



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 10:50 am BST

PDB ID : 6MO0
Title : Structure of dengue virus protease with an allosteric Inhibitor that blocks replication
Authors : Lin, Y.-L.; Nie, S.; Hua, Y.; Wu, J.; Wu, F.; Huo, T.; Yao, Y.; Song, Y.
Deposited on : 2018-10-03
Resolution : 2.70 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
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

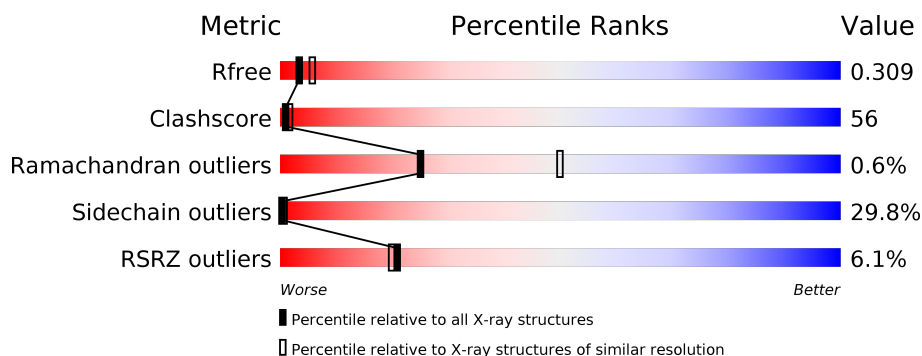
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	247	<div> <div>3%</div> <div>23%</div> <div>33%</div> <div>10%</div> <div>34%</div> </div>
1	B	247	<div> <div>5%</div> <div>23%</div> <div>32%</div> <div>11%</div> <div>34%</div> </div>

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
2	JVJ	A	1201	-	-	-	X
2	JVJ	B	1201	-	-	-	X

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2578 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 FLAVIVIRUS_NS2B/Peptidase S7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	164	Total	C	N	O	S	0	0	0
			1256	796	217	240	3			
1	B	164	Total	C	N	O	S	0	0	0
			1256	796	217	240	3			

There are 32 discrepancies between the modelled and reference sequences:

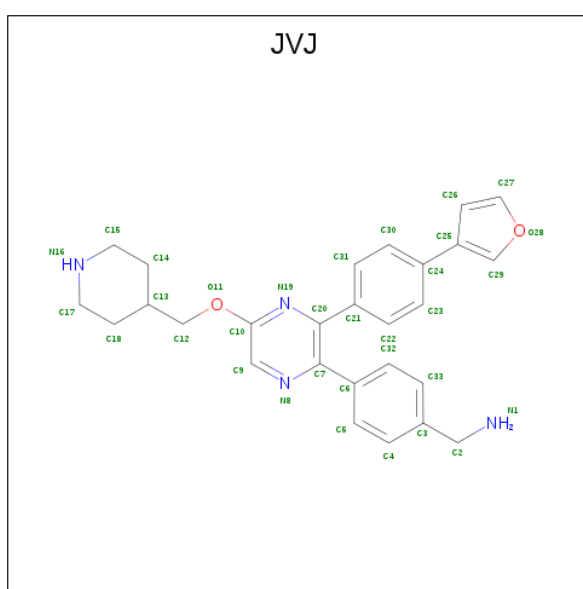
Chain	Residue	Modelled	Actual	Comment	Reference
A	43	GLY	-	expression tag	UNP A0A0B4L2Y4
A	44	SER	-	expression tag	UNP A0A0B4L2Y4
A	45	HIS	-	expression tag	UNP A0A0B4L2Y4
A	46	MET	-	expression tag	UNP A0A0B4L2Y4
A	47	LEU	-	expression tag	UNP A0A0B4L2Y4
A	48	MET	-	expression tag	UNP A0A0B4L2Y4
A	992	GLY	-	linker	UNP A0A0B4L2Y4
A	993	GLY	-	linker	UNP A0A0B4L2Y4
A	994	GLY	-	linker	UNP A0A0B4L2Y4
A	995	GLY	-	linker	UNP A0A0B4L2Y4
A	996	SER	-	linker	UNP A0A0B4L2Y4
A	997	GLY	-	linker	UNP A0A0B4L2Y4
A	998	GLY	-	linker	UNP A0A0B4L2Y4
A	999	GLY	-	linker	UNP A0A0B4L2Y4
A	1000	GLY	-	linker	UNP A0A0B4L2Y4
A	1167	ASN	GLN	conflict	UNP Q91H74
B	43	GLY	-	expression tag	UNP A0A0B4L2Y4
B	44	SER	-	expression tag	UNP A0A0B4L2Y4
B	45	HIS	-	expression tag	UNP A0A0B4L2Y4
B	46	MET	-	expression tag	UNP A0A0B4L2Y4
B	47	LEU	-	expression tag	UNP A0A0B4L2Y4
B	48	MET	-	expression tag	UNP A0A0B4L2Y4
B	992	GLY	-	linker	UNP A0A0B4L2Y4
B	993	GLY	-	linker	UNP A0A0B4L2Y4
B	994	GLY	-	linker	UNP A0A0B4L2Y4

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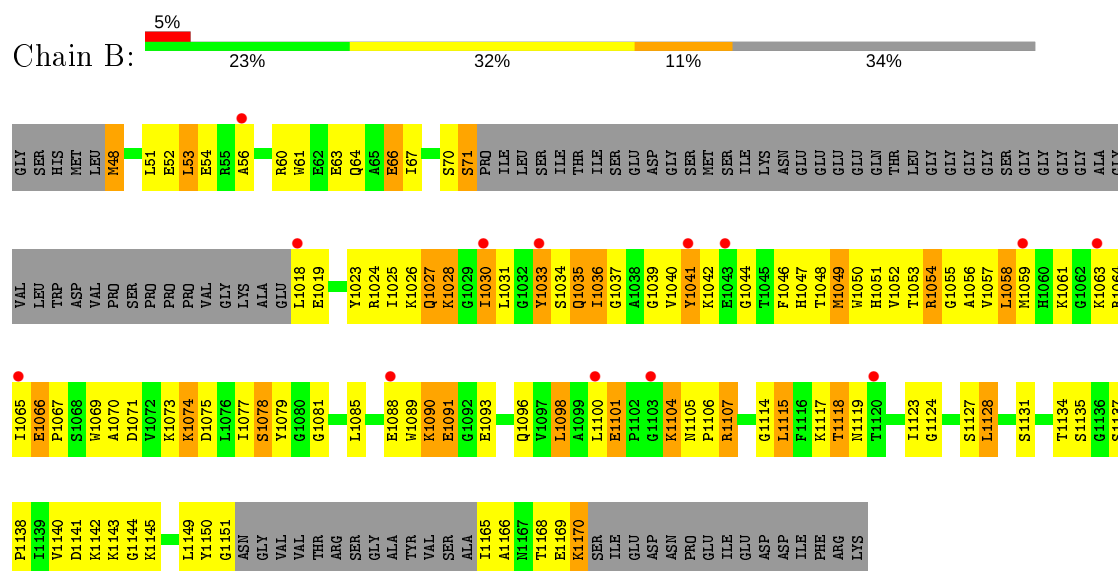
Chain	Residue	Modelled	Actual	Comment	Reference
B	995	GLY	-	linker	UNP A0A0B4L2Y4
B	996	SER	-	linker	UNP A0A0B4L2Y4
B	997	GLY	-	linker	UNP A0A0B4L2Y4
B	998	GLY	-	linker	UNP A0A0B4L2Y4
B	999	GLY	-	linker	UNP A0A0B4L2Y4
B	1000	GLY	-	linker	UNP A0A0B4L2Y4
B	1167	ASN	GLN	conflict	UNP Q91H74

