



Full wwPDB X-ray Structure Validation Report ⓘ

May 28, 2020 – 07:54 pm BST

PDB ID : 1JL5
Title : Novel Molecular Architecture of YopM-a Leucine-rich Effector Protein from *Yersinia pestis*
Authors : Evdokimov, A.G.; Anderson, D.E.; Routzahn, K.M.; Waugh, D.S.
Deposited on : 2001-07-15
Resolution : 2.10 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

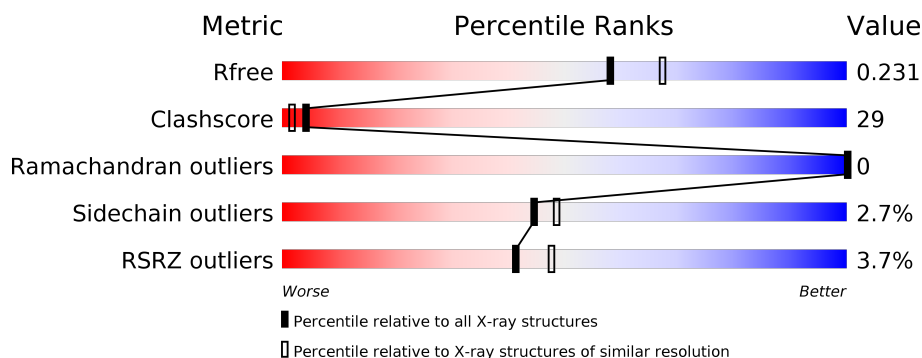
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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))
R_{free}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	454	<div> <div>3%</div> <div>37%</div> <div>40%</div> <div>•</div> <div>22%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3137 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 outer protein YopM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	353	Total	C	N	O	S	0	0	0
			2793	1773	457	558	5			

There are 45 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1385	ASN	-	SEE REMARK 999	UNP P17778
A	1386	SER	-	SEE REMARK 999	UNP P17778
A	1387	HIS	-	SEE REMARK 999	UNP P17778
A	1388	LEU	-	SEE REMARK 999	UNP P17778
A	1389	ALA	-	SEE REMARK 999	UNP P17778
A	1390	GLU	-	SEE REMARK 999	UNP P17778
A	1391	VAL	-	SEE REMARK 999	UNP P17778
A	1392	PRO	-	SEE REMARK 999	UNP P17778
A	1393	GLU	-	SEE REMARK 999	UNP P17778
A	1394	LEU	-	SEE REMARK 999	UNP P17778
A	1395	PRO	-	SEE REMARK 999	UNP P17778
A	1396	GLN	-	SEE REMARK 999	UNP P17778
A	1397	ASN	-	SEE REMARK 999	UNP P17778
A	1398	LEU	-	SEE REMARK 999	UNP P17778
A	1399	LYS	-	SEE REMARK 999	UNP P17778
A	1400	GLN	-	SEE REMARK 999	UNP P17778
A	1401	LEU	-	SEE REMARK 999	UNP P17778
A	1402	HIS	-	SEE REMARK 999	UNP P17778
A	1403	VAL	-	SEE REMARK 999	UNP P17778
A	1404	GLU	-	SEE REMARK 999	UNP P17778
A	1405	THR	-	SEE REMARK 999	UNP P17778
A	1406	ASN	-	SEE REMARK 999	UNP P17778
A	1407	PRO	-	SEE REMARK 999	UNP P17778
A	1408	LEU	-	SEE REMARK 999	UNP P17778
A	1409	ARG	-	SEE REMARK 999	UNP P17778
A	1410	GLU	-	SEE REMARK 999	UNP P17778
A	1411	PHE	-	SEE REMARK 999	UNP P17778

