



wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 17, 2024 – 02:24 pm BST

PDB ID : 6ZP8
Title : Yeast 20S proteasome in complex with glidobactin-like natural product HB335
Authors : Zhao, L.; Le Chapelain, C.; Brachmann, A.O.; Kaiser, M.; Groll, M.; Bode, H.B.
Deposited on : 2020-07-08
Resolution : 3.00 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

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>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.002 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.38.2

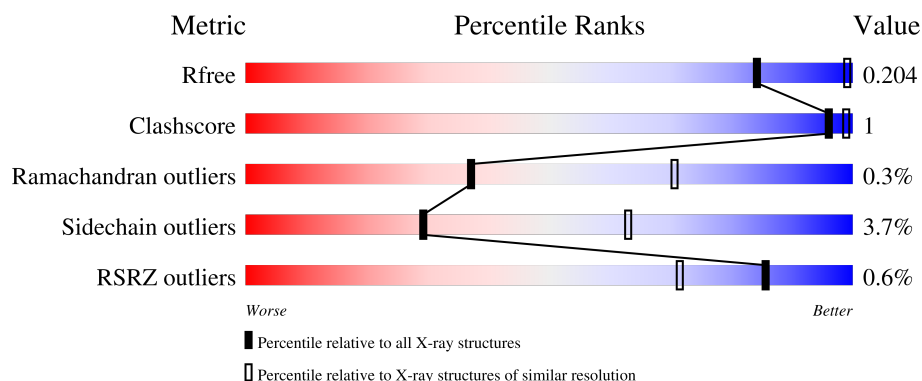
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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}	164625	2511 (3.00-3.00)
Clashscore	180529	2866 (3.00-3.00)
Ramachandran outliers	177936	2778 (3.00-3.00)
Sidechain outliers	177891	2781 (3.00-3.00)
RSRZ outliers	164620	2523 (3.00-3.00)

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 ≥ 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	<div> <div style="width: 98%;"></div> <div>98%</div> </div>
1	O	250	<div> <div style="width: 98%;"></div> <div>98%</div> </div>
2	B	258	<div> <div style="width: 88%;"></div> <div>88%</div> <div style="width: 6%;"></div> <div>6%</div> <div style="width: 5%;"></div> <div>5%</div> </div>
2	P	258	<div> <div style="width: 88%;"></div> <div>88%</div> <div style="width: 6%;"></div> <div>6%</div> <div style="width: 5%;"></div> <div>5%</div> </div>
3	C	254	<div> <div style="width: 86%;"></div> <div>86%</div> <div style="width: 7%;"></div> <div>7%</div> <div style="width: 6%;"></div> <div>6%</div> </div>

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Mol	Chain	Length	Quality of chain
3	Q	254	 2% 87% 7% • 6%
4	D	260	 % 86% • • 10%
4	R	260	 84% 6% 10%
5	E	234	 93% 5% •
5	S	234	 % 93% 6% •
6	F	288	 % 80% • 16%
6	T	288	 % 80% • 16%
7	G	252	 90% 6% •
7	U	252	 90% 6% •
8	H	232	 93% • • •
8	V	232	 % 92% • • •
9	I	205	 94% 5%
9	W	205	 95% 5%
10	J	198	 % 93% 5% • •
10	X	198	 % 93% 5% • •
11	K	212	 92% 7% •
11	Y	212	 % 93% 6%
12	L	222	 % 96% •
12	Z	222	 95% 5%
13	M	246	 90% • 7%
13	a	246	 91% • 6%
14	N	196	 95% • •
14	b	196	 % 97% •

2 Entry composition

There are 18 unique types of molecules in this entry. The entry contains 49792 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
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
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
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
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
5	E	231	Total	C	N	O	S	0	0	0
			1773	1114	307	348	4			
5	S	231	Total	C	N	O	S	0	0	0
			1773	1114	307	348	4			

- Molecule 6 is a protein called Probable proteasome subunit alpha type-7.

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

- Molecule 7 is a protein called Proteasome subunit alpha type-1.

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

- Molecule 8 is a protein called Proteasome subunit beta type-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	226	Total	C	N	O	S	0	0	0
			1719	1082	298	332	7			
8	V	226	Total	C	N	O	S	0	0	0
			1719	1082	298	332	7			

- Molecule 9 is a protein called Proteasome subunit beta type-3.

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

- Molecule 10 is a protein called Proteasome subunit beta type-4.

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

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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	229	Total	C	N	O	S	0	0	0
			1790	1133	306	344	7			
13	a	232	Total	C	N	O	S	0	0	0
			1815	1148	311	349	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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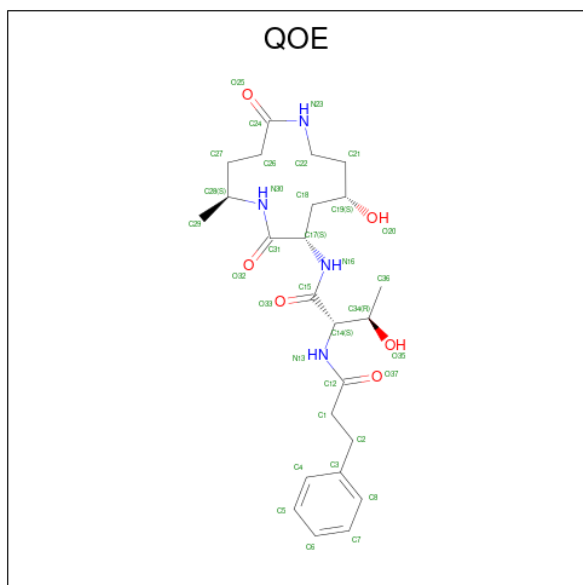
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	K	1	Total	Mg	0	0
			1	1		
15	N	1	Total	Mg	0	0
			1	1		
15	V	1	Total	Mg	0	0
			1	1		
15	W	1	Total	Mg	0	0
			1	1		
15	Y	1	Total	Mg	0	0
			1	1		
15	Z	1	Total	Mg	0	0
			1	1		

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

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

- Molecule 17 is (2 {S},3 {R})- {N}-[(5 {S},8 {S},10 {S})-5-methyl-10-oxidanyl-2,7-bis(oxidanylidene)-1,6-diazacyclododec-8-yl]-3-oxidanyl-2-(3-phenylpropanoylamino)butanamide (three-letter code: QOE) (formula: C₂₄H₃₆N₄O₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
17	H	1	Total	C	N	O	0	0
			34	24	4	6		
17	K	1	Total	C	N	O	0	0
			34	24	4	6		
17	V	1	Total	C	N	O	0	0
			34	24	4	6		
17	Y	1	Total	C	N	O	0	0
			34	24	4	6		

