



# wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 17, 2023 – 03:03 AM EDT

PDB ID : 4QWX  
Title : yCP in complex with the epoxyketone inhibitor ONX 0914  
Authors : Huber, E.M.; Heinemeyer, W.; Groll, M.  
Deposited on : 2014-07-17  
Resolution : 2.90 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

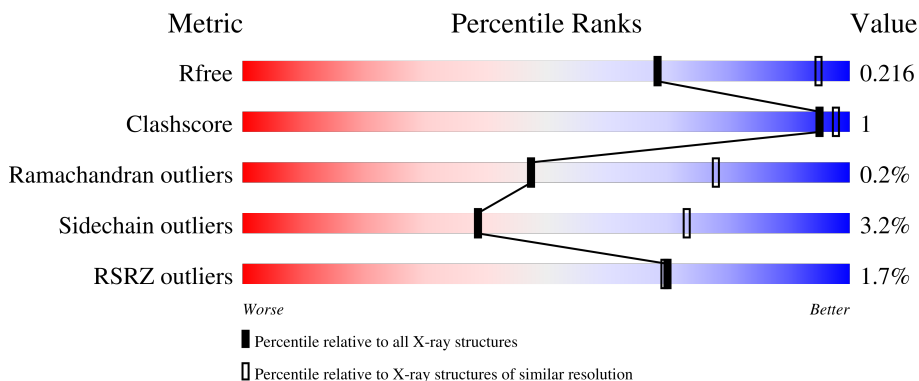
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	250	 98%
1	O	250	 98%
2	B	258	 90% 5%
2	P	258	 89% 5% 5%
3	C	254	 89% 5% 6%

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Mol	Chain	Length	Quality of chain
3	Q	254	 6% 89% 5% • 6%
4	D	260	 2% 86% • 10%
4	R	260	 2% 86% • 10%
5	E	234	 2% 94% 5% •
5	S	234	 3% 93% 6% •
6	F	288	 % 81% • 16%
6	T	288	 % 81% • 16%
7	G	252	 2% 90% 5% •
7	U	252	 % 90% 6% •
8	H	232	 % 91% • •
8	V	232	 % 92% • •
9	I	205	 94% 6%
9	W	205	 95% 5%
10	J	198	 % 93% • • •
10	X	198	 % 92% • • •
11	K	212	 93% 6% •
11	Y	212	 92% 8%
12	L	222	 96% •
12	Z	222	 95% 5%
13	M	246	 2% 91% • 5%
13	a	246	 % 92% • 5%
14	N	196	 % 95% 5%
14	b	196	 % 99% •

## 2 Entry composition

There are 20 unique types of molecules in this entry. The entry contains 49814 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 PROTEASOME SUBUNIT ALPHA TYPE-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	250	Total	C	N	O	S	0	0	0
			1915	1219	315	377	4			
1	O	250	Total	C	N	O	S	0	0	0
			1915	1219	315	377	4			

- Molecule 2 is a protein called PROTEASOME SUBUNIT ALPHA TYPE-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	244	Total	C	N	O	S	0	0	0
			1904	1201	321	379	3			
2	P	244	Total	C	N	O	S	0	0	0
			1904	1201	321	379	3			

- Molecule 3 is a protein called PROTEASOME SUBUNIT ALPHA TYPE-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	240	Total	C	N	O	S	0	0	0
			1881	1176	329	372	4			
3	Q	240	Total	C	N	O	S	0	0	0
			1881	1176	329	372	4			

- Molecule 4 is a protein called PROTEASOME SUBUNIT ALPHA TYPE-5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	D	235	Total	C	N	O	S	0	0	0
			1813	1136	304	366	7			
4	R	235	Total	C	N	O	S	0	0	0
			1813	1136	304	366	7			

- Molecule 5 is a protein called PROTEASOME SUBUNIT ALPHA TYPE-6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
5	E	231	Total 1773	C 1114	N 307	O 348	S 4	0	0	0
5	S	231	Total 1773	C 1114	N 307	O 348	S 4	0	0	0

- Molecule 6 is a protein called PROBABLE PROTEASOME SUBUNIT ALPHA TYPE-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
6	F	243	Total 1892	C 1203	N 329	O 356	S 4	0	0	0
6	T	243	Total 1892	C 1203	N 329	O 356	S 4	0	0	0

- Molecule 7 is a protein called PROTEASOME SUBUNIT ALPHA TYPE-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
7	G	241	Total 1907	C 1214	N 320	O 365	S 8	0	0	0
7	U	241	Total 1907	C 1214	N 320	O 365	S 8	0	0	0

- Molecule 8 is a protein called PROTEASOME SUBUNIT BETA TYPE-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
8	H	222	Total 1684	C 1061	N 293	O 323	S 7	0	0	0
8	V	222	Total 1684	C 1061	N 293	O 323	S 7	0	0	0

- Molecule 9 is a protein called PROTEASOME SUBUNIT BETA TYPE-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
9	I	204	Total 1581	C 1010	N 258	O 305	S 8	0	0	0
9	W	204	Total 1581	C 1010	N 258	O 305	S 8	0	0	0

- Molecule 10 is a protein called PROTEASOME SUBUNIT BETA TYPE-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
10	J	195	Total 1561	C 992	N 264	O 299	S 6	0	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	X	195	Total	C	N	O	S	0	0	0
			1561	992	264	299	6			

- Molecule 11 is a protein called PROTEASOME SUBUNIT BETA TYPE-5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	212	Total	C	N	O	S	0	0	0
			1644	1045	280	312	7			
11	Y	212	Total	C	N	O	S	0	0	0
			1644	1045	280	312	7			

- Molecule 12 is a protein called PROTEASOME SUBUNIT BETA TYPE-6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	222	Total	C	N	O	S	0	0	0
			1757	1115	303	335	4			
12	Z	222	Total	C	N	O	S	0	0	0
			1757	1115	303	335	4			

- Molecule 13 is a protein called PROTEASOME SUBUNIT BETA TYPE-7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	M	233	Total	C	N	O	S	0	0	0
			1824	1154	312	351	7			
13	a	233	Total	C	N	O	S	0	0	0
			1824	1154	312	351	7			

- Molecule 14 is a protein called PROTEASOME SUBUNIT BETA TYPE-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	N	196	Total	C	N	O	S	0	0	0
			1512	955	250	300	7			
14	b	196	Total	C	N	O	S	0	0	0
			1512	955	250	300	7			

- Molecule 15 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	G	1	Total	Mg	0	0
			1	1		
15	I	1	Total	Mg	0	0
			1	1		

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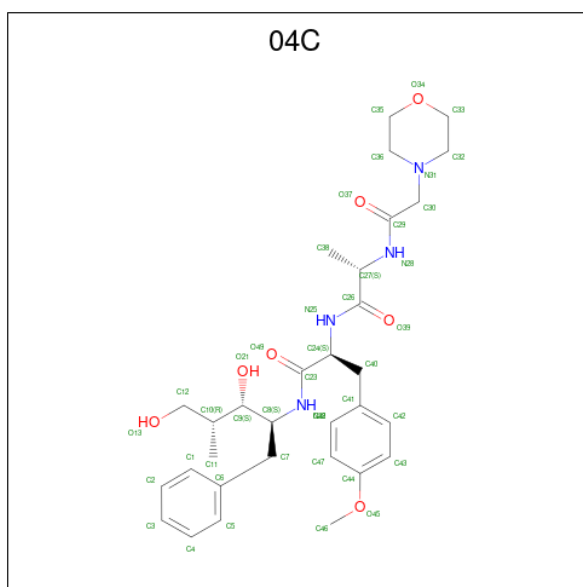
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
15	K	1	Total Mg 1 1	0	0
15	N	1	Total Mg 1 1	0	0
15	V	1	Total Mg 1 1	0	0
15	X	1	Total Mg 1 1	0	0
15	Y	1	Total Mg 1 1	0	0
15	Z	1	Total Mg 1 1	0	0

- Molecule 16 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
16	G	1	Total Cl 1 1	0	0
16	U	1	Total Cl 1 1	0	0

- Molecule 17 is 1,2,4-trideoxy-4-methyl-2-{[N-(morpholin-4-ylacetyl)-L-alanyl-O-methyl-L-tyrosyl]amino}-1-phenyl-D-xylitol (three-letter code: 04C) (formula: C<sub>31</sub>H<sub>44</sub>N<sub>4</sub>O<sub>7</sub>).



