



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2023 – 01:01 AM EDT

PDB ID : 7MBI  
Title : Structure of SARS-CoV2 3CL protease covalently bound to peptidomimetic inhibitor  
Authors : Khan, M.B.; Lu, J.; Young, H.S.; Lemieux, M.J.  
Deposited on : 2021-03-31  
Resolution : 2.15 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 2.35.1  
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) : 2.35.1

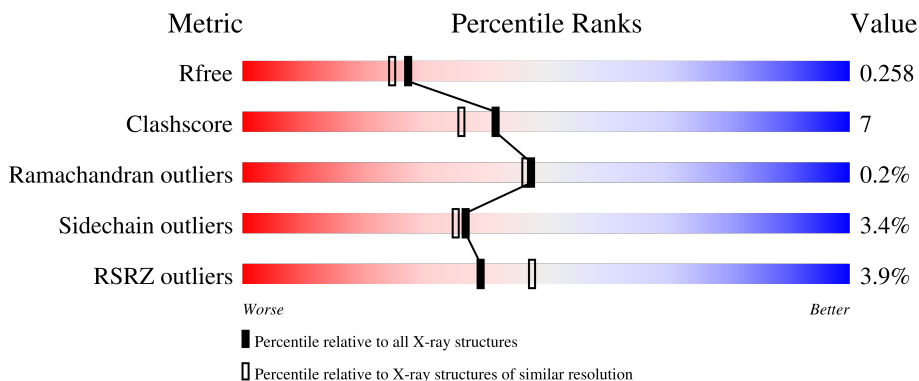
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.15 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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 of 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	306	 2% (red), 87% (green), 11% (yellow), .. (grey)
1	B	306	 5% (red), 82% (green), 16% (yellow), . (grey)
1	C	306	 6% (red), 79% (green), 19% (yellow), .. (grey)
1	D	306	 4% (red), 80% (green), 17% (yellow), .. (grey)

## 2 Entry composition [i](#)

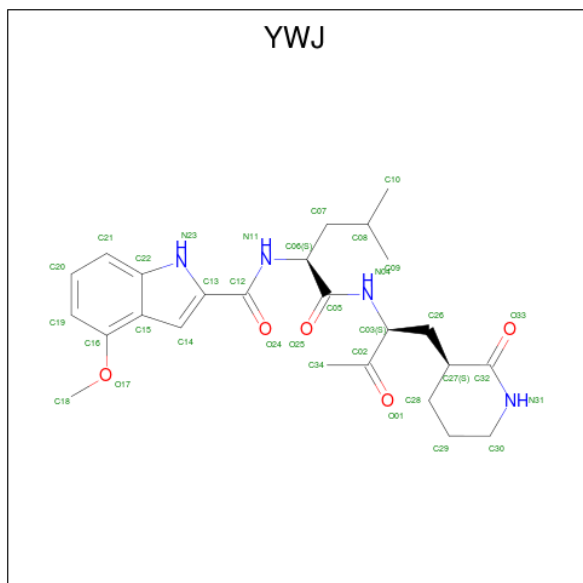
There are 4 unique types of molecules in this entry. The entry contains 19089 atoms, of which 9353 are hydrogens and 0 are deuteriums.

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 3C-like proteinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
1	A	303	Total 4638	C 1483	H 2295	N 398	O 439	S 23	0	1	0
1	B	306	Total 4669	C 1501	H 2299	N 402	O 444	S 23	8	1	0
1	C	305	Total 4667	C 1496	H 2306	N 400	O 442	S 23	0	1	0
1	D	306	Total 4688	C 1501	H 2318	N 402	O 444	S 23	0	1	0

- Molecule 2 is 4-methoxy-N-[(2S)-4-methyl-1-oxo-1-((2S)-3-oxo-1-[(3S)-2-oxopiperidin-3-yl]butan-2-yl)amino)pentan-2-yl]-1H-indole-2-carboxamide (three-letter code: YWJ) (formula: C<sub>25</sub>H<sub>34</sub>N<sub>4</sub>O<sub>5</sub>) (labeled as "Ligand of Interest" by depositor).



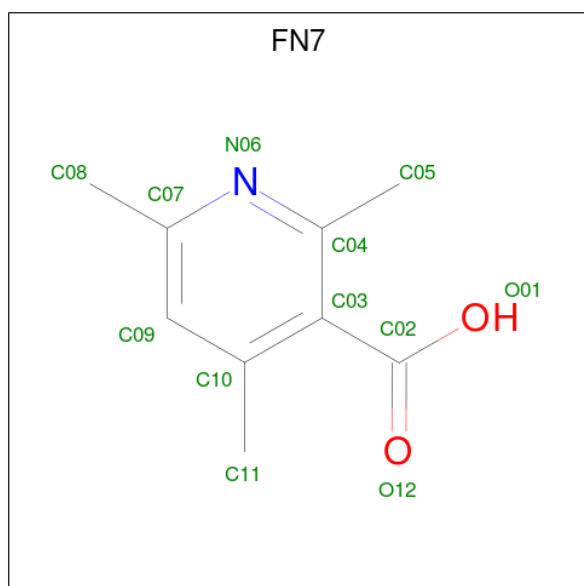
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	H	N	O		
2	A	1	Total 65	C 25	H 31	N 4	O 5	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	H	N	O		
2	B	1	Total	C	H	N	O	0	0
			65	25	31	4	5		
2	C	1	Total	C	H	N	O	0	0
			66	25	32	4	5		
2	D	1	Total	C	H	N	O	0	0
			65	25	31	4	5		

- Molecule 3 is 2,4,6-trimethylpyridine-3-carboxylic acid (three-letter code: FN7) (formula:  $C_9H_{11}NO_2$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	H	N	O		
3	B	1	Total	C	H	N	O	0	0
			22	9	10	1	2		

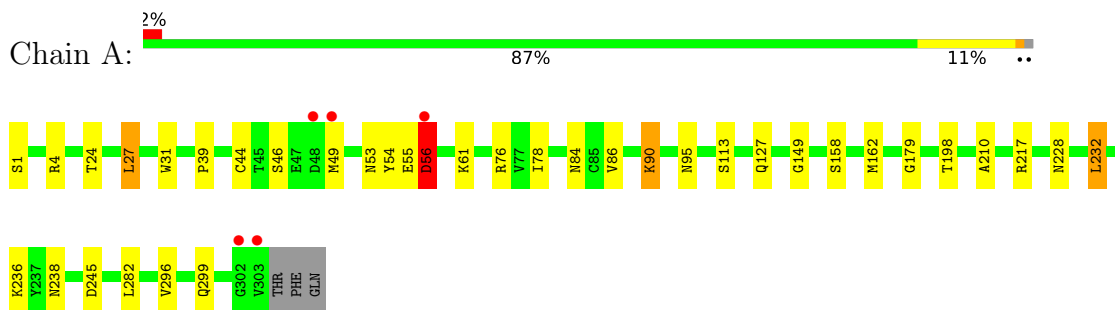
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	56	Total	O	0	0
			56	56		
4	B	22	Total	O	0	0
			22	22		
4	C	40	Total	O	0	0
			40	40		
4	D	26	Total	O	0	0
			26	26		

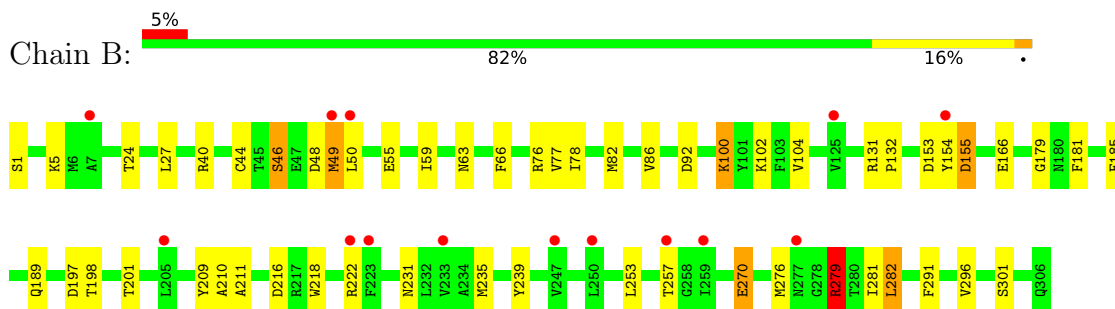
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