- Molecule 18 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
18	A	10	Total	O	0	0
			10	10		
18	B	17	Total	O	0	0
			17	17		
18	C	12	Total	O	0	0
			12	12		
18	D	5	Total	O	0	0
			5	5		
18	E	9	Total	O	0	0
			9	9		
18	F	14	Total	O	0	0
			14	14		
18	G	7	Total	O	0	0
			7	7		
18	H	13	Total	O	0	0
			13	13		
18	I	5	Total	O	0	0
			5	5		
18	J	13	Total	O	0	0
			13	13		
18	K	6	Total	O	0	0
			6	6		
18	L	21	Total	O	0	0
			21	21		
18	M	18	Total	O	0	0
			18	18		
18	N	11	Total	O	0	0
			11	11		
18	O	5	Total	O	0	0
			5	5		
18	P	14	Total	O	0	0
			14	14		

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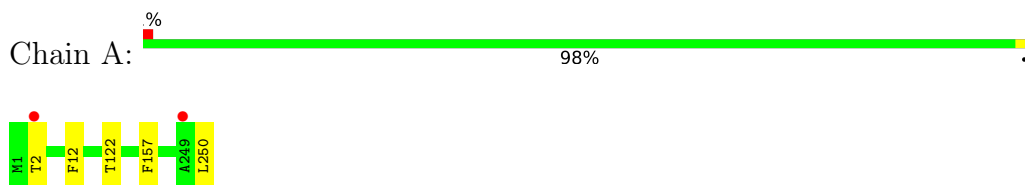
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
18	Q	11	Total 11	O 11	0	0
18	R	10	Total 10	O 10	0	0
18	S	7	Total 7	O 7	0	0
18	T	11	Total 11	O 11	0	0
18	U	15	Total 15	O 15	0	0
18	V	16	Total 16	O 16	0	0
18	W	7	Total 7	O 7	0	0
18	X	12	Total 12	O 12	0	0
18	Y	9	Total 9	O 9	0	0
18	Z	18	Total 18	O 18	0	0
18	a	14	Total 14	O 14	0	0
18	b	13	Total 13	O 13	0	0

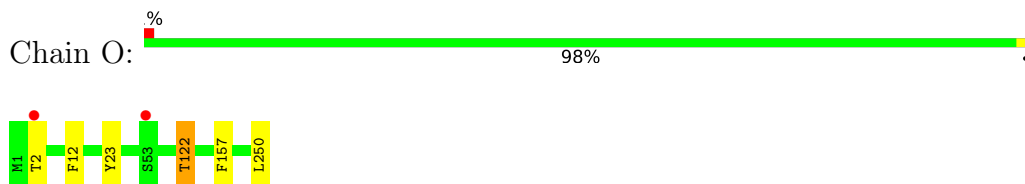
3 Residue-property plots [i](#)

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.

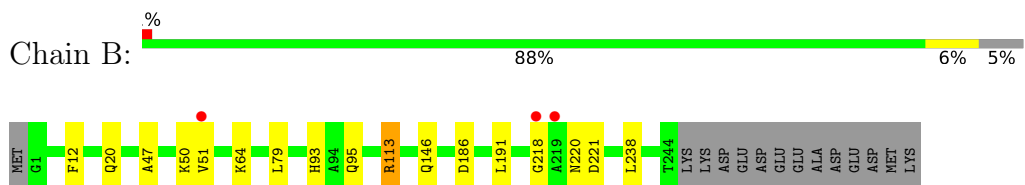
- Molecule 1: Proteasome subunit alpha type-2



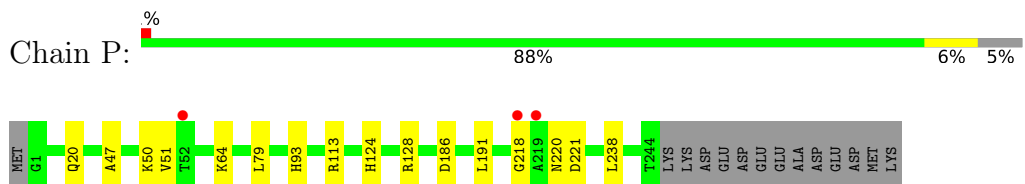
- Molecule 1: Proteasome subunit alpha type-2



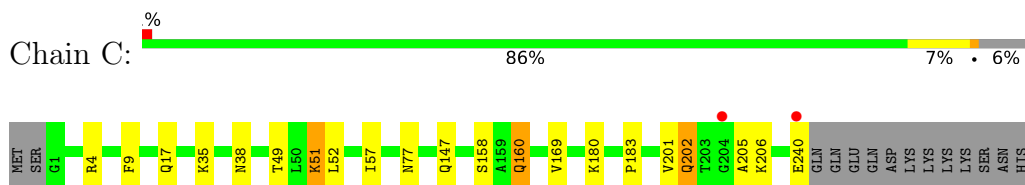
- Molecule 2: Proteasome subunit alpha type-3



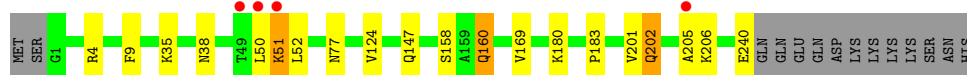
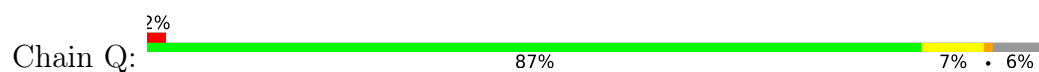
- Molecule 2: Proteasome subunit alpha type-3



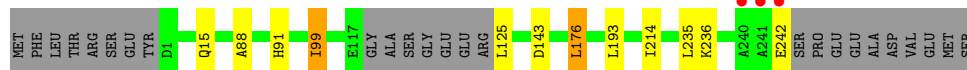
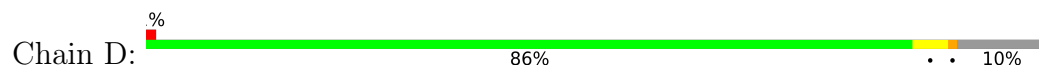
- Molecule 3: Proteasome subunit alpha type-4



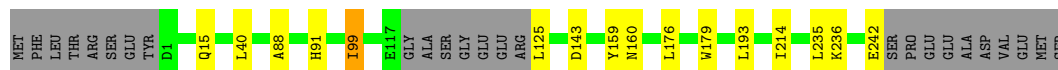
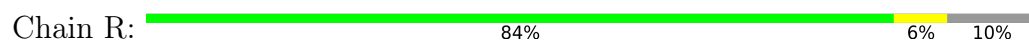
- Molecule 3: Proteasome subunit alpha type-4



- Molecule 4: Proteasome subunit alpha type-5



- Molecule 4: Proteasome subunit alpha type-5



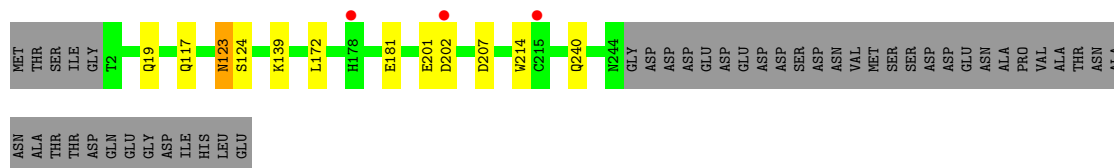
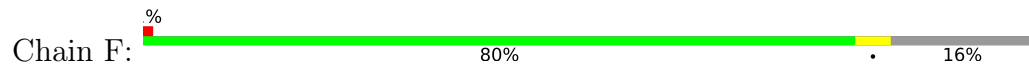
- Molecule 5: Proteasome subunit alpha type-6



