



wwPDB EM Map/Model Validation Report ⓘ

Jul 19, 2016 – 02:31 AM EDT

PDB ID : 3JB5
EMDB ID: : EMD-6398
Title : Capsid Structure of the Propionibacterium acnes Bacteriophage ATCC_Clear
Authors : Chiou, J.; Zhang, X.; Marinelli, L.J.; Modlin, R.L.; Zhou, Z.H.
Deposited on : 2015-07-23
Resolution : 3.70 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

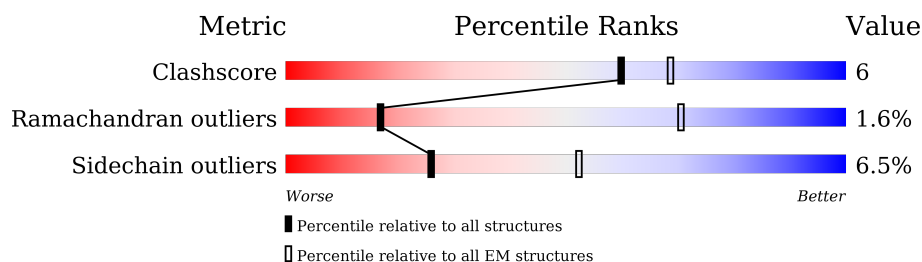
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.70 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	315	81% 13% • 5%
1	B	315	78% 15% • 5%
1	C	315	70% 23% • 5%
1	D	315	78% 15% • 5%
1	E	315	78% 15% • 5%
1	F	315	78% 15% • 5%
1	G	315	70% 17% • 12%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 15232 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 major capsid protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	300	Total	C	N	O	S	0	0
			2200	1398	374	423	5		
1	B	300	Total	C	N	O	S	0	0
			2200	1398	374	423	5		
1	C	300	Total	C	N	O	S	0	0
			2200	1398	374	423	5		
1	D	300	Total	C	N	O	S	0	0
			2200	1398	374	423	5		
1	E	300	Total	C	N	O	S	0	0
			2200	1398	374	423	5		
1	F	300	Total	C	N	O	S	0	0
			2200	1398	374	423	5		
1	G	276	Total	C	N	O	S	0	0
			2032	1289	348	391	4		

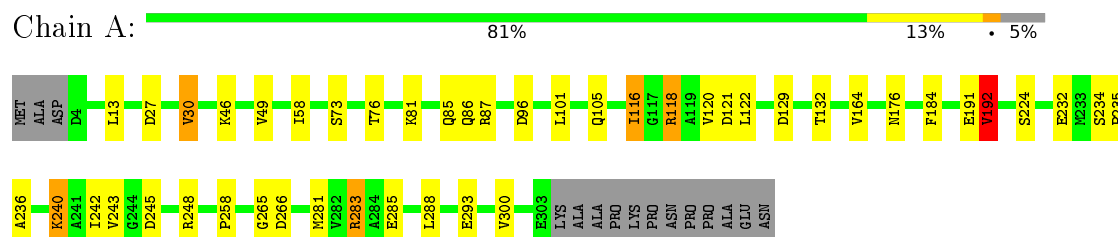
There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	8	VAL	ALA	CONFLICT	UNP A4K473
B	8	VAL	ALA	CONFLICT	UNP A4K473
C	8	VAL	ALA	CONFLICT	UNP A4K473
D	8	VAL	ALA	CONFLICT	UNP A4K473
E	8	VAL	ALA	CONFLICT	UNP A4K473
F	8	VAL	ALA	CONFLICT	UNP A4K473
G	8	VAL	ALA	CONFLICT	UNP A4K473

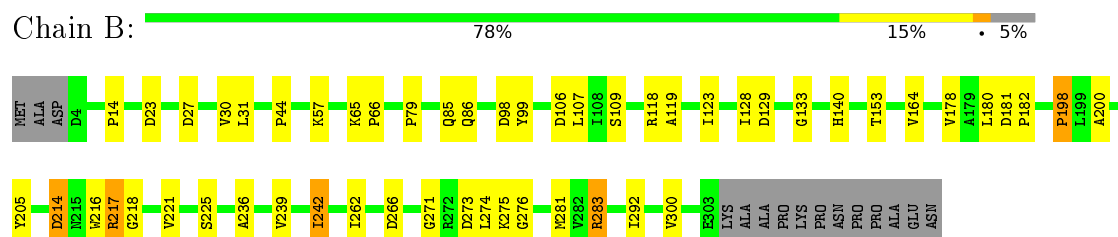
3 Residue-property plots

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.