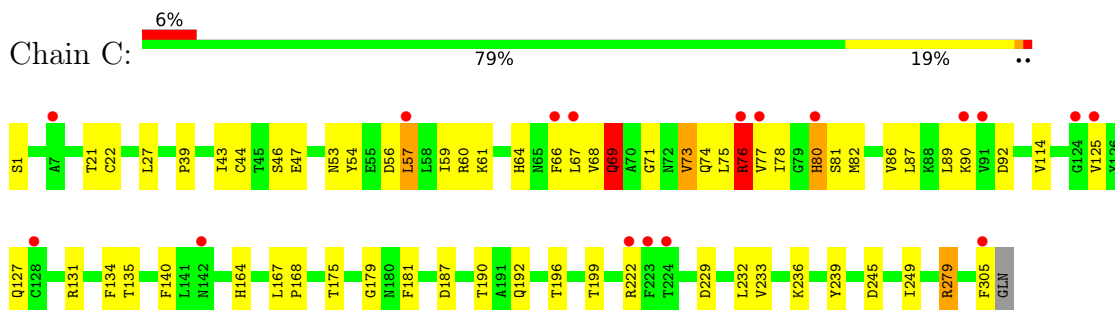
- Molecule 1: 3C-like proteinase



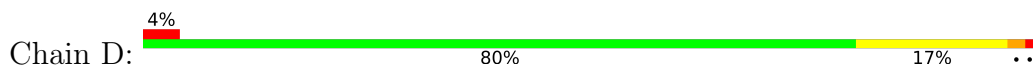
- Molecule 1: 3C-like proteinase

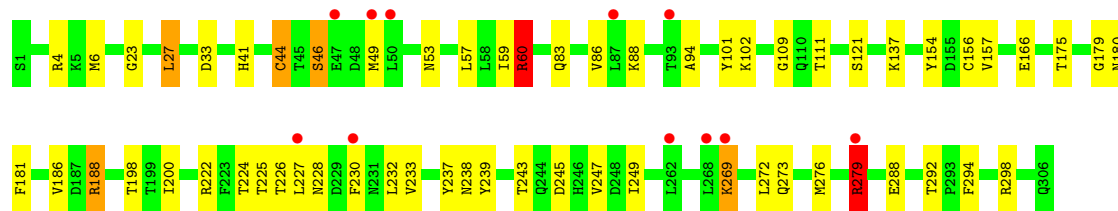


- Molecule 1: 3C-like proteinase



- Molecule 1: 3C-like proteinase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.02Å 123.43Å 139.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.14 – 2.15 39.46 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.7 (37.14-2.15) 99.2 (39.46-1.95)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.09 (at 1.95Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, $R_{free}$	0.204 , 0.261 0.204 , 0.258	Depositor DCC
$R_{free}$ test set	4615 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	46.8	Xtrriage
Anisotropy	0.284	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 48.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	19089	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	68.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.64% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: YWJ, FN7

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	0.68	1/2398 (0.0%)	0.97	6/3259 (0.2%)
1	B	0.74	3/2426 (0.1%)	0.93	11/3297 (0.3%)
1	C	0.69	2/2417 (0.1%)	0.92	9/3285 (0.3%)
1	D	0.67	3/2426 (0.1%)	0.94	13/3297 (0.4%)
All	All	0.70	9/9667 (0.1%)	0.94	39/13138 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	3
1	C	0	2
1	D	0	2
All	All	0	9

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	49	MET	CG-SD	-13.79	1.45	1.81
1	D	269	LYS	CB-CG	-9.81	1.26	1.52
1	C	90	LYS	CD-CE	7.37	1.69	1.51
1	C	47	GLU	CB-CG	6.95	1.65	1.52
1	B	270	GLU	CB-CG	6.00	1.63	1.52
1	A	56	ASP	CB-CG	5.63	1.63	1.51
1	D	44	CYS	CB-SG	-5.09	1.73	1.81
1	B	44	CYS	CB-SG	-5.08	1.73	1.81
1	D	269	LYS	CE-NZ	5.08	1.61	1.49



All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	232	LEU	CB-CG-CD2	15.25	136.92	111.00
1	A	56	ASP	CB-CG-OD1	-15.05	104.75	118.30
1	B	49	MET	CA-CB-CG	14.41	137.81	113.30
1	A	56	ASP	CB-CG-OD2	13.36	130.33	118.30
1	D	279	ARG	CG-CD-NE	12.57	138.20	111.80
1	C	57	LEU	CB-CG-CD1	11.85	131.15	111.00
1	D	188	ARG	CD-NE-CZ	-11.65	107.30	123.60
1	D	269	LYS	CA-CB-CG	11.23	138.11	113.40
1	B	279	ARG	CB-CG-CD	10.71	139.45	111.60
1	D	188	ARG	CB-CG-CD	-10.49	84.33	111.60
1	C	76	ARG	CG-CD-NE	10.43	133.69	111.80
1	B	49	MET	CB-CA-C	-10.15	90.10	110.40
1	B	279	ARG	CB-CA-C	-9.56	91.29	110.40
1	A	56	ASP	CB-CA-C	9.34	129.09	110.40
1	C	57	LEU	CB-CG-CD2	-9.22	95.33	111.00
1	D	188	ARG	CA-CB-CG	8.48	132.06	113.40
1	D	279	ARG	CD-NE-CZ	-8.32	111.94	123.60
1	C	90	LYS	CD-CE-NZ	-8.24	92.75	111.70
1	C	57	LEU	CA-CB-CG	8.05	133.81	115.30
1	A	56	ASP	N-CA-CB	-8.04	96.12	110.60
1	B	270	GLU	CB-CA-C	-7.84	94.72	110.40
1	B	49	MET	N-CA-CB	7.67	124.41	110.60
1	C	90	LYS	CG-CD-CE	-7.65	88.94	111.90
1	B	279	ARG	NE-CZ-NH1	7.65	124.13	120.30
1	D	188	ARG	CG-CD-NE	7.40	127.34	111.80
1	B	279	ARG	N-CA-CB	7.30	123.73	110.60
1	C	90	LYS	CA-CB-CG	6.99	128.78	113.40
1	D	188	ARG	NE-CZ-NH1	-6.98	116.81	120.30
1	B	270	GLU	CA-CB-CG	6.46	127.62	113.40
1	B	279	ARG	CD-NE-CZ	-6.45	114.57	123.60
1	D	279	ARG	CA-CB-CG	-6.21	99.73	113.40
1	C	46	SER	N-CA-CB	6.16	119.74	110.50
1	A	232	LEU	CA-CB-CG	-6.04	101.40	115.30
1	D	279	ARG	NE-CZ-NH2	-5.94	117.33	120.30
1	D	6	MET	CG-SD-CE	5.67	109.27	100.20
1	D	269	LYS	N-CA-CB	5.59	120.66	110.60
1	C	69	GLN	CA-CB-CG	-5.37	101.59	113.40
1	B	50	LEU	CA-CB-CG	5.29	127.47	115.30
1	D	188	ARG	CB-CA-C	-5.10	100.20	110.40

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	55	GLU	Peptide
1	A	56	ASP	Mainchain
1	B	270	GLU	Sidechain
1	B	279	ARG	Sidechain
1	B	48	ASP	Peptide
1	C	76	ARG	Sidechain
1	C	80	HIS	Sidechain
1	D	279	ARG	Sidechain
1	D	60	ARG	Sidechain

## 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	2343	2295	2294	22	8
1	B	2370	2299	2318	32	0
1	C	2361	2306	2310	44	2
1	D	2370	2318	2318	36	5
2	A	34	31	0	0	0
2	B	34	31	0	0	0
2	C	34	32	0	0	0
2	D	34	31	0	1	0
3	B	12	10	0	1	3
4	A	56	0	0	4	0
4	B	22	0	0	0	0
4	C	40	0	0	3	0
4	D	26	0	0	0	0
All	All	9736	9353	9240	131	9

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 7.