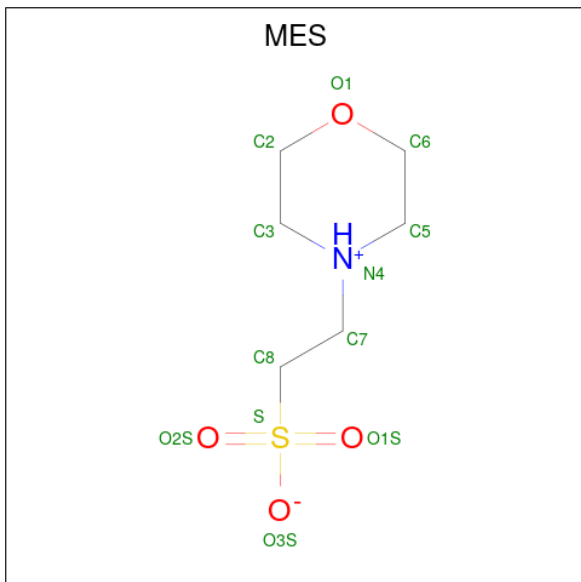
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
17	H	1	Total C N O 42 31 4 7	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
17	K	1	Total 42	C 31	N 4	O 7	0	0
17	N	1	Total 42	C 31	N 4	O 7	0	0
17	V	1	Total 42	C 31	N 4	O 7	0	0
17	Y	1	Total 42	C 31	N 4	O 7	0	0
17	b	1	Total 42	C 31	N 4	O 7	0	0

- Molecule 18 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C<sub>6</sub>H<sub>13</sub>NO<sub>4</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
18	H	1	Total 12	C 6	N 1	O 4	S 1	0	0
18	K	1	Total 12	C 6	N 1	O 4	S 1	0	0
18	V	1	Total 12	C 6	N 1	O 4	S 1	0	0
18	Y	1	Total 12	C 6	N 1	O 4	S 1	0	0

- Molecule 19 is SODIUM ION (three-letter code: NA) (formula: Na).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
19	K	1	Total Na 1 1	0	0
19	Y	1	Total Na 1 1	0	0

- Molecule 20 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
20	A	4	Total O 4 4	0	0
20	B	14	Total O 14 14	0	0
20	C	5	Total O 5 5	0	0
20	D	5	Total O 5 5	0	0
20	E	4	Total O 4 4	0	0
20	F	5	Total O 5 5	0	0
20	G	5	Total O 5 5	0	0
20	H	17	Total O 17 17	0	0
20	I	3	Total O 3 3	0	0
20	J	10	Total O 10 10	0	0
20	K	12	Total O 12 12	0	0
20	L	9	Total O 9 9	0	0
20	M	8	Total O 8 8	0	0
20	N	4	Total O 4 4	0	0
20	O	2	Total O 2 2	0	0
20	P	8	Total O 8 8	0	0
20	Q	7	Total O 7 7	0	0
20	R	5	Total O 5 5	0	0

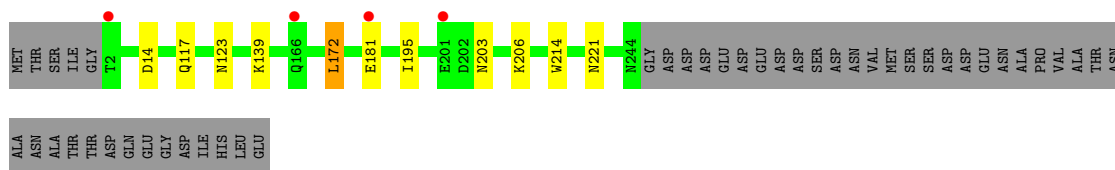
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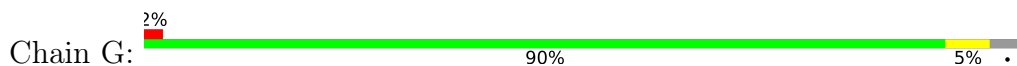
<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>	<b>ZeroOcc</b>	<b>AltConf</b>
20	S	4	Total O 4 4	0	0
20	T	4	Total O 4 4	0	0
20	U	6	Total O 6 6	0	0
20	V	12	Total O 12 12	0	0
20	W	9	Total O 9 9	0	0
20	X	9	Total O 9 9	0	0
20	Y	8	Total O 8 8	0	0
20	Z	11	Total O 11 11	0	0
20	a	11	Total O 11 11	0	0
20	b	5	Total O 5 5	0	0



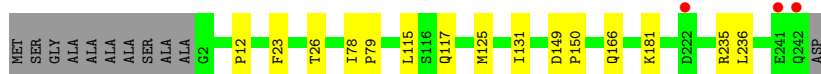
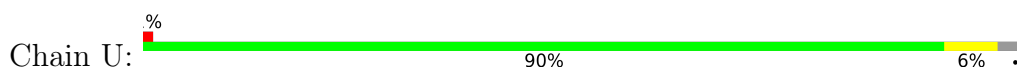




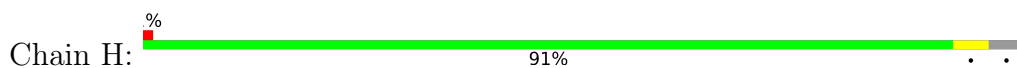
- Molecule 7: PROTEASOME SUBUNIT ALPHA TYPE-1

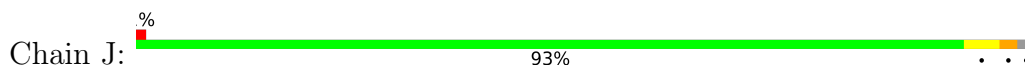


- Molecule 7: PROTEASOME SUBUNIT ALPHA TYPE-1

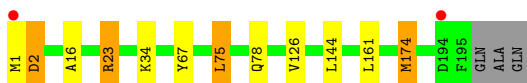
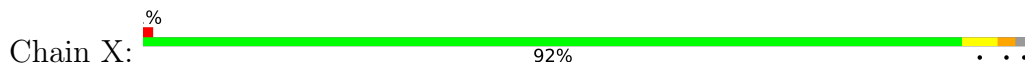


- Molecule 8: PROTEASOME SUBUNIT BETA TYPE-2





- Molecule 10: PROTEASOME SUBUNIT BETA TYPE-4



- Molecule 11: PROTEASOME SUBUNIT BETA TYPE-5



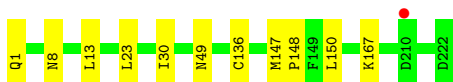
- Molecule 11: PROTEASOME SUBUNIT BETA TYPE-5



- Molecule 12: PROTEASOME SUBUNIT BETA TYPE-6



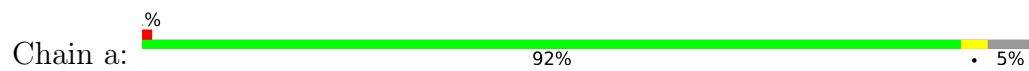
- Molecule 12: PROTEASOME SUBUNIT BETA TYPE-6



- Molecule 13: PROTEASOME SUBUNIT BETA TYPE-7



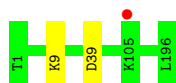
- Molecule 13: PROTEASOME SUBUNIT BETA TYPE-7



- Molecule 14: PROTEASOME SUBUNIT BETA TYPE-1



- Molecule 14: PROTEASOME SUBUNIT BETA TYPE-1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	135.60Å 300.05Å 145.62Å 90.00° 112.77° 90.00°	Depositor
Resolution (Å)	15.00 – 2.90 15.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	95.9 (15.00-2.90) 96.0 (15.00-2.90)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.41 (at 2.91Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.172 , 0.212 0.177 , 0.216	Depositor DCC
$R_{free}$ test set	11257 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	66.5	Xtrriage
Anisotropy	0.080	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 46.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	49814	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	75.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.35% 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

Bond lengths and bond angles in the following residue types are not validated in this section: CL, NA, MES, MG, 04C

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	0.27	0/1952	0.47	0/2642
1	O	0.27	0/1952	0.47	0/2642
2	B	0.27	0/1934	0.49	0/2618
2	P	0.27	0/1934	0.49	0/2618
3	C	0.28	0/1910	0.50	0/2586
3	Q	0.27	0/1910	0.50	0/2586
4	D	0.27	0/1837	0.47	0/2475
4	R	0.27	0/1837	0.47	0/2475
5	E	0.27	0/1800	0.47	0/2433
5	S	0.27	0/1800	0.47	0/2433
6	F	0.28	0/1932	0.45	0/2609
6	T	0.27	0/1932	0.45	0/2609
7	G	0.28	0/1945	0.47	0/2634
7	U	0.27	0/1945	0.47	0/2634
8	H	0.26	0/1715	0.48	1/2326 (0.0%)
8	V	0.25	0/1715	0.48	1/2326 (0.0%)
9	I	0.27	0/1611	0.48	0/2174
9	W	0.27	0/1611	0.48	0/2174
10	J	0.26	0/1589	0.49	0/2142
10	X	0.27	0/1589	0.49	0/2142
11	K	0.28	0/1681	0.51	0/2274
11	Y	0.39	0/1681	0.54	1/2274 (0.0%)
12	L	0.27	0/1795	0.47	0/2420
12	Z	0.27	0/1795	0.47	0/2420
13	M	0.27	0/1855	0.51	0/2514
13	a	0.27	0/1855	0.51	0/2514
14	N	0.26	0/1541	0.48	0/2087
14	b	0.26	0/1541	0.48	0/2087
All	All	0.27	0/50194	0.48	3/67868 (0.0%)

