



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 08:15 pm BST

PDB ID : 2WWS
Title : Physalis Mottle Virus: Natural Empty Capsid
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Deposited on : 2009-10-27
Resolution : 3.90 Å(reported)

This is a Full wwPDB X-ray Structure 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/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

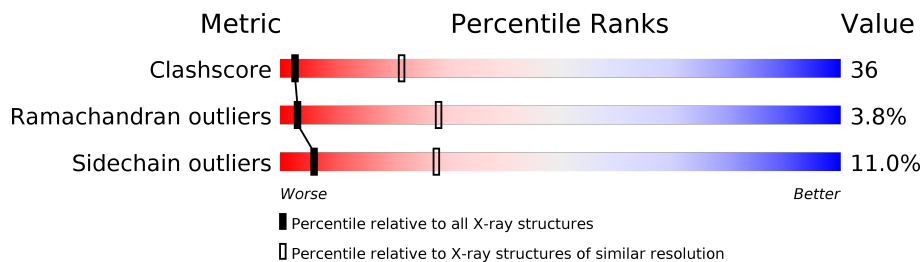
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

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)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	1004 (4.12-3.68)
Ramachandran outliers	138981	1021 (4.14-3.66)
Sidechain outliers	138945	1014 (4.14-3.66)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower 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	188	
1	B	188	
1	C	188	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3976 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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.

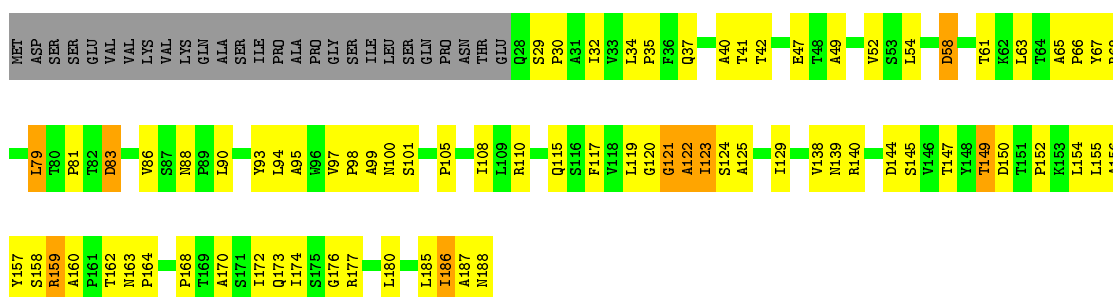
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	161	Total 1210	C 774	N 203	O 230	S 3	0	0	0
1	B	185	Total 1383	C 882	N 232	O 266	S 3	0	0	0
1	C	185	Total 1383	C 882	N 232	O 266	S 3	0	0	0

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 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.

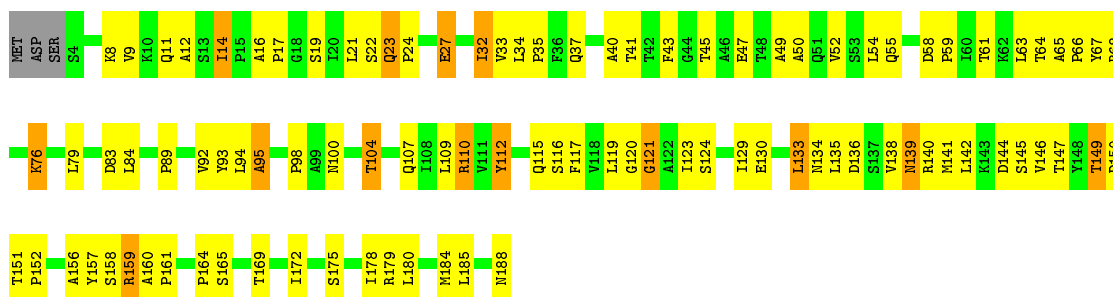
- Molecule 1: COAT PROTEIN

Chain A: 