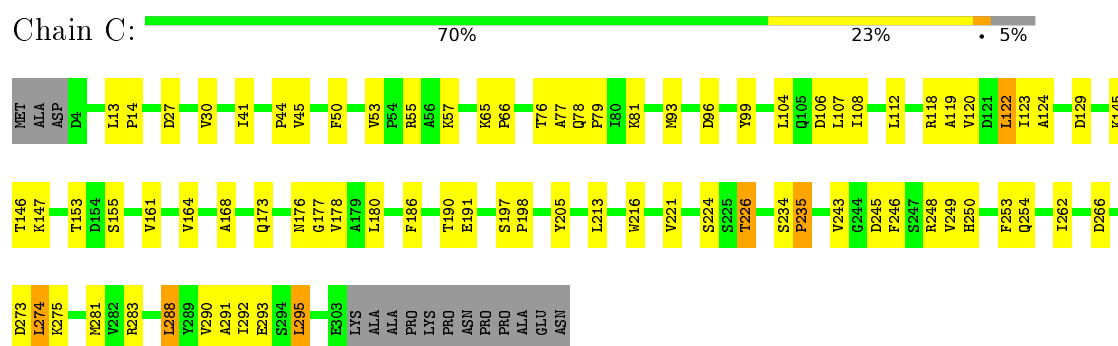
- Molecule 1: major capsid protein



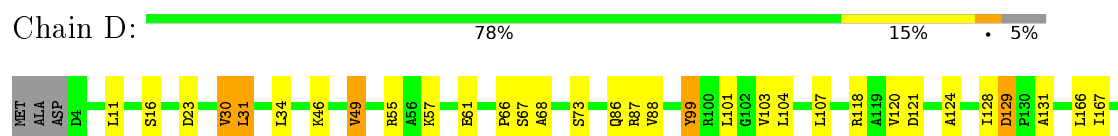
- Molecule 1: major capsid protein



- Molecule 1: major capsid protein



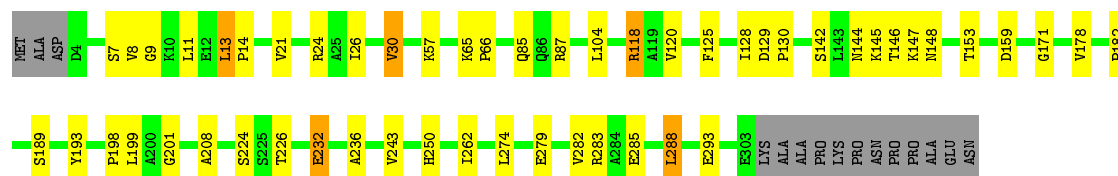
- Molecule 1: major capsid protein





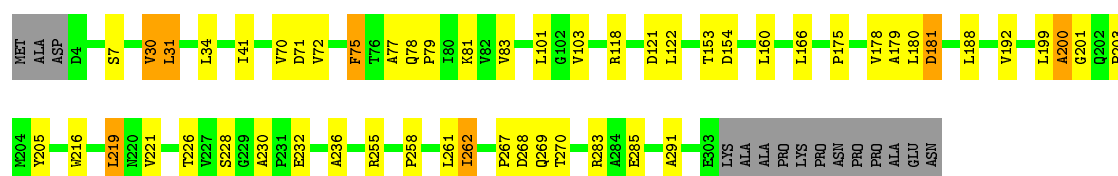
- Molecule 1: major capsid protein

Chain E: 78% 15% 5%



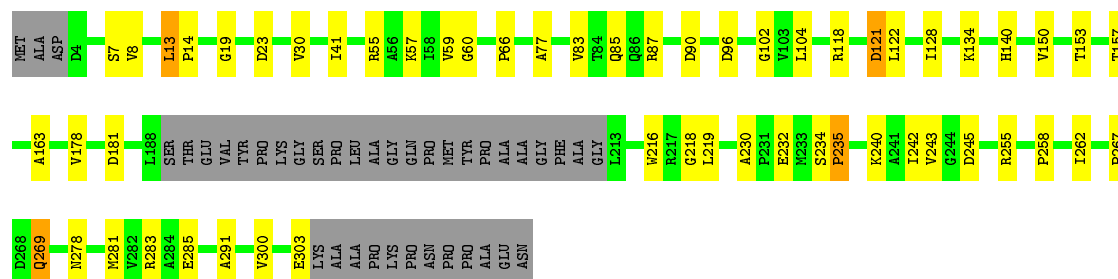
- Molecule 1: major capsid protein

Chain F: 78% 15% 5%



- Molecule 1: major capsid protein

Chain G: 70% 17% 12%



4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, I	Depositor
Number of particles used	27504	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 ($e^-/\text{\AA}^2$)	25	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2270	Depositor
Magnification	38462	Depositor
Image detector	GATAN K2 (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	0.21	0/2246	0.40	0/3057
1	B	0.21	0/2246	0.39	0/3057
1	C	0.21	0/2246	0.41	0/3057
1	D	0.21	0/2246	0.38	0/3057
1	E	0.21	0/2246	0.40	0/3057
1	F	0.21	0/2246	0.39	0/3057
1	G	0.21	0/2070	0.39	0/2814
All	All	0.21	0/15546	0.40	0/21156

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	2200	0	2199	21	0
1	B	2200	0	2199	27	0
1	C	2200	0	2199	44	0
1	D	2200	0	2199	29	0
1	E	2200	0	2199	27	0
1	F	2200	0	2199	24	0
1	G	2032	0	2038	25	0
All	All	15232	0	15232	178	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 6.