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
11	K	0	1
11	Y	0	1
All	All	0	2

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	Y	73	ARG	NE-CZ-NH1	5.89	123.25	120.30
8	H	1	THR	N-CA-C	5.10	124.77	111.00
8	V	1	THR	N-CA-C	5.04	124.62	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
11	K	48	GLY	Peptide
11	Y	48	GLY	Peptide

## 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	1915	0	1929	1	0
1	O	1915	0	1929	2	0
2	B	1904	0	1904	4	0
2	P	1904	0	1904	6	0
3	C	1881	0	1895	5	0
3	Q	1881	0	1895	7	0
4	D	1813	0	1797	3	0
4	R	1813	0	1797	3	0
5	E	1773	0	1775	3	0
5	S	1773	0	1775	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	F	1892	0	1883	1	0
6	T	1892	0	1883	1	0
7	G	1907	0	1901	3	0
7	U	1907	0	1901	5	0
8	H	1684	0	1685	2	0
8	V	1684	0	1685	1	0
9	I	1581	0	1574	6	0
9	W	1581	0	1574	4	0
10	J	1561	0	1569	7	0
10	X	1561	0	1569	7	0
11	K	1644	0	1592	9	0
11	Y	1644	0	1592	12	0
12	L	1757	0	1711	2	0
12	Z	1757	0	1711	3	0
13	M	1824	0	1832	2	0
13	a	1824	0	1832	0	0
14	N	1512	0	1478	5	0
14	b	1512	0	1478	0	0
15	G	1	0	0	0	0
15	I	1	0	0	0	0
15	K	1	0	0	0	0
15	N	1	0	0	0	0
15	V	1	0	0	0	0
15	X	1	0	0	0	0
15	Y	1	0	0	0	0
15	Z	1	0	0	0	0
16	G	1	0	0	0	0
16	U	1	0	0	0	0
17	H	42	0	42	3	0
17	K	42	0	42	0	0
17	N	42	0	42	5	0
17	V	42	0	42	3	0
17	Y	42	0	42	0	0
17	b	42	0	42	0	0
18	H	12	0	13	0	0
18	K	12	0	13	1	0
18	V	12	0	13	0	0
18	Y	12	0	13	1	0
19	K	1	0	0	0	0
19	Y	1	0	0	0	0
20	A	4	0	0	0	0
20	B	14	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
20	C	5	0	0	0	0
20	D	5	0	0	0	0
20	E	4	0	0	0	0
20	F	5	0	0	0	0
20	G	5	0	0	0	0
20	H	17	0	0	0	0
20	I	3	0	0	0	0
20	J	10	0	0	0	0
20	K	12	0	0	0	0
20	L	9	0	0	0	0
20	M	8	0	0	0	0
20	N	4	0	0	1	0
20	O	2	0	0	0	0
20	P	8	0	0	0	0
20	Q	7	0	0	0	0
20	R	5	0	0	0	0
20	S	4	0	0	0	0
20	T	4	0	0	0	0
20	U	6	0	0	0	0
20	V	12	0	0	0	0
20	W	9	0	0	0	0
20	X	9	0	0	0	0
20	Y	8	0	0	0	0
20	Z	11	0	0	0	0
20	a	11	0	0	0	0
20	b	5	0	0	0	0
All	All	49814	0	49354	105	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 1.

The worst 5 of 105 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)
11:K:73:ARG:HH11	11:K:73:ARG:HG3	1.21	1.02
10:J:1:MET:O	10:J:2:ASP:HB2	1.85	0.74
10:X:1:MET:O	10:X:2:ASP:HB2	1.87	0.74
11:K:73:ARG:HG3	11:K:73:ARG:NH1	1.96	0.71
12:Z:13:LEU:HD13	12:Z:150:LEU:HD21	1.80	0.64

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	248/250 (99%)	241 (97%)	7 (3%)	0	100	100
1	O	248/250 (99%)	241 (97%)	7 (3%)	0	100	100
2	B	242/258 (94%)	233 (96%)	7 (3%)	2 (1%)	19	51
2	P	242/258 (94%)	233 (96%)	7 (3%)	2 (1%)	19	51
3	C	238/254 (94%)	230 (97%)	6 (2%)	2 (1%)	19	51
3	Q	238/254 (94%)	230 (97%)	6 (2%)	2 (1%)	19	51
4	D	231/260 (89%)	226 (98%)	5 (2%)	0	100	100
4	R	231/260 (89%)	226 (98%)	5 (2%)	0	100	100
5	E	229/234 (98%)	223 (97%)	6 (3%)	0	100	100
5	S	229/234 (98%)	223 (97%)	6 (3%)	0	100	100
6	F	241/288 (84%)	234 (97%)	7 (3%)	0	100	100
6	T	241/288 (84%)	234 (97%)	7 (3%)	0	100	100
7	G	239/252 (95%)	235 (98%)	4 (2%)	0	100	100
7	U	239/252 (95%)	234 (98%)	5 (2%)	0	100	100
8	H	220/232 (95%)	216 (98%)	4 (2%)	0	100	100
8	V	220/232 (95%)	216 (98%)	4 (2%)	0	100	100
9	I	202/205 (98%)	195 (96%)	7 (4%)	0	100	100
9	W	202/205 (98%)	195 (96%)	7 (4%)	0	100	100
10	J	193/198 (98%)	189 (98%)	3 (2%)	1 (0%)	29	61
10	X	193/198 (98%)	189 (98%)	3 (2%)	1 (0%)	29	61
11	K	210/212 (99%)	207 (99%)	3 (1%)	0	100	100
11	Y	210/212 (99%)	207 (99%)	3 (1%)	0	100	100
12	L	220/222 (99%)	216 (98%)	4 (2%)	0	100	100
12	Z	220/222 (99%)	216 (98%)	4 (2%)	0	100	100
13	M	231/246 (94%)	222 (96%)	9 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
13	a	231/246 (94%)	222 (96%)	9 (4%)	0	100	100
14	N	194/196 (99%)	188 (97%)	6 (3%)	0	100	100
14	b	194/196 (99%)	188 (97%)	6 (3%)	0	100	100
All	All	6276/6614 (95%)	6109 (97%)	157 (2%)	10 (0%)	47	78

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	51	VAL
3	C	202	GLN
10	J	2	ASP
2	P	51	VAL
3	Q	202	GLN

### 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	209/209 (100%)	206 (99%)	3 (1%)	67	89
1	O	209/209 (100%)	206 (99%)	3 (1%)	67	89
2	B	203/216 (94%)	196 (97%)	7 (3%)	37	71
2	P	203/216 (94%)	196 (97%)	7 (3%)	37	71
3	C	212/226 (94%)	204 (96%)	8 (4%)	33	67
3	Q	212/226 (94%)	204 (96%)	8 (4%)	33	67
4	D	194/215 (90%)	186 (96%)	8 (4%)	30	64
4	R	194/215 (90%)	186 (96%)	8 (4%)	30	64
5	E	190/193 (98%)	182 (96%)	8 (4%)	30	63
5	S	190/193 (98%)	182 (96%)	8 (4%)	30	63
6	F	201/239 (84%)	191 (95%)	10 (5%)	24	57
6	T	201/239 (84%)	191 (95%)	10 (5%)	24	57
7	G	206/210 (98%)	199 (97%)	7 (3%)	37	71

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
7	U	206/210 (98%)	199 (97%)	7 (3%)	37	71
8	H	181/190 (95%)	175 (97%)	6 (3%)	38	72
8	V	181/190 (95%)	176 (97%)	5 (3%)	43	76
9	I	172/173 (99%)	169 (98%)	3 (2%)	60	86
9	W	172/173 (99%)	169 (98%)	3 (2%)	60	86
10	J	173/175 (99%)	168 (97%)	5 (3%)	42	76
10	X	173/175 (99%)	168 (97%)	5 (3%)	42	76
11	K	169/169 (100%)	162 (96%)	7 (4%)	30	64
11	Y	169/169 (100%)	162 (96%)	7 (4%)	30	64
12	L	185/185 (100%)	180 (97%)	5 (3%)	44	77
12	Z	185/185 (100%)	180 (97%)	5 (3%)	44	77
13	M	199/208 (96%)	192 (96%)	7 (4%)	36	70
13	a	199/208 (96%)	192 (96%)	7 (4%)	36	70
14	N	162/162 (100%)	160 (99%)	2 (1%)	71	91
14	b	162/162 (100%)	160 (99%)	2 (1%)	71	91
All	All	5312/5540 (96%)	5141 (97%)	171 (3%)	39	73

5 of 171 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
5	S	54	GLU
9	W	171	LEU
5	S	202	ASP
7	U	115	LEU
11	Y	4	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 76 such sidechains are listed below:

Mol	Chain	Res	Type
6	T	117	GLN
13	a	18	ASN
6	T	191	GLN
10	X	55	GLN
13	a	213	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 22 ligands modelled in this entry, 12 are monoatomic - leaving 10 for Mogul analysis.

