



# Full wwPDB X-ray Structure Validation Report ⓘ

May 12, 2020 – 11:33 pm BST

PDB ID : 6JKY  
Title : Crystal structure of MvcA-UBE2N-Ub complex from Legionella pneumophila  
Authors : Ouyang, S.Y.; Guan, H.  
Deposited on : 2019-03-03  
Resolution : 2.45 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](mailto: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

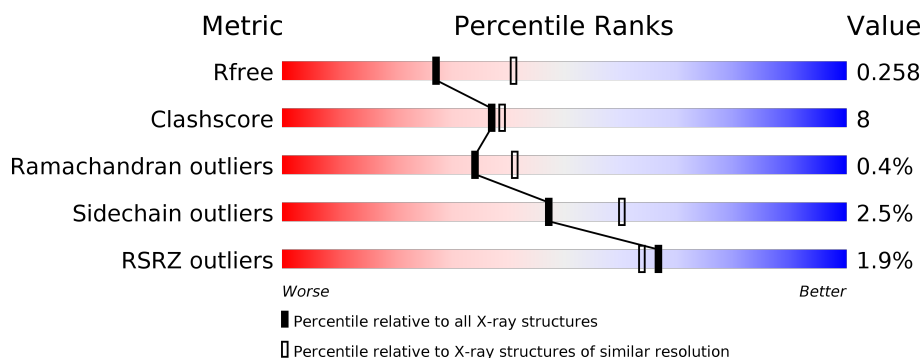
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.45 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	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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  $\geq 3$ , 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	385	<div> <div>82%</div> <div>18%</div> </div>
1	D	385	<div> <div>6%</div> <div>79%</div> <div>19%</div> <div>• •</div> </div>
2	B	152	<div> <div>6%</div> <div>79%</div> <div>16%</div> <div>• •</div> </div>
2	E	152	<div> <div>6%</div> <div>66%</div> <div>30%</div> <div>• •</div> </div>
3	C	79	<div> <div>84%</div> <div>13%</div> <div>•</div> </div>
3	F	79	<div> <div>84%</div> <div>10%</div> <div>5%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9776 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 Mvca.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	384	Total	C	N	O	S	0	0	0
			3091	1967	514	596	14			
1	D	381	Total	C	N	O	S	0	0	0
			3069	1954	511	590	14			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	11	ALA	-	expression tag	UNP Q5ZTL3
A	12	SER	-	expression tag	UNP Q5ZTL3
A	83	ALA	CYS	engineered mutation	UNP Q5ZTL3
D	11	ALA	-	expression tag	UNP Q5ZTL3
D	12	SER	-	expression tag	UNP Q5ZTL3
D	83	ALA	CYS	engineered mutation	UNP Q5ZTL3

- Molecule 2 is a protein called Ubiquitin-conjugating enzyme E2 N.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	149	Total	C	N	O	S	0	0	0
			1179	757	203	215	4			
2	E	148	Total	C	N	O	S	0	0	0
			1174	754	202	214	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	94	ALA	LYS	engineered mutation	UNP P61088
E	94	ALA	LYS	engineered mutation	UNP P61088

- Molecule 3 is a protein called Ub.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	76	Total	C	N	O	S	0	0	0
			602	379	104	118	1			
3	F	75	Total	C	N	O	S	0	0	0
			598	377	103	117	1			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	CYS	-	expression tag	UNP A0A3P7ZMV6
C	-1	GLY	-	expression tag	UNP A0A3P7ZMV6
C	0	SER	-	expression tag	UNP A0A3P7ZMV6
F	-2	CYS	-	expression tag	UNP A0A3P7ZMV6
F	-1	GLY	-	expression tag	UNP A0A3P7ZMV6
F	0	SER	-	expression tag	UNP A0A3P7ZMV6

