



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 20, 2020 – 06:31 PM BST

PDB ID : 3KRD
Title : Crystal Structure of Mycobacterium Tuberculosis Proteasome in complex with
Fellutamide B
Authors : Li, D.; Li, H.
Deposited on : 2009-11-18
Resolution : 2.50 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 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	:	2.13.1
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)	:	2.13.1

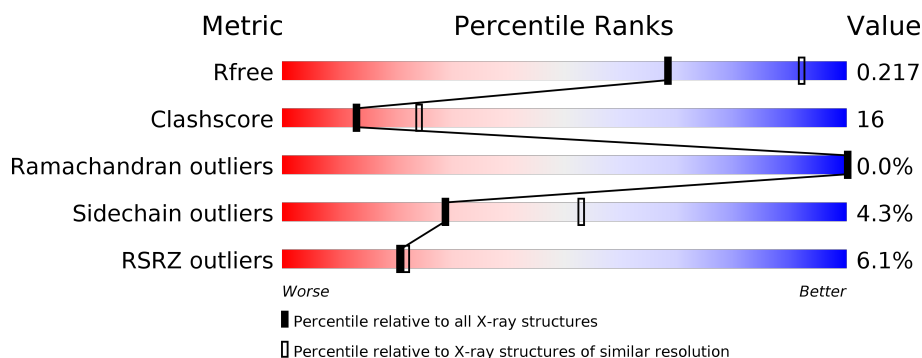
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	1	248	<div> <div>7%</div> <div>52%</div> <div>31%</div> <div>•</div> <div>13%</div> </div>
1	A	248	<div> <div>12%</div> <div>54%</div> <div>31%</div> <div>•</div> <div>14%</div> </div>
1	B	248	<div> <div>6%</div> <div>58%</div> <div>27%</div> <div>•</div> <div>14%</div> </div>
1	D	248	<div> <div>20%</div> <div>45%</div> <div>37%</div> <div>•</div> <div>14%</div> </div>
1	F	248	<div> <div>5%</div> <div>54%</div> <div>31%</div> <div>•</div> <div>13%</div> </div>
1	I	248	<div> <div>8%</div> <div>57%</div> <div>28%</div> <div>•</div> <div>13%</div> </div>

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Mol	Chain	Length	Quality of chain
1	K	248	
1	M	248	
1	O	248	
1	Q	248	
1	S	248	
1	U	248	
1	W	248	
1	Y	248	
2	2	240	
2	C	240	
2	E	240	
2	G	240	
2	H	240	
2	J	240	
2	L	240	
2	N	240	
2	P	240	
2	R	240	
2	T	240	
2	V	240	
2	X	240	
2	Z	240	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 48828 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	0	0	0
			1650	1033	302	312	3			
1	B	213	Total	C	N	O	S	0	0	0
			1642	1027	301	311	3			
1	D	213	Total	C	N	O	S	0	0	0
			1646	1031	301	311	3			
1	F	216	Total	C	N	O	S	0	0	0
			1661	1040	304	314	3			
1	I	215	Total	C	N	O	S	0	0	0
			1654	1035	303	313	3			
1	K	215	Total	C	N	O	S	0	0	0
			1654	1035	303	313	3			
1	M	214	Total	C	N	O	S	0	0	0
			1650	1033	302	312	3			
1	O	214	Total	C	N	O	S	0	0	0
			1650	1033	302	312	3			
1	Q	214	Total	C	N	O	S	0	0	0
			1650	1033	302	312	3			
1	S	215	Total	C	N	O	S	0	0	0
			1657	1038	303	313	3			
1	U	214	Total	C	N	O	S	0	0	0
			1650	1033	302	312	3			
1	W	215	Total	C	N	O	S	0	0	0
			1654	1035	303	313	3			
1	Y	216	Total	C	N	O	S	0	0	0
			1661	1040	304	314	3			
1	1	215	Total	C	N	O	S	0	0	0
			1654	1035	303	313	3			