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
17	04C	N	201	14	44,44,44	1.44	3 (6%)	56,58,58	1.34	8 (14%)
17	04C	H	301	8	44,44,44	1.20	4 (9%)	56,58,58	1.23	7 (12%)
18	MES	V	303	-	12,12,12	2.24	1 (8%)	14,16,16	1.37	3 (21%)
17	04C	V	301	8	44,44,44	1.19	4 (9%)	56,58,58	1.22	6 (10%)
17	04C	b	201	14	44,44,44	1.39	3 (6%)	56,58,58	1.29	8 (14%)
18	MES	K	304	-	12,12,12	2.39	1 (8%)	14,16,16	1.40	2 (14%)
17	04C	Y	301	11,19	44,44,44	1.30	4 (9%)	56,58,58	1.18	8 (14%)
17	04C	K	301	11	44,44,44	1.28	3 (6%)	56,58,58	1.17	9 (16%)
18	MES	H	302	-	12,12,12	2.21	1 (8%)	14,16,16	1.50	2 (14%)
18	MES	Y	304	-	12,12,12	2.33	1 (8%)	14,16,16	1.40	2 (14%)

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
17	04C	N	201	14	-	16/44/52/52	0/3/3/3
17	04C	H	301	8	-	11/44/52/52	0/3/3/3
18	MES	V	303	-	-	0/6/14/14	0/1/1/1
17	04C	V	301	8	-	11/44/52/52	0/3/3/3
17	04C	b	201	14	-	16/44/52/52	0/3/3/3
18	MES	K	304	-	-	0/6/14/14	0/1/1/1
17	04C	Y	301	11,19	-	13/44/52/52	0/3/3/3
17	04C	K	301	11	-	12/44/52/52	0/3/3/3
18	MES	H	302	-	-	3/6/14/14	0/1/1/1
18	MES	Y	304	-	-	0/6/14/14	0/1/1/1

The worst 5 of 25 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
18	K	304	MES	C8-S	-7.99	1.66	1.77
18	Y	304	MES	C8-S	-7.76	1.66	1.77
18	V	303	MES	C8-S	-7.47	1.66	1.77
18	H	302	MES	C8-S	-7.40	1.67	1.77
17	N	201	04C	C10-C9	5.23	1.62	1.53

The worst 5 of 55 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	N	201	04C	C11-C10-C12	-5.53	102.61	109.88
17	b	201	04C	C11-C10-C12	-4.94	103.39	109.88
18	H	302	MES	O2S-S-C8	3.90	111.62	106.92
17	V	301	04C	O34-C33-C32	-3.81	103.39	111.80
17	H	301	04C	O34-C33-C32	-3.81	103.41	111.80

There are no chirality outliers.

5 of 82 torsion outliers are listed below:

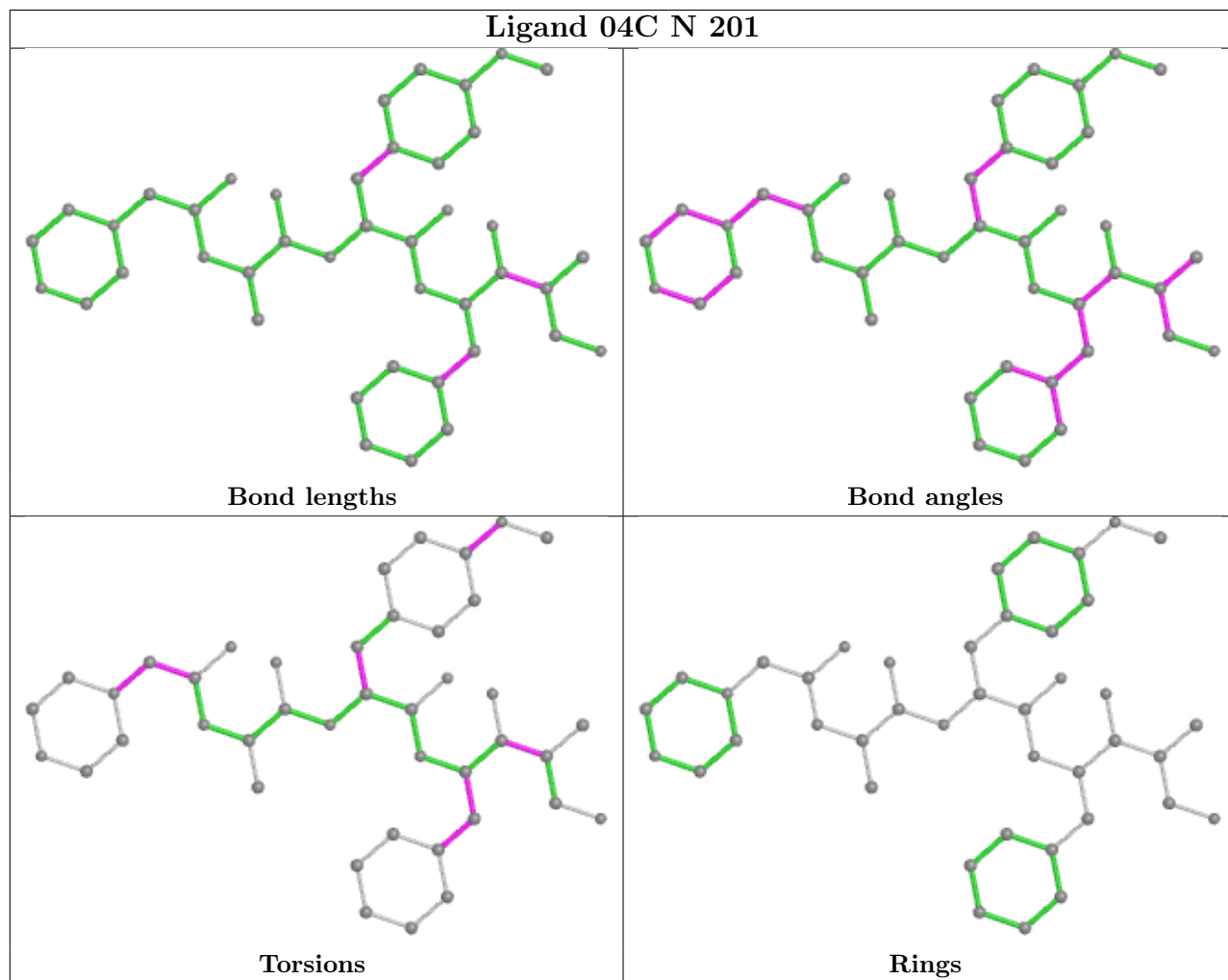
Mol	Chain	Res	Type	Atoms
17	H	301	04C	C11-C10-C9-C8
17	K	301	04C	C11-C10-C9-C8
17	K	301	04C	C12-C10-C9-C8
17	N	201	04C	C11-C10-C9-C8
17	N	201	04C	C12-C10-C9-C8