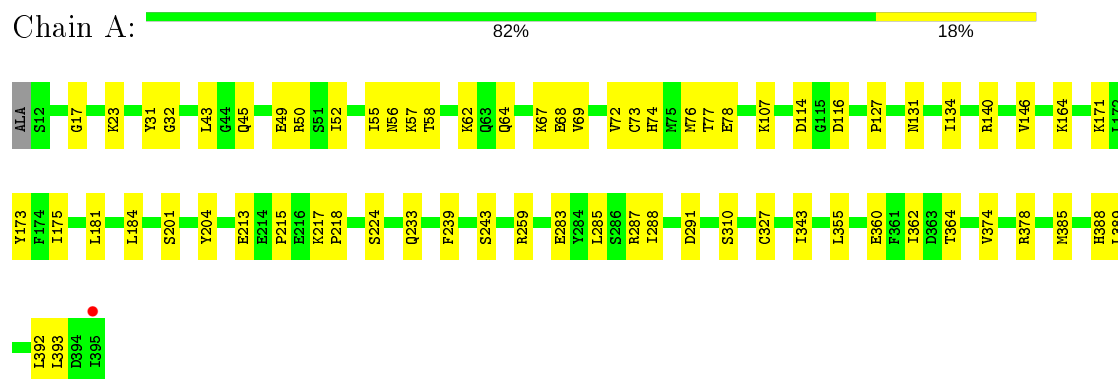
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	39	Total	O	0	0
			39	39		
4	B	5	Total	O	0	0
			5	5		
4	C	5	Total	O	0	0
			5	5		
4	D	10	Total	O	0	0
			10	10		
4	E	3	Total	O	0	0
			3	3		
4	F	1	Total	O	0	0
			1	1		

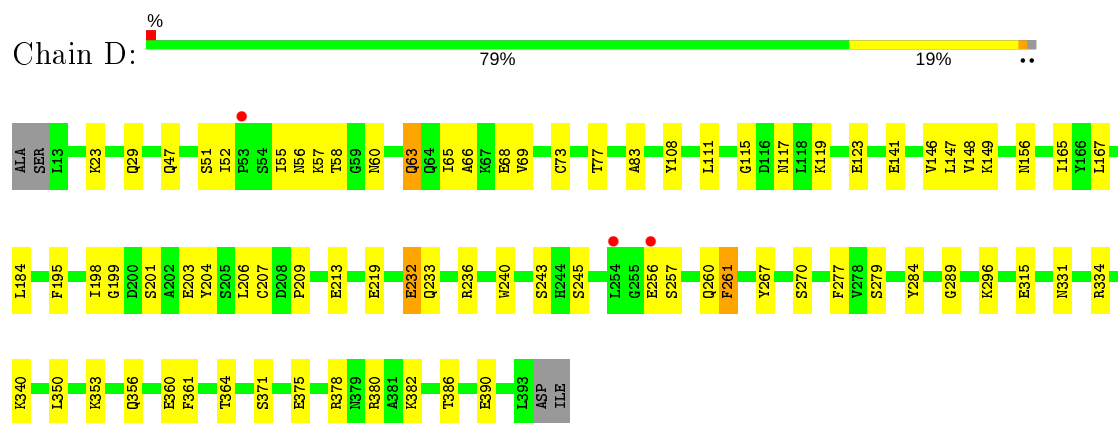
### 3 Residue-property plots [i](#)

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.