- Molecule 2 is a protein called Proteasome subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	222	Total 1638	C 1027	N 282	O 324	S 5	0	0	0
2	E	222	Total 1638	C 1027	N 282	O 324	S 5	0	0	0
2	G	222	Total 1638	C 1027	N 282	O 324	S 5	0	0	0
2	H	222	Total 1638	C 1027	N 282	O 324	S 5	0	0	0
2	J	222	Total 1638	C 1027	N 282	O 324	S 5	0	0	0
2	L	222	Total 1638	C 1027	N 282	O 324	S 5	0	0	0
2	N	222	Total 1638	C 1027	N 282	O 324	S 5	0	0	0
2	P	222	Total 1638	C 1027	N 282	O 324	S 5	0	0	0
2	R	229	Total 1683	C 1054	N 289	O 335	S 5	0	0	0
2	T	222	Total 1638	C 1027	N 282	O 324	S 5	0	0	0
2	V	229	Total 1683	C 1054	N 289	O 335	S 5	0	0	0
2	X	222	Total 1638	C 1027	N 282	O 324	S 5	0	0	0
2	Z	222	Total 1638	C 1027	N 282	O 324	S 5	0	0	0
2	2	222	Total 1638	C 1027	N 282	O 324	S 5	0	0	0

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	535	HIS	-	EXPRESSION TAG	UNP O33245
C	536	HIS	-	EXPRESSION TAG	UNP O33245
C	537	HIS	-	EXPRESSION TAG	UNP O33245
C	538	HIS	-	EXPRESSION TAG	UNP O33245
C	539	HIS	-	EXPRESSION TAG	UNP O33245
C	540	HIS	-	EXPRESSION TAG	UNP O33245
E	535	HIS	-	EXPRESSION TAG	UNP O33245
E	536	HIS	-	EXPRESSION TAG	UNP O33245
E	537	HIS	-	EXPRESSION TAG	UNP O33245
E	538	HIS	-	EXPRESSION TAG	UNP O33245
E	539	HIS	-	EXPRESSION TAG	UNP O33245
E	540	HIS	-	EXPRESSION TAG	UNP O33245

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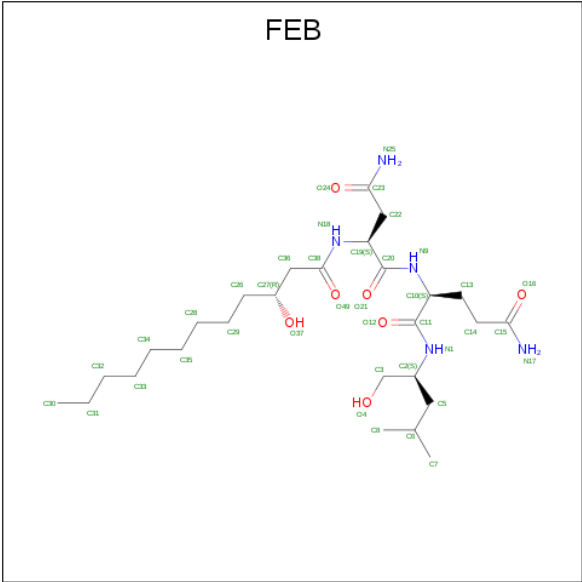
Chain	Residue	Modelled	Actual	Comment	Reference
G	535	HIS	-	EXPRESSION TAG	UNP O33245
G	536	HIS	-	EXPRESSION TAG	UNP O33245
G	537	HIS	-	EXPRESSION TAG	UNP O33245
G	538	HIS	-	EXPRESSION TAG	UNP O33245
G	539	HIS	-	EXPRESSION TAG	UNP O33245
G	540	HIS	-	EXPRESSION TAG	UNP O33245
H	535	HIS	-	EXPRESSION TAG	UNP O33245
H	536	HIS	-	EXPRESSION TAG	UNP O33245
H	537	HIS	-	EXPRESSION TAG	UNP O33245
H	538	HIS	-	EXPRESSION TAG	UNP O33245
H	539	HIS	-	EXPRESSION TAG	UNP O33245
H	540	HIS	-	EXPRESSION TAG	UNP O33245
J	535	HIS	-	EXPRESSION TAG	UNP O33245
J	536	HIS	-	EXPRESSION TAG	UNP O33245
J	537	HIS	-	EXPRESSION TAG	UNP O33245
J	538	HIS	-	EXPRESSION TAG	UNP O33245
J	539	HIS	-	EXPRESSION TAG	UNP O33245
J	540	HIS	-	EXPRESSION TAG	UNP O33245
L	535	HIS	-	EXPRESSION TAG	UNP O33245
L	536	HIS	-	EXPRESSION TAG	UNP O33245
L	537	HIS	-	EXPRESSION TAG	UNP O33245
L	538	HIS	-	EXPRESSION TAG	UNP O33245
L	539	HIS	-	EXPRESSION TAG	UNP O33245
L	540	HIS	-	EXPRESSION TAG	UNP O33245
N	535	HIS	-	EXPRESSION TAG	UNP O33245
N	536	HIS	-	EXPRESSION TAG	UNP O33245
N	537	HIS	-	EXPRESSION TAG	UNP O33245
N	538	HIS	-	EXPRESSION TAG	UNP O33245
N	539	HIS	-	EXPRESSION TAG	UNP O33245
N	540	HIS	-	EXPRESSION TAG	UNP O33245
P	535	HIS	-	EXPRESSION TAG	UNP O33245
P	536	HIS	-	EXPRESSION TAG	UNP O33245
P	537	HIS	-	EXPRESSION TAG	UNP O33245
P	538	HIS	-	EXPRESSION TAG	UNP O33245
P	539	HIS	-	EXPRESSION TAG	UNP O33245
P	540	HIS	-	EXPRESSION TAG	UNP O33245
R	535	HIS	-	EXPRESSION TAG	UNP O33245
R	536	HIS	-	EXPRESSION TAG	UNP O33245
R	537	HIS	-	EXPRESSION TAG	UNP O33245
R	538	HIS	-	EXPRESSION TAG	UNP O33245
R	539	HIS	-	EXPRESSION TAG	UNP O33245
R	540	HIS	-	EXPRESSION TAG	UNP O33245