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1412	PRO	-	SEE REMARK 999	UNP P17778
A	1413	ASP	-	SEE REMARK 999	UNP P17778
A	1414	ILE	-	SEE REMARK 999	UNP P17778
A	1415	PRO	-	SEE REMARK 999	UNP P17778
A	1416	GLU	-	SEE REMARK 999	UNP P17778
A	1417	SER	-	SEE REMARK 999	UNP P17778
A	1418	VAL	-	SEE REMARK 999	UNP P17778
A	1419	GLU	-	SEE REMARK 999	UNP P17778
A	1420	ASP	-	SEE REMARK 999	UNP P17778
A	1421	LEU	-	SEE REMARK 999	UNP P17778
A	1422	ARG	-	SEE REMARK 999	UNP P17778
A	1423	MET	-	SEE REMARK 999	UNP P17778
A	1449	HIS	-	EXPRESSION TAG	UNP P17778
A	1450	HIS	-	EXPRESSION TAG	UNP P17778
A	1451	HIS	-	EXPRESSION TAG	UNP P17778
A	1452	HIS	-	EXPRESSION TAG	UNP P17778
A	1453	HIS	-	EXPRESSION TAG	UNP P17778
A	1454	HIS	-	EXPRESSION TAG	UNP P17778

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	5	Total Ca 5 5	0	0

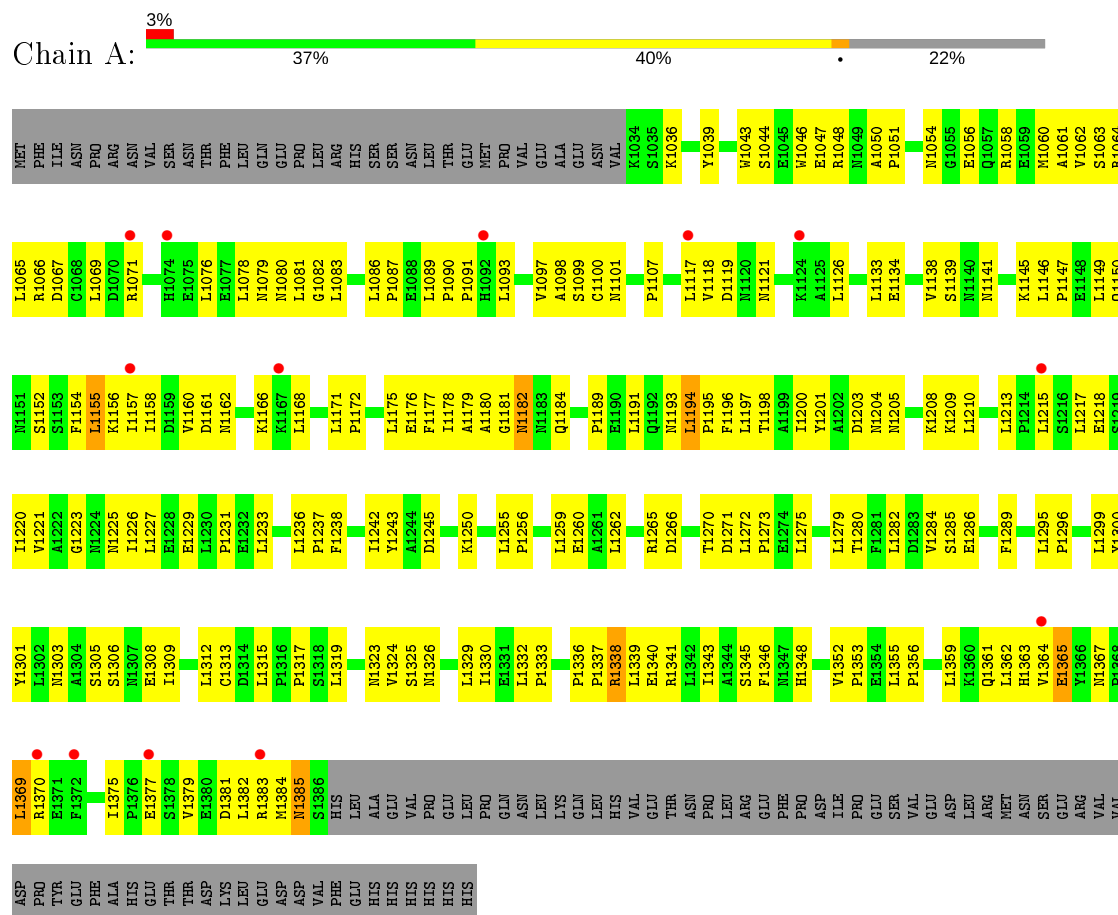
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	339	Total O 339 339	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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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: outer protein YopM



4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	149.43 Å 149.43 Å 191.18 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	70.89 – 2.10 70.89 – 2.02	Depositor EDS
% Data completeness (in resolution range)	92.7 (70.89-2.10) 91.9 (70.89-2.02)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.50 (at 2.03 Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.220 , 0.240 0.221 , 0.231	Depositor DCC
R_{free} test set	3416 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	23.6	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 45.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3137	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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](#)

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

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	0.41	0/2853	0.72	0/3903

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	2793	0	2784	164	0
2	A	5	0	0	0	0
3	A	339	0	0	9	0
All	All	3137	0	2784	164	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 29.