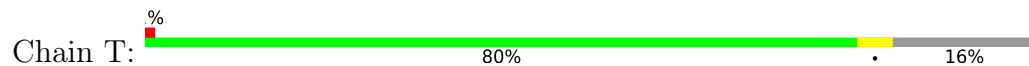
- Molecule 5: Proteasome subunit alpha type-6

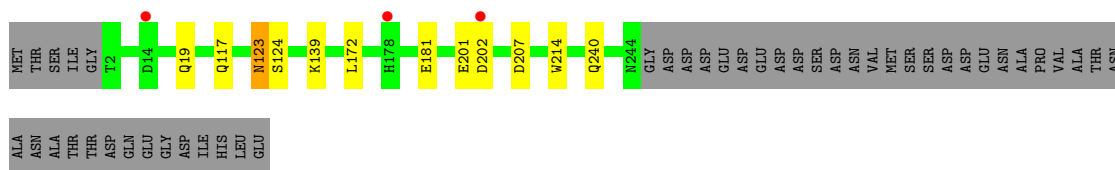


- Molecule 6: Probable proteasome subunit alpha type-7




- Molecule 6: Probable proteasome subunit alpha type-7





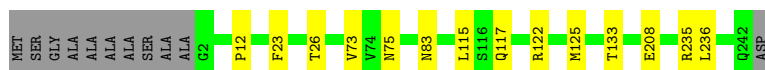
- Molecule 7: Proteasome subunit alpha type-1

Chain G:  90% 6% .



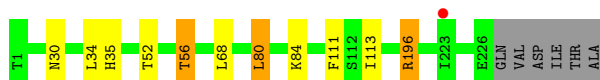
- Molecule 7: Proteasome subunit alpha type-1

Chain U:  90% 6% .



- Molecule 8: Proteasome subunit beta type-2

Chain H:  93% . . .



- Molecule 8: Proteasome subunit beta type-2

Chain V:  92% . . .



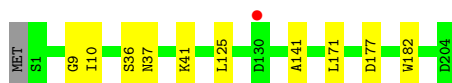
- Molecule 9: Proteasome subunit beta type-3

Chain I:  94% 5%

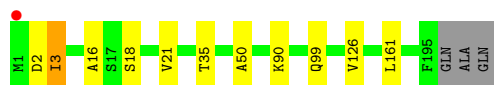
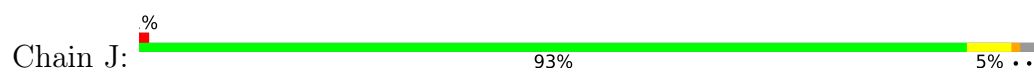


- Molecule 9: Proteasome subunit beta type-3

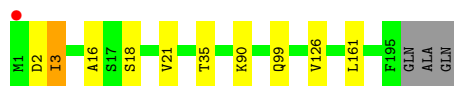
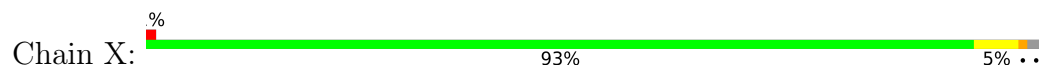
Chain W:  95% 5%



- Molecule 10: Proteasome subunit beta type-4



- Molecule 10: Proteasome subunit beta type-4



- Molecule 11: Proteasome subunit beta type-5



- Molecule 11: Proteasome subunit beta type-5



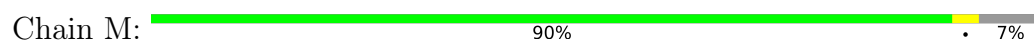
- Molecule 12: Proteasome subunit beta type-6



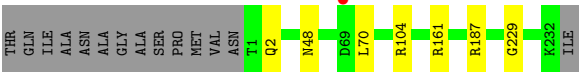
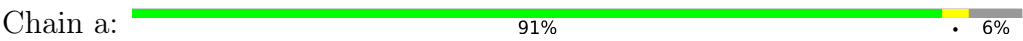
- Molecule 12: Proteasome subunit beta type-6



- Molecule 13: Proteasome subunit beta type-7



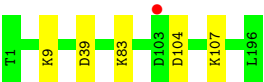
● Molecule 13: Proteasome subunit beta type-7



● 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, α , β , γ	136.16Å 300.12Å 144.83Å 90.00° 113.11° 90.00°	Depositor
Resolution (Å)	15.00 – 3.00 15.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	98.4 (15.00-3.00) 98.4 (15.00-3.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.167 , 0.200 0.173 , 0.204	Depositor DCC
R_{free} test set	10458 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	60.6	Xtriage
Anisotropy	0.274	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 41.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	49792	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.25% 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: CL, QOE, MG

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	0.36	0/1952	0.57	0/2642
1	O	0.36	0/1952	0.58	0/2642
2	B	0.37	0/1934	0.62	0/2618
2	P	0.36	0/1934	0.62	0/2618
3	C	0.36	0/1910	0.64	0/2586
3	Q	0.36	0/1910	0.63	0/2586
4	D	0.35	0/1837	0.59	0/2475
4	R	0.35	0/1837	0.59	0/2475
5	E	0.36	0/1800	0.59	0/2433
5	S	0.35	0/1800	0.59	0/2433
6	F	0.36	0/1932	0.56	0/2609
6	T	0.35	0/1932	0.56	0/2609
7	G	0.36	0/1945	0.58	0/2634
7	U	0.36	0/1945	0.57	0/2634
8	H	0.34	0/1750	0.60	1/2373 (0.0%)
8	V	0.34	0/1750	0.61	1/2373 (0.0%)
9	I	0.36	0/1611	0.59	0/2174
9	W	0.35	0/1611	0.59	0/2174
10	J	0.35	0/1589	0.61	0/2142
10	X	0.35	0/1589	0.62	0/2142
11	K	0.35	0/1681	0.62	1/2274 (0.0%)
11	Y	0.34	0/1681	0.63	1/2274 (0.0%)
12	L	0.36	0/1795	0.60	0/2420
12	Z	0.36	0/1795	0.59	0/2420
13	M	0.37	0/1821	0.64	0/2470
13	a	0.36	0/1846	0.64	0/2503
14	N	0.34	0/1541	0.60	1/2087 (0.0%)
14	b	0.34	0/1541	0.60	0/2087
All	All	0.35	0/50221	0.60	5/67907 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	V	196	ARG	NE-CZ-NH1	5.39	122.99	120.30
8	H	196	ARG	NE-CZ-NH1	5.32	122.96	120.30
11	Y	4	LEU	CA-CB-CG	5.24	127.34	115.30
11	K	4	LEU	CA-CB-CG	5.13	127.10	115.30
14	N	36	ARG	NE-CZ-NH2	-5.03	117.78	120.30