- Molecule 2 is 1-(4-{3-[4-(furan-3-yl)phenyl]-5-[(piperidin-4-yl)methoxy]pyrazin-2-yl}phenyl) methanamine (three-letter code: JVJ) (formula: C₂₇H₂₈N₄O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			33	27	4	2		
2	B	1	Total	C	N	O	0	0
			33	27	4	2		

● Molecule 1: FLAVIVIRUS NS2B/Peptidase S7



4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	109.11Å 109.11Å 75.85Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.13 – 2.70 40.10 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.7 (40.13-2.70) 99.8 (40.10-2.70)	Depositor EDS
R_{merge}	0.40	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.256 , 0.311 0.277 , 0.309	Depositor DCC
R_{free} test set	709 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	56.9	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 49.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.037 for -h,-k,l	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	2578	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% 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

5.1 Standard geometry

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

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	0.71	0/1280	0.84	0/1721
1	B	0.74	0/1280	0.83	1/1721 (0.1%)
All	All	0.72	0/2560	0.84	1/3442 (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	B	1066	GLU	CB-CA-C	5.57	121.55	110.40

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	1256	0	1252	140	3
1	B	1256	0	1252	149	3
2	A	33	0	0	3	0
2	B	33	0	0	5	0
All	All	2578	0	2504	283	6

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 56.