All (178) 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:F:181:ASP:HB2	1:F:228:SER:HB3	1.70	0.72
1:B:271:GLY:O	1:G:269:GLN:NE2	2.24	0.71
1:C:77:ALA:HB2	1:C:291:ALA:HB2	1.76	0.67
1:B:65:LYS:HD3	1:C:283:ARG:HH12	1.58	0.67
1:E:57:LYS:HD3	1:E:66:PRO:HG3	1.77	0.67
1:F:205:TYR:HH	1:F:216:TRP:HE1	1.43	0.66
1:C:57:LYS:HD3	1:C:66:PRO:HG3	1.77	0.66
1:C:205:TYR:HH	1:C:216:TRP:HE1	1.43	0.66
1:C:120:VAL:HG22	1:C:288:LEU:HG	1.78	0.66
1:A:242:ILE:HD13	1:A:300:VAL:HG12	1.78	0.65
1:G:118:ARG:NH1	1:G:232:GLU:OE2	2.31	0.63
1:B:57:LYS:HD3	1:B:66:PRO:HG3	1.81	0.63
1:C:176:ASN:ND2	1:C:245:ASP:OD1	2.32	0.63
1:C:178:VAL:HG13	1:C:243:VAL:HG22	1.80	0.62
1:A:30:VAL:HG22	1:A:224:SER:HB3	1.82	0.62
1:F:199:LEU:O	1:F:201:GLY:N	2.31	0.62
1:E:283:ARG:NH1	1:E:285:GLU:OE1	2.32	0.62
1:G:57:LYS:HD3	1:G:66:PRO:HG3	1.82	0.62
1:E:13:LEU:HD23	1:E:14:PRO:HD2	1.82	0.62
1:B:274:LEU:HD21	1:B:281:MET:HB2	1.82	0.60
1:C:248:ARG:NH1	1:C:293:GLU:OE1	2.35	0.60
1:D:124:ALA:HB2	1:D:290:VAL:HG11	1.83	0.59
1:C:234:SER:HB2	1:C:235:PRO:HD3	1.83	0.59
1:G:13:LEU:HD23	1:G:14:PRO:HD2	1.84	0.59
1:E:120:VAL:HG22	1:E:288:LEU:HG	1.84	0.59
1:C:122:LEU:HD13	1:C:226:THR:HG22	1.83	0.59
1:C:106:ASP:O	1:C:108:ILE:N	2.36	0.58
1:A:248:ARG:NH2	1:B:23:ASP:O	2.35	0.58
1:C:178:VAL:HB	1:C:221:VAL:HG12	1.85	0.58
1:D:176:ASN:ND2	1:D:245:ASP:OD1	2.36	0.58
1:D:262:ILE:HB	1:D:281:MET:HB3	1.85	0.58
1:D:274:LEU:HD21	1:D:281:MET:HB2	1.86	0.57
1:C:65:LYS:HD3	1:D:283:ARG:HH12	1.68	0.57
1:D:171:GLY:HA3	1:E:182:PRO:HB2	1.87	0.56
1:C:249:VAL:HG22	1:C:292:ILE:HG22	1.86	0.56
1:D:167:ILE:HD13	1:D:175:PRO:HB3	1.87	0.56
1:A:118:ARG:NH1	1:A:232:GLU:OE1	2.39	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:262:ILE:HB	1:B:281:MET:HB3	1.88	0.56
1:G:283:ARG:NH1	1:G:285:GLU:OE1	2.38	0.56
1:D:176:ASN:HA	1:D:219:LEU:HD22	1.87	0.56
1:B:128:ILE:HD12	1:B:133:GLY:HA2	1.87	0.55
1:E:87:ARG:NH1	1:E:279:GLU:OE2	2.38	0.55
1:A:46:LYS:HG2	1:A:76:THR:HG22	1.88	0.55
1:C:197:SER:OG	1:D:191:GLU:OE2	2.20	0.55
1:B:242:ILE:HG22	1:B:300:VAL:HG22	1.90	0.55
1:B:27:ASP:OD2	1:B:118:ARG:NH1	2.41	0.54
1:G:234:SER:HB2	1:G:235:PRO:HD3	1.91	0.53
1:B:178:VAL:HB	1:B:221:VAL:HG22	1.91	0.53
1:C:262:ILE:HD11	1:C:283:ARG:HE	1.73	0.53
1:D:120:VAL:HB	1:D:288:LEU:HD22	1.90	0.53
1:A:176:ASN:ND2	1:A:245:ASP:O	2.41	0.53
1:B:273:ASP:OD2	1:B:276:GLY:N	2.42	0.52
1:F:118:ARG:NH2	1:F:232:GLU:OE1	2.41	0.52
1:F:77:ALA:HB2	1:F:291:ALA:HB2	1.91	0.52
1:A:58:ILE:HD11	1:B:123:ILE:HD11	1.92	0.52
1:A:49:VAL:HG23	1:A:73:SER:HB3	1.90	0.52
1:A:176:ASN:ND2	1:A:245:ASP:OD1	2.42	0.51
1:A:248:ARG:HG3	1:A:293:GLU:HB3	1.91	0.51
1:C:274:LEU:HD21	1:C:281:MET:HB2	1.91	0.51
1:G:85:GLN:HE22	1:G:267:PRO:HB2	1.75	0.51
1:A:266:ASP:N	1:A:266:ASP:OD1	2.43	0.51