All (131) 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:61:LYS:NZ	4:A:503:HOH:O	2.00	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:SER:OG	4:A:501:HOH:O	1.87	0.90
1:B:49:MET:HG2	1:B:189:GLN:HG3	1.54	0.87
1:D:186:VAL:HG23	1:D:188:ARG:HG3	1.57	0.86
1:A:84:ASN:OD1	4:A:502:HOH:O	1.99	0.79
1:C:199:THR:HG21	1:C:239:TYR:CZ	2.18	0.79
1:B:211:ALA:HA	1:B:282:LEU:HD21	1.65	0.78
1:C:279:ARG:HG2	1:C:279:ARG:HH21	1.50	0.76
1:A:127:GLN:NE2	4:A:504:HOH:O	2.13	0.75
1:B:49:MET:HG2	1:B:189:GLN:CG	2.16	0.74
1:A:53:ASN:OD1	1:A:56:ASP:HB3	1.88	0.71
1:D:245:ASP:O	1:D:249:ILE:HD12	1.92	0.70
1:D:186:VAL:CG2	1:D:188:ARG:HG3	2.20	0.70
1:C:43:ILE:HD12	1:C:57:LEU:HB3	1.76	0.67
1:C:67:LEU:HD11	1:C:74:GLN:NE2	2.10	0.66
1:D:60:ARG:HG3	1:D:60:ARG:HH21	1.60	0.65
1:D:233:VAL:HB	1:D:269:LYS:HE2	1.77	0.65
1:A:53:ASN:OD1	1:A:56:ASP:CB	2.45	0.65
1:B:253:LEU:O	1:B:257:THR:HG23	1.97	0.64
1:A:198:THR:HG22	1:A:238:ASN:HD21	1.63	0.63
1:C:67:LEU:HD11	1:C:74:GLN:HE22	1.61	0.63
1:D:109:GLY:HA2	1:D:200:ILE:HD13	1.81	0.62
1:C:21:THR:HB	1:C:67:LEU:HD23	1.82	0.62
1:C:68:VAL:HG23	1:C:75:LEU:HB2	1.81	0.61
1:B:153:ASP:O	1:B:155:ASP:N	2.34	0.61
1:B:55:GLU:O	1:B:59:ILE:HG13	2.01	0.60
1:A:232:LEU:HD12	1:A:232:LEU:N	2.16	0.60
1:C:305:PHE:O	4:C:501:HOH:O	2.15	0.60
1:D:102:LYS:HG3	1:D:156:CYS:SG	2.42	0.60
1:C:86:VAL:HG23	1:C:179:GLY:HA2	1.84	0.59
1:A:27:LEU:HD22	1:A:39:PRO:HG2	1.83	0.59
1:D:188:ARG:HH11	1:D:188:ARG:CG	2.13	0.59
1:B:1:SER:N	1:C:140:PHE:O	2.35	0.59
1:C:44:CYS:HG	1:C:54:TYR:HE1	1.51	0.58
1:C:76:ARG:HD3	1:C:78:ILE:HG22	1.84	0.58
1:B:209:TYR:HD2	1:B:257:THR:HG21	1.67	0.57
1:C:56:ASP:O	1:C:59:ILE:HG13	2.04	0.57
1:A:228:ASN:O	1:A:232:LEU:HD13	2.06	0.56
1:C:87:LEU:HD21	1:C:89:LEU:HD21	1.87	0.56
1:A:46:SER:O	1:A:49:MET:HG3	2.05	0.55
1:C:229:ASP:HA	1:C:232:LEU:HD12	1.87	0.55
1:B:218:TRP:CZ2	1:B:279:ARG:HG3	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:222:ARG:HA	1:D:222:ARG:HE	1.72	0.55
1:A:39:PRO:HD3	1:A:162:MET:SD	2.47	0.55
1:A:86:VAL:HG13	1:A:179:GLY:HA2	1.89	0.54
1:B:218:TRP:CE2	1:B:279:ARG:HG3	2.41	0.54
1:C:245:ASP:O	1:C:249:ILE:HG13	2.07	0.54
1:C:199:THR:HG21	1:C:239:TYR:CE2	2.42	0.54
1:D:53:ASN:O	1:D:57:LEU:HD13	2.08	0.54
1:C:66:PHE:HB2	1:C:77:VAL:HG21	1.89	0.53
1:C:134:PHE:O	4:C:502:HOH:O	2.18	0.52
1:B:76:ARG:HB3	1:B:92:ASP:OD2	2.10	0.52
1:D:86:VAL:HG13	1:D:179:GLY:HA2	1.92	0.52
1:C:175:THR:HG22	1:C:181:PHE:HA	1.92	0.52
1:A:232:LEU:N	1:A:232:LEU:CD1	2.73	0.51
1:B:66:PHE:HB2	1:B:77:VAL:HG21	1.91	0.51
1:C:39:PRO:HB3	1:C:164:HIS:NE2	2.25	0.51
1:C:233:VAL:HA	1:C:236:LYS:HE3	1.93	0.51
1:A:76:ARG:HG2	1:A:76:ARG:HH11	1.76	0.51
1:C:127:GLN:HG2	4:C:514:HOH:O	2.12	0.50
1:A:76:ARG:HH11	1:A:76:ARG:CG	2.23	0.50
1:C:69:GLN:OE1	1:C:71:GLY:O	2.29	0.50
1:D:166:GLU:O	2:D:401:YWJ:N23	2.45	0.50
1:B:100:LYS:HE2	1:B:155:ASP:OD2	2.11	0.50
1:B:46:SER:HA	1:B:49:MET:SD	2.53	0.48
1:C:53:ASN:O	1:C:57:LEU:HD23	2.12	0.48
1:C:59:ILE:HD12	1:C:60:ARG:N	2.28	0.48
1:D:243:THR:O	1:D:247:VAL:HG23	2.13	0.48
1:D:230:PHE:HA	1:D:269:LYS:NZ	2.29	0.48
1:C:67:LEU:CD1	1:C:74:GLN:HE22	2.25	0.47
1:D:225:THR:HG22	1:D:226:THR:O	2.14	0.47
1:C:190:THR:O	1:C:192:GLN:HG3	2.14	0.47
1:B:102:LYS:HD2	1:B:104:VAL:CG1	2.45	0.47
1:B:218:TRP:CZ3	1:B:279:ARG:NH2	2.84	0.46
1:D:188:ARG:HG3	1:D:188:ARG:HH11	1.79	0.46
1:C:279:ARG:HG2	1:C:279:ARG:NH2	2.26	0.46
1:D:198:THR:HG22	1:D:238:ASN:OD1	2.16	0.46
1:C:187:ASP:OD1	1:C:187:ASP:N	2.49	0.46
1:B:132:PRO:HD2	1:B:197:ASP:OD1	2.16	0.46
1:D:228:ASN:O	1:D:232:LEU:HD22	2.16	0.46
1:B:40:ARG:HH11	1:B:82:MET:CE	2.29	0.46
1:D:225:THR:CG2	1:D:226:THR:N	2.79	0.45
1:C:69:GLN:HG3	1:C:74:GLN:HG2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:46:SER:HA	1:D:49:MET:SD	2.56	0.45
1:D:228:ASN:O	1:D:232:LEU:CD2	2.64	0.45
1:C:232:LEU:O	1:C:236:LYS:HE3	2.16	0.45
1:D:233:VAL:HB	1:D:269:LYS:CE	2.46	0.45
1:C:114:VAL:O	1:C:125:VAL:HA	2.17	0.45
1:D:225:THR:HG22	1:D:226:THR:N	2.32	0.45
1:D:111:THR:HG23	1:D:292:THR:HG23	1.99	0.44
1:C:67:LEU:CD1	1:C:74:GLN:NE2	2.80	0.44
1:D:175:THR:HG22	1:D:181:PHE:HA	2.00	0.44
1:D:239:TYR:CZ	1:D:272:LEU:HD21	2.53	0.44
1:C:167:LEU:HB3	1:C:168:PRO:HD2	2.00	0.44
1:B:231:ASN:O	1:B:235:MET:HG2	2.17	0.44
1:C:53:ASN:OD1	1:C:56:ASP:OD1	2.36	0.44
1:D:269:LYS:O	1:D:273:GLN:HG3	2.18	0.44
1:C:68:VAL:HG22	1:C:75:LEU:O	2.18	0.43
1:B:166:GLU:OE1	1:C:1:SER:OG	2.34	0.43
1:C:69:GLN:HG2	1:C:73:VAL:O	2.18	0.43
1:B:86:VAL:HG23	1:B:179:GLY:HA2	2.01	0.43
1:C:54:TYR:HA	1:C:57:LEU:HB2	2.01	0.43
1:A:210:ALA:HB2	1:A:296:VAL:HG13	2.00	0.42
1:D:101:TYR:HA	1:D:157:VAL:O	2.19	0.42
1:B:24:THR:O	3:B:402:FN7:C03	2.67	0.42
1:C:86:VAL:HG23	1:C:179:GLY:CA	2.49	0.42
1:D:33:ASP:O	1:D:94:ALA:HA	2.19	0.42
1:D:294:PHE:CE2	1:D:298:ARG:HG3	2.55	0.42
1:A:282:LEU:HD23	1:A:282:LEU:HA	1.92	0.42
1:B:181:PHE:CD1	1:B:185:PHE:HB2	2.55	0.42
1:B:63:ASN:ND2	1:B:78:ILE:C	2.74	0.41
1:D:27:LEU:HD23	1:D:27:LEU:C	2.40	0.41
1:D:237:TYR:OH	1:D:273:GLN:HA	2.21	0.41
1:B:5:LYS:HG2	1:B:291:PHE:CZ	2.55	0.41
1:B:235:MET:HE2	1:D:180:ASN:HB3	2.02	0.41
1:C:131:ARG:HG2	1:C:135:THR:O	2.21	0.41
1:D:41:HIS:O	1:D:44:CYS:HB2	2.21	0.41
1:D:83:GLN:OE1	1:D:88:LYS:HD2	2.21	0.41
1:A:78:ILE:HG13	1:A:90:LYS:HG3	2.02	0.41
1:B:210:ALA:HB2	1:B:296:VAL:HG13	2.03	0.41
1:A:31:TRP:CD2	1:A:95:ASN:HB2	2.56	0.41
1:D:276:MET:O	1:D:279:ARG:HB2	2.20	0.41
1:A:44:CYS:SG	1:A:54:TYR:CE1	3.14	0.41
1:B:276:MET:CE	1:B:281:ILE:HD12	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:43:ILE:CD1	1:C:57:LEU:HB3	2.48	0.41
1:B:63:ASN:HB3	1:B:77:VAL:O	2.21	0.40
1:B:66:PHE:CB	1:B:77:VAL:HG21	2.51	0.40
1:B:76:ARG:N	1:B:92:ASP:OD2	2.54	0.40
1:B:201:THR:OG1	1:B:239:TYR:HB3	2.20	0.40
1:A:113:SER:O	1:A:149:GLY:HA2	2.21	0.40
1:C:68:VAL:HG23	1:C:68:VAL:O	2.22	0.40