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	1915	0	1929	1	0
1	O	1915	0	1929	3	0
2	B	1904	0	1904	7	0
2	P	1904	0	1904	5	0
3	C	1881	0	1895	7	0
3	Q	1881	0	1895	6	0
4	D	1813	0	1797	5	0
4	R	1813	0	1797	6	0
5	E	1773	0	1775	4	0
5	S	1773	0	1775	3	0
6	F	1892	0	1883	3	0
6	T	1892	0	1883	3	0
7	G	1907	0	1901	3	0
7	U	1907	0	1901	3	0
8	H	1719	0	1718	6	0
8	V	1719	0	1718	8	0
9	I	1581	0	1574	7	0
9	W	1581	0	1574	6	0
10	J	1561	0	1569	5	0
10	X	1561	0	1569	4	0
11	K	1644	0	1594	5	0
11	Y	1644	0	1594	3	0
12	L	1757	0	1711	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
12	Z	1757	0	1711	4	0
13	M	1790	0	1793	2	0
13	a	1815	0	1821	0	0
14	N	1512	0	1481	4	0
14	b	1512	0	1481	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	W	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	34	0	0	2	0
17	K	34	0	0	0	0
17	V	34	0	0	3	0
17	Y	34	0	0	1	0
18	A	10	0	0	0	0
18	B	17	0	0	2	0
18	C	12	0	0	0	0
18	D	5	0	0	0	0
18	E	9	0	0	1	0
18	F	14	0	0	0	0
18	G	7	0	0	0	0
18	H	13	0	0	0	0
18	I	5	0	0	0	0
18	J	13	0	0	0	0
18	K	6	0	0	0	0
18	L	21	0	0	0	0
18	M	18	0	0	0	0
18	N	11	0	0	0	0
18	O	5	0	0	0	0
18	P	14	0	0	1	0
18	Q	11	0	0	0	0
18	R	10	0	0	0	0
18	S	7	0	0	0	0
18	T	11	0	0	0	0
18	U	15	0	0	0	0
18	V	16	0	0	0	0
18	W	7	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
18	X	12	0	0	0	0
18	Y	9	0	0	0	0
18	Z	18	0	0	0	0
18	a	14	0	0	0	0
18	b	13	0	0	0	0
All	All	49792	0	49076	95	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 95 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)
17:H:901:QOE:C5	9:I:125:LEU:HD21	2.12	0.80
8:H:52:THR:O	8:H:56:THR:OG1	2.09	0.71
17:H:901:QOE:C4	9:I:125:LEU:HD21	2.21	0.70
8:V:52:THR:O	8:V:56:THR:OG1	2.11	0.68
11:Y:1:THR:O	11:Y:130:GLY:HA3	1.97	0.65

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	248/250 (99%)	241 (97%)	6 (2%)	1 (0%)	30 66
1	O	248/250 (99%)	239 (96%)	8 (3%)	1 (0%)	30 66
2	B	242/258 (94%)	234 (97%)	4 (2%)	4 (2%)	7 33
2	P	242/258 (94%)	234 (97%)	4 (2%)	4 (2%)	7 33
3	C	238/254 (94%)	233 (98%)	2 (1%)	3 (1%)	10 39

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	Q	238/254 (94%)	233 (98%)	2 (1%)	3 (1%)	10	39
4	D	231/260 (89%)	227 (98%)	4 (2%)	0	100	100
4	R	231/260 (89%)	227 (98%)	4 (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%)	238 (99%)	3 (1%)	0	100	100
6	T	241/288 (84%)	238 (99%)	3 (1%)	0	100	100
7	G	239/252 (95%)	237 (99%)	2 (1%)	0	100	100
7	U	239/252 (95%)	238 (100%)	1 (0%)	0	100	100
8	H	224/232 (97%)	219 (98%)	5 (2%)	0	100	100
8	V	224/232 (97%)	218 (97%)	6 (3%)	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%)	4 (2%)	0	100	100
10	X	193/198 (98%)	189 (98%)	4 (2%)	0	100	100
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	227/246 (92%)	220 (97%)	7 (3%)	0	100	100
13	a	230/246 (94%)	223 (97%)	6 (3%)	1 (0%)	30	66
14	N	194/196 (99%)	190 (98%)	4 (2%)	0	100	100
14	b	194/196 (99%)	190 (98%)	4 (2%)	0	100	100
All	All	6279/6614 (95%)	6139 (98%)	123 (2%)	17 (0%)	37	70

5 of 17 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	THR
2	B	51	VAL
2	B	221	ASP
3	C	202	GLN
1	O	2	THR

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	209/209 (100%)	206 (99%)	3 (1%)	62	83
1	O	209/209 (100%)	206 (99%)	3 (1%)	62	83
2	B	203/216 (94%)	197 (97%)	6 (3%)	36	69
2	P	203/216 (94%)	197 (97%)	6 (3%)	36	69
3	C	212/226 (94%)	201 (95%)	11 (5%)	19	52
3	Q	212/226 (94%)	201 (95%)	11 (5%)	19	52
4	D	194/215 (90%)	185 (95%)	9 (5%)	23	56
4	R	194/215 (90%)	184 (95%)	10 (5%)	19	52
5	E	190/193 (98%)	181 (95%)	9 (5%)	22	56
5	S	190/193 (98%)	181 (95%)	9 (5%)	22	56
6	F	201/239 (84%)	191 (95%)	10 (5%)	20	53
6	T	201/239 (84%)	191 (95%)	10 (5%)	20	53
7	G	206/210 (98%)	197 (96%)	9 (4%)	24	58
7	U	206/210 (98%)	197 (96%)	9 (4%)	24	58
8	H	185/190 (97%)	179 (97%)	6 (3%)	34	67
8	V	185/190 (97%)	179 (97%)	6 (3%)	34	67
9	I	172/173 (99%)	169 (98%)	3 (2%)	56	81
9	W	172/173 (99%)	169 (98%)	3 (2%)	56	81
10	J	173/175 (99%)	168 (97%)	5 (3%)	37	70
10	X	173/175 (99%)	168 (97%)	5 (3%)	37	70
11	K	169/169 (100%)	160 (95%)	9 (5%)	19	51
11	Y	169/169 (100%)	160 (95%)	9 (5%)	19	51
12	L	185/185 (100%)	178 (96%)	7 (4%)	28	62
12	Z	185/185 (100%)	178 (96%)	7 (4%)	28	62
13	M	195/208 (94%)	189 (97%)	6 (3%)	35	68
13	a	198/208 (95%)	192 (97%)	6 (3%)	36	69

