG	2.53	0.49
2:2:186:ILE:HA	3:3:49:VAL:HG11	1.94	0.49
3:3:156:TRP:CD1	3:3:164:CYS:HB2	2.47	0.49
3:3:182:ASP:OD1	3:3:183:GLU:N	2.46	0.49
2:2:214:ASN:OD1	2:2:215:MET:N	2.46	0.48
1:1:213:MET:HE2	5:1:301:PLM:H92	1.95	0.48
1:1:182:PHE:CZ	1:1:184:SER:HB3	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:186:ILE:HD11	3:3:99:LEU:HD22	1.94	0.48
1:1:35:GLU:HA	2:2:191:TRP:HB2	1.95	0.48
3:3:72:VAL:HG21	3:3:210:ILE:HD12	1.95	0.48
4:4:29:ASN:OD1	4:4:30:ILE:N	2.46	0.48
2:2:126:CYS:SG	2:2:204:MET:HG3	2.54	0.48
4:4:6:SER:HB2	4:4:27:TYR:CE1	2.49	0.48
3:3:104:LEU:HD11	3:3:220:PHE:HZ	1.78	0.48
1:1:143:TYR:CD1	5:1:301:PLM:HD2	2.49	0.47
3:3:82:VAL:HG12	3:3:83:PHE:HD2	1.78	0.47
2:2:183:ASN:O	2:2:186:ILE:HD12	2.14	0.47
2:2:122:LEU:HB2	2:2:192:ILE:HB	1.94	0.47
3:3:161:GLN:HE22	4:4:68:LEU:H	1.60	0.47
1:1:201:ARG:HH22	2:2:216:PHE:HD1	1.62	0.47
1:1:22:PRO:N	1:1:49:GLN:HE21	2.12	0.47
1:1:272:ARG:HG2	3:3:57:ASN:HB2	1.97	0.47
1:1:143:TYR:HB2	5:1:301:PLM:HF2	1.96	0.47
4:4:33:TYR:HD2	4:4:38:SER:HB3	1.79	0.47
2:2:99:HIS:CD2	2:2:254:TYR:HB3	2.50	0.46
3:3:172:SER:HB2	3:3:177:ARG:NH2	2.27	0.46
4:4:3:ALA:HA	4:4:30:ILE:HG12	1.96	0.46
1:1:184:SER:HB2	1:1:189:TYR:CE1	2.50	0.46
2:2:30:ASN:O	2:2:190:GLN:NE2	2.43	0.46
3:3:91:TYR:CG	3:3:181:SER:HB3	2.50	0.46
1:1:143:TYR:CB	5:1:301:PLM:HF2	2.46	0.46
2:2:59:ALA:HB1	2:2:91:LEU:HD12	1.97	0.46
1:1:116:LEU:CD1	5:1:301:PLM:CC	2.79	0.45
3:3:156:TRP:CE2	3:3:164:CYS:HB2	2.51	0.45
1:1:28:ILE:HD12	1:1:31:LEU:HB2	1.98	0.45
2:2:73:LYS:HG2	2:2:231:LEU:O	2.17	0.45
1:1:169:TRP:CH2	1:1:171:GLU:HA	2.51	0.45
2:2:149:GLY:C	2:2:151:THR:H	2.19	0.45
2:2:98:TYR:CE2	2:2:260:ALA:HB2	2.52	0.45
1:1:75:TYR:CE2	1:1:90:TRP:CD1	3.05	0.45
1:1:74:VAL:O	1:1:97:ALA:HB2	2.17	0.45
2:2:71:TRP:HD1	2:2:78:TRP:HH2	1.65	0.45
1:1:78:GLU:O	1:1:85:LYS:HE3	2.16	0.45
3:3:177:ARG:NH1	3:3:187:GLY:HA2	2.32	0.45
1:1:109:TYR:OH	2:2:129:GLU:OE2	2.32	0.45
3:3:98:THR:HG22	3:3:99:LEU:N	2.32	0.44
1:1:93:THR:HG22	1:1:211:ASN:ND2	2.33	0.44
2:2:181:VAL:HG13	3:3:68:TYR:HE2	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:72:ALA:O	1:1:236:TYR:HA	2.17	0.44
2:2:83:PRO:HG2	2:2:220:ASN:H	1.82	0.44
1:1:280:ASN:HD21	3:3:93:SER:H	1.66	0.43
2:2:185:THR:OG1	3:3:50:ASP:O	2.31	0.43
2:2:117:PHE:CE1	3:3:125:MET:HG3	2.49	0.43
1:1:202:ASN:OD1	1:1:203:GLY:N	2.43	0.43
2:2:186:ILE:HD11	3:3:99:LEU:HD13	1.99	0.43
3:3:170:TRP:HH2	3:3:177:ARG:HG3	1.83	0.43
1:1:114:LEU:HD11	5:1:301:PLM:C8	2.44	0.43
2:2:57:ASP:OD2	2:2:58:VAL:HG23	2.17	0.43
1:1:81:ASN:OD1	1:1:223:ALA:HA	2.18	0.43
1:1:169:TRP:NE1	1:1:176:PRO:HD3	2.34	0.43
2:2:171:VAL:HG23	2:2:174:ASN:HB2	2.01	0.43
1:1:280:ASN:HD21	3:3:93:SER:N	2.16	0.43
2:2:82:LEU:HA	2:2:82:LEU:HD23	1.84	0.43
1:1:106:PHE:CE1	1:1:250:PRO:HG3	2.54	0.43
1:1:34:ALA:N	3:3:163:SER:OG	2.36	0.42
3:3:83:PHE:HE1	3:3:85:PHE:HE1	1.66	0.42
2:2:109:HIS:CE1	2:2:111:GLN:HE21	2.38	0.42
2:2:123:LEU:HG	2:2:125:VAL:HG13	2.01	0.42
1:1:75:TYR:CE2	1:1:77:THR:HG23	2.54	0.42
1:1:123:THR:HG22	1:1:124:GLN:O	2.20	0.42
2:2:63:PHE:CD1	2:2:248:ALA:HB2	2.54	0.42
3:3:53:VAL:HG21	3:3:214:VAL:HG23	2.02	0.42
1:1:120:ILE:HD11	1:1:141:ILE:HD11	2.02	0.41
1:1:143:TYR:HB2	5:1:301:PLM:CF	2.50	0.41
2:2:172:VAL:O	2:2:175:ALA:N	2.49	0.41
1:1:59:ARG:HD3	3:3:44:MET:HG3	2.03	0.41
1:1:251:ARG:HD3	1:1:255:TYR:CZ	2.56	0.41
2:2:63:PHE:HD1	2:2:246:THR:HG22	1.86	0.41
3:3:53:VAL:O	3:3:55:VAL:N	2.53	0.41
3:3:43:LEU:HA	3:3:43:LEU:HD23	1.92	0.41
1:1:193:TYR:N	1:1:207:ILE:HG22	2.36	0.41
2:2:71:TRP:HD1	2:2:78:TRP:CH2	2.39	0.41
1:1:278:MET:HE2	3:3:189:PHE:HD1	1.86	0.41
1:1:61:GLU:HA	1:1:66:ASN:HD22	1.85	0.40
3:3:83:PHE:HE1	3:3:85:PHE:CE1	2.40	0.40
3:3:108:THR:HG22	3:3:109:HIS:CD2	2.57	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	1	268/281 (95%)	260 (97%)	8 (3%)	0	100 100
2	2	254/263 (97%)	243 (96%)	10 (4%)	1 (0%)	39 79
3	3	236/238 (99%)	229 (97%)	7 (3%)	0	100 100
4	4	52/68 (76%)	49 (94%)	2 (4%)	1 (2%)	10 53
All	All	810/850 (95%)	781 (96%)	27 (3%)	2 (0%)	56 86