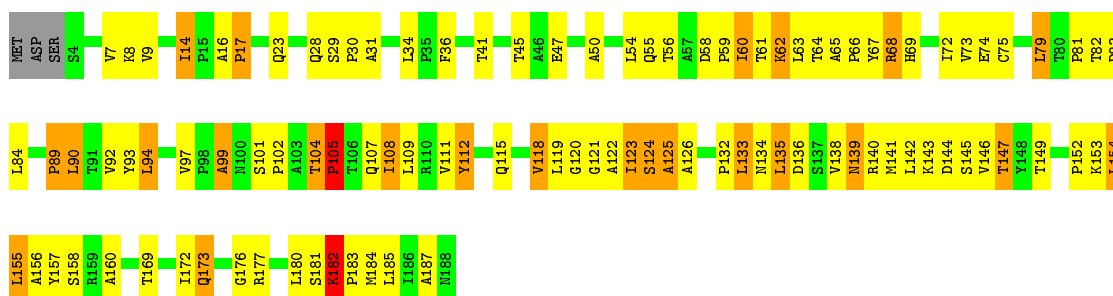
- Molecule 1: COAT PROTEIN

Chain B: 



- Molecule 1: COAT PROTEIN

Chain C: 



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	325.76Å 325.76Å 738.94Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.90 19.97 – 3.90	Depositor EDS
% Data completeness (in resolution range)	79.4 (20.00-3.90) 63.9 (19.97-3.90)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.70 (at 3.94Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.230 , 0.242 0.546 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	59.7	Xtrriage
Anisotropy	0.237	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	1.33 , -10.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.408 for h,-h-k,-l	Xtrriage
F_o, F_c correlation	0.10	EDS
Total number of atoms	3976	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.64% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	0.49	0/1236	0.73	1/1695 (0.1%)
1	B	0.51	0/1412	0.69	0/1936
1	C	0.53	0/1412	0.72	0/1936
All	All	0.51	0/4060	0.72	1/5567 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	121	GLY	N-CA-C	-5.71	98.84	113.10

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	1210	0	1254	78	0
1	B	1383	0	1435	99	0
1	C	1383	0	1435	127	0
All	All	3976	0	4124	294	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 36.