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
14	N	162/162 (100%)	157 (97%)	5 (3%)	35	68
14	b	162/162 (100%)	157 (97%)	5 (3%)	35	68
All	All	5315/5540 (96%)	5118 (96%)	197 (4%)	29	63

5 of 197 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	Q	180	LYS
6	T	202	ASP
4	R	99	ILE
5	S	55	LEU
7	U	117	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 125 such sidechains are listed below:

Mol	Chain	Res	Type
13	M	102	GLN
11	Y	176	ASN
3	Q	147	GLN
11	Y	85	ASN
13	a	18	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 10 are monoatomic - leaving 4 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	QOE	H	901	8	35,35,35	2.22	6 (17%)	44,46,46	1.15	4 (9%)
17	QOE	V	901	8	35,35,35	2.17	5 (14%)	44,46,46	1.35	7 (15%)
17	QOE	K	901	11	35,35,35	2.31	6 (17%)	44,46,46	1.26	3 (6%)
17	QOE	Y	901	11	35,35,35	2.18	4 (11%)	44,46,46	1.35	5 (11%)

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	QOE	H	901	8	-	5/44/44/44	0/1/2/2
17	QOE	V	901	8	-	5/44/44/44	0/1/2/2
17	QOE	K	901	11	-	15/44/44/44	0/1/2/2
17	QOE	Y	901	11	-	7/44/44/44	0/1/2/2

The worst 5 of 21 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
17	K	901	QOE	C26-C24	-8.29	1.35	1.51
17	H	901	QOE	C26-C24	-8.16	1.35	1.51
17	V	901	QOE	C26-C24	-8.13	1.36	1.51
17	Y	901	QOE	C17-C31	-7.13	1.34	1.52
17	Y	901	QOE	C26-C24	-7.10	1.37	1.51

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	Y	901	QOE	C26-C27-C28	-4.21	107.54	114.02
17	V	901	QOE	C26-C27-C28	-3.81	108.16	114.02
17	K	901	QOE	C22-N23-C24	3.34	129.03	122.84
17	K	901	QOE	C26-C27-C28	-3.20	109.10	114.02
17	V	901	QOE	C15-C14-N13	-3.04	102.08	110.36

There are no chirality outliers.

5 of 32 torsion outliers are listed below:

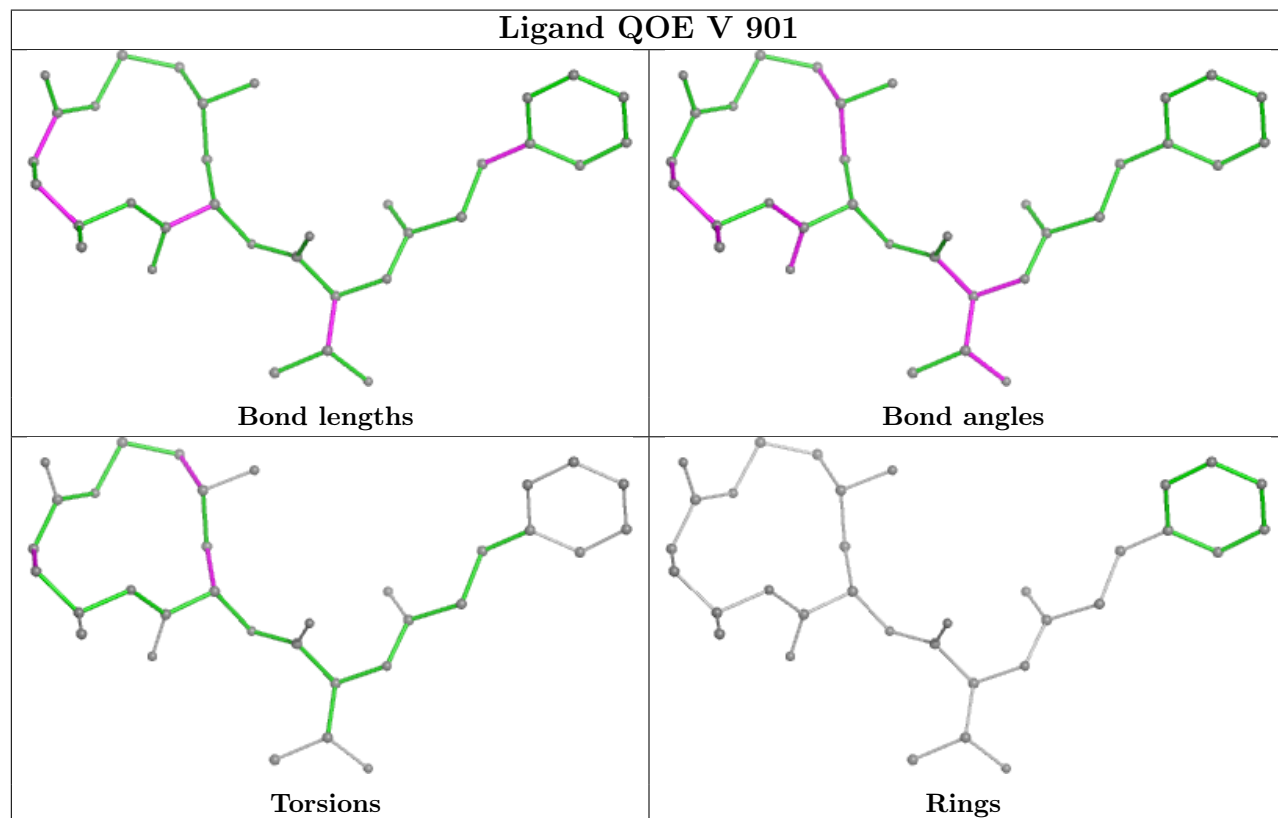
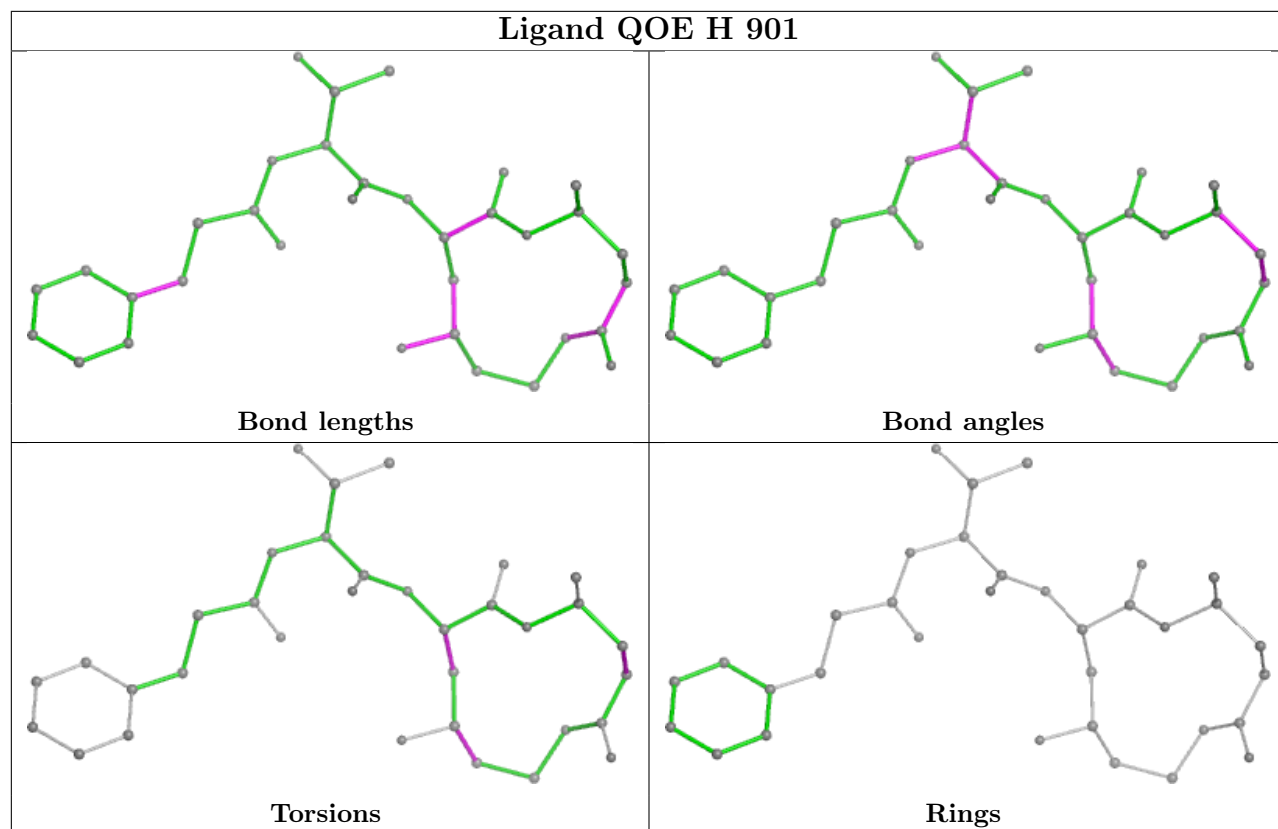
Mol	Chain	Res	Type	Atoms
17	H	901	QOE	N16-C17-C18-C19
17	H	901	QOE	C24-C26-C27-C28
17	K	901	QOE	N16-C17-C18-C19
17	K	901	QOE	O20-C19-C21-C22
17	K	901	QOE	C18-C19-C21-C22