1:D:178:VAL:N	1:D:220:ASN:O	2.43	0.51
1:G:85:GLN:HB3	1:G:281:MET:HE3	1.93	0.51
1:D:49:VAL:HG23	1:D:73:SER:HB3	1.92	0.51
1:B:27:ASP:OD2	1:B:225:SER:OG	2.28	0.50
1:E:262:ILE:HD11	1:E:283:ARG:HG2	1.92	0.50
1:E:189:SER:HB2	1:E:208:ALA:HB1	1.93	0.50
1:G:30:VAL:N	1:G:121:ASP:OD2	2.44	0.50
1:C:13:LEU:HD12	1:C:14:PRO:HD2	1.93	0.50
1:D:118:ARG:HH22	1:D:230:ALA:HB1	1.76	0.50
1:C:155:SER:OG	1:C:191:GLU:OE2	2.27	0.50
1:D:242:ILE:HG12	1:D:300:VAL:HG12	1.93	0.50
1:C:44:PRO:HA	1:C:79:PRO:HD3	1.94	0.50
1:D:57:LYS:HD3	1:D:66:PRO:HG3	1.94	0.50
1:E:85:GLN:OE1	1:E:283:ARG:NH2	2.45	0.49
1:D:30:VAL:N	1:D:121:ASP:OD2	2.43	0.49
1:B:214:ASP:N	1:B:214:ASP:OD1	2.44	0.49
1:C:253:PHE:CD1	1:C:288:LEU:HB3	2.47	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:GLN:OE1	1:B:283:ARG:NH2	2.46	0.49
1:G:121:ASP:N	1:G:121:ASP:OD1	2.46	0.49
1:E:9:GLY:C	1:E:11:LEU:H	2.16	0.49
1:F:268:ASP:OD1	1:F:270:THR:OG1	2.27	0.48
1:B:98:ASP:HB3	1:G:41:ILE:HG23	1.96	0.48
1:E:178:VAL:HG22	1:E:243:VAL:HG13	1.96	0.48
1:F:192:VAL:HG12	1:F:203:PRO:HA	1.94	0.48
1:F:71:ASP:OD1	1:F:71:ASP:N	2.46	0.48
1:A:85:GLN:HG2	1:A:283:ARG:HD2	1.96	0.47
1:G:7:SER:OG	1:G:8:VAL:N	2.48	0.47
1:D:87:ARG:NH2	1:D:272:ARG:HB2	2.30	0.47
1:F:262:ILE:HD11	1:F:283:ARG:HD2	1.97	0.47
1:B:216:TRP:O	1:B:218:GLY:N	2.46	0.47
1:E:250:HIS:NE2	1:E:293:GLU:OE2	2.37	0.47
1:E:30:VAL:HG22	1:E:224:SER:HB3	1.96	0.47
1:A:164:VAL:HG22	1:A:243:VAL:HG11	1.95	0.47
1:D:67:SER:OG	1:D:68:ALA:N	2.48	0.47
1:C:213:LEU:H	1:C:213:LEU:HD22	1.80	0.46
1:F:34:LEU:HD11	1:F:179:ALA:HB2	1.97	0.46
1:C:45:VAL:HG23	1:D:16:SER:HB2	1.96	0.46
1:F:153:THR:OG1	1:F:154:ASP:N	2.46	0.46
1:F:101:LEU:HB3	1:F:103:VAL:HG23	1.97	0.46
1:E:198:PRO:HD3	1:F:200:ALA:HB3	1.98	0.46
1:C:93:MET:SD	1:C:104:LEU:HD22	2.56	0.46
1:G:87:ARG:HB3	1:G:281:MET:SD	2.55	0.46
1:C:253:PHE:HD1	1:C:288:LEU:HB3	1.81	0.46
1:D:55:ARG:HG2	1:E:232:GLU:HG3	1.97	0.46
1:A:120:VAL:HB	1:A:288:LEU:HD22	1.98	0.45
1:B:85:GLN:HG2	1:B:283:ARG:HB2	1.98	0.45
1:B:266:ASP:OD1	1:B:266:ASP:N	2.49	0.45
1:G:150:VAL:HG21	1:G:163:ALA:HA	1.97	0.45
1:G:262:ILE:HB	1:G:281:MET:HB3	1.97	0.45
1:G:83:VAL:HG22	1:G:285:GLU:HG2	1.97	0.45
1:G:77:ALA:HB2	1:G:291:ALA:HB2	1.97	0.45
1:B:205:TYR:HH	1:B:216:TRP:HE1	1.63	0.45
1:C:168:ALA:HB1	1:D:182:PRO:HB2	1.98	0.45
1:E:118:ARG:HD2	1:E:226:THR:HG22	1.98	0.45
1:C:119:ALA:O	1:C:123:ILE:HG12	2.16	0.45
1:C:146:THR:OG1	1:C:147:LYS:N	2.49	0.45
1:F:121:ASP:OD2	1:F:226:THR:OG1	2.21	0.45
1:G:90:ASP:HB2	1:G:278:ASN:O	2.16	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:SER:HB3	1:A:235:PRO:HD3	1.98	0.45
1:C:295:LEU:H	1:C:295:LEU:HD13	1.82	0.44
1:B:181:ASP:OD1	1:B:182:PRO:HD2	2.17	0.44
1:A:184:PHE:HB2	1:A:240:LYS:HG3	1.99	0.44