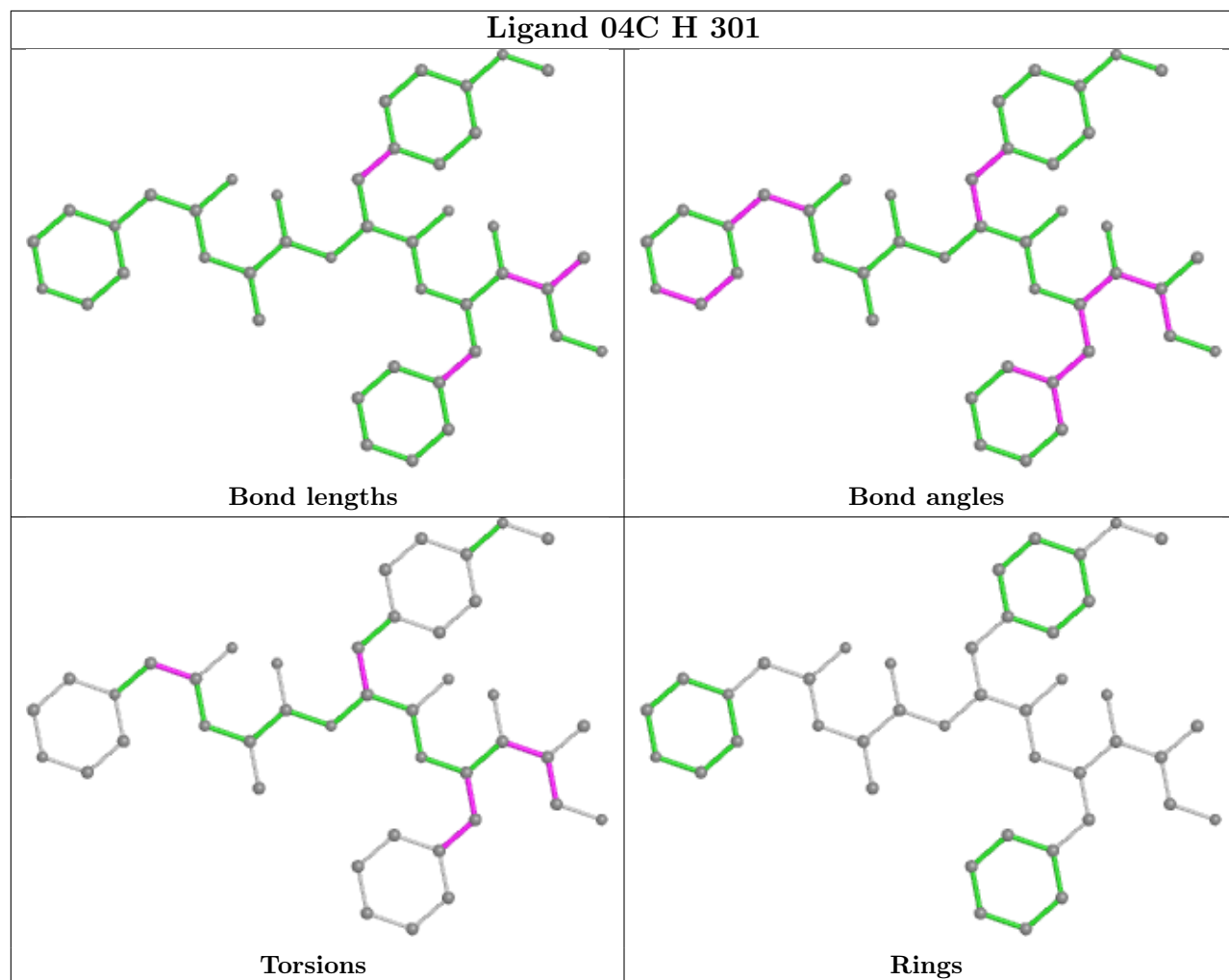
There are no ring outliers.

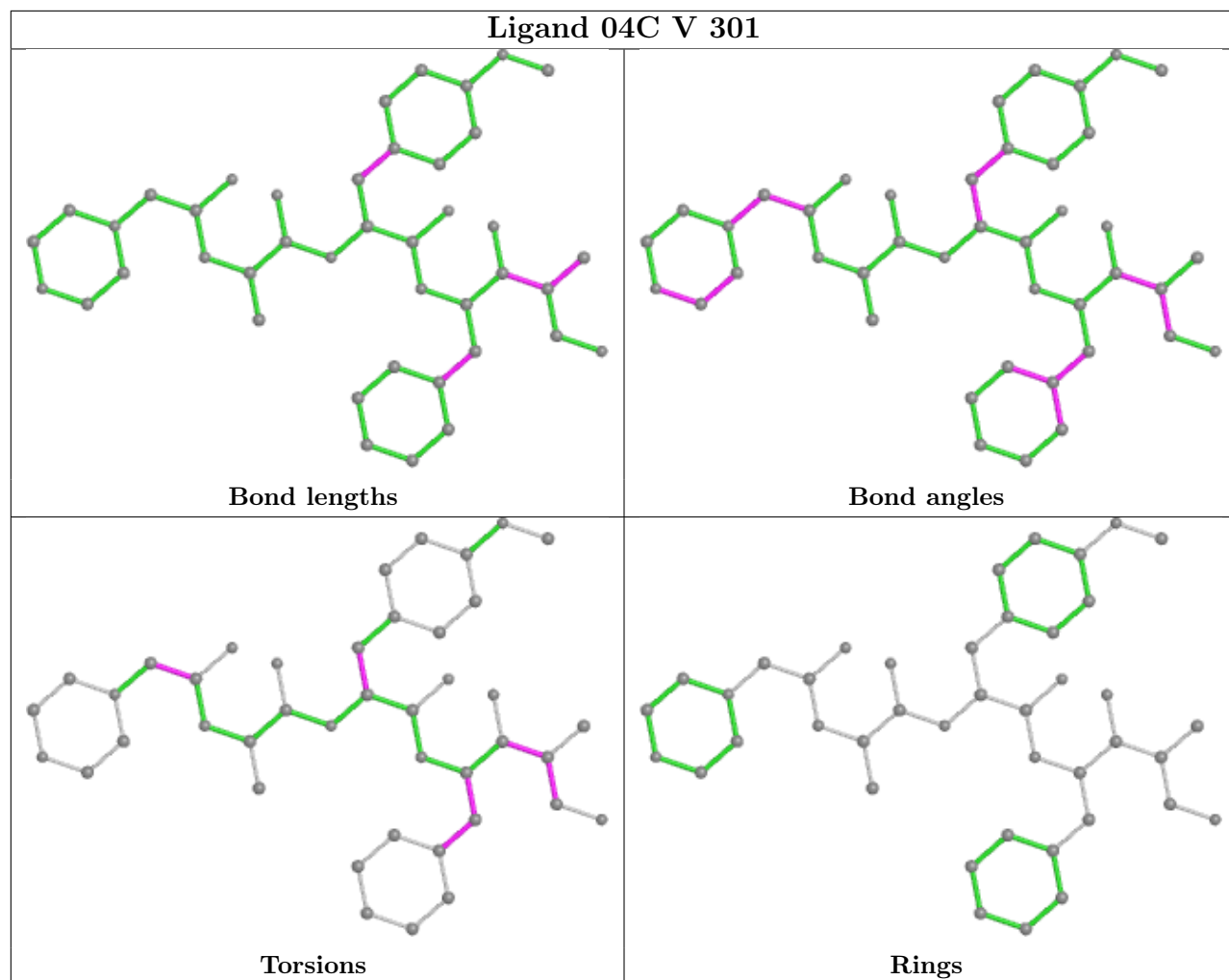
5 monomers are involved in 13 short contacts:

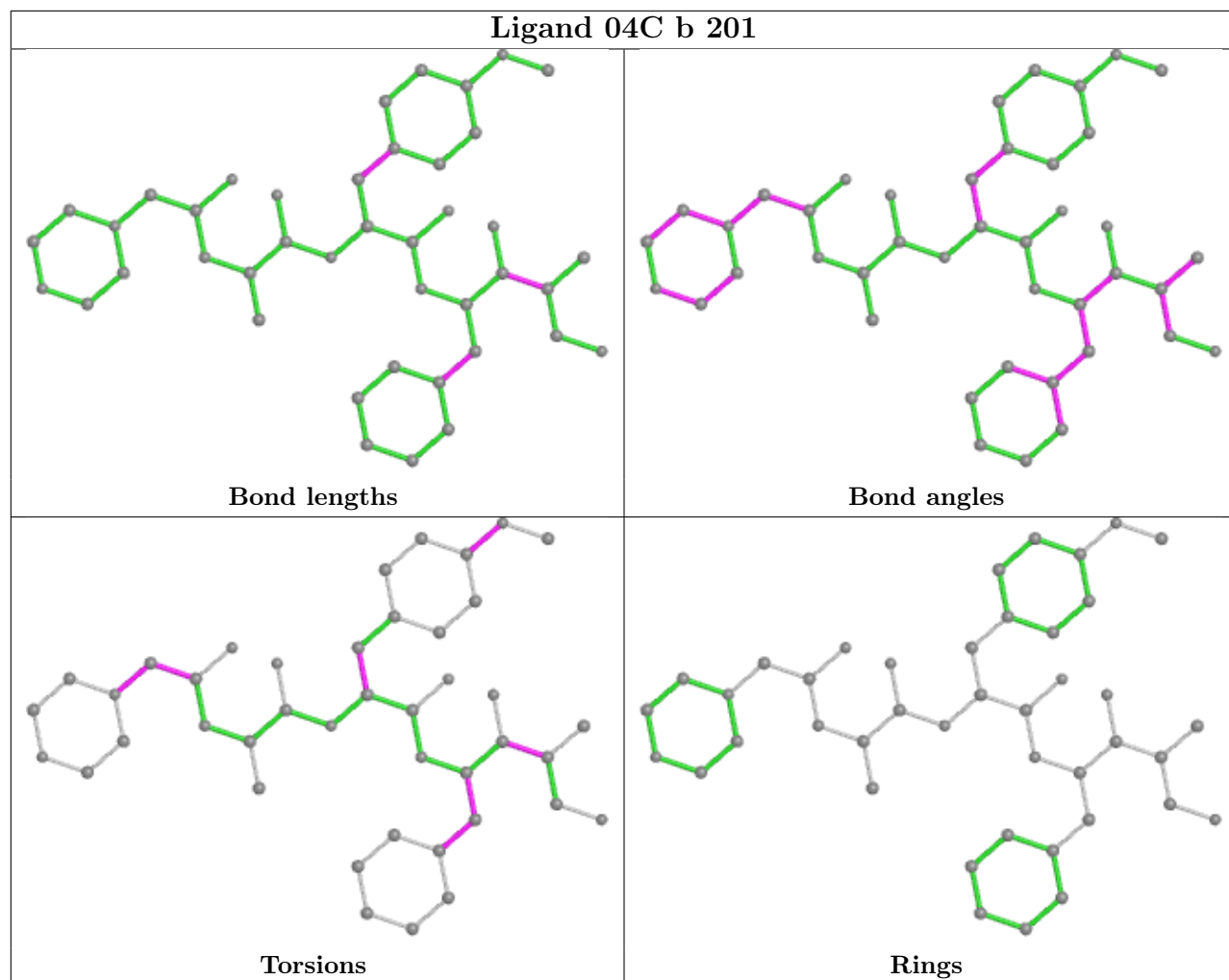
Mol	Chain	Res	Type	Clashes	Symm-Clashes
17	N	201	04C	5	0
17	H	301	04C	3	0
17	V	301	04C	3	0
18	K	304	MES	1	0
18	Y	304	MES	1	0