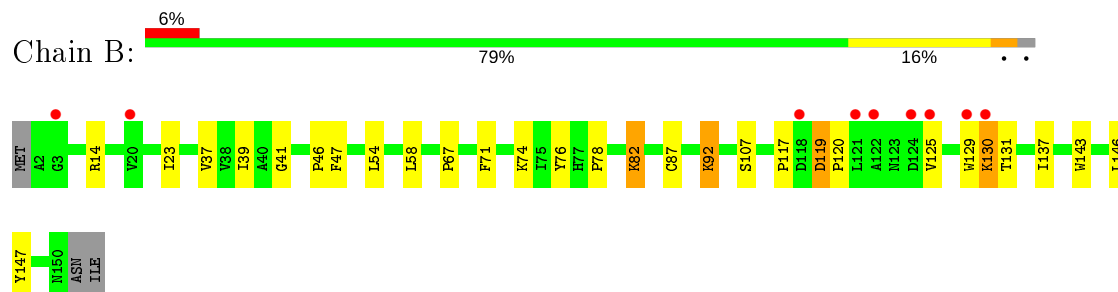
#### • Molecule 1: MvCA



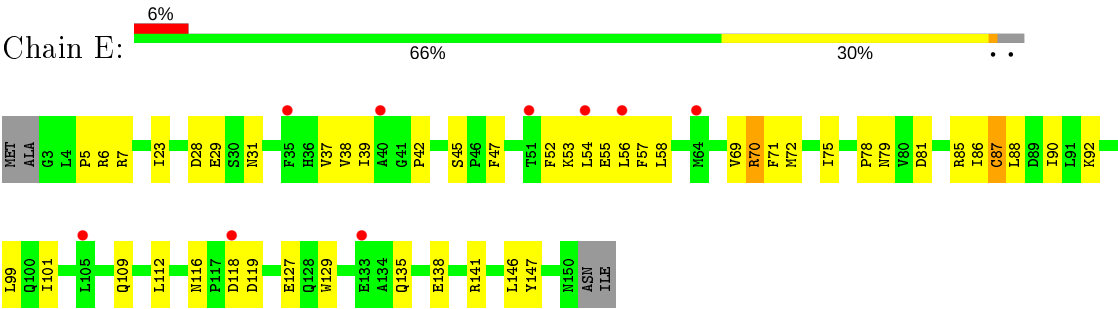
#### • Molecule 1: MvCA



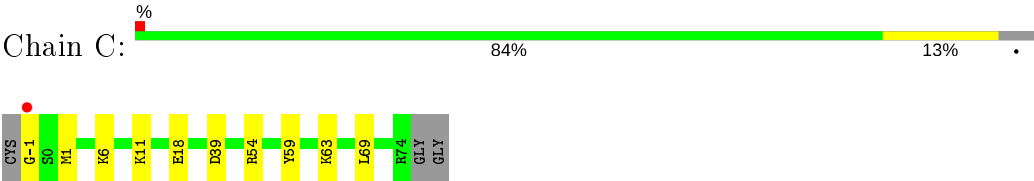
#### • Molecule 2: Ubiquitin-conjugating enzyme E2 N



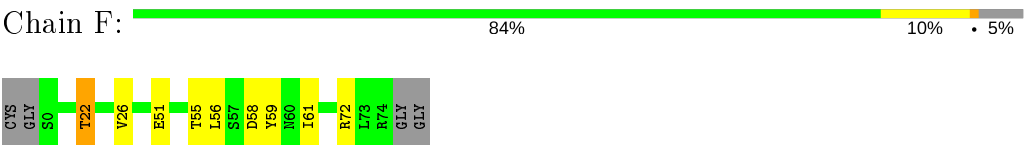
#### • Molecule 2: Ubiquitin-conjugating enzyme E2 N



● Molecule 3: Ub



● Molecule 3: Ub



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	104.33Å 107.61Å 263.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	65.14 – 2.45 65.14 – 2.45	Depositor EDS
% Data completeness (in resolution range)	96.4 (65.14-2.45) 96.4 (65.14-2.45)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.21 (at 2.45Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, $R_{free}$	0.209 , 0.258 0.208 , 0.258	Depositor DCC
$R_{free}$ test set	2639 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	58.5	Xtriage
Anisotropy	0.627	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 46.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.067 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	9776	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.28% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

### 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	0.26	0/3157	0.43	0/4258
1	D	0.26	0/3135	0.46	0/4228
2	B	0.33	0/1209	0.51	1/1648 (0.1%)
2	E	0.33	0/1204	0.61	1/1641 (0.1%)
3	C	0.32	0/608	0.48	0/817
3	F	0.41	0/604	0.52	0/812
All	All	0.29	0/9917	0.49	2/13404 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	92	LYS	CD-CE-NZ	6.95	127.69	111.70
2	E	92	LYS	CB-CG-CD	5.41	125.66	111.60

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	3091	0	3053	38	0
1	D	3069	0	3033	51	0
2	B	1179	0	1185	20	0
2	E	1174	0	1180	37	0
3	C	602	0	629	6	0

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	598	0	626	4	0
4	A	39	0	0	0	0
4	B	5	0	0	0	0
4	C	5	0	0	0	0
4	D	10	0	0	1	0
4	E	3	0	0	0	0
4	F	1	0	0	0	0
All	All	9776	0	9706	152	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 8.