All (9) 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:24:THR:O	3:B:402:FN7:C03[4_456]	1.42	0.78
1:A:4:ARG:NH1	1:D:137:LYS:O[1_455]	1.62	0.58
1:C:196:THR:OG1	1:D:23:GLY:O[3_655]	1.80	0.40
1:A:127:GLN:O	1:D:4:ARG:NH2[1_455]	1.81	0.39
1:A:4:ARG:HH12	1:D:137:LYS:O[1_455]	1.29	0.31
1:A:24:THR:O	3:B:402:FN7:C04[4_456]	1.99	0.21
1:A:24:THR:O	3:B:402:FN7:C10[4_456]	2.01	0.19
1:A:127:GLN:O	1:D:4:ARG:HH21[1_455]	1.47	0.13
1:A:217:ARG:HH22	1:C:245:ASP:OD1[2_564]	1.50	0.10

## 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	302/306 (99%)	295 (98%)	7 (2%)	0	100 100
1	B	305/306 (100%)	294 (96%)	10 (3%)	1 (0%)	41 37
1	C	304/306 (99%)	298 (98%)	6 (2%)	0	100 100
1	D	305/306 (100%)	299 (98%)	5 (2%)	1 (0%)	41 37

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1216/1224 (99%)	1186 (98%)	28 (2%)	2 (0%)	47 46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	154	TYR
1	D	154	TYR

### 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	261/263 (99%)	255 (98%)	6 (2%)	50 53
1	B	264/263 (100%)	254 (96%)	10 (4%)	33 31
1	C	263/263 (100%)	251 (95%)	12 (5%)	27 23
1	D	264/263 (100%)	256 (97%)	8 (3%)	41 40
All	All	1052/1052 (100%)	1016 (97%)	36 (3%)	37 35

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	SER
1	A	27	LEU
1	A	90	LYS
1	A	236	LYS
1	A	245	ASP
1	A	299	GLN
1	B	27	LEU
1	B	46	SER
1	B	100	LYS
1	B	131	ARG
1	B	155	ASP
1	B	198	THR
1	B	216	ASP

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Mol	Chain	Res	Type
1	B	222	ARG
1	B	282	LEU
1	B	301	SER
1	C	22	CYS
1	C	27	LEU
1	C	61	LYS
1	C	64	HIS
1	C	69	GLN
1	C	73	VAL
1	C	80	HIS
1	C	81	SER
1	C	82	MET
1	C	92	ASP
1	C	222	ARG
1	C	279	ARG
1	D	27	LEU
1	D	46	SER
1	D	59	ILE
1	D	60	ARG
1	D	121	SER
1	D	224	THR
1	D	227	LEU
1	D	288	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	110	GLN
1	A	238	ASN
1	A	277	ASN
1	B	80	HIS
1	B	228	ASN
1	B	256	GLN
1	C	41	HIS
1	C	53	ASN
1	C	63	ASN
1	C	72	ASN
1	C	74	GLN
1	C	80	HIS



### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

5 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	YWJ	C	401	1	34,36,36	2.98	11 (32%)	39,50,50	1.60	8 (20%)
2	YWJ	B	401	1	34,36,36	2.96	10 (29%)	39,50,50	1.73	9 (23%)
2	YWJ	D	401	1	34,36,36	2.87	10 (29%)	39,50,50	1.50	9 (23%)
3	FN7	B	402	-	12,12,12	1.74	3 (25%)	16,17,17	1.48	3 (18%)
2	YWJ	A	401	1	34,36,36	2.94	9 (26%)	39,50,50	1.43	5 (12%)

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	YWJ	C	401	1	-	8/27/41/41	0/3/3/3
2	YWJ	B	401	1	-	5/27/41/41	0/3/3/3
2	YWJ	D	401	1	-	2/27/41/41	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	FN7	B	402	-	-	0/4/4/4	0/1/1/1
2	YWJ	A	401	1	-	4/27/41/41	0/3/3/3

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	YWJ	C32-N31	12.32	1.55	1.33
2	D	401	YWJ	C32-N31	11.57	1.53	1.33
2	A	401	YWJ	C32-N31	11.15	1.53	1.33
2	B	401	YWJ	C32-N31	11.04	1.52	1.33
2	A	401	YWJ	C27-C32	7.48	1.61	1.50
2	B	401	YWJ	C27-C32	7.14	1.60	1.50
2	C	401	YWJ	C27-C32	6.09	1.59	1.50
2	B	401	YWJ	C05-N04	5.58	1.46	1.34
2	D	401	YWJ	C27-C32	5.52	1.58	1.50
2	A	401	YWJ	C05-N04	5.39	1.45	1.34
2	A	401	YWJ	C12-N11	5.11	1.45	1.34
2	B	401	YWJ	C12-N11	4.84	1.44	1.34
2	C	401	YWJ	C12-N11	4.72	1.44	1.34
2	C	401	YWJ	C05-N04	4.65	1.44	1.34
2	D	401	YWJ	C12-N11	4.62	1.44	1.34
2	D	401	YWJ	C05-N04	3.99	1.42	1.34
2	C	401	YWJ	C28-C27	-3.71	1.43	1.53
2	A	401	YWJ	C28-C27	-3.66	1.43	1.53
2	D	401	YWJ	C28-C27	-3.59	1.43	1.53
2	D	401	YWJ	C30-N31	3.55	1.54	1.46
2	B	401	YWJ	C28-C27	-3.42	1.44	1.53
2	D	401	YWJ	C29-C28	-3.31	1.44	1.53
2	C	401	YWJ	C30-N31	3.14	1.53	1.46
3	B	402	FN7	O12-C02	3.04	1.31	1.22
2	B	401	YWJ	O17-C16	2.87	1.41	1.36
2	A	401	YWJ	C26-C27	2.83	1.60	1.53
2	A	401	YWJ	C30-N31	2.77	1.52	1.46
2	C	401	YWJ	C29-C28	-2.74	1.46	1.53
3	B	402	FN7	C03-C10	-2.70	1.37	1.40
2	D	401	YWJ	O24-C12	-2.66	1.17	1.23
2	B	401	YWJ	O24-C12	-2.64	1.17	1.23
2	D	401	YWJ	O17-C16	2.62	1.41	1.36
2	B	401	YWJ	C29-C28	-2.60	1.46	1.53
2	B	401	YWJ	C26-C27	2.49	1.59	1.53
2	C	401	YWJ	O25-C05	-2.46	1.18	1.23
2	A	401	YWJ	C29-C28	-2.38	1.46	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	FN7	C03-C04	-2.31	1.37	1.41
2	C	401	YWJ	O24-C12	-2.17	1.18	1.23
2	D	401	YWJ	O25-C05	-2.13	1.19	1.23
2	B	401	YWJ	C30-N31	2.09	1.51	1.46
2	A	401	YWJ	O17-C16	2.05	1.40	1.36
2	C	401	YWJ	C15-C22	-2.04	1.37	1.42
2	C	401	YWJ	C26-C27	2.03	1.58	1.53

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	YWJ	O17-C16-C15	4.92	119.27	114.46
2	A	401	YWJ	O17-C16-C15	4.34	118.70	114.46
2	D	401	YWJ	C30-N31-C32	-3.68	115.13	125.79
2	C	401	YWJ	C30-N31-C32	-3.59	115.37	125.79
2	C	401	YWJ	C18-O17-C16	-3.58	112.73	117.75
2	B	401	YWJ	C18-O17-C16	-3.55	112.78	117.75
2	B	401	YWJ	C13-C12-N11	3.51	121.72	115.20
2	A	401	YWJ	C18-O17-C16	-3.38	113.02	117.75
2	C	401	YWJ	C29-C30-N31	3.18	117.22	110.24
2	A	401	YWJ	C30-N31-C32	-3.15	116.65	125.79
2	C	401	YWJ	C13-C12-N11	3.15	121.05	115.20
2	C	401	YWJ	C08-C07-C06	-3.15	106.77	115.43
3	B	402	FN7	C05-C04-C03	-3.13	118.29	122.27
2	D	401	YWJ	C26-C03-N04	-2.96	103.76	110.58
2	B	401	YWJ	C30-N31-C32	-2.79	117.70	125.79
2	C	401	YWJ	C13-N23-C22	2.77	110.22	104.45
2	B	401	YWJ	O25-C05-N04	-2.76	117.82	122.93
2	A	401	YWJ	C13-N23-C22	2.70	110.09	104.45
2	D	401	YWJ	C29-C30-N31	2.69	116.14	110.24
2	B	401	YWJ	O33-C32-N31	-2.69	118.64	122.38
2	B	401	YWJ	C13-N23-C22	2.67	110.02	104.45
2	B	401	YWJ	C26-C03-N04	-2.65	104.47	110.58
2	D	401	YWJ	O25-C05-N04	-2.63	118.05	122.93
2	D	401	YWJ	C18-O17-C16	-2.63	114.06	117.75
2	D	401	YWJ	C28-C27-C32	2.60	119.58	112.01
2	A	401	YWJ	O33-C32-N31	-2.55	118.82	122.38
2	C	401	YWJ	C14-C15-C22	2.53	108.48	106.27
3	B	402	FN7	C05-C04-N06	2.46	120.30	116.49
2	D	401	YWJ	C13-C12-N11	2.37	119.61	115.20
2	D	401	YWJ	C08-C07-C06	-2.28	109.15	115.43
2	C	401	YWJ	O24-C12-C13	-2.17	116.35	121.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	YWJ	O24-C12-C13	-2.08	116.54	121.08
3	B	402	FN7	C04-C03-C02	-2.04	116.78	119.92
2	D	401	YWJ	O25-C05-C06	2.01	124.69	120.45