All (164) 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:1158:ILE:HD13	1:A:1175:LEU:HD13	1.42	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1312:LEU:HD11	1:A:1324:VAL:HG11	1.51	0.93
1:A:1155:LEU:HD13	1:A:1158:ILE:HD11	1.57	0.87
1:A:1158:ILE:HD13	1:A:1175:LEU:CD1	2.08	0.83
1:A:1146:LEU:HD11	1:A:1160:VAL:HG11	1.64	0.78
1:A:1179:ALA:HA	1:A:1201:TYR:HB2	1.66	0.77
1:A:1191:LEU:HA	1:A:1194:LEU:HD22	1.65	0.77
1:A:1044:SER:O	1:A:1048:ARG:HG3	1.87	0.74
1:A:1266:ASP:HA	1:A:1286:GLU:O	1.88	0.73
1:A:1339:LEU:HB3	1:A:1356:PRO:HG3	1.68	0.73
1:A:1080:ASN:HA	1:A:1100:CYS:O	1.89	0.72
1:A:1345:SER:HB3	1:A:1365:GLU:HG3	1.69	0.72
1:A:1191:LEU:HD13	1:A:1200:ILE:HD13	1.72	0.71
1:A:1367:ASN:HB3	1:A:1369:LEU:HD13	1.73	0.70
1:A:1086:LEU:HB2	1:A:1107:PRO:HD3	1.74	0.69
1:A:1319:LEU:HB3	1:A:1336:PRO:HG3	1.75	0.69
1:A:1175:LEU:HD21	1:A:1178:ILE:HB	1.77	0.67
1:A:1301:TYR:HB2	3:A:321:HOH:O	1.94	0.67
1:A:1364:VAL:HG12	1:A:1364:VAL:O	1.95	0.66
1:A:1325:SER:HB2	1:A:1343:ILE:HG22	1.79	0.65
1:A:1330:ILE:HA	1:A:1348:HIS:O	1.97	0.65
1:A:1217:LEU:HD21	1:A:1220:ILE:HB	1.77	0.64
1:A:1155:LEU:HD22	1:A:1158:ILE:CD1	2.28	0.64
1:A:1079:ASN:OD1	1:A:1097:VAL:HG12	1.98	0.64
1:A:1050:ALA:HB2	1:A:1058:ARG:HD2	1.80	0.63
1:A:1218:GLU:HA	1:A:1238:PHE:O	1.98	0.63
1:A:1272:LEU:HD11	1:A:1284:VAL:HG11	1.81	0.63
1:A:1355:LEU:HD11	1:A:1362:LEU:HD22	1.80	0.63
1:A:1134:GLU:HA	1:A:1154:PHE:O	1.99	0.62
1:A:1171:LEU:HD13	1:A:1175:LEU:HD23	1.80	0.62
1:A:1056:GLU:HG2	1:A:1082:GLY:O	2.00	0.62
1:A:1221:VAL:HA	1:A:1243:TYR:HB2	1.82	0.61
1:A:1171:LEU:HD13	1:A:1175:LEU:CD2	2.30	0.61
1:A:1168:LEU:HB2	1:A:1189:PRO:HD3	1.83	0.61
1:A:1060:MET:O	1:A:1064:ARG:HG2	2.00	0.61
1:A:1210:LEU:HB2	1:A:1231:PRO:HD3	1.81	0.61
1:A:1060:MET:HE2	1:A:1064:ARG:HG2	1.84	0.60
1:A:1326:ASN:HA	1:A:1346:PHE:O	2.02	0.59
1:A:1039:TYR:CZ	1:A:1091:PRO:HD3	2.37	0.58
1:A:1043:TRP:HH2	1:A:1069:LEU:HB2	1.68	0.58
1:A:1250:LYS:N	1:A:1250:LYS:HD2	2.19	0.57
1:A:1062:VAL:O	1:A:1066:ARG:HG3	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1101:ASN:HB2	1:A:1121:ASN:OD1	2.06	0.56
1:A:1194:LEU:HD23	1:A:1197:LEU:HD22	1.88	0.56
1:A:1047:GLU:HB2	1:A:1062:VAL:HG21	1.88	0.56
1:A:1205:ASN:HB2	1:A:1225:ASN:OD1	2.07	0.55
1:A:1337:PRO:HA	1:A:1356:PRO:HB3	1.88	0.55
1:A:1370:ARG:HH11	1:A:1370:ARG:HG2	1.71	0.55
1:A:1126:LEU:HB2	1:A:1147:PRO:HD3	1.88	0.55
1:A:1155:LEU:HD22	1:A:1158:ILE:HD12	1.89	0.55
1:A:1330:ILE:HG22	1:A:1348:HIS:HB3	1.90	0.54
1:A:1312:LEU:HB2	1:A:1333:PRO:HD3	1.90	0.54
1:A:1279:LEU:HD21	1:A:1282:LEU:HD13	1.90	0.54
1:A:1213:LEU:HB2	1:A:1236:LEU:HD11	1.89	0.53