All (152) 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)
2:E:56:LEU:HG	2:E:69:VAL:HG12	1.57	0.87
1:A:78:GLU:OE1	1:A:378:ARG:NH2	2.12	0.83
1:A:52:ILE:HD11	1:A:393:LEU:HD21	1.69	0.75
1:D:69:VAL:HG11	1:D:386:THR:HG22	1.67	0.74
1:A:134:ILE:HG12	1:A:343:ILE:CD1	2.19	0.72
2:B:46:PRO:HG3	2:B:137:ILE:HG23	1.71	0.72
1:D:47:GLN:NE2	1:D:51:SER:OG	2.23	0.71
2:E:146:LEU:HD12	2:E:147:TYR:CE2	2.24	0.71
2:E:79:ASN:ND2	2:E:119:ASP:OD2	2.23	0.71
2:E:146:LEU:HD12	2:E:147:TYR:CZ	2.27	0.70
2:E:58:LEU:HD21	2:E:101:ILE:HD11	1.75	0.69
1:D:199:GLY:O	1:D:204:TYR:OH	2.13	0.67
2:B:82:LYS:H	2:B:82:LYS:HD3	1.59	0.67
1:D:184:LEU:O	1:D:233:GLN:NE2	2.27	0.67
2:E:135:GLN:OE1	2:E:135:GLN:N	2.28	0.66
1:D:233:GLN:OE1	1:D:236:ARG:NH1	2.29	0.65
1:D:371:SER:O	1:D:375:GLU:HG2	1.96	0.65
1:D:65:ILE:O	1:D:69:VAL:HG13	1.97	0.65
1:A:217:LYS:HG2	1:A:218:PRO:HD2	1.78	0.64
2:E:55:GLU:OE2	2:E:70:ARG:NE	2.29	0.64
2:E:23:ILE:HG23	2:E:39:ILE:HG13	1.80	0.64
1:A:55:ILE:HD13	1:A:64:GLN:HB3	1.79	0.62
2:E:86:ILE:HG22	2:E:87:CYS:H	1.64	0.62
2:B:130:LYS:O	2:B:130:LYS:HD3	2.00	0.62
2:E:29:GLU:OE1	2:E:29:GLU:N	2.31	0.61
1:D:52:ILE:O	1:D:55:ILE:HG13	2.00	0.61

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:ILE:HG12	1:A:343:ILE:HD13	1.81	0.61
2:B:119:ASP:HB3	2:B:120:PRO:HD2	1.81	0.61
2:B:119:ASP:HB3	2:B:120:PRO:CD	2.30	0.61
1:D:111:LEU:HD11	1:D:115:GLY:HA2	1.82	0.60
1:D:261:PHE:HE2	1:D:289:GLY:C	2.05	0.60
1:A:134:ILE:HG12	1:A:343:ILE:HD11	1.83	0.60
1:A:50:ARG:NH2	3:C:69:LEU:O	2.34	0.59
2:B:74:LYS:HE3	2:B:143:TRP:CZ3	2.37	0.59
1:D:256:GLU:O	1:D:256:GLU:HG2	2.03	0.59
2:E:56:LEU:HG	2:E:69:VAL:CG1	2.30	0.59
1:D:141:GLU:HG3	1:D:240:TRP:CZ3	2.38	0.58
1:D:66:ALA:HA	1:D:69:VAL:HG22	1.85	0.58
1:A:57:LYS:HE2	1:A:393:LEU:HD22	1.85	0.58
1:A:127:PRO:HB3	1:A:259:ARG:HD3	1.85	0.58
2:E:138:GLU:OE1	2:E:141:ARG:NH2	2.36	0.58
2:B:117:PRO:HB3	2:B:129:TRP:HB3	1.85	0.57
1:D:198:ILE:HG12	1:D:219:GLU:HG3	1.87	0.56
2:E:54:LEU:HD21	2:E:71:PHE:CZ	2.40	0.56
1:D:60:ASN:OD1	1:D:63:GLN:NE2	2.37	0.56
1:D:83:ALA:HB1	1:D:245:SER:O	2.04	0.56
2:E:42:PRO:O	2:E:45:SER:HB3	2.06	0.56
1:A:140:ARG:HG3	1:A:243:SER:HB2	1.87	0.55
1:D:296:LYS:NZ	1:D:315:GLU:OE2	2.36	0.55
2:B:54:LEU:HD13	2:B:71:PHE:HA	1.88	0.54
1:D:361:PHE:O	1:D:364:THR:HG22	2.07	0.54
1:D:260:GLN:O	1:D:261:PHE:CD2	2.61	0.54
1:D:209:PRO:O	1:D:213:GLU:HG2	2.08	0.54
1:A:175:ILE:HG12	1:A:181:LEU:HG	1.90	0.54
1:A:283:GLU:OE2	1:A:287:ARG:NH2	2.41	0.54
1:D:108:TYR:HB3	1:D:340:LYS:HB3	1.90	0.54
1:A:184:LEU:O	1:A:233:GLN:NE2	2.35	0.54
2:E:38:VAL:HA	2:E:52:PHE:O	2.08	0.53
2:B:146:LEU:HB3	2:B:147:TYR:CD1	2.43	0.53
2:E:87:CYS:O	2:E:88:LEU:HG	2.07	0.53
3:C:1:MET:HB2	3:C:63:LYS:HB3	1.90	0.53
1:A:68:GLU:O	1:A:72:VAL:HG22	2.09	0.53
1:A:23:LYS:HB3	1:A:355:LEU:HD13	1.91	0.53
2:B:76:TYR:HB2	2:B:143:TRP:CE3	2.44	0.53
1:D:261:PHE:CE2	1:D:289:GLY:C	2.82	0.53
3:F:51:GLU:HG2	3:F:59:TYR:OH	2.09	0.52
2:E:37:VAL:O	2:E:53:LYS:HA	2.10	0.52