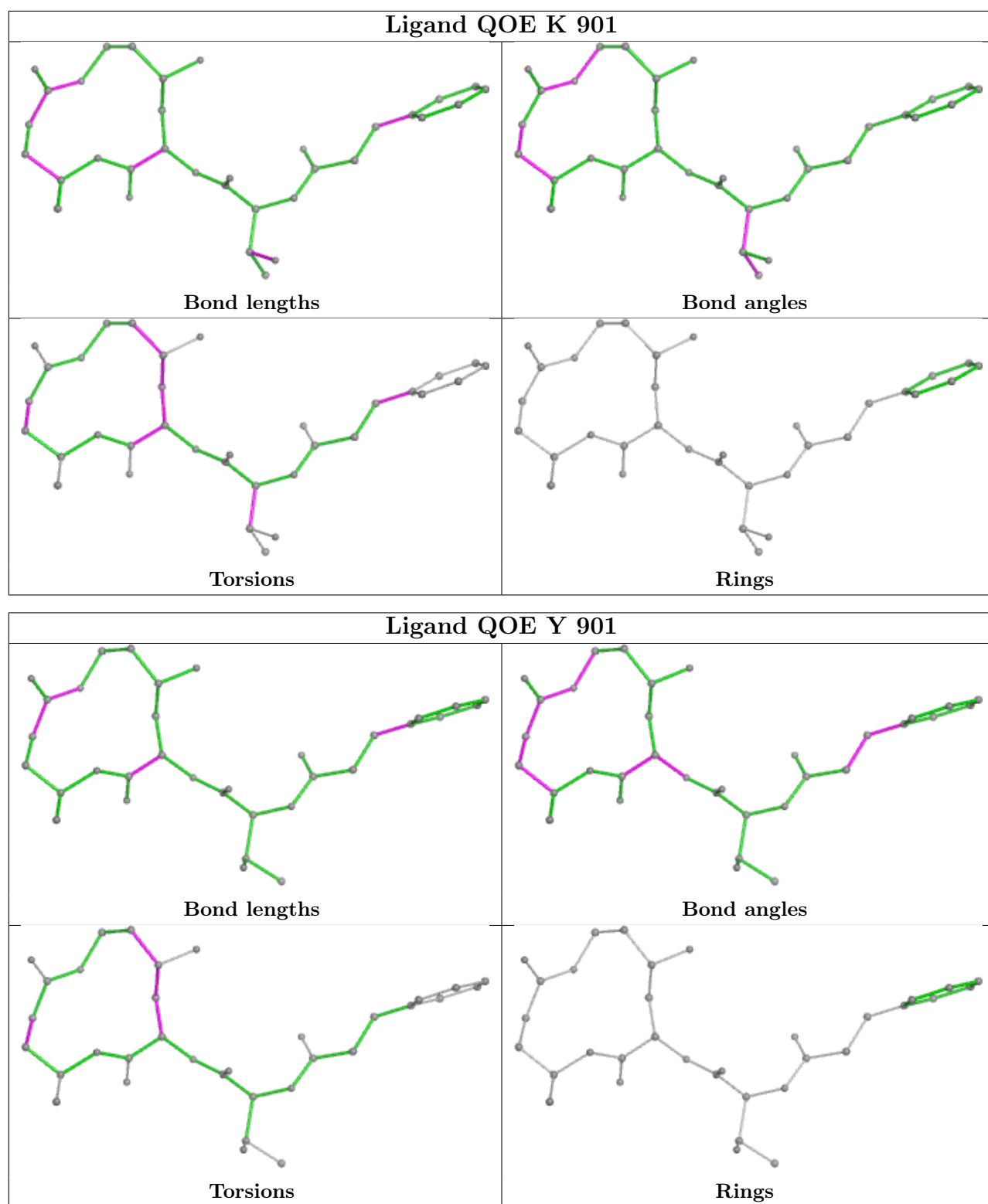
There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
17	H	901	QOE	2	0
17	V	901	QOE	3	0
17	Y	901	QOE	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 ⓘ