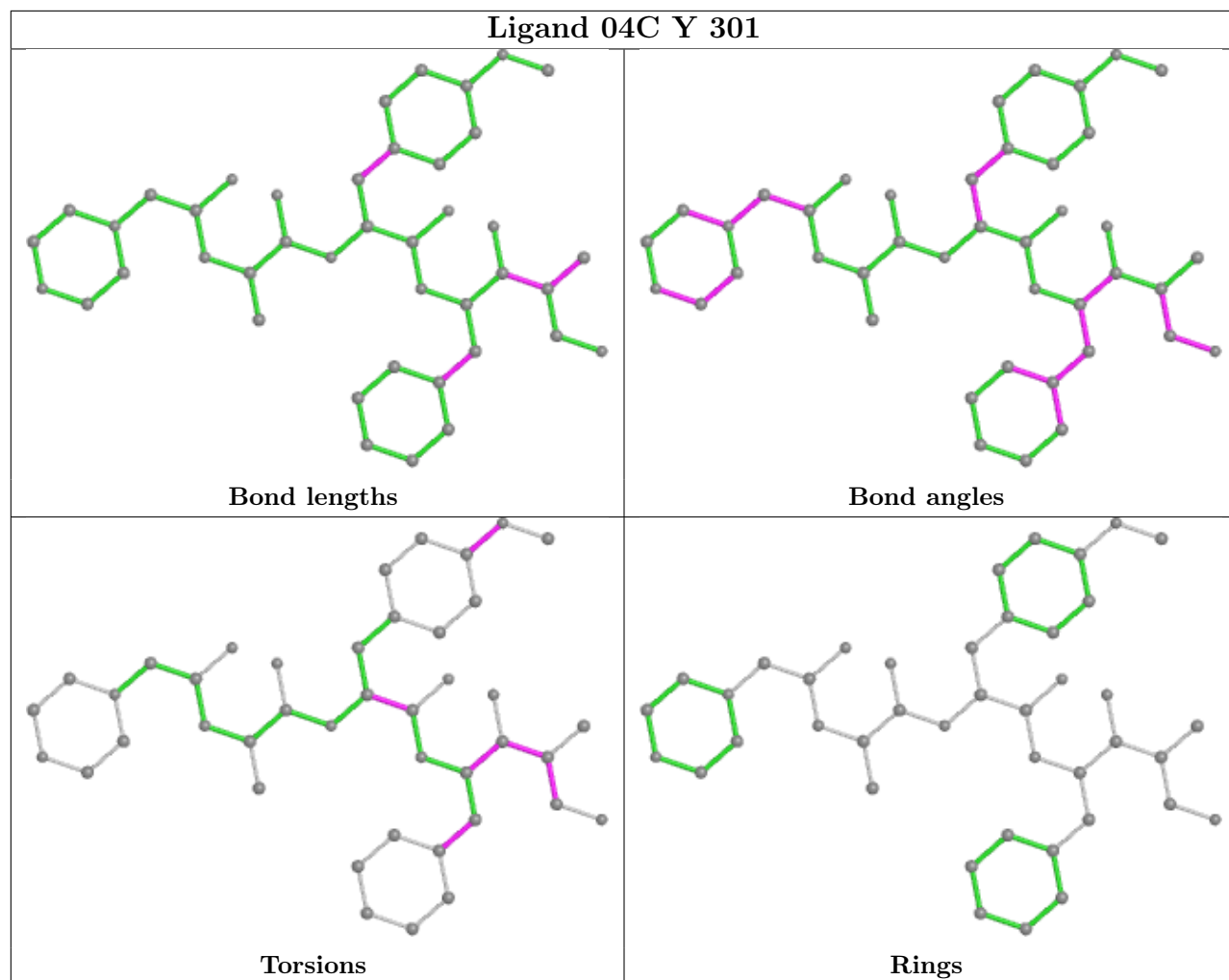
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