1:C:147:LYS:HE3	1:C:147:LYS:HB2	1.88	0.44
1:G:242:ILE:HG12	1:G:300:VAL:HG12	1.99	0.44
1:G:128:ILE:HG13	1:G:134:LYS:O	2.17	0.44
1:G:245:ASP:OD1	1:G:245:ASP:N	2.49	0.44
1:B:119:ALA:O	1:B:123:ILE:HG12	2.18	0.44
1:F:267:PRO:C	1:F:269:GLN:H	2.21	0.44
1:A:81:LYS:NZ	1:A:285:GLU:OE2	2.31	0.44
1:B:198:PRO:O	1:B:200:ALA:N	2.45	0.44
1:C:250:HIS:NE2	1:C:293:GLU:OE2	2.50	0.43
1:C:118:ARG:O	1:C:122:LEU:HB2	2.18	0.43
1:F:175:PRO:O	1:F:219:LEU:HD22	2.18	0.43
1:B:44:PRO:HA	1:B:79:PRO:HD3	2.00	0.43
1:C:266:ASP:OD1	1:C:266:ASP:N	2.52	0.43
1:E:129:ASP:OD1	1:E:129:ASP:N	2.48	0.43
1:E:7:SER:OG	1:E:8:VAL:N	2.51	0.43
1:A:27:ASP:N	1:A:27:ASP:OD1	2.48	0.43
1:F:30:VAL:HB	1:F:31:LEU:H	1.62	0.43
1:F:83:VAL:HG12	1:F:285:GLU:HG3	2.01	0.43
1:C:55:ARG:HB3	1:D:131:ALA:HA	2.01	0.43
1:E:142:SER:HB3	1:E:145:LYS:HG2	2.01	0.43
1:D:101:LEU:HB3	1:D:103:VAL:HG23	2.01	0.42
1:E:65:LYS:HB3	1:F:83:VAL:HG21	2.01	0.42
1:F:160:LEU:HD11	1:F:178:VAL:HG21	2.00	0.42
1:G:178:VAL:HG22	1:G:243:VAL:HG13	2.01	0.42
1:A:191:GLU:O	1:A:192:VAL:HG22	2.18	0.42
1:F:78:GLN:HA	1:F:79:PRO:HD3	1.90	0.42
1:C:79:PRO:HG3	1:C:254:GLN:HE22	1.84	0.42
1:D:234:SER:HB2	1:D:235:PRO:HD3	2.02	0.42
1:G:240:LYS:HE2	1:G:303:GLU:HG2	2.01	0.42
1:B:216:TRP:CD2	1:B:217:ARG:HG2	2.55	0.42
1:C:124:ALA:HA	1:C:290:VAL:HG11	2.02	0.41
1:C:78:GLN:HA	1:C:79:PRO:HD3	1.82	0.41
1:E:146:THR:HG22	1:E:148:ASN:H	1.84	0.41
1:F:178:VAL:HG13	1:F:221:VAL:HG22	2.01	0.41
1:C:198:PRO:HG3	1:E:201:GLY:H	1.86	0.41
1:A:116:ILE:H	1:A:116:ILE:HD13	1.84	0.41
1:C:30:VAL:HG13	1:C:224:SER:HB2	2.01	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:59:VAL:HG12	1:G:60:GLY:O	2.21	0.41
1:G:96:ASP:OD1	1:G:102:GLY:N	2.52	0.41
1:A:87:ARG:HD2	1:A:281:MET:SD	2.60	0.41
1:C:161:VAL:HA	1:C:164:VAL:HG12	2.03	0.41
1:F:75:PHE:HD1	1:F:75:PHE:H	1.68	0.41
1:B:164:VAL:HG13	1:C:186:PHE:CZ	2.56	0.41
1:D:61:GLU:OE2	1:F:7:SER:OG	2.37	0.41
1:D:49:VAL:HA	1:E:21:VAL:O	2.21	0.41
1:B:239:VAL:HG13	1:B:300:VAL:HG13	2.03	0.40
1:E:125:PHE:O	1:E:144:ASN:HB2	2.21	0.40
1:C:177:GLY:HA3	1:C:246:PHE:HD2	1.86	0.40
1:C:50:PHE:CD1	1:D:107:LEU:HD13	2.56	0.40
1:E:148:ASN:N	1:E:148:ASN:OD1	2.54	0.40
1:D:129:ASP:N	1:D:129:ASP:OD1	2.55	0.40
1:E:193:TYR:HB2	1:E:199:LEU:O	2.21	0.40
1:C:41:ILE:H	1:C:41:ILE:HG13	1.63	0.40
1:D:30:VAL:HB	1:D:31:LEU:H	1.66	0.40
1:E:129:ASP:HA	1:E:130:PRO:HD3	1.89	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	298/315 (95%)	266 (89%)	28 (9%)	4 (1%)	15	62
1	B	298/315 (95%)	272 (91%)	21 (7%)	5 (2%)	11	57
1	C	298/315 (95%)	274 (92%)	20 (7%)	4 (1%)	15	62
1	D	298/315 (95%)	272 (91%)	21 (7%)	5 (2%)	11	57
1	E	298/315 (95%)	278 (93%)	17 (6%)	3 (1%)	19	66
1	F	298/315 (95%)	275 (92%)	17 (6%)	6 (2%)	9	55