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	250/250 (100%)	-0.53	2 (0%) 82 66	44, 58, 93, 135	0
1	O	250/250 (100%)	-0.47	2 (0%) 82 66	46, 64, 107, 135	0
2	B	244/258 (94%)	-0.33	3 (1%) 76 56	43, 62, 107, 163	0
2	P	244/258 (94%)	-0.21	3 (1%) 76 56	49, 67, 109, 159	0
3	C	240/254 (94%)	-0.37	2 (0%) 82 66	43, 67, 127, 148	0
3	Q	240/254 (94%)	-0.01	4 (1%) 69 47	52, 79, 158, 182	0
4	D	235/260 (90%)	-0.32	3 (1%) 74 54	49, 69, 102, 140	0
4	R	235/260 (90%)	-0.19	0 100 100	50, 74, 114, 147	0
5	E	231/234 (98%)	-0.22	0 100 100	52, 72, 111, 152	0
5	S	231/234 (98%)	0.11	2 (0%) 81 63	52, 80, 128, 169	0
6	F	243/288 (84%)	-0.32	3 (1%) 76 56	42, 63, 114, 140	0
6	T	243/288 (84%)	-0.04	3 (1%) 76 56	42, 71, 128, 159	0
7	G	241/252 (95%)	-0.49	0 100 100	38, 58, 92, 138	0
7	U	241/252 (95%)	-0.39	0 100 100	44, 60, 95, 139	0
8	H	226/232 (97%)	-0.44	1 (0%) 89 77	39, 55, 94, 159	0
8	V	226/232 (97%)	-0.42	2 (0%) 81 63	44, 57, 94, 176	0
9	I	204/205 (99%)	-0.63	0 100 100	40, 56, 86, 111	0
9	W	204/205 (99%)	-0.62	1 (0%) 87 75	42, 56, 90, 113	0
10	J	195/198 (98%)	-0.52	1 (0%) 87 75	40, 57, 81, 133	0
10	X	195/198 (98%)	-0.51	1 (0%) 87 75	42, 59, 83, 148	0
11	K	212/212 (100%)	-0.35	1 (0%) 87 75	42, 59, 87, 108	0
11	Y	212/212 (100%)	-0.41	2 (0%) 81 63	46, 59, 89, 108	0
12	L	222/222 (100%)	-0.40	2 (0%) 81 63	42, 58, 96, 127	0
12	Z	222/222 (100%)	-0.48	1 (0%) 87 75	42, 57, 95, 126	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	229/246 (93%)	-0.53	0 100 100	38, 57, 82, 98	0
13	a	232/246 (94%)	-0.58	1 (0%) 89 77	38, 56, 81, 93	0
14	N	196/196 (100%)	-0.60	0 100 100	39, 51, 82, 105	0
14	b	196/196 (100%)	-0.59	1 (0%) 87 75	39, 52, 81, 117	0
All	All	6339/6614 (95%)	-0.38	41 (0%) 85 71	38, 62, 108, 182	0

The worst 5 of 41 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	Q	50	LEU	5.4
10	X	1	MET	5.1
8	V	223	ILE	5.0
8	H	223	ILE	4.3
3	Q	49	THR	4.2

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, 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.

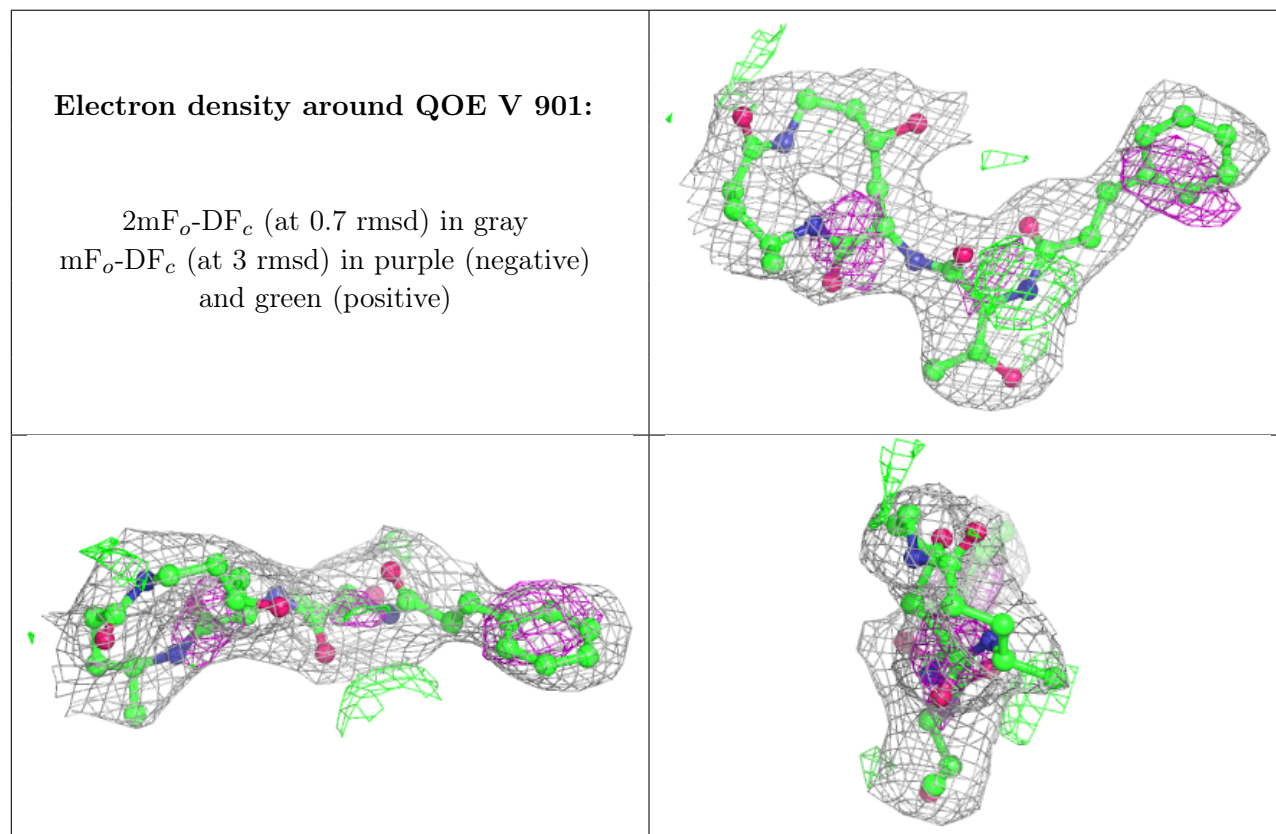
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
15	MG	K	902	1/1	0.82	0.13	68,68,68,68	0
15	MG	Y	902	1/1	0.88	0.09	68,68,68,68	0
15	MG	I	301	1/1	0.89	0.24	77,77,77,77	0
17	QOE	V	901	34/34	0.89	0.11	41,50,60,63	0
17	QOE	K	901	34/34	0.90	0.10	49,54,59,60	0
17	QOE	H	901	34/34	0.90	0.10	42,51,60,61	0
15	MG	V	902	1/1	0.91	0.12	78,78,78,78	0