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:87:CYS:SG	2:E:88:LEU:N	2.83	0.51
1:D:60:ASN:HB3	1:D:63:GLN:HG2	1.92	0.51
1:A:56:ASN:OD1	1:A:58:THR:OG1	2.22	0.51
1:D:146:VAL:HG13	1:D:203:GLU:HB3	1.91	0.51
3:C:-1:GLY:N	3:C:18:GLU:OE1	2.38	0.51
2:B:37:VAL:HB	2:B:54:LEU:HB2	1.93	0.51
1:D:147:LEU:HB2	1:D:206:LEU:HD11	1.92	0.50
1:D:51:SER:HB3	1:D:68:GLU:OE1	2.12	0.50
2:E:146:LEU:CD1	2:E:147:TYR:CZ	2.95	0.50
1:D:77:THR:HB	1:D:378:ARG:HD3	1.93	0.50
2:E:47:PHE:CE1	2:E:75:ILE:HD11	2.47	0.50
1:A:45:GLN:O	1:A:49:GLU:HG3	2.12	0.50
1:D:284:TYR:OH	4:D:401:HOH:O	2.19	0.49
2:E:54:LEU:HD21	2:E:71:PHE:CE2	2.47	0.49
2:B:82:LYS:H	2:B:82:LYS:CD	2.24	0.49
2:E:81:ASP:OD1	2:E:85:ARG:HG3	2.13	0.49
1:D:356:GLN:O	1:D:360:GLU:HG3	2.13	0.49
1:D:382:LYS:O	1:D:386:THR:HG23	2.13	0.48
1:A:285:LEU:O	1:A:288:ILE:HG12	2.13	0.48
1:A:52:ILE:HD11	1:A:393:LEU:CD2	2.40	0.48
2:E:6:ARG:HG3	2:E:6:ARG:O	2.12	0.48
1:A:114:ASP:N	1:A:114:ASP:OD1	2.47	0.48
1:D:119:LYS:O	1:D:123:GLU:HG2	2.13	0.48
1:D:55:ILE:HD13	1:D:57:LYS:HG2	1.95	0.48
1:A:74:HIS:O	1:A:74:HIS:CG	2.67	0.47
2:E:70:ARG:HH22	2:E:72:MET:HE1	1.80	0.47
2:E:28:ASP:HB3	2:E:31:ASN:O	2.15	0.47
1:A:171:LYS:HG2	1:A:173:TYR:CZ	2.49	0.47
3:C:54:ARG:NH1	3:C:59:TYR:OH	2.48	0.47
2:B:41:GLY:HA3	2:B:47:PHE:O	2.14	0.47
2:B:78:PRO:HA	2:B:125:VAL:HG13	1.97	0.47
1:D:267:TYR:HB3	1:D:270:SER:HB3	1.96	0.47
1:D:117:ASN:OD1	1:D:117:ASN:N	2.48	0.47
2:E:23:ILE:HD11	2:E:109:GLN:CD	2.35	0.47
1:A:43:LEU:HG	1:A:76:MET:HG2	1.97	0.46
1:D:386:THR:O	1:D:390:GLU:HG3	2.16	0.46
1:D:331:ASN:OD1	2:E:90:ILE:HD12	2.16	0.46
1:D:261:PHE:HE2	1:D:289:GLY:CA	2.29	0.46
2:E:52:PHE:CE1	2:E:75:ILE:HD13	2.51	0.45
1:D:165:ILE:HD12	1:D:206:LEU:HD21	1.98	0.45
1:D:69:VAL:O	1:D:73:CYS:N	2.46	0.45