All (294) 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:B:89:PRO:HA	1:B:120:GLY:HA3	1.17	1.16
1:C:139:ASN:HD22	1:C:140:ARG:N	1.58	1.01
1:C:182:LYS:HE2	1:C:182:LYS:C	1.81	1.01
1:C:68:ARG:HG2	1:C:182:LYS:HZ2	1.28	0.98
1:C:31:ALA:HB2	1:C:181:SER:HB3	1.45	0.96
1:C:146:VAL:HG12	1:C:147:THR:H	1.31	0.94
1:A:138:VAL:HG21	1:A:152:PRO:HG3	1.50	0.93
1:B:11:GLN:HG2	1:B:12:ALA:H	1.33	0.91
1:A:159:ARG:HH11	1:A:159:ARG:HB3	1.34	0.89
1:C:68:ARG:HG2	1:C:182:LYS:NZ	1.90	0.87
1:B:139:ASN:HD22	1:B:140:ARG:N	1.73	0.86
1:B:32:ILE:HD13	1:B:33:VAL:H	1.40	0.83
1:B:139:ASN:HB2	1:B:150:ASP:HB2	1.60	0.82
1:A:94:LEU:HD21	1:A:172:ILE:HD13	1.61	0.80
1:C:139:ASN:ND2	1:C:141:MET:H	1.81	0.79
1:B:12:ALA:HB1	1:B:112:TYR:CE2	2.18	0.79
1:C:59:PRO:C	1:C:61:THR:H	1.85	0.79
1:C:146:VAL:HG12	1:C:147:THR:N	1.97	0.78
1:B:159:ARG:HB2	1:B:159:ARG:HH11	1.47	0.78
1:C:90:LEU:HD21	1:C:119:LEU:HB2	1.65	0.78
1:C:72:ILE:HD11	1:C:75:CYS:HB2	1.67	0.77
1:A:97:VAL:HG12	1:A:99:ALA:H	1.51	0.76
1:A:52:VAL:CG2	1:A:154:LEU:HB3	2.16	0.75
1:A:40:ALA:HB1	1:A:52:VAL:HG11	1.68	0.75
1:B:49:ALA:HB2	1:B:157:TYR:HD2	1.52	0.74
1:B:24:PRO:HA	1:B:27:GLU:CG	2.17	0.74
1:B:43:PHE:CE1	1:B:172:ILE:HD13	2.23	0.73
1:C:68:ARG:CG	1:C:182:LYS:HZ2	2.01	0.73
1:B:32:ILE:HD13	1:B:33:VAL:N	2.04	0.73
1:C:14:ILE:HD13	1:C:14:ILE:H	1.54	0.72
1:B:89:PRO:HA	1:B:120:GLY:CA	2.10	0.72
1:C:108:ILE:O	1:C:111:VAL:HG22	1.89	0.71
1:A:41:THR:HG22	1:A:42:THR:N	2.04	0.71
1:B:180:LEU:H	1:B:180:LEU:HD22	1.56	0.70
1:C:31:ALA:CB	1:C:181:SER:HB3	2.21	0.70
1:A:97:VAL:CG1	1:A:101:SER:HB3	2.21	0.70
1:C:182:LYS:HE2	1:C:182:LYS:O	1.91	0.70
1:C:64:THR:HG22	1:C:180:LEU:HD21	1.73	0.70
1:A:129:ILE:HD12	1:A:129:ILE:N	2.06	0.69
1:A:47:GLU:HB2	1:A:160:ALA:HB2	1.77	0.67
1:C:138:VAL:HG21	1:C:152:PRO:HB3	1.75	0.67
1:B:24:PRO:HA	1:B:27:GLU:HG3	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:PRO:HB3	1:A:119:LEU:HD13	1.77	0.67
1:C:139:ASN:HD22	1:C:140:ARG:H	1.40	0.67
1:B:159:ARG:HB2	1:B:159:ARG:NH1	2.10	0.66
1:B:37:GLN:HG3	1:B:175:SER:HB3	1.76	0.66
1:A:95:ALA:HB2	1:A:108:ILE:HG23	1.77	0.66
1:A:186:ILE:H	1:A:186:ILE:HD13	1.60	0.66
1:B:8:LYS:HG2	1:B:130:GLU:HB3	1.77	0.66
1:B:180:LEU:N	1:B:180:LEU:HD22	2.10	0.66
1:B:43:PHE:HE1	1:B:172:ILE:HD13	1.61	0.66
1:A:100:ASN:HD21	1:A:149:THR:HB	1.61	0.66
1:C:54:LEU:HD12	1:C:60:ILE:HD13	1.77	0.66
1:C:82:THR:HG22	1:C:84:LEU:H	1.59	0.66
1:A:97:VAL:HG11	1:A:101:SER:HB3	1.77	0.66
1:A:121:GLY:O	1:A:123:ILE:HG13	1.95	0.65
1:B:49:ALA:HB2	1:B:157:TYR:CD2	2.31	0.65
1:A:83:ASP:O	1:A:86:VAL:HG22	1.97	0.65