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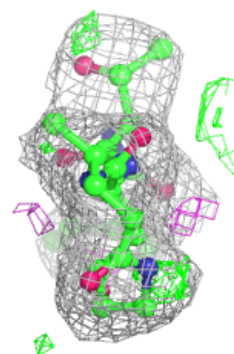
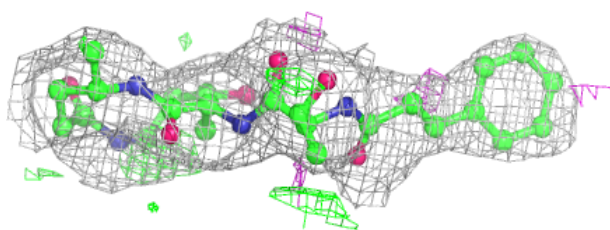
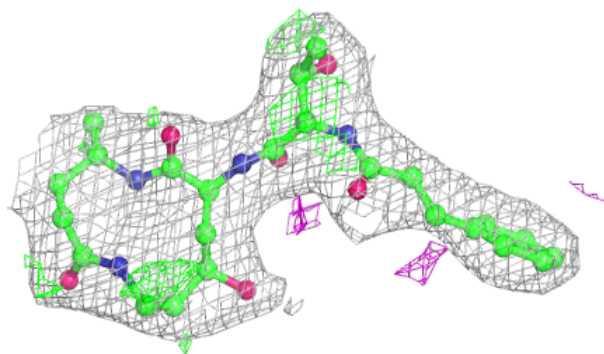
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
15	MG	W	301	1/1	0.92	0.19	70,70,70,70	0
17	QOE	Y	901	34/34	0.92	0.09	48,54,58,63	0
15	MG	Z	301	1/1	0.94	0.16	56,56,56,56	0
16	CL	U	301	1/1	0.94	0.14	54,54,54,54	0
15	MG	N	201	1/1	0.98	0.06	36,36,36,36	0
15	MG	G	301	1/1	0.98	0.05	48,48,48,48	0
16	CL	G	302	1/1	0.99	0.07	42,42,42,42	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.

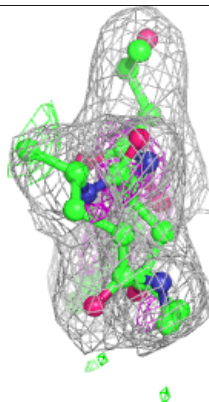
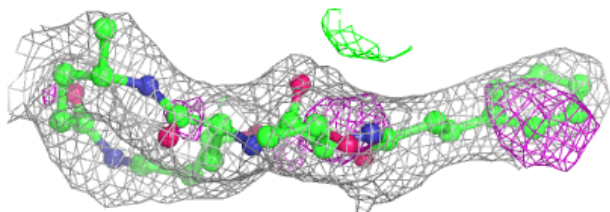
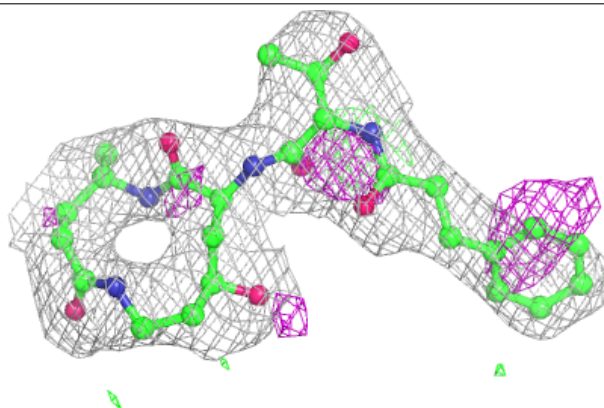


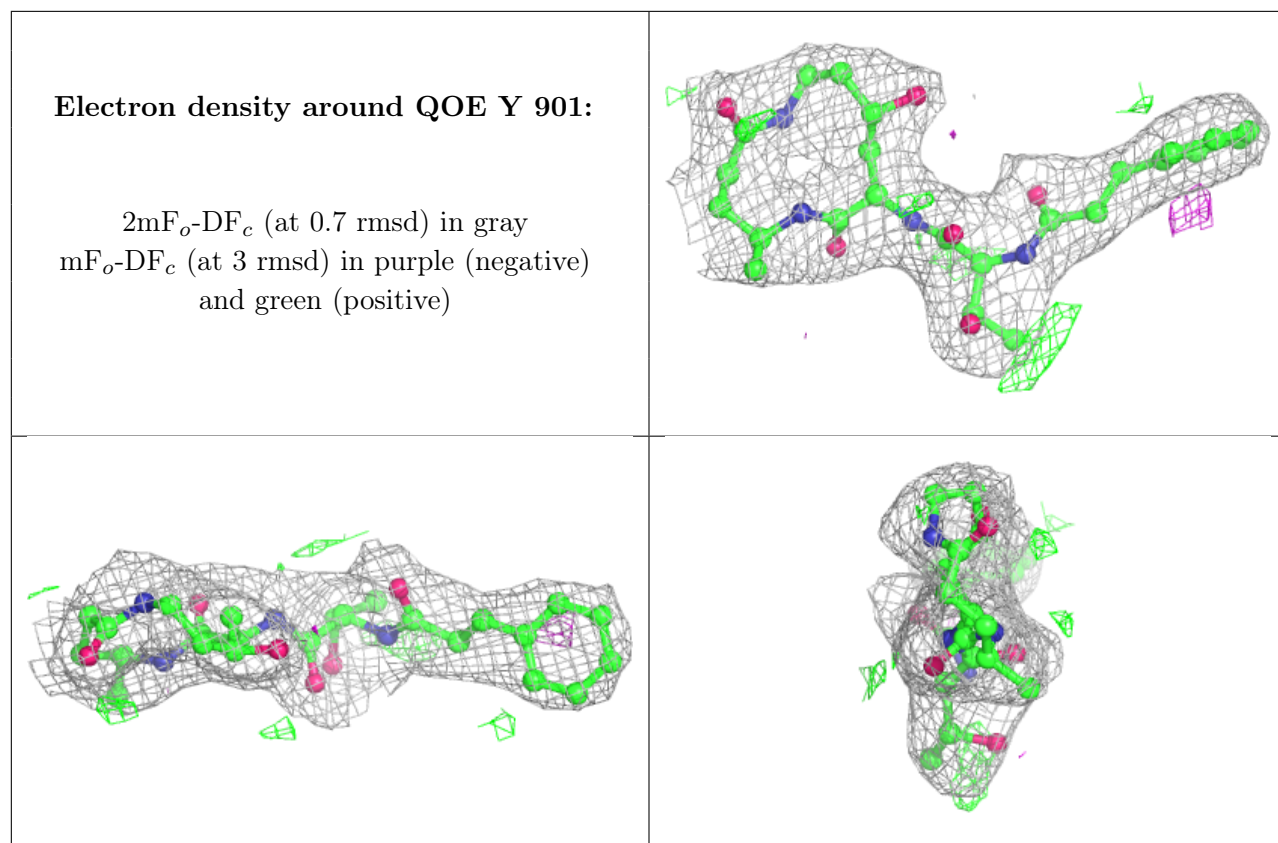
Electron density around QOE K 901:

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

**Electron density around QOE H 901:**

$2mF_o-DF_c$ (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.