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	YWJ	C15-C16-O17-C18
2	A	401	YWJ	C19-C16-O17-C18
2	B	401	YWJ	C19-C16-O17-C18
2	D	401	YWJ	C34-C02-C03-N04
2	B	401	YWJ	C15-C16-O17-C18
2	C	401	YWJ	C06-C07-C08-C09
2	B	401	YWJ	C06-C07-C08-C09
2	C	401	YWJ	C06-C07-C08-C10
2	C	401	YWJ	C19-C16-O17-C18
2	C	401	YWJ	C15-C16-O17-C18
2	B	401	YWJ	C06-C07-C08-C10
2	A	401	YWJ	C34-C02-C03-C26
2	A	401	YWJ	O01-C02-C03-C26
2	C	401	YWJ	O01-C02-C03-N04
2	D	401	YWJ	C03-C26-C27-C28
2	B	401	YWJ	C34-C02-C03-N04
2	C	401	YWJ	C34-C02-C03-N04
2	C	401	YWJ	C34-C02-C03-C26
2	C	401	YWJ	O01-C02-C03-C26

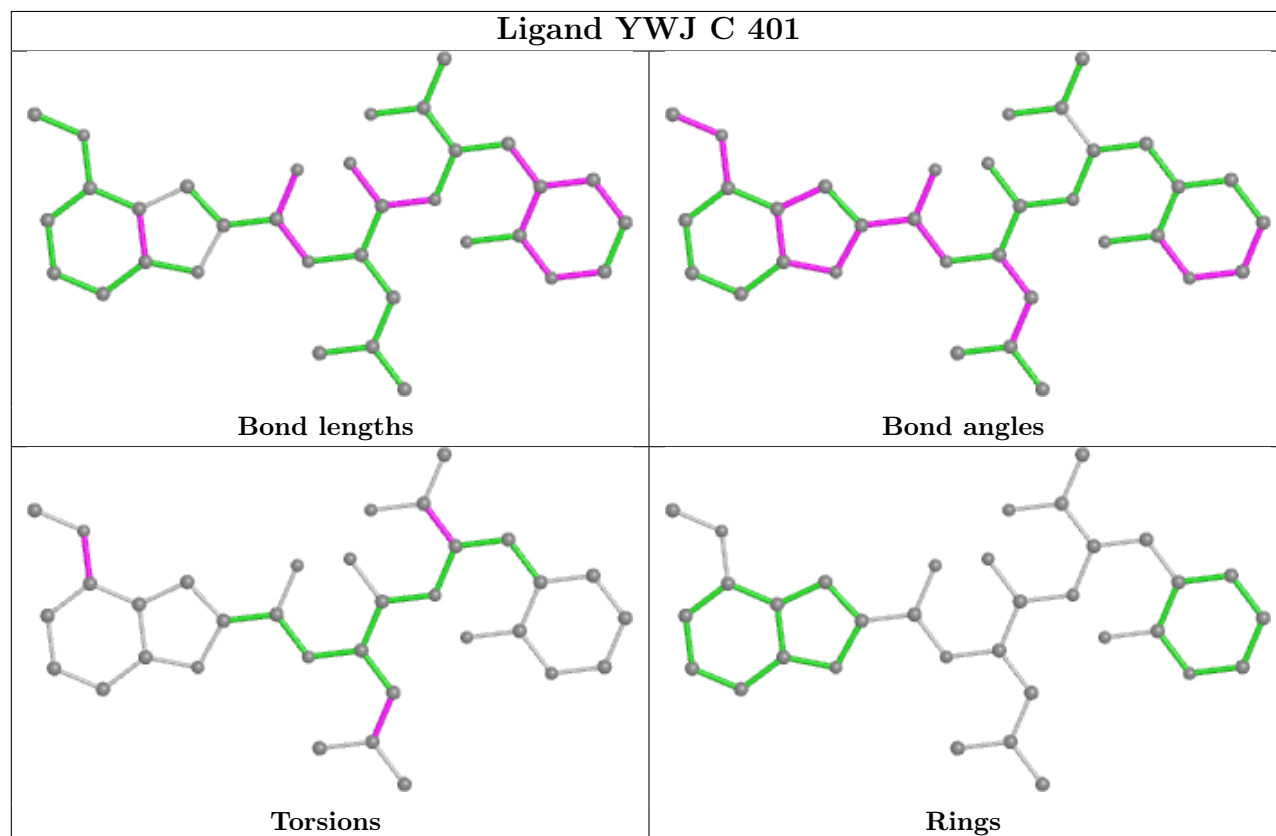
There are no ring outliers.

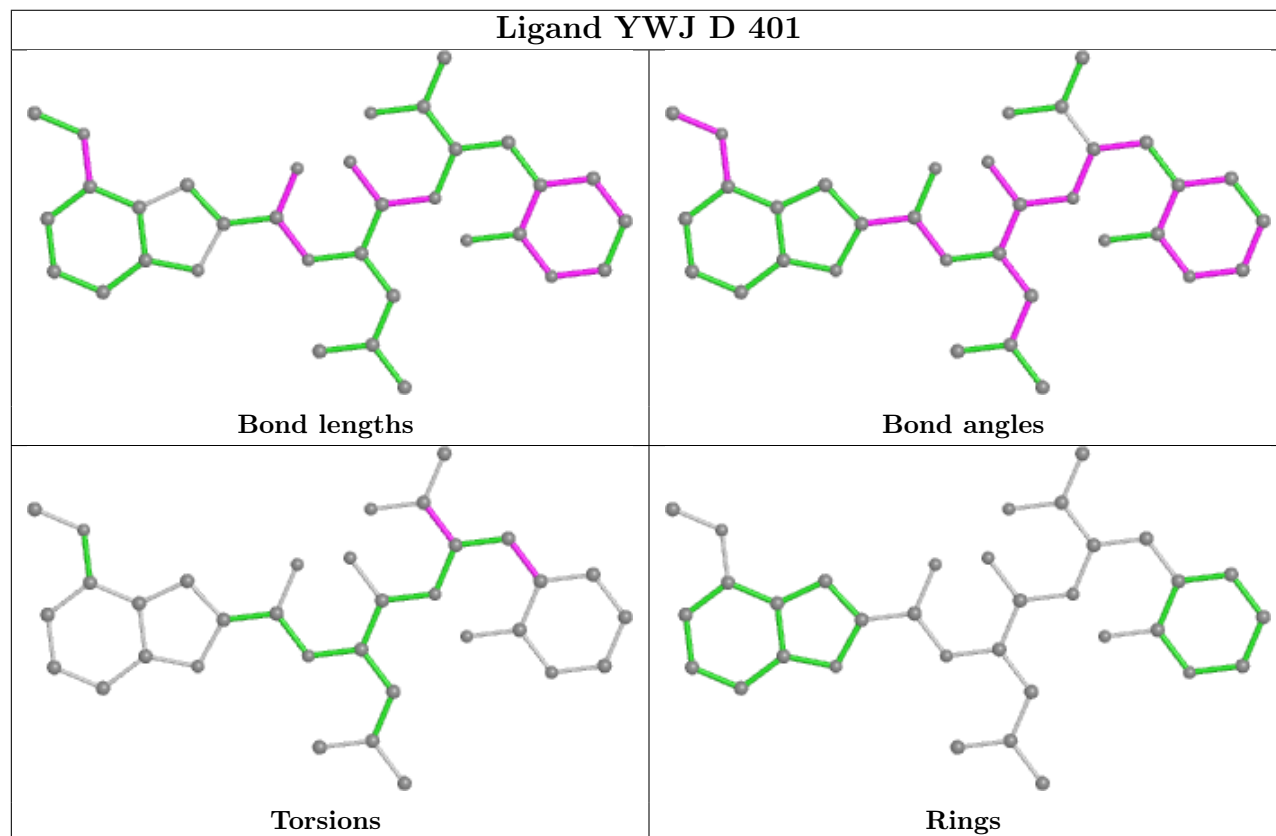
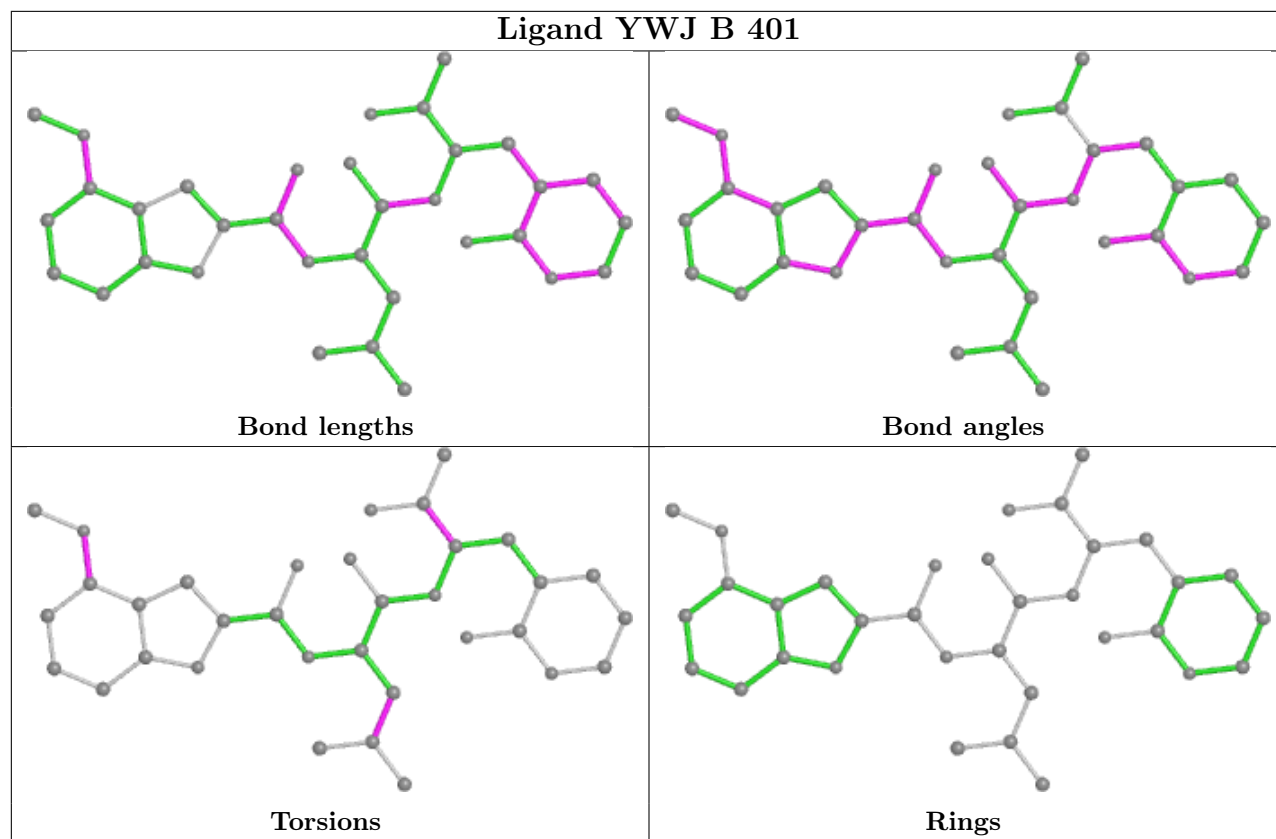
2 monomers are involved in 5 short contacts:

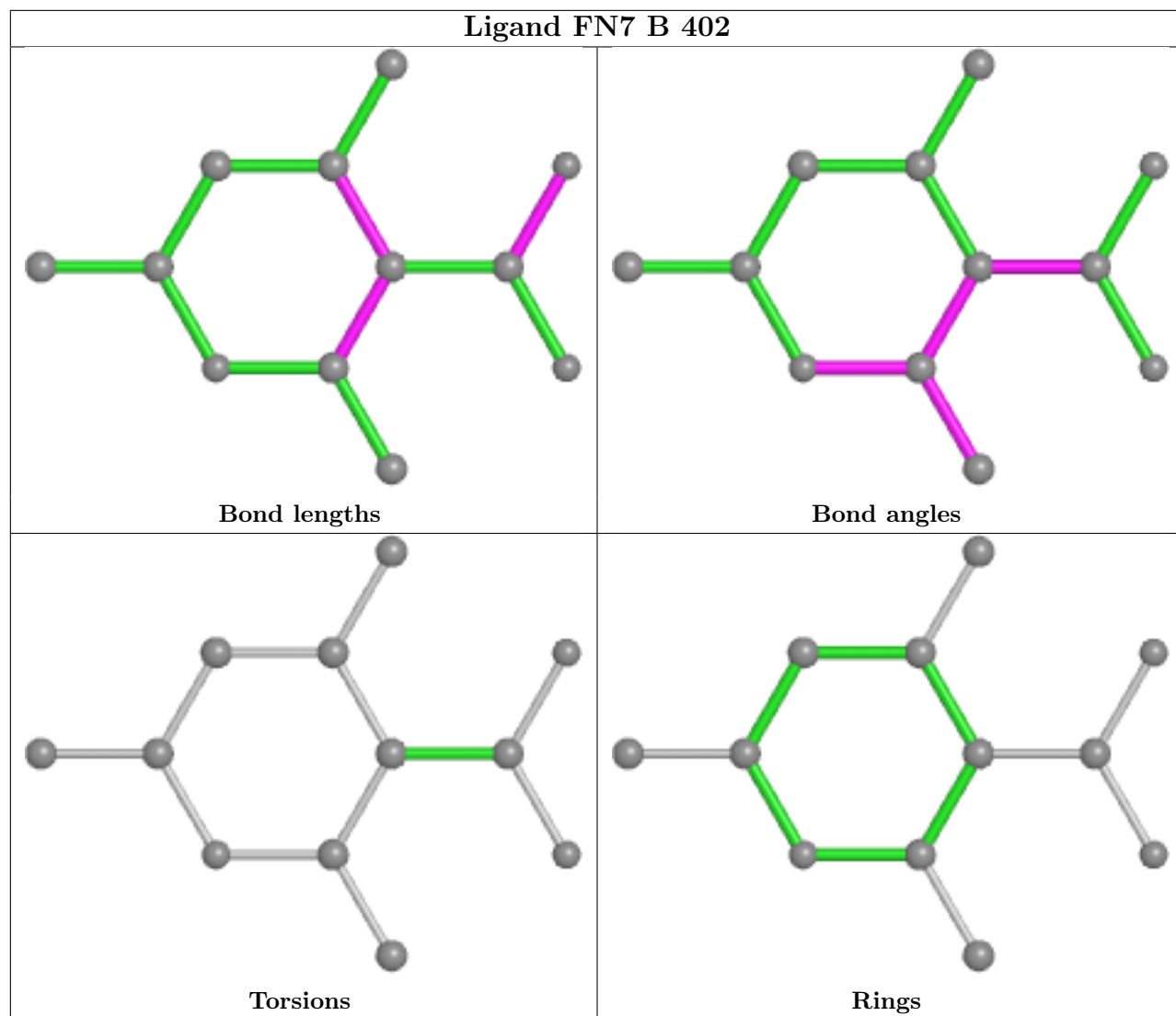
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	401	YWJ	1	0
3	B	402	FN7	1	3