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:GLU:O	1:A:364:THR:HG23	2.17	0.45
1:A:69:VAL:O	1:A:73:CYS:N	2.50	0.45
1:A:388:HIS:CE1	1:A:392:LEU:HD11	2.52	0.44
1:A:201:SER:HB2	1:A:204:TYR:CZ	2.52	0.44
2:E:70:ARG:CZ	2:E:72:MET:SD	3.06	0.44
1:D:232:GLU:HG3	1:D:232:GLU:O	2.17	0.44
1:D:65:ILE:HD12	1:D:66:ALA:N	2.33	0.44
1:A:146:VAL:HG22	1:A:164:LYS:HG2	2.00	0.44
2:B:78:PRO:HD3	2:B:129:TRP:NE1	2.33	0.44
1:D:334:ARG:CZ	2:E:99:LEU:HD21	2.48	0.44
1:D:149:LYS:NZ	1:D:199:GLY:HA2	2.33	0.44
1:D:141:GLU:HG3	1:D:240:TRP:CH2	2.53	0.43
2:E:116:ASN:O	2:E:119:ASP:HB3	2.19	0.43
1:A:62:LYS:HA	1:A:62:LYS:HD2	1.95	0.43
2:B:23:ILE:HG23	2:B:39:ILE:HG13	2.01	0.43
2:E:71:PHE:HZ	2:E:112:LEU:HD11	1.83	0.43
2:E:70:ARG:NH1	2:E:72:MET:SD	2.92	0.43
1:D:350:LEU:HD12	1:D:353:LYS:HE2	2.00	0.43
2:B:58:LEU:HD23	2:B:67:PRO:HB3	2.01	0.43
3:F:22:THR:O	3:F:26:VAL:HG23	2.19	0.43
2:E:23:ILE:HD11	2:E:109:GLN:NE2	2.34	0.42
2:E:78:PRO:HD3	2:E:129:TRP:NE1	2.34	0.42
1:D:69:VAL:O	1:D:73:CYS:HB3	2.20	0.42
3:C:6:LYS:HA	3:C:11:LYS:O	2.20	0.42
1:A:374:VAL:HG12	1:A:378:ARG:HE	1.84	0.42
1:D:56:ASN:C	1:D:58:THR:H	2.23	0.42
2:B:39:ILE:HD12	2:B:54:LEU:HD23	2.01	0.42
1:D:29:GLN:HG3	1:D:277:PHE:HB3	2.02	0.42
3:F:56:LEU:O	3:F:61:ILE:HG12	2.20	0.42
1:D:148:VAL:HG22	1:D:203:GLU:HG2	2.02	0.41
1:A:74:HIS:HA	1:A:77:THR:HG22	2.03	0.41
1:A:389:LEU:O	1:A:393:LEU:HG	2.21	0.41
1:A:31:TYR:CD2	1:A:362:ILE:HD13	2.56	0.41
1:D:201:SER:HB2	1:D:204:TYR:CZ	2.56	0.41
1:A:213:GLU:C	1:A:215:PRO:HD3	2.42	0.41
1:A:69:VAL:HA	1:A:385:MET:HE3	2.01	0.41
2:B:78:PRO:HD3	2:B:129:TRP:CD1	2.56	0.41
2:B:92:LYS:HG2	3:C:39:ASP:HB3	2.03	0.41
3:F:55:THR:HG22	3:F:58:ASP:CG	2.41	0.41
1:A:17:GLY:HA3	1:A:131:ASN:OD1	2.21	0.40
1:D:167:LEU:HD21	1:D:195:PHE:HA	2.02	0.40