All (283) 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:1018:LEU:HD23	1:A:1060:HIS:CD2	1.58	1.39
1:A:1112:LYS:HE2	1:A:1113:PRO:O	1.35	1.26
1:B:1053:THR:HG21	1:B:1079:TYR:OH	1.37	1.22
1:A:1018:LEU:O	1:A:1023:TYR:OH	1.64	1.15
1:A:1101:GLU:N	1:A:1101:GLU:OE2	1.82	1.12
1:A:1091:GLU:HB2	1:A:1168:THR:CG2	1.85	1.06
1:A:1091:GLU:HB2	1:A:1168:THR:HG23	1.34	1.05
1:B:1026:LYS:NZ	1:B:1035:GLN:HB3	1.72	1.04
1:B:1070:ALA:HB2	1:B:1077:ILE:HD12	1.41	1.03
1:B:1026:LYS:HZ3	1:B:1035:GLN:HB3	1.19	1.02
1:A:1018:LEU:HD23	1:A:1060:HIS:NE2	1.73	1.02
1:B:1065:ILE:CG2	1:B:1079:TYR:HB3	1.89	1.02
1:A:51:LEU:HD21	1:A:1025:ILE:CG2	1.90	1.01
1:A:1105:ASN:OD1	1:A:1106:PRO:HD2	1.60	1.01
1:A:1030:ILE:HG23	1:A:1031:LEU:HD13	1.40	0.99
1:B:56:ALA:HB3	1:B:1100:LEU:HD21	1.38	0.98
1:B:1053:THR:CG2	1:B:1079:TYR:OH	2.13	0.96
1:B:1105:ASN:OD1	1:B:1106:PRO:HD2	1.66	0.96
1:A:1026:LYS:HE3	1:A:1035:GLN:HB2	1.50	0.94
1:A:1116:PHE:HE2	1:A:1125:ALA:HB2	1.31	0.93
1:B:1124:GLY:O	1:B:1166:ALA:HB3	1.67	0.93
1:B:1090:LYS:NZ	1:B:1093:GLU:OE1	2.01	0.92
1:A:1056:ALA:O	1:A:1079:TYR:OH	1.88	0.91
1:B:1057:VAL:HG21	1:B:1064:ARG:NE	1.84	0.91
1:A:55:ARG:NH1	1:A:1020:ASP:O	2.03	0.90
1:B:56:ALA:CB	1:B:1100:LEU:HD21	2.01	0.90
1:A:1151:GLY:C	1:A:1165:ILE:HD12	1.91	0.90
1:B:1019:GLU:O	1:B:1023:TYR:OH	1.89	0.90
1:B:1118:THR:HG21	1:B:1123:ILE:HG12	1.52	0.89
1:B:1091:GLU:HB2	1:B:1168:THR:CG2	2.02	0.89
1:A:1018:LEU:CD2	1:A:1060:HIS:CD2	2.53	0.89
1:A:1116:PHE:CE2	1:A:1125:ALA:HB2	2.08	0.88
1:A:55:ARG:HB3	1:B:64:GLN:NE2	1.88	0.88
1:A:1101:GLU:HG2	1:A:1104:LYS:HB2	1.57	0.87
1:A:1027:GLN:NE2	1:A:1027:GLN:O	2.08	0.86
1:A:65:ALA:O	1:A:66:GLU:HB2	1.72	0.86
1:B:1101:GLU:OE1	1:B:1101:GLU:N	2.07	0.86
1:A:50:ASP:OD2	1:A:1028:LYS:HE3	1.76	0.84
1:A:51:LEU:CD2	1:A:1025:ILE:HG23	2.08	0.84
1:A:1037:GLY:HA3	1:A:1052:VAL:HG11	1.60	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:LEU:CD2	1:A:1025:ILE:CG2	2.56	0.83
1:A:51:LEU:HD22	1:A:1058:LEU:HD11	1.59	0.83
1:A:68:SER:OG	1:A:69:GLY:N	2.13	0.82
1:B:1091:GLU:HB2	1:B:1168:THR:HG23	1.61	0.80
1:A:1097:VAL:CG1	1:A:1109:VAL:HG13	2.12	0.80
1:B:1057:VAL:HG21	1:B:1064:ARG:HE	1.45	0.80
1:A:1101:GLU:HG2	1:A:1104:LYS:CB	2.13	0.79
1:A:1070:ALA:HB2	1:A:1077:ILE:HD12	1.64	0.79
1:A:1061:LYS:HG3	1:A:1062:GLY:H	1.48	0.78
1:A:1101:GLU:CD	1:A:1101:GLU:H	1.86	0.78
1:B:1049:MET:HE3	1:B:1151:GLY:HA3	1.65	0.78
1:B:1048:THR:HG23	1:B:1077:ILE:CG2	2.12	0.78
1:B:1070:ALA:HB2	1:B:1077:ILE:CD1	2.13	0.78
1:A:1097:VAL:HG12	1:A:1109:VAL:HG13	1.64	0.78
1:A:1101:GLU:HG3	1:A:1104:LYS:HE2	1.66	0.78
1:B:1026:LYS:HZ2	1:B:1035:GLN:N	1.80	0.78
1:B:1070:ALA:CB	1:B:1077:ILE:HD12	2.14	0.78
1:B:1036:ILE:HD12	1:B:1036:ILE:H	1.49	0.77
1:B:1090:LYS:O	1:B:1091:GLU:HB3	1.83	0.76
1:A:1053:THR:CG2	1:A:1079:TYR:OH	2.34	0.76
1:A:1090:LYS:HD3	1:A:1093:GLU:HB2	1.68	0.76
1:A:55:ARG:HD3	1:B:64:GLN:HE22	1.49	0.76
1:A:1088:GLU:HG2	1:A:1088:GLU:O	1.84	0.76
1:A:1031:LEU:HD13	1:A:1031:LEU:N	2.02	0.75
1:A:1053:THR:HG21	1:A:1079:TYR:OH	1.86	0.75
1:A:1030:ILE:HG23	1:A:1031:LEU:CD1	2.17	0.75
1:A:1116:PHE:HE2	1:A:1125:ALA:CB	1.99	0.75
1:A:1037:GLY:HA3	1:A:1052:VAL:CG1	2.16	0.75
1:B:56:ALA:CB	1:B:1100:LEU:CD2	2.64	0.75
1:A:55:ARG:HB3	1:B:64:GLN:HE21	1.51	0.74
1:B:1151:GLY:O	1:B:1165:ILE:HG13	1.87	0.74
1:A:1091:GLU:HB2	1:A:1168:THR:HG21	1.70	0.73
1:B:1067:PRO:HG3	1:B:1079:TYR:CE2	2.23	0.73
1:A:1018:LEU:CD2	1:A:1060:HIS:NE2	2.51	0.73
1:B:1040:VAL:CG2	1:B:1138:PRO:HG3	2.19	0.73
1:B:1053:THR:HG22	1:B:1055:GLY:H	1.54	0.72
1:B:51:LEU:HB3	1:B:1058:LEU:HD12	1.71	0.72
1:A:1047:HIS:HE1	1:A:1084:LYS:H	1.37	0.72
1:B:1101:GLU:CD	1:B:1101:GLU:H	1.90	0.72
1:B:63:GLU:N	1:B:63:GLU:OE2	2.21	0.72
1:B:66:GLU:HG2	1:B:70:SER:OG	1.88	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1031:LEU:H	1:A:1031:LEU:HD13	1.54	0.71
1:A:1025:ILE:HG22	1:A:1036:ILE:HD11	1.70	0.71
1:A:1090:LYS:HD3	1:A:1093:GLU:CB	2.20	0.71
1:B:1026:LYS:HD2	1:B:1035:GLN:HB2	1.72	0.71
1:B:1065:ILE:HG22	1:B:1079:TYR:HB3	1.71	0.71
1:A:1099:ALA:C	1:A:1100:LEU:HD12	2.12	0.71
1:B:1050:TRP:CE3	1:B:1077:ILE:HD13	2.26	0.71
1:A:1071:ASP:OD2	1:A:1074:LYS:HE3	1.91	0.71
1:A:51:LEU:HD22	1:A:1058:LEU:CD1	2.21	0.70
1:A:1112:LYS:CE	1:A:1113:PRO:O	2.28	0.70
1:B:1151:GLY:O	1:B:1165:ILE:CD1	2.41	0.69
1:B:1101:GLU:OE2	1:B:1107:ARG:CZ	2.41	0.69