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	272/315 (86%)	251 (92%)	15 (6%)	6 (2%)	8	52
All	All	2060/2205 (93%)	1888 (92%)	139 (7%)	33 (2%)	17	58

All (33) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	192	VAL
1	A	258	PRO
1	B	198	PRO
1	C	107	LEU
1	F	30	VAL
1	B	14	PRO
1	B	140	HIS
1	D	99	TYR
1	E	171	GLY
1	F	262	ILE
1	A	236	ALA
1	B	217	ARG
1	B	236	ALA
1	C	173	GLN
1	D	23	ASP
1	D	236	ALA
1	E	24	ARG
1	E	236	ALA
1	F	200	ALA
1	F	236	ALA
1	G	218	GLY
1	C	53	VAL
1	D	30	VAL
1	D	230	ALA
1	F	258	PRO
1	G	19	GLY
1	G	235	PRO
1	G	269	GLN
1	C	235	PRO
1	F	230	ALA
1	A	265	GLY
1	G	230	ALA
1	G	258	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	230/241 (95%)	215 (94%)	15 (6%)	21	64
1	B	230/241 (95%)	215 (94%)	15 (6%)	21	64
1	C	230/241 (95%)	212 (92%)	18 (8%)	16	56
1	D	230/241 (95%)	215 (94%)	15 (6%)	21	64
1	E	230/241 (95%)	217 (94%)	13 (6%)	25	68
1	F	230/241 (95%)	216 (94%)	14 (6%)	23	65
1	G	214/241 (89%)	201 (94%)	13 (6%)	23	65
All	All	1594/1687 (94%)	1491 (94%)	103 (6%)	26	64