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:56:LEU:HD23	2:E:57:PHE:N	2.36	0.40
1:A:310:SER:HB3	1:A:327:CYS:HB3	2.04	0.40
2:E:58:LEU:CD2	2:E:101:ILE:HD11	2.49	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	382/385 (99%)	372 (97%)	9 (2%)	1 (0%)	41	49
1	D	379/385 (98%)	354 (93%)	25 (7%)	0	100	100
2	B	147/152 (97%)	140 (95%)	4 (3%)	3 (2%)	7	5
2	E	146/152 (96%)	136 (93%)	9 (6%)	1 (1%)	22	25
3	C	74/79 (94%)	73 (99%)	1 (1%)	0	100	100
3	F	73/79 (92%)	72 (99%)	1 (1%)	0	100	100
All	All	1201/1232 (98%)	1147 (96%)	49 (4%)	5 (0%)	34	41

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	119	ASP
2	B	87	CYS
2	B	131	THR
1	A	32	GLY
2	E	5	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 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	344/344 (100%)	338 (98%)	6 (2%)	60	73
1	D	341/344 (99%)	331 (97%)	10 (3%)	42	53
2	B	125/128 (98%)	121 (97%)	4 (3%)	39	50
2	E	125/128 (98%)	120 (96%)	5 (4%)	31	41
3	C	69/70 (99%)	69 (100%)	0	100	100
3	F	69/70 (99%)	67 (97%)	2 (3%)	42	53
All	All	1073/1084 (99%)	1046 (98%)	27 (2%)	47	60

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	LYS
1	A	107	LYS
1	A	116	ASP
1	A	224	SER
1	A	239	PHE
1	A	291	ASP
2	B	14	ARG
2	B	82	LYS
2	B	107	SER
2	B	130	LYS
1	D	23	LYS
1	D	63	GLN
1	D	156	ASN
1	D	207	CYS
1	D	232	GLU
1	D	243	SER
1	D	257	SER
1	D	261	PHE
1	D	279	SER
1	D	380	ARG
2	E	7	ARG
2	E	70	ARG

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
2	E	87	CYS
2	E	118	ASP
2	E	127	GLU
3	F	22	THR
3	F	72	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	47	GLN
1	D	156	ASN
2	E	116	ASN
2	E	123	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](#)

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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	384/385 (99%)	-0.18	1 (0%) 94 94	41, 62, 100, 153	0
1	D	381/385 (98%)	-0.06	3 (0%) 86 86	49, 73, 126, 164	0
2	B	149/152 (98%)	0.39	9 (6%) 21 18	52, 95, 143, 183	0
2	E	148/152 (97%)	0.42	9 (6%) 21 18	60, 100, 139, 179	0
3	C	76/79 (96%)	-0.23	1 (1%) 77 76	43, 66, 94, 138	0
3	F	75/79 (94%)	-0.03	0 100 100	51, 79, 109, 129	0
All	All	1213/1232 (98%)	0.01	23 (1%) 66 64	41, 76, 127, 183	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	121	LEU	6.8
2	B	122	ALA	6.1
1	D	254	LEU	4.6
2	B	124	ASP	4.4
2	B	125	VAL	3.8
2	E	56	LEU	3.5
2	E	118	ASP	3.2
2	B	3	GLY	3.1
3	C	-1	GLY	3.0
2	E	105	LEU	3.0
2	E	35	PHE	2.7
2	B	129	TRP	2.6
2	E	133	GLU	2.5
2	E	54	LEU	2.5
1	D	256	GLU	2.4
2	B	118	ASP	2.3
2	B	130	LYS	2.3
1	A	395	ILE	2.2
1	D	53	PRO	2.1

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
2	E	40	ALA	2.1
2	E	51	THR	2.1
2	E	64	MET	2.0
2	B	20	VAL	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 [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.