1:A:1270:THR:HG23	1:A:1271:ASP:OD2	2.08	0.53
1:A:1315:LEU:HB2	1:A:1336:PRO:HD3	1.90	0.53
1:A:1166:LYS:CD	1:A:1184:GLN:HE21	2.21	0.53
1:A:1209:LYS:HG2	3:A:241:HOH:O	2.09	0.53
1:A:1138:VAL:CG2	1:A:1160:VAL:HG12	2.38	0.53
1:A:1279:LEU:HD23	1:A:1296:PRO:HG2	1.91	0.52
1:A:1176:GLU:HA	1:A:1196:PHE:O	2.09	0.52
1:A:1385:ASN:C	1:A:1385:ASN:HD22	2.13	0.52
1:A:1215:LEU:HD12	1:A:1237:PRO:HD2	1.91	0.52
1:A:1210:LEU:HD13	1:A:1220:ILE:CD1	2.40	0.52
1:A:1352:VAL:CG2	1:A:1364:VAL:HG11	2.40	0.52
1:A:1046:TRP:CE2	1:A:1087:PRO:HD3	2.45	0.52
1:A:1265:ARG:NH2	1:A:1286:GLU:OE2	2.43	0.52
1:A:1126:LEU:HD11	1:A:1138:VAL:HG11	1.92	0.52
1:A:1367:ASN:CB	1:A:1369:LEU:HD13	2.39	0.52
1:A:1155:LEU:HD22	1:A:1158:ILE:HD11	1.91	0.51
1:A:1060:MET:O	1:A:1060:MET:HE2	2.10	0.51
1:A:1146:LEU:CD1	1:A:1160:VAL:HG11	2.38	0.51
1:A:1098:ALA:HB3	1:A:1118:VAL:HG12	1.93	0.51
1:A:1284:VAL:HB	1:A:1289:PHE:CE2	2.45	0.51
1:A:1191:LEU:HD13	1:A:1200:ILE:CD1	2.39	0.50
1:A:1355:LEU:CD1	1:A:1362:LEU:HD22	2.40	0.50
1:A:1139:SER:HA	1:A:1161:ASP:O	2.12	0.50
1:A:1161:ASP:HB2	3:A:23:HOH:O	2.12	0.50
1:A:1210:LEU:HD13	1:A:1220:ILE:HD11	1.92	0.50
1:A:1119:ASP:HA	1:A:1139:SER:O	2.10	0.50
1:A:1117:LEU:HD13	3:A:252:HOH:O	2.11	0.50
1:A:1036:LYS:HA	1:A:1069:LEU:HD21	1.94	0.50
1:A:1051:PRO:HD2	1:A:1056:GLU:OE2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1301:TYR:CZ	1:A:1303:ASN:HB2	2.47	0.49
1:A:1280:THR:HA	1:A:1299:LEU:HA	1.93	0.49
1:A:1060:MET:HE3	1:A:1063:SER:HB3	1.93	0.49
1:A:1081:LEU:HB2	1:A:1083:LEU:HG	1.93	0.49
1:A:1198:THR:HB	1:A:1218:GLU:HG3	1.95	0.49
1:A:1152:SER:CB	1:A:1155:LEU:HB2	2.42	0.48
1:A:1343:ILE:N	1:A:1343:ILE:HD12	2.27	0.48
1:A:1158:ILE:O	1:A:1178:ILE:HA	2.13	0.48
1:A:1177:PHE:C	1:A:1177:PHE:CD1	2.87	0.48
1:A:1336:PRO:HG2	1:A:1339:LEU:HB2	1.95	0.48
1:A:1229:GLU:HG2	3:A:241:HOH:O	2.14	0.48
1:A:1168:LEU:HD11	1:A:1180:ALA:CB	2.44	0.47
1:A:1152:SER:HB2	1:A:1155:LEU:HB2	1.96	0.47
1:A:1355:LEU:HD22	1:A:1359:LEU:HD22	1.96	0.47
1:A:1259:LEU:HD21	1:A:1262:LEU:HB2	1.96	0.47
1:A:1166:LYS:HD3	1:A:1184:GLN:NE2	2.29	0.47
1:A:1223:GLY:HA2	1:A:1245:ASP:O	2.13	0.47
1:A:1345:SER:CB	1:A:1365:GLU:HG3	2.43	0.47
1:A:1121:ASN:HB2	1:A:1141:ASN:OD1	2.15	0.47
1:A:1157:ILE:C	1:A:1158:ILE:HD12	2.35	0.46
1:A:1272:LEU:HG	1:A:1289:PHE:CZ	2.50	0.46
1:A:1157:ILE:HG23	1:A:1177:PHE:CD1	2.51	0.46
1:A:1155:LEU:CD1	1:A:1158:ILE:HD11	2.38	0.46
1:A:1166:LYS:HD3	1:A:1184:GLN:HE21	1.80	0.46
1:A:1295:LEU:HG	1:A:1313:CYS:SG	2.56	0.46
1:A:1145:LYS:HE3	3:A:332:HOH:O	2.15	0.46
1:A:1279:LEU:HB3	1:A:1296:PRO:HG2	1.98	0.46