1:A:1040:VAL:HG21	1:A:1138:PRO:HB3	1.76	0.68
1:A:1067:PRO:HG3	1:A:1079:TYR:CE2	2.29	0.68
1:B:1049:MET:CE	1:B:1151:GLY:HA3	2.23	0.68
1:B:1165:ILE:HD11	2:B:1201:JVJ:C33	2.24	0.67
1:B:1065:ILE:CG2	1:B:1079:TYR:CB	2.71	0.67
1:A:1053:THR:HG21	1:A:1079:TYR:CE1	2.30	0.66
1:B:1091:GLU:HB2	1:B:1168:THR:HG21	1.76	0.66
1:A:1078:SER:OG	1:A:1081:GLY:O	2.14	0.66
1:A:1057:VAL:CG2	1:A:1064:ARG:HG3	2.27	0.65
1:B:1091:GLU:O	1:B:1091:GLU:HG2	1.96	0.64
1:B:1057:VAL:CG2	1:B:1064:ARG:NE	2.60	0.64
1:A:1090:LYS:O	1:A:1091:GLU:HB3	1.96	0.64
1:B:1026:LYS:NZ	1:B:1035:GLN:CB	2.56	0.64
1:A:51:LEU:HD21	1:A:1025:ILE:HG22	1.78	0.64
1:B:1089:TRP:HH2	1:B:1115:LEU:HD11	1.62	0.64
1:B:1041:TYR:CD2	1:B:1046:PHE:HD1	2.16	0.63
1:A:1040:VAL:CG2	1:A:1138:PRO:HB3	2.29	0.63
1:A:1061:LYS:HG3	1:A:1062:GLY:N	2.14	0.63
1:A:1057:VAL:HG21	1:A:1064:ARG:HG3	1.81	0.62
1:A:1031:LEU:HD22	1:A:1031:LEU:H	1.62	0.62
1:B:1030:ILE:N	1:B:1030:ILE:HD12	2.14	0.62
1:A:55:ARG:HD3	1:B:64:GLN:NE2	2.14	0.62
1:A:1065:ILE:HG22	1:A:1079:TYR:CD2	2.35	0.61
1:A:1053:THR:HG22	1:A:1055:GLY:H	1.65	0.61
1:B:1054:ARG:CG	1:B:1054:ARG:HH11	2.14	0.61
1:B:1040:VAL:HG23	1:B:1138:PRO:HG3	1.84	0.60
1:B:1030:ILE:HD12	1:B:1030:ILE:H	1.67	0.60
1:B:1048:THR:HG23	1:B:1077:ILE:HG21	1.83	0.60
1:B:1151:GLY:O	1:B:1165:ILE:CG1	2.50	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1025:ILE:N	1:A:1025:ILE:HD12	2.18	0.59
1:B:1105:ASN:OD1	1:B:1106:PRO:CD	2.45	0.59
1:A:55:ARG:CD	1:B:64:GLN:HE22	2.15	0.59
1:A:1053:THR:HG21	1:A:1079:TYR:CZ	2.38	0.59
1:A:1047:HIS:CE1	1:A:1084:LYS:H	2.19	0.58
1:B:1050:TRP:HE3	1:B:1077:ILE:HD13	1.68	0.58
1:B:1089:TRP:CH2	1:B:1115:LEU:HD11	2.38	0.58
1:A:1127:SER:O	1:A:1128:LEU:HD23	2.03	0.58
1:A:1067:PRO:HG3	1:A:1079:TYR:CZ	2.39	0.58
1:B:56:ALA:HB3	1:B:1100:LEU:CD2	2.22	0.57
1:A:1033:TYR:N	1:A:1033:TYR:CD1	2.72	0.57
1:A:1098:LEU:HB3	1:A:1138:PRO:HG2	1.87	0.57
1:A:1033:TYR:N	1:A:1033:TYR:HD1	2.02	0.56
1:B:1048:THR:CG2	1:B:1077:ILE:CG2	2.83	0.56
1:B:1090:LYS:O	1:B:1091:GLU:CB	2.53	0.56
1:A:1151:GLY:C	1:A:1165:ILE:CD1	2.71	0.56
1:A:1036:ILE:HG12	1:A:1052:VAL:HG13	1.87	0.56
1:A:1042:LYS:HE2	1:A:1085:LEU:CD2	2.35	0.56
1:A:1093:GLU:HG3	1:A:1094:GLU:N	2.21	0.56
1:B:1054:ARG:NH1	1:B:1054:ARG:HG2	2.21	0.56
1:B:1026:LYS:HE3	1:B:1034:SER:C	2.26	0.55
1:B:1037:GLY:HA3	1:B:1052:VAL:CG1	2.36	0.55
1:A:1058:LEU:HD22	1:A:1079:TYR:CZ	2.41	0.55
1:B:1026:LYS:HZ2	1:B:1035:GLN:HB3	1.70	0.55
1:B:1028:LYS:HG3	1:B:1033:TYR:HB3	1.88	0.55
1:B:1065:ILE:HG22	1:B:1079:TYR:CD2	2.41	0.55
1:A:1036:ILE:H	1:A:1036:ILE:HD13	1.72	0.55
1:A:1042:LYS:HE2	1:A:1085:LEU:HD23	1.87	0.55
1:A:1099:ALA:O	1:A:1100:LEU:HD12	2.07	0.55
1:A:1025:ILE:HD13	1:A:1037:GLY:C	2.28	0.55
1:B:1118:THR:CG2	1:B:1123:ILE:HG12	2.32	0.55
1:A:1065:ILE:HG22	1:A:1079:TYR:HD2	1.71	0.54
1:A:1067:PRO:HA	1:A:1079:TYR:HA	1.89	0.54
1:A:55:ARG:CB	1:B:64:GLN:NE2	2.67	0.54
1:A:1057:VAL:HG23	1:A:1065:ILE:O	2.08	0.54
1:B:1018:LEU:CD2	1:B:1044:GLY:HA2	2.38	0.54
1:A:1025:ILE:CG2	1:A:1036:ILE:HD11	2.38	0.53
2:A:1201:JVJ:C6	2:A:1201:JVJ:C31	2.85	0.53
1:B:1025:ILE:HD12	1:B:1025:ILE:N	2.24	0.53
1:B:1041:TYR:CD2	1:B:1046:PHE:CD1	2.96	0.53
1:B:1115:LEU:CD1	1:B:1115:LEU:N	2.71	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:LEU:HB3	1:A:1058:LEU:HD12	1.91	0.53
1:B:1026:LYS:HD2	1:B:1035:GLN:CB	2.39	0.53
1:B:1065:ILE:HG21	1:B:1079:TYR:CB	2.39	0.52
1:A:1091:GLU:HG2	1:A:1091:GLU:O	2.09	0.52
1:B:1141:ASP:OD2	1:B:1145:LYS:HE3	2.10	0.52
1:B:1053:THR:HG22	1:B:1055:GLY:N	2.21	0.52
1:B:1067:PRO:HG3	1:B:1079:TYR:HE2	1.74	0.52
1:A:1061:LYS:CG	1:A:1062:GLY:H	2.17	0.52
1:B:1058:LEU:HB3	1:B:1065:ILE:HD12	1.91	0.52
1:B:1065:ILE:HG21	1:B:1079:TYR:HB3	1.85	0.52
1:A:1030:ILE:CG2	1:A:1031:LEU:HD13	2.28	0.52
1:B:1037:GLY:HA3	1:B:1052:VAL:HG11	1.92	0.51
1:A:1090:LYS:HG3	1:A:1092:GLY:H	1.75	0.51
1:B:1042:LYS:HD3	1:B:1085:LEU:CD2	2.40	0.51
1:A:1053:THR:HG21	1:A:1079:TYR:HE1	1.76	0.51
1:B:1025:ILE:HD11	1:B:1046:PHE:HZ	1.75	0.51
1:B:1028:LYS:HA	1:B:1033:TYR:HB3	1.92	0.51
1:B:1026:LYS:CD	1:B:1035:GLN:HB2	2.40	0.51
1:B:1098:LEU:HD12	1:B:1098:LEU:C	2.31	0.51
1:A:1050:TRP:O	1:A:1054:ARG:N	2.41	0.51
1:B:1070:ALA:CB	1:B:1077:ILE:CD1	2.83	0.51
1:B:1054:ARG:HH11	1:B:1054:ARG:HG2	1.74	0.51
1:B:1065:ILE:HG23	1:B:1079:TYR:HB3	1.88	0.51
1:B:1141:ASP:OD2	1:B:1145:LYS:HB2	2.11	0.51