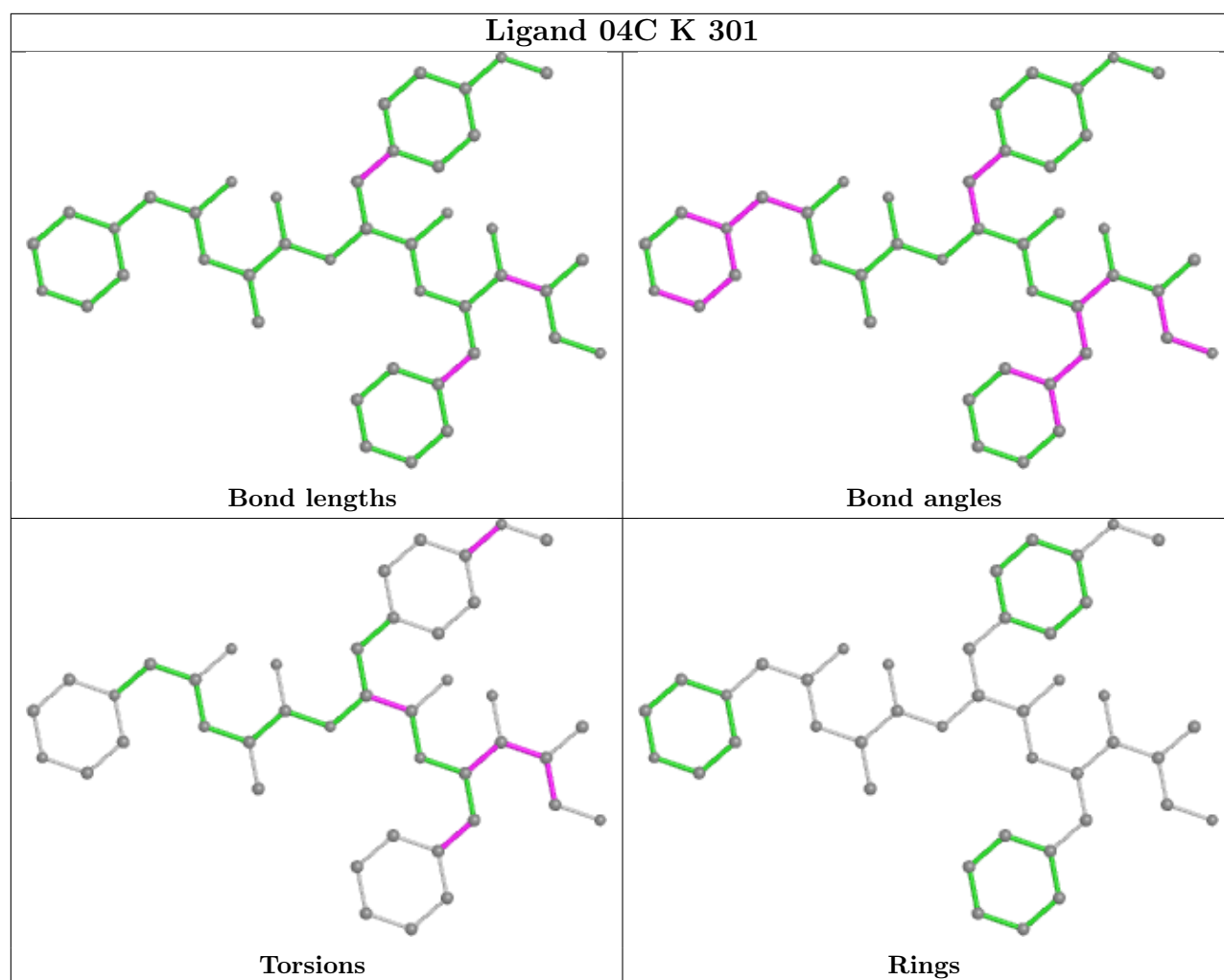












## 5.7 Other polymers [\(i\)](#)

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 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	250/250 (100%)	-0.48	3 (1%) 79 79	53, 66, 108, 152	0
1	O	250/250 (100%)	-0.45	5 (2%) 65 63	51, 72, 112, 154	0
2	B	244/258 (94%)	-0.42	8 (3%) 46 41	49, 69, 118, 174	0
2	P	244/258 (94%)	-0.41	7 (2%) 51 47	51, 73, 118, 160	0
3	C	240/254 (94%)	-0.30	11 (4%) 32 29	48, 74, 136, 163	0
3	Q	240/254 (94%)	-0.03	15 (6%) 20 16	59, 91, 172, 208	0
4	D	235/260 (90%)	-0.46	4 (1%) 70 69	51, 77, 110, 151	0
4	R	235/260 (90%)	-0.39	4 (1%) 70 69	58, 78, 118, 170	0
5	E	231/234 (98%)	-0.29	5 (2%) 62 59	58, 85, 122, 170	0
5	S	231/234 (98%)	-0.14	8 (3%) 44 38	58, 90, 147, 184	0
6	F	243/288 (84%)	-0.49	4 (1%) 72 71	55, 80, 127, 162	0
6	T	243/288 (84%)	-0.36	4 (1%) 72 71	56, 84, 143, 175	0
7	G	241/252 (95%)	-0.48	5 (2%) 63 61	49, 71, 111, 155	0
7	U	241/252 (95%)	-0.50	3 (1%) 79 79	51, 70, 106, 140	0
8	H	222/232 (95%)	-0.61	2 (0%) 84 84	48, 65, 93, 117	0
8	V	222/232 (95%)	-0.58	2 (0%) 84 84	48, 67, 95, 123	0
9	I	204/205 (99%)	-0.76	1 (0%) 91 91	44, 59, 88, 119	0
9	W	204/205 (99%)	-0.73	0 100 100	43, 60, 89, 119	0
10	J	195/198 (98%)	-0.65	2 (1%) 82 82	42, 58, 88, 131	0
10	X	195/198 (98%)	-0.62	2 (1%) 82 82	46, 61, 92, 144	0
11	K	212/212 (100%)	-0.69	0 100 100	40, 61, 87, 107	0
11	Y	212/212 (100%)	-0.68	0 100 100	32, 60, 88, 115	0
12	L	222/222 (100%)	-0.65	0 100 100	44, 65, 94, 111	0
12	Z	222/222 (100%)	-0.70	1 (0%) 91 91	43, 63, 92, 109	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
13	M	233/246 (94%)	-0.56	5 (2%) 63 61	44, 69, 99, 124	0
13	a	233/246 (94%)	-0.58	3 (1%) 77 77	43, 66, 95, 127	0
14	N	196/196 (100%)	-0.67	1 (0%) 91 91	44, 64, 93, 120	0
14	b	196/196 (100%)	-0.65	1 (0%) 91 91	45, 63, 95, 116	0
All	All	6336/6614 (95%)	-0.50	106 (1%) 70 69	32, 70, 119, 208	0

The worst 5 of 106 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	Q	49	THR	6.9
3	C	206	LYS	6.1
4	R	241	ALA	5.3
1	A	1	MET	5.2
3	Q	206	LYS	5.1

## 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(Å <sup>2</sup> )	Q<0.9
17	04C	b	201	42/42	0.89	0.36	47,63,83,90	0
17	04C	N	201	42/42	0.91	0.37	45,62,76,81	0
17	04C	H	301	42/42	0.94	0.16	41,50,75,77	0
17	04C	V	301	42/42	0.95	0.15	44,50,84,86	0
17	04C	K	301	42/42	0.96	0.13	34,49,61,62	0
15	MG	G	301	1/1	0.96	0.09	67,67,67,67	0
18	MES	V	303	12/12	0.96	0.25	64,71,86,91	0

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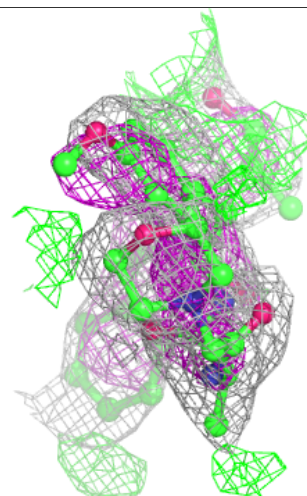
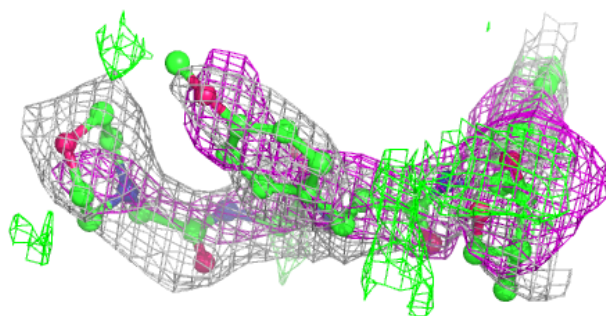
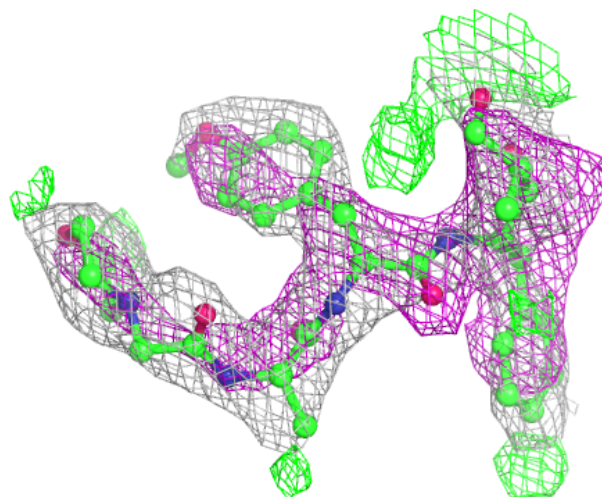
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
15	MG	Z	301	1/1	0.97	0.15	77,77,77,77	0
18	MES	H	302	12/12	0.97	0.21	66,70,74,75	0
17	04C	Y	301	42/42	0.97	0.14	29,44,63,63	0
15	MG	N	202	1/1	0.98	0.17	45,45,45,45	0
15	MG	V	302	1/1	0.98	0.13	75,75,75,75	0
15	MG	X	201	1/1	0.98	0.32	29,29,29,29	0
15	MG	Y	302	1/1	0.98	0.08	63,63,63,63	0
15	MG	I	301	1/1	0.98	0.17	61,61,61,61	0
16	CL	U	301	1/1	0.98	0.17	52,52,52,52	0
15	MG	K	302	1/1	0.98	0.07	62,62,62,62	0
18	MES	Y	304	12/12	0.98	0.17	44,45,57,57	0
19	NA	Y	303	1/1	0.98	0.11	24,24,24,24	0
18	MES	K	304	12/12	0.99	0.19	42,47,54,58	0
19	NA	K	303	1/1	0.99	0.17	32,32,32,32	0
16	CL	G	302	1/1	0.99	0.12	52,52,52,52	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.