All (103) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	30	VAL
1	A	86	GLN
1	A	96	ASP
1	A	101	LEU
1	A	105	GLN
1	A	116	ILE
1	A	118	ARG
1	A	121	ASP
1	A	122	LEU
1	A	129	ASP
1	A	132	THR
1	A	192	VAL
1	A	240	LYS
1	A	283	ARG
1	B	30	VAL
1	B	31	LEU
1	B	86	GLN
1	B	99	TYR
1	B	106	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	107	LEU
1	B	109	SER
1	B	129	ASP
1	B	153	THR
1	B	180	LEU
1	B	214	ASP
1	B	242	ILE
1	B	275	LYS
1	B	283	ARG
1	B	292	ILE
1	C	27	ASP
1	C	76	THR
1	C	81	LYS
1	C	96	ASP
1	C	99	TYR
1	C	112	LEU
1	C	122	LEU
1	C	129	ASP
1	C	145	LYS
1	C	153	THR
1	C	180	LEU
1	C	190	THR
1	C	226	THR
1	C	273	ASP
1	C	274	LEU
1	C	275	LYS
1	C	288	LEU
1	C	295	LEU
1	D	11	LEU
1	D	31	LEU
1	D	34	LEU
1	D	46	LYS
1	D	49	VAL
1	D	86	GLN
1	D	88	VAL
1	D	99	TYR
1	D	104	LEU
1	D	128	ILE
1	D	129	ASP
1	D	166	LEU
1	D	176	ASN
1	D	281	MET

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	287	VAL
1	E	13	LEU
1	E	26	ILE
1	E	30	VAL
1	E	104	LEU
1	E	118	ARG
1	E	128	ILE
1	E	147	LYS
1	E	153	THR
1	E	159	ASP
1	E	232	GLU
1	E	274	LEU
1	E	282	VAL
1	E	288	LEU
1	F	31	LEU
1	F	41	ILE
1	F	70	VAL
1	F	72	VAL
1	F	75	PHE
1	F	81	LYS
1	F	122	LEU
1	F	166	LEU
1	F	180	LEU
1	F	181	ASP
1	F	188	LEU
1	F	219	LEU
1	F	255	ARG
1	F	261	LEU
1	G	13	LEU
1	G	23	ASP
1	G	55	ARG
1	G	104	LEU
1	G	121	ASP
1	G	122	LEU
1	G	140	HIS
1	G	153	THR
1	G	157	THR
1	G	181	ASP
1	G	216	TRP
1	G	219	LEU
1	G	255	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such

sidechains are listed below:

Mol	Chain	Res	Type
1	D	176	ASN
1	G	85	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates 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.