1:A:1370:ARG:NH1	1:A:1370:ARG:HG2	2.31	0.46
1:A:1182:ASN:HB2	3:A:132:HOH:O	2.16	0.45
1:A:1356:PRO:HB2	3:A:75:HOH:O	2.17	0.45
1:A:1233:LEU:HD13	1:A:1242:ILE:HD13	1.98	0.45
1:A:1352:VAL:HG22	1:A:1364:VAL:HG11	1.97	0.45
1:A:1308:GLU:OE1	1:A:1308:GLU:HA	2.16	0.45
1:A:1341:ARG:NH1	1:A:1361:GLN:OE1	2.49	0.45
1:A:1306:SER:HA	1:A:1326:ASN:O	2.17	0.45
1:A:1275:LEU:HB3	1:A:1296:PRO:HG3	1.98	0.44
1:A:1182:ASN:HA	1:A:1204:ASN:O	2.17	0.44
1:A:1233:LEU:HB3	1:A:1256:PRO:HG3	1.99	0.44
1:A:1149:LEU:O	1:A:1172:PRO:HB3	2.17	0.44
1:A:1233:LEU:HD13	1:A:1242:ILE:CD1	2.48	0.44
1:A:1305:SER:HA	1:A:1325:SER:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1076:LEU:N	1:A:1093:LEU:HD11	2.32	0.44
1:A:1309:ILE:HG22	1:A:1329:LEU:HD21	2.00	0.44
1:A:1208:LYS:HA	1:A:1226:ILE:O	2.18	0.44
1:A:1265:ARG:HG3	1:A:1285:SER:OG	2.18	0.43
1:A:1332:LEU:HB2	1:A:1353:PRO:CD	2.48	0.43
1:A:1157:ILE:N	1:A:1157:ILE:HD12	2.32	0.43
1:A:1061:ALA:HA	1:A:1078:LEU:HD22	2.00	0.43
1:A:1046:TRP:CG	1:A:1087:PRO:HB3	2.54	0.43
1:A:1150:GLN:NE2	3:A:340:HOH:O	2.51	0.43
1:A:1286:GLU:HA	1:A:1306:SER:O	2.19	0.43
1:A:1067:ASP:OD1	1:A:1071:ARG:NH1	2.51	0.43
1:A:1161:ASP:OD1	1:A:1162:ASN:HB2	2.18	0.43
1:A:1062:VAL:HA	1:A:1065:LEU:HD12	2.01	0.43
1:A:1043:TRP:CH2	1:A:1069:LEU:HB2	2.50	0.43
1:A:1181:GLY:HA2	1:A:1203:ASP:O	2.19	0.43
1:A:1089:LEU:O	1:A:1090:PRO:C	2.57	0.42
1:A:1175:LEU:HG	1:A:1197:LEU:HD13	2.00	0.42
1:A:1303:ASN:HA	1:A:1323:ASN:HB3	2.01	0.42
1:A:1058:ARG:O	1:A:1062:VAL:HG23	2.20	0.42
1:A:1166:LYS:HD2	1:A:1184:GLN:HE21	1.85	0.42
1:A:1193:ASN:O	1:A:1195:PRO:HD3	2.20	0.42
1:A:1086:LEU:HB2	1:A:1107:PRO:CD	2.47	0.42
1:A:1156:LYS:HB2	1:A:1157:ILE:HD12	2.01	0.42
1:A:1340:GLU:C	1:A:1359:LEU:HD12	2.40	0.41
1:A:1168:LEU:HD13	1:A:1178:ILE:HD11	2.02	0.41
1:A:1317:PRO:O	1:A:1338:ARG:HB2	2.20	0.41
1:A:1162:ASN:HA	1:A:1182:ASN:O	2.19	0.41
1:A:1260:GLU:C	1:A:1279:LEU:HD12	2.41	0.41
1:A:1332:LEU:HB2	1:A:1353:PRO:CG	2.51	0.41
1:A:1361:GLN:HA	1:A:1381:ASP:O	2.21	0.41
1:A:1079:ASN:HA	1:A:1099:SER:O	2.21	0.41
1:A:1208:LYS:O	1:A:1227:LEU:HA	2.20	0.41
1:A:1382:LEU:HD21	1:A:1384:MET:HE1	2.03	0.41
1:A:1213:LEU:HB2	1:A:1236:LEU:CD1	2.51	0.41
1:A:1275:LEU:CB	1:A:1296:PRO:HG3	2.51	0.41
1:A:1280:THR:HB	1:A:1300:TYR:CD2	2.56	0.41
1:A:1133:LEU:HA	1:A:1133:LEU:HD12	1.90	0.40
1:A:1255:LEU:HG	1:A:1273:PRO:HG2	2.02	0.40
1:A:1363:HIS:CD2	1:A:1383:ARG:HG3	2.55	0.40
1:A:1375:ILE:HG23	1:A:1379:VAL:HG21	2.04	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 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	351/454 (77%)	324 (92%)	27 (8%)	0	100	100