1:C:9:VAL:HG12	1:C:115:GLN:HE22	1.62	0.64
1:A:105:PRO:O	1:A:108:ILE:HG12	1.97	0.64
1:C:47:GLU:HB2	1:C:160:ALA:HB2	1.80	0.63
1:C:90:LEU:CD2	1:C:119:LEU:HB2	2.28	0.63
1:B:185:LEU:H	1:B:185:LEU:HD12	1.63	0.63
1:C:60:ILE:HD12	1:C:142:LEU:HD13	1.80	0.63
1:B:139:ASN:HD22	1:B:139:ASN:C	1.99	0.63
1:B:139:ASN:ND2	1:B:141:MET:H	1.96	0.63
1:C:139:ASN:HD22	1:C:139:ASN:C	1.98	0.63
1:B:185:LEU:N	1:B:185:LEU:HD12	2.14	0.63
1:B:150:ASP:OD2	1:C:145:SER:HB3	1.99	0.63
1:C:59:PRO:O	1:C:61:THR:N	2.32	0.62
1:A:108:ILE:HD11	1:A:155:LEU:HB3	1.82	0.62
1:C:79:LEU:HD22	1:C:119:LEU:HD11	1.81	0.62
1:A:41:THR:CG2	1:A:42:THR:N	2.63	0.62
1:A:98:PRO:HA	1:A:138:VAL:HG12	1.80	0.62
1:C:68:ARG:HG2	1:C:182:LYS:CD	2.30	0.61
1:A:34:LEU:HD23	1:A:34:LEU:H	1.64	0.61
1:B:104:THR:HG23	1:B:107:GLN:OE1	2.01	0.61
1:C:72:ILE:HD11	1:C:75:CYS:CB	2.30	0.61
1:B:79:LEU:HD22	1:B:119:LEU:HD21	1.83	0.61
1:C:123:ILE:HD13	1:C:123:ILE:O	2.01	0.60
1:A:129:ILE:HD12	1:A:129:ILE:H	1.65	0.60
1:A:93:TYR:O	1:A:156:ALA:HA	2.02	0.60
1:B:23:GLN:N	1:B:24:PRO:CD	2.64	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:11:GLN:HG2	1:B:12:ALA:N	2.10	0.60
1:C:139:ASN:ND2	1:C:140:ARG:N	2.41	0.60
1:B:55:GLN:HB2	1:B:151:THR:HG21	1.84	0.60
1:C:34:LEU:HD12	1:C:180:LEU:HD12	1.84	0.60
1:C:59:PRO:C	1:C:61:THR:N	2.53	0.60
1:B:138:VAL:HG21	1:B:152:PRO:HB3	1.82	0.59
1:C:16:ALA:N	1:C:17:PRO:HD3	2.18	0.59
1:C:41:THR:HG21	1:C:50:ALA:HB3	1.83	0.59
1:A:54:LEU:HD21	1:A:174:ILE:HG21	1.85	0.59
1:A:159:ARG:NH1	1:A:159:ARG:HB3	2.13	0.58
1:A:97:VAL:HG13	1:A:98:PRO:HD2	1.85	0.58
1:B:117:PHE:CD1	1:B:129:ILE:HD13	2.37	0.58
1:C:155:LEU:H	1:C:155:LEU:HD23	1.67	0.58
1:C:9:VAL:CG1	1:C:115:GLN:HE22	2.16	0.58
1:C:81:PRO:HG3	1:C:119:LEU:HD22	1.85	0.58
1:B:135:LEU:HD23	1:B:135:LEU:H	1.67	0.57
1:C:58:ASP:HB2	1:C:59:PRO:HA	1.85	0.57
1:A:98:PRO:O	1:A:99:ALA:HB3	2.05	0.57
1:B:41:THR:HG21	1:B:50:ALA:HB3	1.86	0.57
1:C:93:TYR:CZ	1:C:109:LEU:HD11	2.40	0.57
1:A:187:ALA:O	1:A:188:ASN:HB2	2.02	0.57
1:B:23:GLN:N	1:B:24:PRO:HD2	2.20	0.57
1:A:122:ALA:C	1:A:124:SER:H	2.08	0.57
1:C:93:TYR:CE2	1:C:109:LEU:HD11	2.40	0.57
1:C:82:THR:HB	1:C:169:THR:O	2.04	0.57
1:B:178:ILE:HG22	1:B:179:ARG:N	2.20	0.57
1:C:120:GLY:H	1:C:125:ALA:HB1	1.68	0.56
1:B:24:PRO:HA	1:B:27:GLU:HG2	1.87	0.56
1:C:146:VAL:CG1	1:C:147:THR:H	2.11	0.56
1:C:93:TYR:O	1:C:156:ALA:HA	2.05	0.56
1:B:67:TYR:CG	1:B:180:LEU:HB3	2.41	0.55
1:B:49:ALA:CB	1:B:157:TYR:HD2	2.17	0.55
1:C:36:PHE:CD2	1:C:60:ILE:HG22	2.41	0.55
1:A:42:THR:OG1	1:A:168:PRO:HB3	2.07	0.55