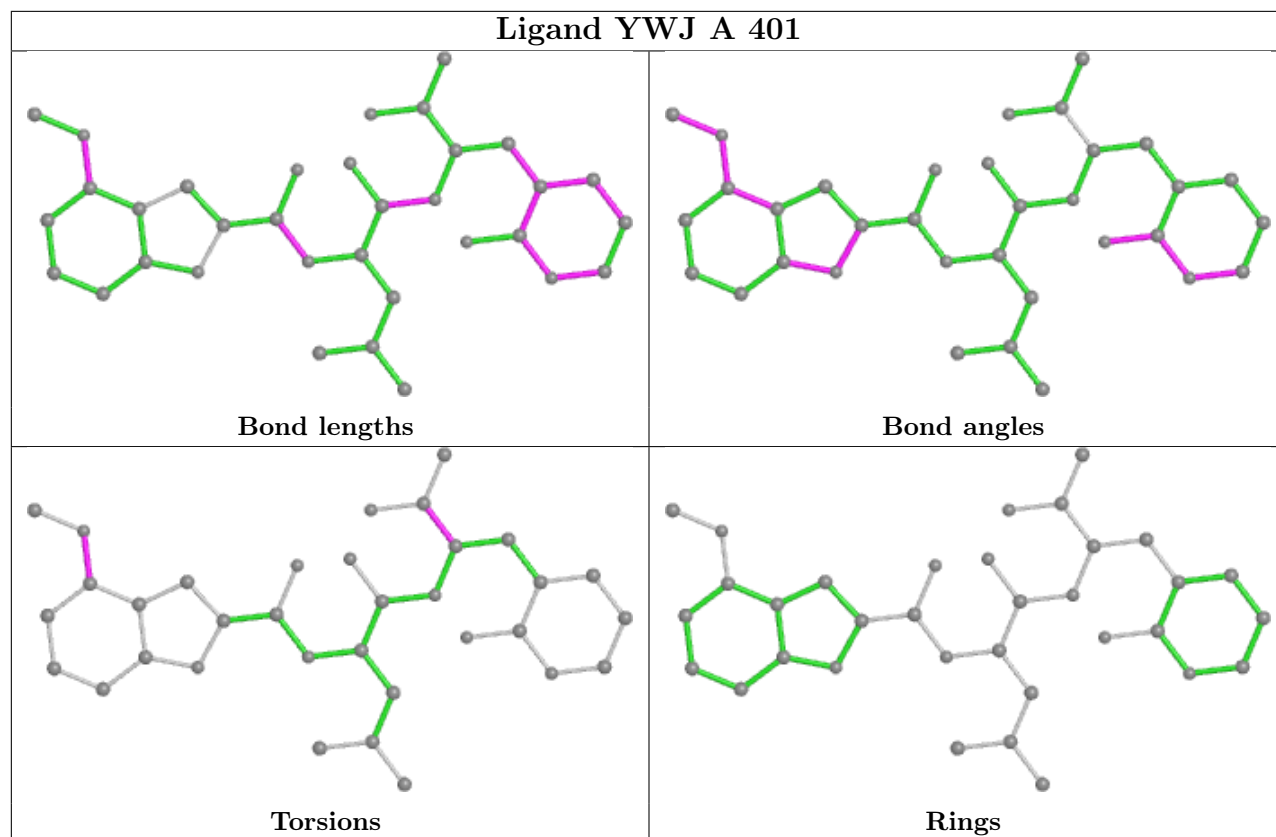
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 [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, 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	303/306 (99%)	0.13	5 (1%) 70 76	40, 50, 78, 142	0
1	B	306/306 (100%)	0.55	14 (4%) 32 42	41, 64, 93, 134	0
1	C	305/306 (99%)	0.50	17 (5%) 24 33	41, 58, 97, 159	0
1	D	306/306 (100%)	0.38	11 (3%) 42 51	43, 61, 91, 133	0
All	All	1220/1224 (99%)	0.39	47 (3%) 39 48	40, 58, 93, 159	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	222	ARG	5.4
1	D	262	LEU	4.8
1	D	49	MET	4.2
1	C	76	ARG	4.2
1	B	49	MET	4.1
1	C	77	VAL	3.8
1	A	48	ASP	3.5
1	B	205	LEU	3.5
1	C	7	ALA	3.4
1	D	227	LEU	3.2
1	C	67	LEU	3.2
1	C	223	PHE	3.2
1	B	154	TYR	3.2
1	B	250	LEU	3.1
1	D	50	LEU	3.1
1	C	66	PHE	3.1
1	C	222	ARG	3.1
1	D	279	ARG	3.0
1	A	303	VAL	3.0
1	C	91	VAL	2.9
1	A	302	GLY	2.9

*Continued on next page...*

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Mol	Chain	Res	Type	RSRZ
1	C	57	LEU	2.8
1	D	47	GLU	2.8
1	B	125	VAL	2.7
1	B	259	ILE	2.7
1	D	269	LYS	2.7
1	B	50	LEU	2.7
1	D	230	PHE	2.6
1	B	277	ASN	2.6
1	B	223	PHE	2.6
1	C	224	THR	2.5
1	C	125	VAL	2.5
1	B	7	ALA	2.4
1	C	305	PHE	2.4
1	D	87	LEU	2.4
1	C	128[A]	CYS	2.3
1	D	93	THR	2.3
1	B	257	THR	2.2
1	C	80	HIS	2.2
1	B	233	VAL	2.2
1	C	142	ASN	2.2
1	C	90	LYS	2.1
1	D	268	LEU	2.1
1	A	49	MET	2.1
1	C	124	GLY	2.1
1	A	56	ASP	2.1
1	B	247	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 monosaccharides in this entry.

## 6.4 Ligands [\(i\)](#)

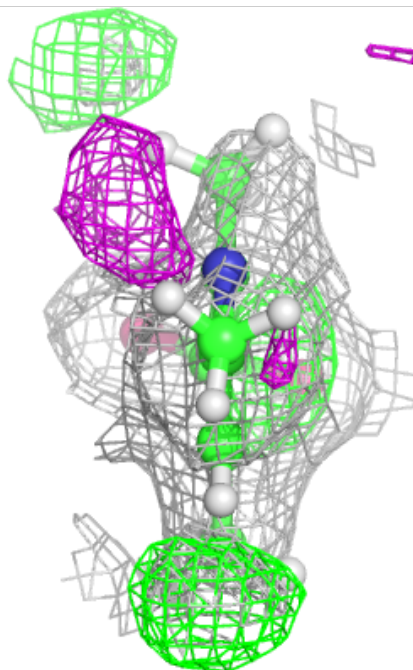
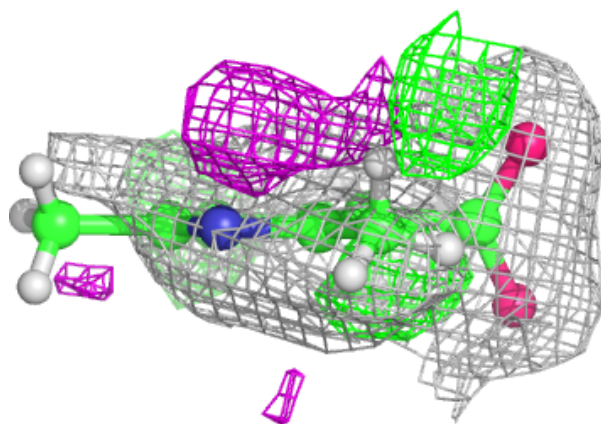
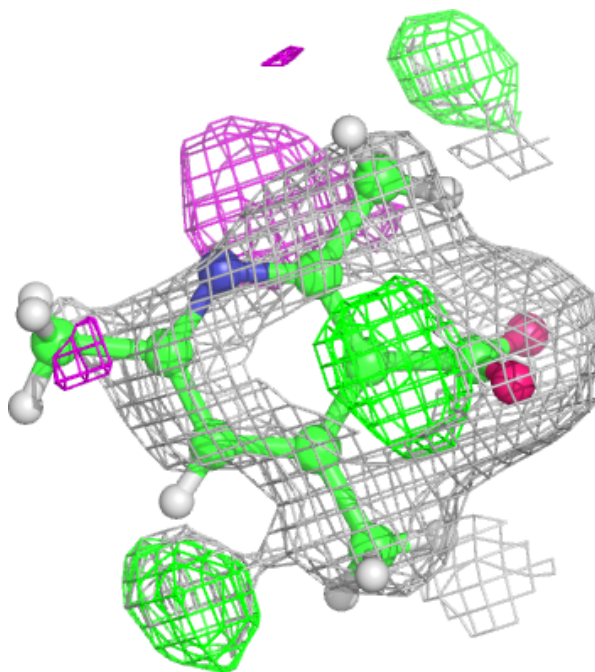
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	FN7	B	402	12/12	0.85	0.22	42,65,88,88	0
2	YWJ	B	401	34/34	0.91	0.15	43,58,70,77	0
2	YWJ	C	401	34/34	0.93	0.18	54,69,90,94	0
2	YWJ	D	401	34/34	0.94	0.13	50,64,81,82	0
2	YWJ	A	401	34/34	0.95	0.11	39,54,69,72	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.

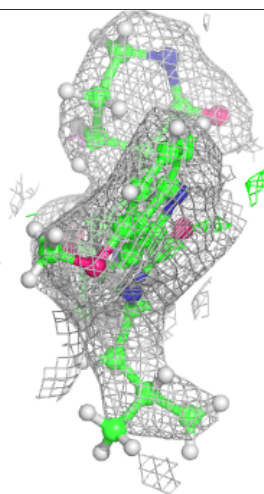
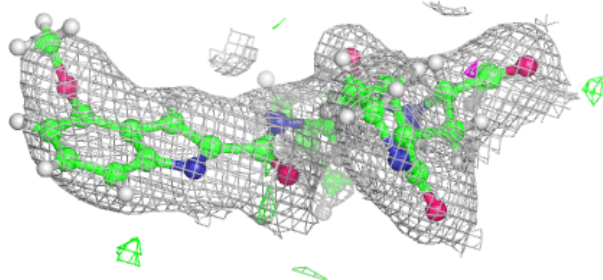
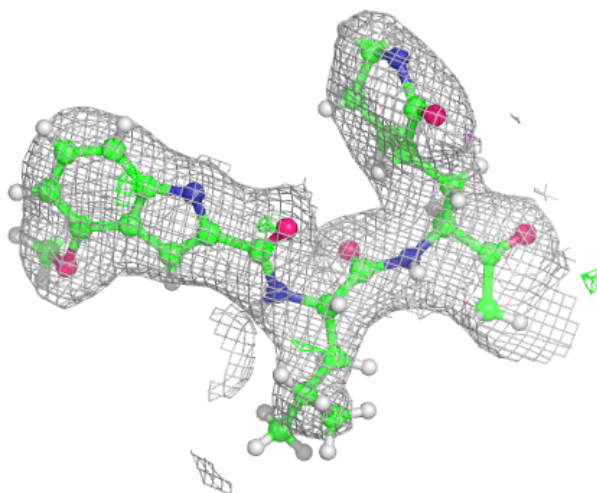
**Electron density around FN7 B 402:**