There are no Ramachandran outliers to report.

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 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	329/427 (77%)	320 (97%)	9 (3%)	44	48

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1054	ASN
1	A	1155	LEU
1	A	1182	ASN
1	A	1194	LEU
1	A	1338	ARG
1	A	1365	GLU
1	A	1369	LEU
1	A	1377	GLU
1	A	1385	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1054	ASN
1	A	1122	ASN
1	A	1182	ASN
1	A	1184	GLN
1	A	1193	ASN
1	A	1326	ASN
1	A	1385	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	353/454 (77%)	0.56	13 (3%)	41 48	16, 30, 42, 53	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1074	HIS	3.1
1	A	1370	ARG	2.8
1	A	1167	LYS	2.4
1	A	1215	LEU	2.4
1	A	1364	VAL	2.3
1	A	1117	LEU	2.3
1	A	1377	GLU	2.2
1	A	1071	ARG	2.2
1	A	1124	LYS	2.0
1	A	1157	ILE	2.0
1	A	1092	HIS	2.0
1	A	1372	PHE	2.0
1	A	1383	ARG	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	CA	A	2005	1/1	0.77	0.24	78,78,78,78	0
2	CA	A	2004	1/1	0.78	0.14	62,62,62,62	0
2	CA	A	2003	1/1	0.95	0.07	70,70,70,70	0
2	CA	A	2001	1/1	0.99	0.13	18,18,18,18	0
2	CA	A	2002	1/1	0.99	0.14	19,19,19,19	0

6.5 Other polymers

There are no such residues in this entry.