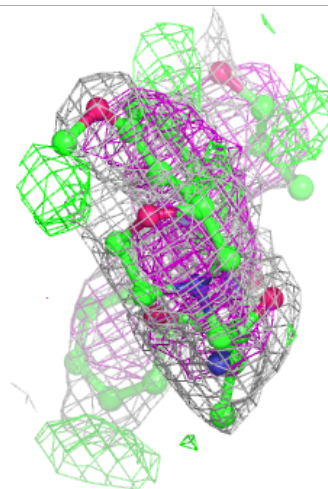
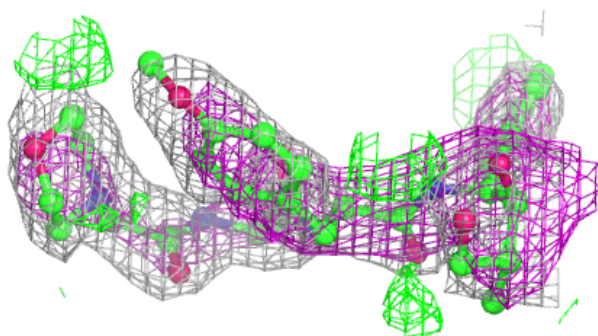
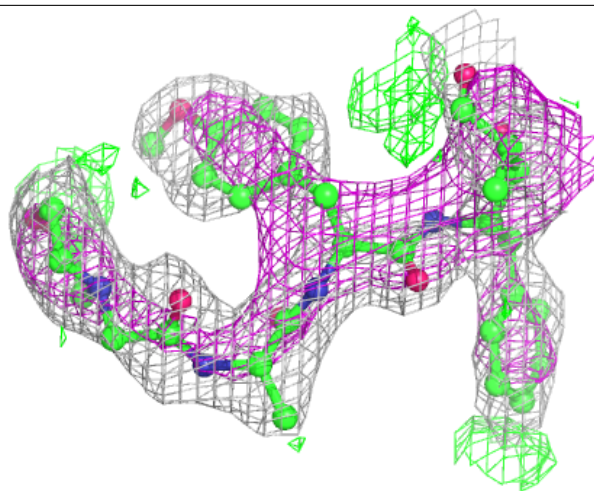
**Electron density around 04C b 201:**

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



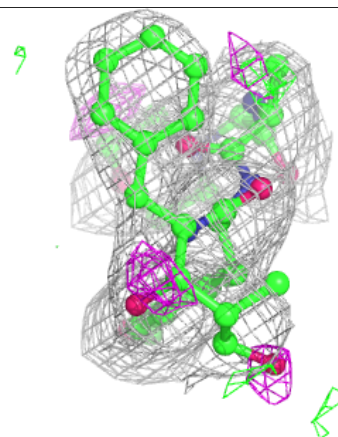
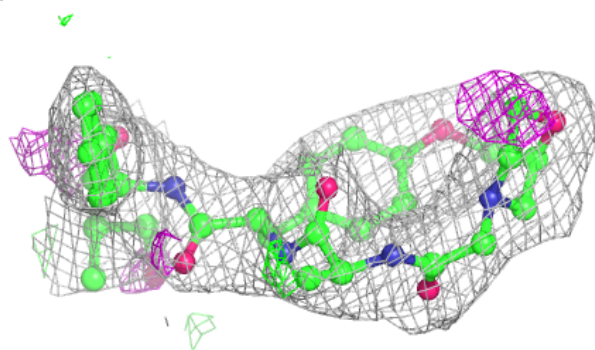
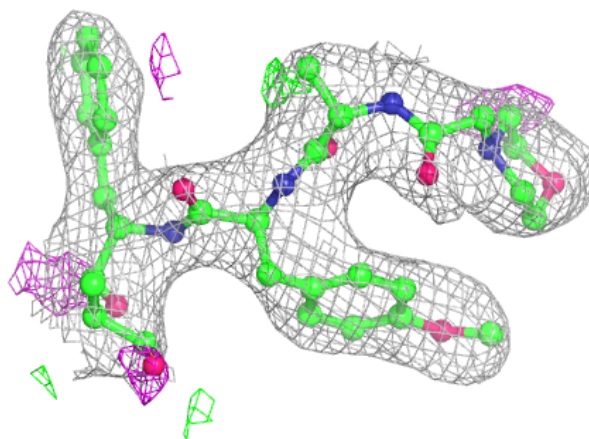
**Electron density around 04C N 201:**

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



**Electron density around 04C H 301:**

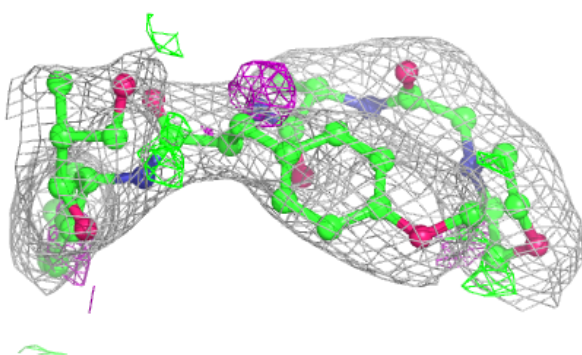
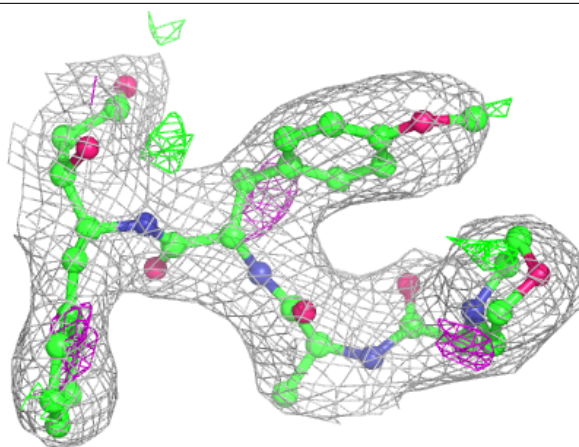
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



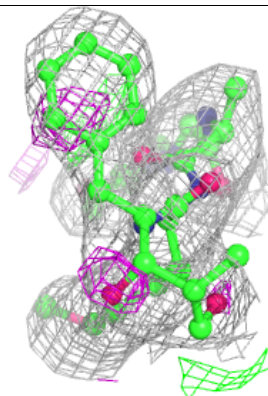
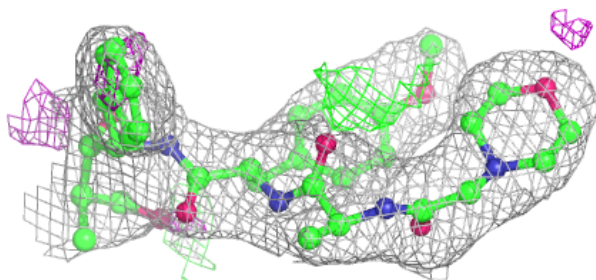
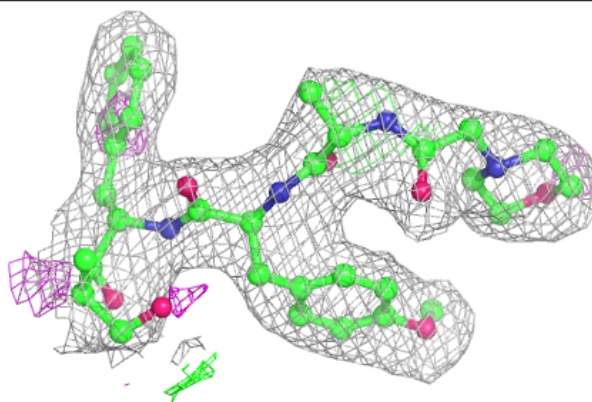


**Electron density around 04C V 301:**

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

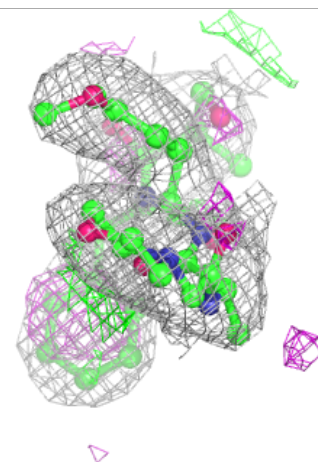
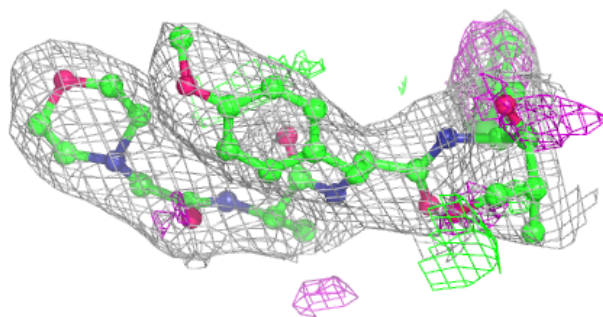
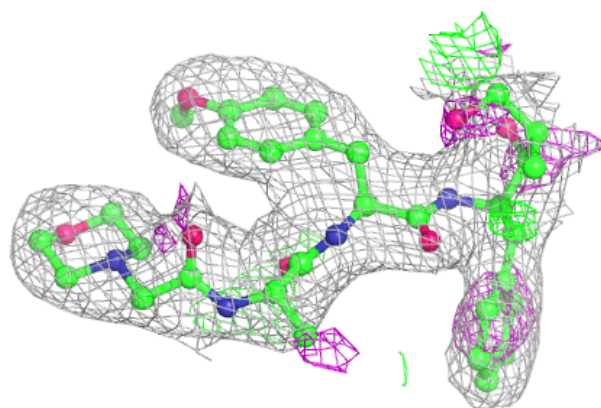
**Electron density around 04C K 301:**

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



**Electron density around 04C Y 301:**

$2mF_o-DF_c$  (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.