1:A:52:VAL:HG22	1:A:154:LEU:HB3	1.88	0.55
1:A:86:VAL:HA	1:A:121:GLY:H	1.72	0.55
1:B:123:ILE:HG13	1:B:124:SER:N	2.21	0.55
1:C:118:VAL:O	1:C:118:VAL:HG22	2.05	0.55
1:C:133:LEU:HD13	1:C:135:LEU:CD1	2.36	0.55
1:C:68:ARG:HG2	1:C:182:LYS:HD3	1.88	0.55
1:B:110:ARG:CB	1:B:110:ARG:HH11	2.19	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:TYR:O	1:B:156:ALA:HA	2.06	0.55
1:C:155:LEU:N	1:C:155:LEU:HD23	2.20	0.55
1:C:73:VAL:HG13	1:C:74:GLU:N	2.22	0.55
1:C:58:ASP:HB3	1:C:61:THR:OG1	2.06	0.55
1:B:120:GLY:O	1:B:121:GLY:C	2.44	0.55
1:B:16:ALA:HA	1:C:182:LYS:HG2	1.88	0.55
1:C:92:VAL:HG12	1:C:93:TYR:N	2.22	0.55
1:A:52:VAL:HG23	1:A:154:LEU:HB3	1.86	0.54
1:C:65:ALA:N	1:C:66:PRO:HD2	2.23	0.54
1:B:161:PRO:HG2	1:B:164:PRO:HG3	1.89	0.54
1:C:105:PRO:O	1:C:108:ILE:HG12	2.08	0.53
1:A:159:ARG:HH11	1:A:159:ARG:CB	2.13	0.53
1:A:186:ILE:N	1:A:186:ILE:HD13	2.23	0.53
1:C:120:GLY:N	1:C:125:ALA:HB1	2.23	0.53
1:C:68:ARG:HA	1:C:143:LYS:HE2	1.90	0.53
1:A:97:VAL:HG13	1:A:101:SER:HB3	1.91	0.52
1:B:119:LEU:HD12	1:B:119:LEU:H	1.74	0.52
1:A:41:THR:CG2	1:A:42:THR:H	2.22	0.52
1:C:73:VAL:CG1	1:C:177:ARG:HB2	2.39	0.52
1:C:104:THR:HG23	1:C:107:GLN:OE1	2.10	0.52
1:C:154:LEU:HD22	1:C:155:LEU:N	2.24	0.52
1:A:34:LEU:HD23	1:A:34:LEU:N	2.25	0.52
1:A:58:ASP:O	1:A:61:THR:HB	2.10	0.52
1:C:139:ASN:C	1:C:139:ASN:ND2	2.63	0.52
1:C:60:ILE:O	1:C:60:ILE:HG13	2.08	0.52
1:B:144:ASP:OD1	1:B:145:SER:N	2.38	0.51
1:C:67:TYR:CD2	1:C:180:LEU:HB3	2.44	0.51
1:B:159:ARG:CB	1:B:159:ARG:HH11	2.20	0.51
1:C:58:ASP:HA	1:C:61:THR:OG1	2.11	0.51
1:B:110:ARG:HH12	1:C:187:ALA:HB1	1.75	0.51
1:B:146:VAL:HG12	1:B:147:THR:N	2.26	0.50
1:A:157:TYR:CG	1:A:158:SER:N	2.79	0.50
1:B:55:GLN:CB	1:B:151:THR:HG21	2.41	0.50
1:A:122:ALA:O	1:A:124:SER:N	2.45	0.50
1:B:40:ALA:HB1	1:B:52:VAL:HG21	1.94	0.50
1:B:123:ILE:HG13	1:B:124:SER:H	1.77	0.49
1:A:108:ILE:CD1	1:A:155:LEU:HB3	2.42	0.49
1:B:178:ILE:CG2	1:B:179:ARG:N	2.76	0.49
1:B:19:SER:OG	1:C:23:GLN:NE2	2.44	0.49
1:B:19:SER:HB2	1:C:145:SER:HB2	1.94	0.49
1:C:14:ILE:N	1:C:14:ILE:HD13	2.23	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:182:LYS:HE2	1:C:183:PRO:N	2.26	0.49
1:A:120:GLY:C	1:A:122:ALA:N	2.61	0.49
1:B:150:ASP:OD1	1:C:68:ARG:HD2	2.13	0.49
1:B:32:ILE:CD1	1:B:33:VAL:N	2.74	0.49
1:B:98:PRO:HB3	1:B:138:VAL:HG12	1.95	0.49
1:C:111:VAL:O	1:C:111:VAL:HG23	2.13	0.49
1:A:129:ILE:H	1:A:129:ILE:CD1	2.26	0.49
1:B:16:ALA:HB1	1:C:68:ARG:NH2	2.28	0.49
1:C:94:LEU:HD11	1:C:172:ILE:HD13	1.95	0.49
1:C:157:TYR:CG	1:C:158:SER:N	2.81	0.48
1:C:29:SER:HB2	1:C:30:PRO:HD2	1.94	0.48