1:B:52:GLU:HA	1:B:1059:MET:O	2.11	0.50
1:A:1033:TYR:H	1:A:1033:TYR:HD1	1.60	0.50
1:A:1048:THR:HG23	1:A:1077:ILE:CG2	2.42	0.50
1:B:1058:LEU:HD22	1:B:1079:TYR:CZ	2.46	0.50
1:B:1018:LEU:HD21	1:B:1044:GLY:HA2	1.93	0.50
1:A:1129:ASP:OD1	1:A:1129:ASP:C	2.50	0.50
1:B:1018:LEU:CD2	1:B:1041:TYR:CE1	2.95	0.50
1:B:66:GLU:CG	1:B:70:SER:OG	2.59	0.50
1:B:1018:LEU:HD23	1:B:1041:TYR:CE1	2.47	0.50
1:A:1025:ILE:HD13	1:A:1038:ALA:N	2.27	0.49
1:B:1169:GLU:O	1:B:1170:LYS:HB3	2.13	0.49
1:B:1051:HIS:ND1	1:B:1075:ASP:OD2	2.40	0.49
1:A:1053:THR:HG23	1:A:1079:TYR:OH	2.13	0.49
1:B:56:ALA:HB1	1:B:1106:PRO:HG3	1.94	0.49
1:B:61:TRP:CE3	1:B:1142:LYS:HA	2.47	0.49
1:A:70:SER:O	1:A:70:SER:OG	2.14	0.49
1:B:1069:TRP:HE3	1:B:1078:SER:HB3	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1057:VAL:HG21	1:B:1064:ARG:CZ	2.40	0.48
1:B:51:LEU:HD23	1:B:1058:LEU:HD11	1.95	0.48
1:A:1097:VAL:CG1	1:A:1109:VAL:CG1	2.90	0.48
1:B:1018:LEU:O	1:B:1018:LEU:HD23	2.13	0.48
1:B:1090:LYS:NZ	1:B:1093:GLU:CD	2.67	0.48
1:A:63:GLU:H	1:A:63:GLU:HG2	1.34	0.47
1:A:1101:GLU:HG2	1:A:1104:LYS:HB3	1.92	0.47
1:A:1058:LEU:HD22	1:A:1079:TYR:CE1	2.50	0.47
1:A:67:ILE:HG13	1:A:1110:GLN:HB3	1.96	0.47
1:A:51:LEU:HD23	1:A:1025:ILE:HG23	1.94	0.47
1:B:1026:LYS:NZ	1:B:1035:GLN:N	2.58	0.47
1:A:1018:LEU:HA	1:A:1060:HIS:CE1	2.49	0.47
1:A:1025:ILE:HG22	1:A:1036:ILE:CD1	2.41	0.47
1:A:1074:LYS:HB3	1:A:1076:LEU:HD12	1.97	0.47
1:B:1042:LYS:HD3	1:B:1085:LEU:HD23	1.96	0.46
1:A:1086:GLU:HG3	1:A:1086:GLU:H	1.58	0.46
1:B:1149:LEU:HB2	2:B:1201:JVJ:C27	2.45	0.46
1:A:1116:PHE:CD2	1:A:1116:PHE:N	2.84	0.46
1:A:51:LEU:HD21	1:A:1025:ILE:HG21	1.90	0.46
1:B:1057:VAL:CG2	1:B:1064:ARG:CZ	2.93	0.46
1:B:1039:GLY:HA3	1:B:1046:PHE:CZ	2.51	0.46
1:A:1036:ILE:H	1:A:1036:ILE:CD1	2.29	0.46
1:A:1054:ARG:HE	1:A:1054:ARG:HB2	1.50	0.46
1:A:1140:VAL:HA	1:A:1145:LYS:O	2.15	0.46
1:B:56:ALA:CB	1:B:1100:LEU:HD23	2.46	0.46
1:A:50:ASP:HB3	1:A:1028:LYS:HB2	1.98	0.45
1:B:1025:ILE:N	1:B:1025:ILE:CD1	2.79	0.45
1:A:1031:LEU:HD22	1:A:1031:LEU:N	2.31	0.45
1:B:1078:SER:OG	1:B:1081:GLY:O	2.35	0.45
1:A:1039:GLY:HA3	1:A:1046:PHE:CZ	2.52	0.45
1:A:1058:LEU:HA	1:A:1058:LEU:HD12	1.78	0.45
1:B:1026:LYS:CD	1:B:1035:GLN:CB	2.93	0.45
1:B:1107:ARG:HH21	1:B:1107:ARG:CG	2.29	0.45
1:B:1127:SER:C	1:B:1128:LEU:HD13	2.36	0.45
1:B:1050:TRP:CZ3	1:B:1077:ILE:HD13	2.51	0.45
1:B:51:LEU:HD22	1:B:1056:ALA:HB3	1.99	0.45
1:A:1026:LYS:HE3	1:A:1035:GLN:CB	2.35	0.45
1:A:55:ARG:HB3	1:B:64:GLN:HE22	1.79	0.45
1:A:1149:LEU:H	2:A:1201:JVJ:C27	2.30	0.44
1:B:1104:LYS:HD2	1:B:1104:LYS:HA	1.49	0.44
1:B:60:ARG:HA	1:B:60:ARG:HD2	1.85	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1067:PRO:HB3	1:A:1077:ILE:HD11	1.98	0.44
1:B:61:TRP:CE2	1:B:1096:GLN:NE2	2.86	0.44
1:B:1030:ILE:CD1	1:B:1030:ILE:H	2.23	0.44
1:A:1100:LEU:HD12	1:A:1100:LEU:N	2.33	0.43
1:B:1107:ARG:HG2	1:B:1107:ARG:NH2	2.33	0.43
1:A:64:GLN:O	1:A:64:GLN:HG3	2.18	0.43
1:B:51:LEU:HD12	1:B:1027:GLN:HB2	1.98	0.43
1:B:1069:TRP:CZ2	1:B:1071:ASP:HB2	2.53	0.43
1:B:51:LEU:CD1	1:B:1027:GLN:HB2	2.47	0.43
1:B:1066:GLU:HA	1:B:1067:PRO:HD3	1.73	0.43
1:B:1114:GLY:C	1:B:1115:LEU:HD12	2.39	0.43
1:B:1074:LYS:O	2:B:1201:JVJ:C31	2.67	0.43
1:A:1018:LEU:CA	1:A:1060:HIS:NE2	2.82	0.43
1:B:1049:MET:HE1	1:B:1150:TYR:O	2.18	0.43
1:B:1091:GLU:CB	1:B:1168:THR:HG23	2.41	0.43
1:A:1047:HIS:HB3	1:A:1083:TRP:CH2	2.53	0.42
1:B:1145:LYS:HB2	1:B:1145:LYS:HE3	1.80	0.42
1:B:1115:LEU:N	1:B:1115:LEU:HD13	2.34	0.42
2:A:1201:JVJ:C32	2:A:1201:JVJ:C31	2.98	0.42
1:B:1165:ILE:CD1	2:B:1201:JVJ:C33	2.97	0.42
1:A:1018:LEU:N	1:A:1060:HIS:CE1	2.88	0.41
1:A:54:GLU:O	1:A:1023:TYR:HA	2.20	0.41
1:B:51:LEU:HB3	1:B:1058:LEU:CD1	2.44	0.41
1:A:1100:LEU:CD1	1:A:1100:LEU:N	2.82	0.41
1:B:53:LEU:HD12	1:B:53:LEU:HA	1.84	0.41
1:A:1053:THR:HG22	1:A:1055:GLY:N	2.30	0.41
1:A:1131:SER:OG	1:A:1132:PRO:HD2	2.20	0.41
1:B:60:ARG:NH2	1:B:1144:GLY:HA3	2.36	0.41
1:A:1122:THR:O	1:A:1168:THR:HB	2.20	0.41
1:B:1114:GLY:C	1:B:1115:LEU:CD1	2.89	0.40
1:B:1137:SER:HA	1:B:1138:PRO:HD3	1.89	0.40
1:B:1165:ILE:HD13	2:B:1201:JVJ:C30	2.51	0.40
1:A:53:LEU:HD13	1:A:1058:LEU:HG	2.04	0.40
1:A:1027:GLN:NE2	1:A:1027:GLN:C	2.73	0.40
1:A:55:ARG:CG	1:B:64:GLN:NE2	2.85	0.40
1:B:1026:LYS:HZ2	1:B:1035:GLN:H	1.66	0.40
1:B:1064:ARG:NH1	1:B:1066:GLU:HB3	2.36	0.40
1:B:1091:GLU:CB	1:B:1168:THR:HG21	2.48	0.40