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



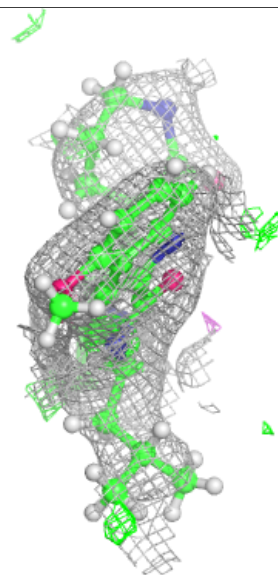
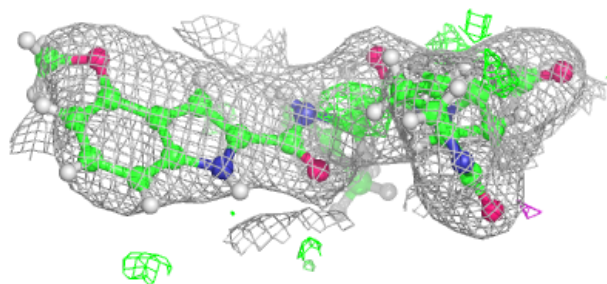
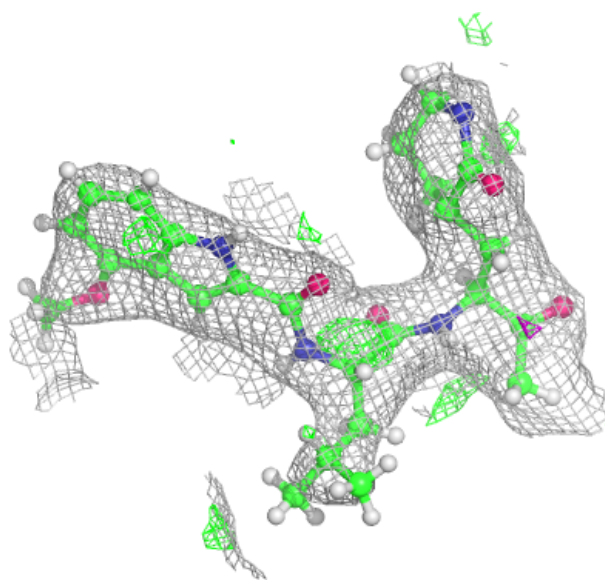
**Electron density around YWJ B 401:**

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



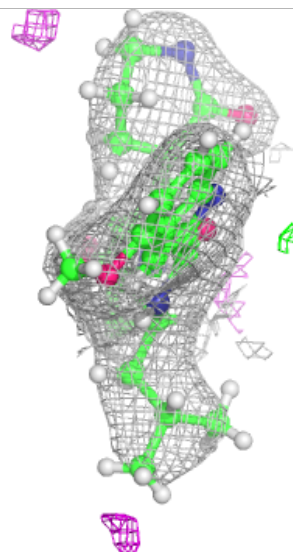
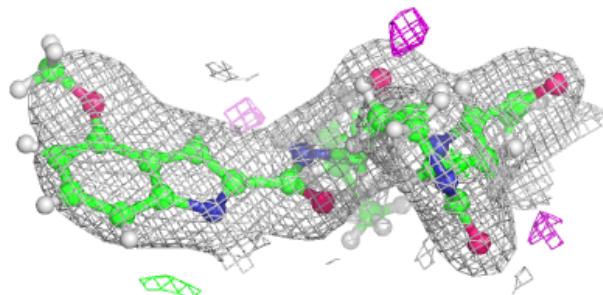
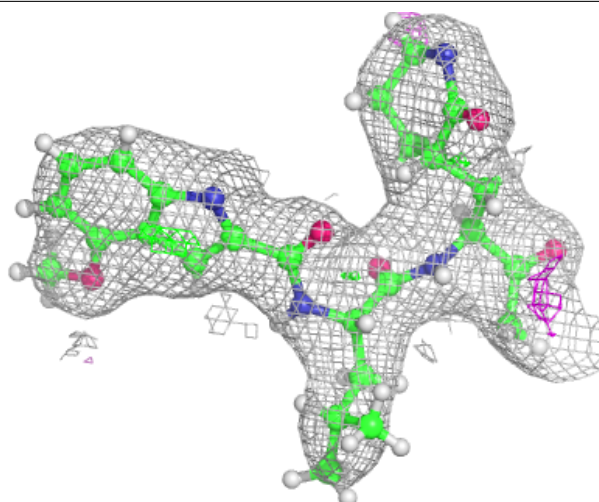
**Electron density around YWJ C 401:**

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

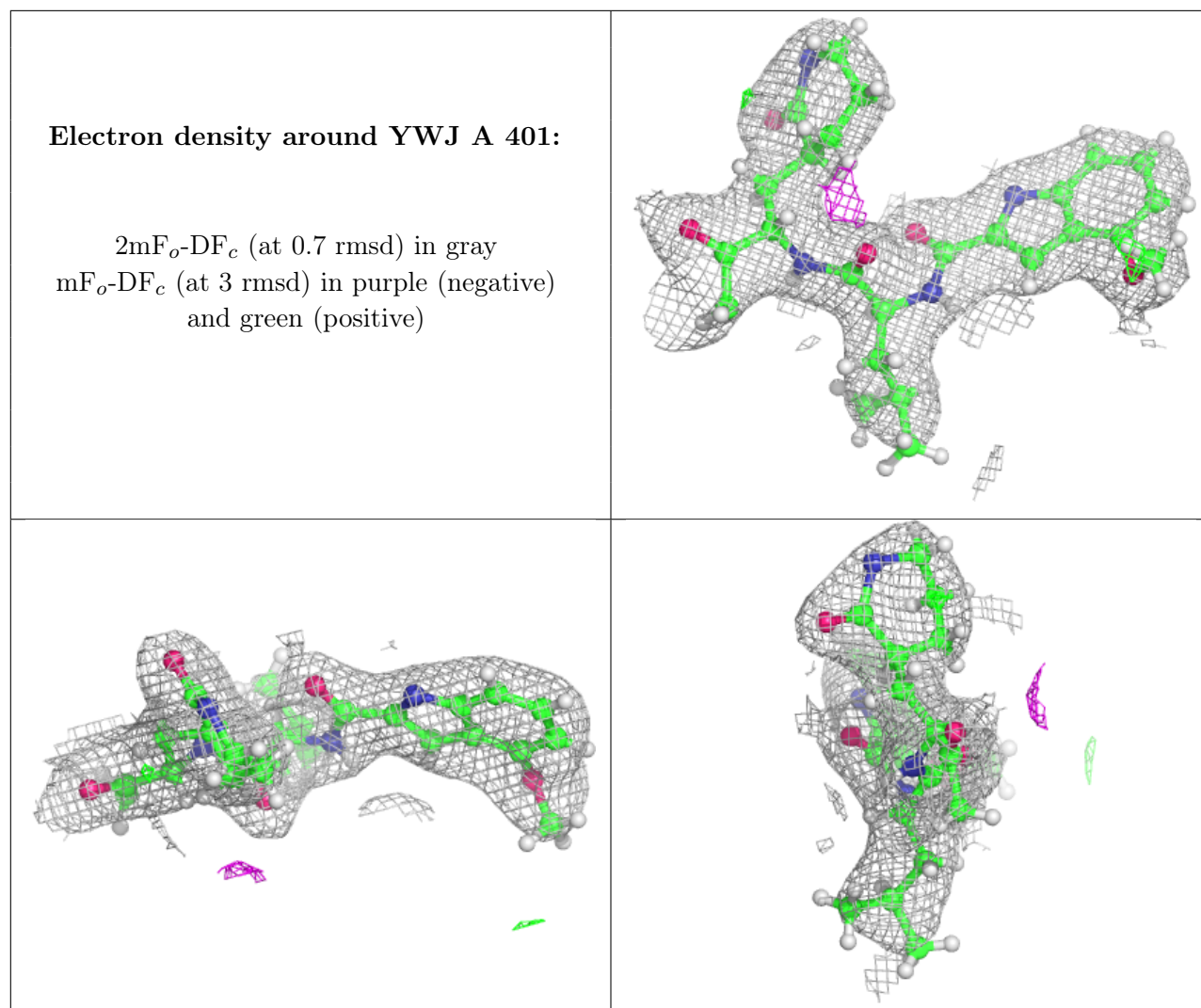


**Electron density around YWJ D 401:**

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







## 6.5 Other polymers [i](#)

There are no such residues in this entry.