1:B:66:PRO:HG2	1:B:67:TYR:CE2	2.48	0.48
1:B:110:ARG:NH1	1:C:187:ALA:HB1	2.28	0.48
1:C:68:ARG:O	1:C:68:ARG:HG3	2.13	0.48
1:C:107:GLN:O	1:C:108:ILE:C	2.50	0.48
1:C:181:SER:O	1:C:182:LYS:O	2.31	0.48
1:B:34:LEU:HD13	1:B:35:PRO:HD2	1.95	0.48
1:B:9:VAL:HG23	1:B:9:VAL:O	2.13	0.48
1:C:133:LEU:HD11	1:C:152:PRO:HB3	1.96	0.48
1:B:100:ASN:HD22	1:B:149:THR:HB	1.79	0.48
1:C:82:THR:HG22	1:C:83:ASP:N	2.28	0.48
1:C:82:THR:CG2	1:C:83:ASP:N	2.76	0.47
1:C:134:ASN:OD1	1:C:136:ASP:HB2	2.13	0.47
1:A:100:ASN:ND2	1:A:149:THR:HB	2.28	0.47
1:A:185:LEU:HD12	1:A:185:LEU:N	2.29	0.47
1:B:12:ALA:HB1	1:B:112:TYR:CD2	2.49	0.47
1:B:79:LEU:HD22	1:B:119:LEU:CD2	2.45	0.47
1:B:93:TYR:CE2	1:B:116:SER:HB2	2.50	0.47
1:C:154:LEU:C	1:C:154:LEU:HD22	2.34	0.47
1:C:61:THR:C	1:C:63:LEU:N	2.68	0.46
1:B:119:LEU:HD12	1:B:119:LEU:N	2.31	0.46
1:A:110:ARG:HD2	1:B:188:ASN:OXT	2.16	0.46
1:C:68:ARG:NE	1:C:182:LYS:HZ2	2.13	0.46
1:C:185:LEU:H	1:C:185:LEU:HD12	1.80	0.46
1:A:97:VAL:HG13	1:A:98:PRO:CD	2.45	0.46
1:C:111:VAL:O	1:C:112:TYR:C	2.53	0.46
1:C:176:GLY:C	1:C:177:ARG:HG2	2.36	0.46
1:A:61:THR:C	1:A:63:LEU:N	2.69	0.46
1:A:41:THR:HG22	1:A:42:THR:H	1.79	0.46
1:B:14:ILE:HG23	1:B:14:ILE:O	2.16	0.46
1:B:58:ASP:N	1:B:59:PRO:HD2	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:182:LYS:CE	1:C:182:LYS:O	2.62	0.46
1:C:72:ILE:HG23	1:C:72:ILE:O	2.16	0.45
1:C:69:HIS:HA	1:C:143:LYS:O	2.16	0.45
1:B:94:LEU:N	1:B:94:LEU:HD12	2.31	0.45
1:A:90:LEU:N	1:A:90:LEU:HD23	2.31	0.45
1:A:68:ARG:O	1:A:68:ARG:HG2	2.17	0.45
1:A:67:TYR:CD2	1:A:180:LEU:HD13	2.51	0.45
1:B:135:LEU:CD2	1:B:135:LEU:H	2.30	0.45
1:C:72:ILE:HG22	1:C:140:ARG:O	2.17	0.45
1:C:8:LYS:HE2	1:C:132:PRO:HG3	1.99	0.45
1:B:65:ALA:HB3	1:B:66:PRO:HD3	1.99	0.44
1:B:94:LEU:O	1:B:95:ALA:HB2	2.17	0.44
1:C:97:VAL:HG22	1:C:155:LEU:CD2	2.47	0.44
1:A:81:PRO:HA	1:A:170:ALA:CB	2.47	0.44
1:B:139:ASN:ND2	1:B:139:ASN:C	2.69	0.44
1:C:185:LEU:HD12	1:C:185:LEU:N	2.32	0.44
1:B:61:THR:O	1:B:63:LEU:N	2.51	0.44
1:C:123:ILE:CG2	1:C:124:SER:N	2.80	0.44
1:C:16:ALA:N	1:C:17:PRO:CD	2.79	0.44
1:A:79:LEU:CD2	1:A:119:LEU:HD11	2.47	0.44
1:B:22:SER:C	1:B:24:PRO:HD2	2.38	0.44
1:A:47:GLU:HB2	1:A:160:ALA:CB	2.45	0.44
1:A:117:PHE:CE2	1:A:129:ILE:HG23	2.53	0.43
1:A:187:ALA:O	1:A:188:ASN:CB	2.66	0.43
1:C:92:VAL:CG1	1:C:93:TYR:N	2.81	0.43
1:B:14:ILE:HD13	1:B:14:ILE:O	2.18	0.43
1:B:16:ALA:N	1:B:17:PRO:CD	2.80	0.43
1:A:97:VAL:HA	1:A:98:PRO:HD3	1.87	0.43
1:C:65:ALA:N	1:C:66:PRO:CD	2.82	0.43
1:B:157:TYR:CG	1:B:158:SER:N	2.86	0.43
1:A:37:GLN:HA	1:A:174:ILE:O	2.19	0.43