All (6) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:MET:N	1:A:71:SER:O[2_445]	0.67	1.53
1:B:48:MET:N	1:B:71:SER:O[3_655]	1.12	1.08
1:B:1143:LYS:CG	1:B:1143:LYS:CG[6_556]	1.75	0.45
1:A:48:MET:N	1:A:71:SER:C[2_445]	1.91	0.29
1:A:48:MET:CA	1:A:71:SER:O[2_445]	2.11	0.09
1:B:48:MET:CA	1:B:71:SER:O[3_655]	2.18	0.02

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	158/247 (64%)	152 (96%)	5 (3%)	1 (1%)	25	50
1	B	158/247 (64%)	153 (97%)	4 (2%)	1 (1%)	25	50
All	All	316/494 (64%)	305 (96%)	9 (3%)	2 (1%)	25	50

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	66	GLU
1	B	1091	GLU

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	129/194 (66%)	92 (71%)	37 (29%)	0	1

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	129/194 (66%)	89 (69%)	40 (31%)	0	0
All	All	258/388 (66%)	181 (70%)	77 (30%)	0	1

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	LEU
1	A	52	GLU
1	A	53	LEU
1	A	60	ARG
1	A	62	GLU
1	A	63	GLU
1	A	64	GLN
1	A	67	ILE
1	A	68	SER
1	A	71	SER
1	A	1018	LEU
1	A	1019	GLU
1	A	1026	LYS
1	A	1028	LYS
1	A	1031	LEU
1	A	1033	TYR
1	A	1036	ILE
1	A	1041	TYR
1	A	1042	LYS
1	A	1054	ARG
1	A	1058	LEU
1	A	1063	LYS
1	A	1064	ARG
1	A	1066	GLU
1	A	1078	SER
1	A	1086	GLU
1	A	1098	LEU
1	A	1101	GLU
1	A	1109	VAL
1	A	1112	LYS
1	A	1117	LYS
1	A	1129	ASP
1	A	1135	SER
1	A	1147	VAL
1	A	1167	ASN
1	A	1169	GLU

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Mol	Chain	Res	Type
1	A	1170	LYS
1	B	48	MET
1	B	53	LEU
1	B	54	GLU
1	B	66	GLU
1	B	67	ILE
1	B	71	SER
1	B	1024	ARG
1	B	1027	GLN
1	B	1028	LYS
1	B	1030	ILE
1	B	1031	LEU
1	B	1033	TYR
1	B	1035	GLN
1	B	1036	ILE
1	B	1041	TYR
1	B	1047	HIS
1	B	1049	MET
1	B	1054	ARG
1	B	1058	LEU
1	B	1061	LYS
1	B	1063	LYS
1	B	1073	LYS
1	B	1074	LYS
1	B	1078	SER
1	B	1088	GLU
1	B	1090	LYS
1	B	1098	LEU
1	B	1101	GLU
1	B	1104	LYS
1	B	1107	ARG
1	B	1115	LEU
1	B	1117	LYS
1	B	1118	THR
1	B	1119	ASN
1	B	1128	LEU
1	B	1131	SER
1	B	1134	THR
1	B	1135	SER
1	B	1140	VAL
1	B	1170	LYS

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

sidechains are listed below:

Mol	Chain	Res	Type
1	A	64	GLN
1	A	1047	HIS
1	B	64	GLN
1	B	1119	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](#)

2 ligands are 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
2	JVJ	B	1201	-	34,37,37	2.64	13 (38%)	43,50,50	1.78	7 (16%)
2	JVJ	A	1201	-	34,37,37	2.41	14 (41%)	43,50,50	2.35	14 (32%)

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
2	JVJ	B	1201	-	-	1/19/27/27	0/5/5/5
2	JVJ	A	1201	-	-	6/19/27/27	0/5/5/5

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1201	JVJ	C29-C25	8.75	1.47	1.37
2	B	1201	JVJ	C21-C20	6.23	1.56	1.49
2	A	1201	JVJ	C20-N19	5.57	1.43	1.34
2	A	1201	JVJ	C29-C25	5.19	1.43	1.37
2	A	1201	JVJ	C30-C24	4.81	1.49	1.39
2	A	1201	JVJ	C30-C31	3.90	1.45	1.38
2	B	1201	JVJ	C20-N19	3.73	1.40	1.34
2	B	1201	JVJ	C30-C31	3.56	1.45	1.38
2	A	1201	JVJ	O11-C10	3.52	1.43	1.36
2	A	1201	JVJ	C22-C21	3.43	1.46	1.39
2	A	1201	JVJ	C21-C20	3.28	1.52	1.49
2	A	1201	JVJ	C6-C7	-3.22	1.45	1.49
2	B	1201	JVJ	C32-C33	3.17	1.44	1.38
2	B	1201	JVJ	C6-C7	-3.08	1.45	1.49
2	B	1201	JVJ	C10-N19	3.01	1.38	1.33
2	B	1201	JVJ	C30-C24	2.97	1.45	1.39
2	A	1201	JVJ	C4-C5	2.96	1.44	1.38
2	B	1201	JVJ	C22-C23	2.89	1.44	1.38
2	B	1201	JVJ	C22-C21	2.77	1.45	1.39
2	A	1201	JVJ	C10-N19	2.49	1.37	1.33
2	B	1201	JVJ	C23-C24	2.43	1.44	1.39
2	B	1201	JVJ	C33-C3	2.43	1.44	1.38
2	A	1201	JVJ	C32-C6	2.33	1.44	1.39
2	B	1201	JVJ	C25-C24	2.31	1.54	1.49
2	A	1201	JVJ	C31-C21	2.29	1.44	1.39
2	A	1201	JVJ	C9-C10	2.29	1.43	1.39
2	A	1201	JVJ	C4-C3	2.26	1.43	1.38