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	2	150	ASP
4	4	11	GLY

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	1	237/244 (97%)	236 (100%)	1 (0%)	93 97
2	2	218/225 (97%)	218 (100%)	0	100 100
3	3	206/206 (100%)	206 (100%)	0	100 100
4	4	48/57 (84%)	48 (100%)	0	100 100
All	All	709/732 (97%)	708 (100%)	1 (0%)	95 97

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

Mol	Chain	Res	Type
1	1	85	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	1	49	GLN
1	1	124	GLN
1	1	132	GLN
1	1	164	ASN
1	1	208	ASN
2	2	95	ASN
2	2	109	HIS
2	2	197	ASN
2	2	255	ASN
3	3	56	GLN
3	3	105	ASN
3	3	153	HIS
3	3	161	GLN

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

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	$\#\|Z\| > 2$
5	PLM	1	301	-	14,17,17	0.66	0	14,17,17	2.07	4 (28%)

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
5	PLM	1	301	-	-	0/13/15/15	0/0/0/0

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	1	301	PLM	C6-C5-C4	-2.30	102.61	114.54
5	1	301	PLM	C3-C2-C1	2.16	124.25	112.58
5	1	301	PLM	C5-C4-C3	2.91	129.63	114.54
5	1	301	PLM	C7-C6-C5	5.54	143.32	114.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 26 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	1	301	PLM	26	0

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.