1:C:14:ILE:CD1	1:C:14:ILE:N	2.82	0.43
1:C:184:MET:HG3	1:C:185:LEU:N	2.33	0.43
1:C:31:ALA:HB1	1:C:180:LEU:O	2.18	0.43
1:C:68:ARG:HG2	1:C:182:LYS:CE	2.49	0.43
1:A:138:VAL:HB	1:A:139:ASN:H	1.61	0.43
1:B:84:LEU:HB3	1:B:169:THR:HG22	2.01	0.43
1:A:98:PRO:O	1:A:150:ASP:HA	2.18	0.43
1:B:120:GLY:O	1:B:121:GLY:O	2.37	0.43
1:C:89:PRO:HB3	1:C:121:GLY:HA3	2.00	0.43
1:B:47:GLU:HB2	1:B:160:ALA:HB2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:92:VAL:O	1:B:93:TYR:HD2	2.02	0.43
1:B:185:LEU:CD1	1:B:185:LEU:H	2.30	0.43
1:C:61:THR:C	1:C:63:LEU:H	2.21	0.43
1:A:49:ALA:HB2	1:A:157:TYR:HD1	1.83	0.42
1:B:110:ARG:HB2	1:B:110:ARG:HH11	1.84	0.42
1:C:59:PRO:HA	1:C:62:LYS:HG3	2.00	0.42
1:C:7:VAL:HG13	1:C:7:VAL:O	2.17	0.42
1:C:125:ALA:O	1:C:126:ALA:HB3	2.18	0.42
1:A:163:ASN:CG	1:A:163:ASN:O	2.58	0.42
1:B:45:THR:HA	1:B:165:SER:O	2.19	0.42
1:C:99:ALA:HB3	1:C:149:THR:O	2.18	0.42
1:A:144:ASP:CG	1:A:145:SER:H	2.22	0.42
1:B:92:VAL:HG12	1:B:93:TYR:N	2.34	0.42
1:B:129:ILE:HG22	1:B:130:GLU:N	2.35	0.42
1:C:73:VAL:CG1	1:C:74:GLU:N	2.82	0.42
1:A:154:LEU:HD12	1:A:155:LEU:N	2.35	0.42
1:A:29:SER:OG	1:A:30:PRO:CD	2.68	0.42
1:C:101:SER:HA	1:C:102:PRO:HD3	1.71	0.42
1:C:133:LEU:HD13	1:C:135:LEU:HD12	2.01	0.42
1:C:54:LEU:O	1:C:56:THR:N	2.52	0.42
1:B:61:THR:C	1:B:63:LEU:N	2.73	0.42
1:C:97:VAL:O	1:C:153:LYS:HB3	2.20	0.42
1:C:173:GLN:HB3	1:C:173:GLN:HE21	1.65	0.42
1:A:140:ARG:H	1:A:140:ARG:HG3	1.65	0.41
1:A:65:ALA:N	1:A:66:PRO:CD	2.83	0.41
1:B:64:THR:HG22	1:B:180:LEU:HD11	2.03	0.41
1:B:66:PRO:HG2	1:B:67:TYR:CD2	2.55	0.41
1:C:144:ASP:CG	1:C:145:SER:H	2.23	0.41
1:C:14:ILE:CD1	1:C:14:ILE:H	2.30	0.41
1:C:58:ASP:CB	1:C:61:THR:OG1	2.69	0.41
1:A:120:GLY:C	1:A:122:ALA:H	2.21	0.41
1:B:76:LYS:HD3	1:B:130:GLU:HG2	2.03	0.41
1:C:157:TYR:CD2	1:C:158:SER:N	2.89	0.41
1:C:74:GLU:HG2	1:C:177:ARG:HG3	2.03	0.41
1:A:67:TYR:CD2	1:A:67:TYR:N	2.89	0.41
1:B:150:ASP:OD2	1:C:145:SER:CB	2.68	0.41
1:B:68:ARG:HG2	1:B:184:MET:HB2	2.02	0.40
1:A:95:ALA:CB	1:A:108:ILE:HG23	2.47	0.40
1:A:154:LEU:CD2	1:A:174:ILE:HD11	2.51	0.40
1:A:176:GLY:C	1:A:177:ARG:HG3	2.42	0.40
1:B:133:LEU:HD22	1:B:134:ASN:N	2.36	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:GLU:HB2	1:A:160:ALA:CA	2.52	0.40
1:A:94:LEU:HD21	1:A:172:ILE:CD1	2.42	0.40
1:C:146:VAL:CG1	1:C:147:THR:N	2.69	0.40
1:B:83:ASP:OD1	1:B:84:LEU:HD12	2.20	0.40
1:A:97:VAL:CG1	1:A:98:PRO:N	2.81	0.40
1:C:84:LEU:HB3	1:C:169:THR:HG22	2.03	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 X-ray entries followed by that with respect to entries of similar resolution.