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1201	JVJ	C9-C10-N19	-8.69	116.72	122.99
2	B	1201	JVJ	C33-C32-C6	-5.98	112.52	121.13
2	A	1201	JVJ	C4-C5-C6	-4.25	115.02	121.13
2	A	1201	JVJ	C5-C6-C32	4.14	125.83	117.59
2	B	1201	JVJ	C5-C6-C32	3.93	125.42	117.59
2	B	1201	JVJ	C17-C18-C13	-3.75	106.33	112.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1201	JVJ	C22-C23-C24	-3.67	115.84	121.13
2	A	1201	JVJ	C21-C20-N19	3.02	119.09	115.09
2	A	1201	JVJ	C29-C25-C24	-2.99	123.50	127.58
2	A	1201	JVJ	C23-C24-C30	2.94	123.44	117.59
2	A	1201	JVJ	C32-C6-C7	-2.90	116.01	120.61
2	A	1201	JVJ	C15-N16-C17	2.88	118.61	110.34
2	B	1201	JVJ	C6-C7-C20	-2.83	120.02	124.20
2	A	1201	JVJ	C30-C24-C25	-2.82	116.47	121.36
2	A	1201	JVJ	C30-C31-C21	-2.76	117.16	121.13
2	A	1201	JVJ	C4-C3-C33	2.73	122.45	118.17
2	B	1201	JVJ	C5-C4-C3	-2.70	117.31	121.03
2	B	1201	JVJ	C32-C6-C7	-2.45	116.72	120.61
2	B	1201	JVJ	C14-C13-C12	-2.42	105.22	111.28
2	A	1201	JVJ	C33-C32-C6	-2.21	117.95	121.13
2	A	1201	JVJ	C32-C33-C3	-2.01	118.26	121.03

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1201	JVJ	C13-C12-O11-C10
2	A	1201	JVJ	C13-C12-O11-C10
2	A	1201	JVJ	O11-C12-C13-C18
2	A	1201	JVJ	C30-C24-C25-C29
2	A	1201	JVJ	O11-C12-C13-C14
2	A	1201	JVJ	C7-C20-C21-C22
2	A	1201	JVJ	C7-C20-C21-C31

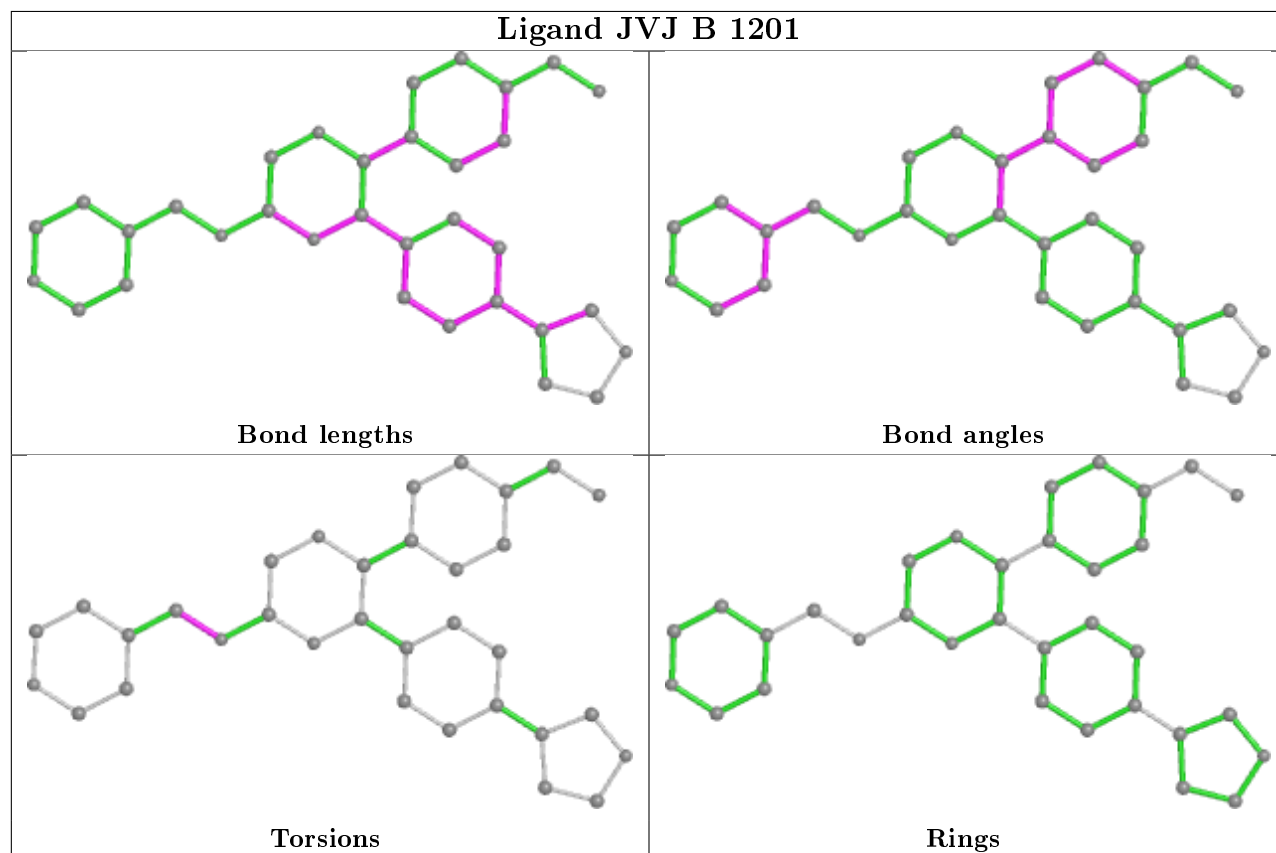
There are no ring outliers.