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	159/188 (85%)	138 (87%)	17 (11%)	4 (2%)	5	36
1	B	183/188 (97%)	154 (84%)	26 (14%)	3 (2%)	9	44
1	C	183/188 (97%)	146 (80%)	24 (13%)	13 (7%)	1	17
All	All	525/564 (93%)	438 (83%)	67 (13%)	20 (4%)	3	28

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	123	ILE
1	A	164	PRO
1	C	55	GLN
1	C	122	ALA
1	C	125	ALA
1	A	122	ALA
1	A	125	ALA
1	B	95	ALA
1	B	121	GLY
1	C	60	ILE
1	C	99	ALA

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Mol	Chain	Res	Type
1	C	135	LEU
1	B	112	TYR
1	C	45	THR
1	C	17	PRO
1	C	108	ILE
1	C	112	TYR
1	C	182	LYS
1	C	105	PRO
1	C	89	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 X-ray entries followed by that with respect to entries of similar resolution.

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	135/159 (85%)	122 (90%)	13 (10%)	8	32
1	B	156/159 (98%)	139 (89%)	17 (11%)	6	28
1	C	156/159 (98%)	137 (88%)	19 (12%)	5	24
All	All	447/477 (94%)	398 (89%)	49 (11%)	6	28

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ILE
1	A	35	PRO
1	A	58	ASP
1	A	79	LEU
1	A	83	ASP
1	A	88	ASN
1	A	115	GLN
1	A	147	THR
1	A	149	THR
1	A	159	ARG
1	A	162	THR
1	A	173	GLN
1	A	186	ILE

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Mol	Chain	Res	Type
1	B	14	ILE
1	B	21	LEU
1	B	23	GLN
1	B	27	GLU
1	B	32	ILE
1	B	54	LEU
1	B	76	LYS
1	B	104	THR
1	B	109	LEU
1	B	110	ARG
1	B	115	GLN
1	B	133	LEU
1	B	136	ASP
1	B	139	ASN
1	B	142	LEU
1	B	149	THR
1	B	159	ARG
1	C	14	ILE
1	C	28	GLN
1	C	62	LYS
1	C	68	ARG
1	C	79	LEU
1	C	90	LEU
1	C	94	LEU
1	C	104	THR
1	C	105	PRO
1	C	118	VAL
1	C	123	ILE
1	C	124	SER
1	C	133	LEU
1	C	139	ASN
1	C	147	THR
1	C	154	LEU
1	C	155	LEU
1	C	173	GLN
1	C	182	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	GLN
1	A	69	HIS

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Mol	Chain	Res	Type
1	A	100	ASN
1	B	69	HIS
1	B	71	GLN
1	B	88	ASN
1	B	115	GLN
1	B	139	ASN
1	B	173	GLN
1	C	23	GLN
1	C	37	GLN
1	C	69	HIS
1	C	115	GLN
1	C	139	ASN
1	C	173	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.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.