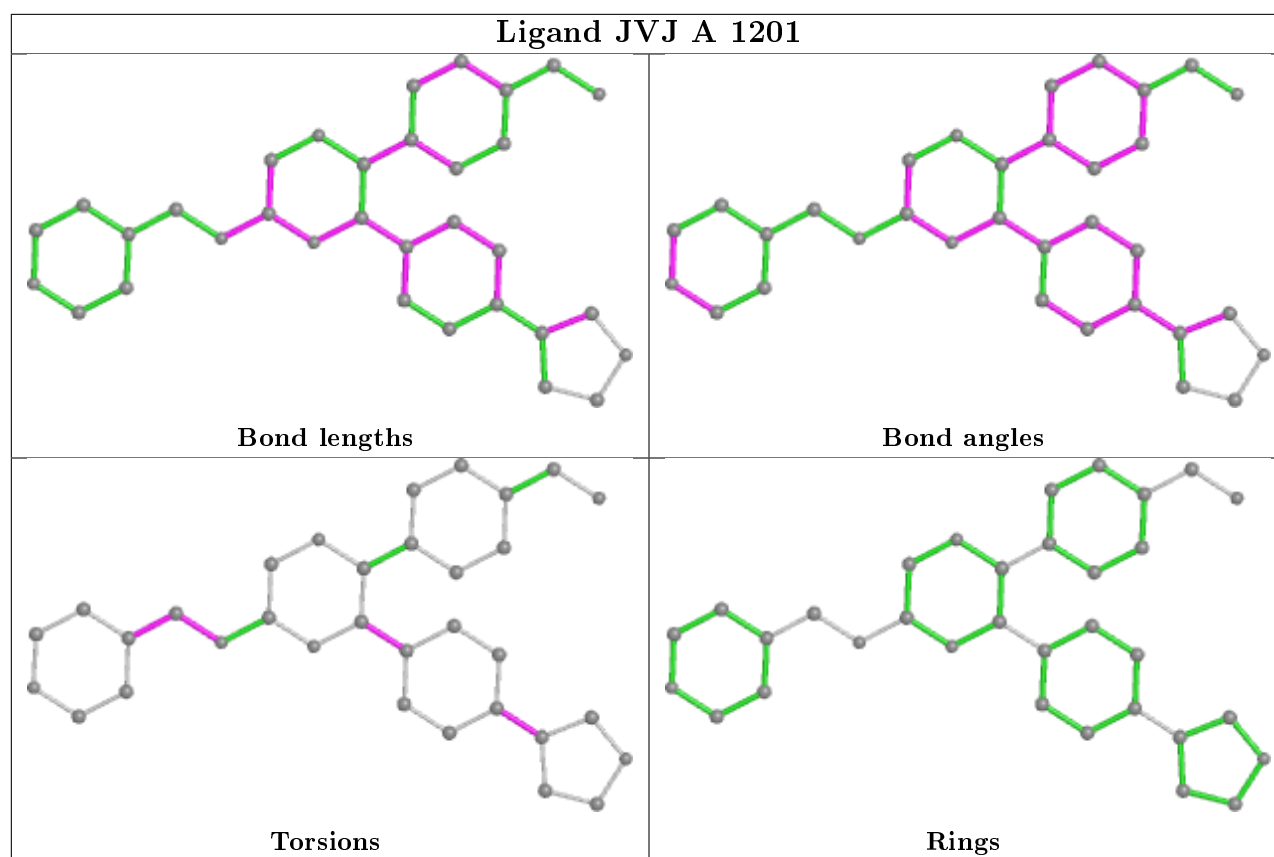
2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1201	JVJ	5	0
2	A	1201	JVJ	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring

in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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, 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	164/247 (66%)	0.39	7 (4%) 35 33	35, 51, 89, 128	0
1	B	164/247 (66%)	0.45	13 (7%) 12 10	33, 50, 99, 115	0
All	All	328/494 (66%)	0.42	20 (6%) 21 20	33, 51, 97, 128	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1169	GLU	7.4
1	B	1018	LEU	4.7
1	B	1030	ILE	4.1
1	A	1033	TYR	3.7
1	A	65	ALA	3.5
1	B	1041	TYR	2.8
1	A	1031	LEU	2.8
1	B	1065	ILE	2.6
1	B	1033	TYR	2.5
1	B	1120	THR	2.5
1	B	1100	LEU	2.4
1	A	1120	THR	2.4
1	B	1103	GLY	2.3
1	B	1063	LYS	2.2
1	A	1057	VAL	2.2
1	B	1043	GLU	2.1
1	A	1091	GLU	2.1
1	B	56	ALA	2.1
1	B	1059	MET	2.1
1	B	1088	GLU	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](#)

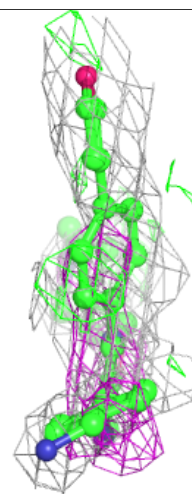
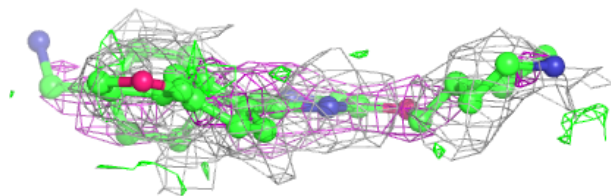
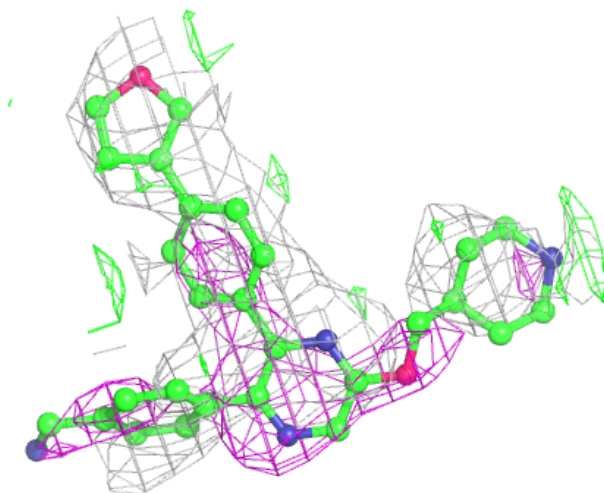
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	JVJ	B	1201	33/33	0.59	0.47	33,46,54,73	0
2	JVJ	A	1201	33/33	0.65	0.49	37,45,52,56	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

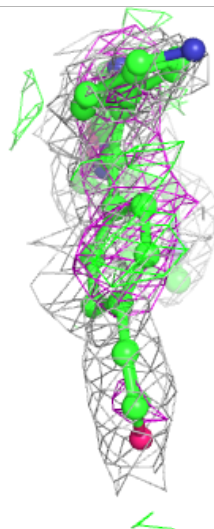
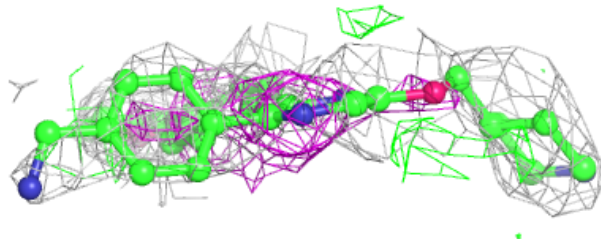
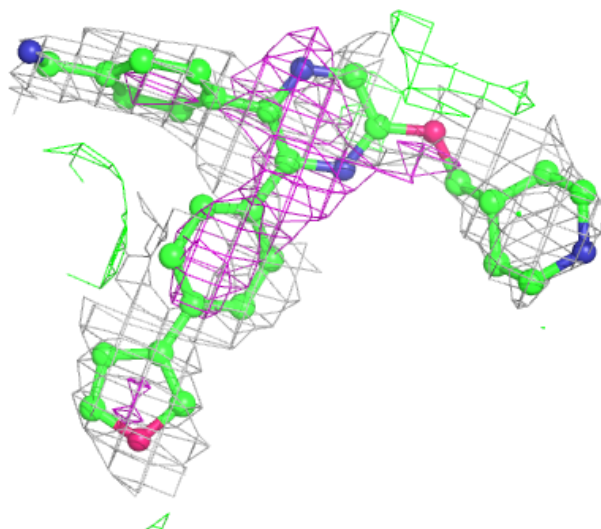
Electron density around JVJ B 1201:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around JVJ A 1201:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers ⓘ

There are no such residues in this entry.