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Chain	Residue	Modelled	Actual	Comment	Reference
T	535	HIS	-	EXPRESSION TAG	UNP O33245
T	536	HIS	-	EXPRESSION TAG	UNP O33245
T	537	HIS	-	EXPRESSION TAG	UNP O33245
T	538	HIS	-	EXPRESSION TAG	UNP O33245
T	539	HIS	-	EXPRESSION TAG	UNP O33245
T	540	HIS	-	EXPRESSION TAG	UNP O33245
V	535	HIS	-	EXPRESSION TAG	UNP O33245
V	536	HIS	-	EXPRESSION TAG	UNP O33245
V	537	HIS	-	EXPRESSION TAG	UNP O33245
V	538	HIS	-	EXPRESSION TAG	UNP O33245
V	539	HIS	-	EXPRESSION TAG	UNP O33245
V	540	HIS	-	EXPRESSION TAG	UNP O33245
X	535	HIS	-	EXPRESSION TAG	UNP O33245
X	536	HIS	-	EXPRESSION TAG	UNP O33245
X	537	HIS	-	EXPRESSION TAG	UNP O33245
X	538	HIS	-	EXPRESSION TAG	UNP O33245
X	539	HIS	-	EXPRESSION TAG	UNP O33245
X	540	HIS	-	EXPRESSION TAG	UNP O33245
Z	535	HIS	-	EXPRESSION TAG	UNP O33245
Z	536	HIS	-	EXPRESSION TAG	UNP O33245
Z	537	HIS	-	EXPRESSION TAG	UNP O33245
Z	538	HIS	-	EXPRESSION TAG	UNP O33245
Z	539	HIS	-	EXPRESSION TAG	UNP O33245
Z	540	HIS	-	EXPRESSION TAG	UNP O33245
2	535	HIS	-	EXPRESSION TAG	UNP O33245
2	536	HIS	-	EXPRESSION TAG	UNP O33245
2	537	HIS	-	EXPRESSION TAG	UNP O33245
2	538	HIS	-	EXPRESSION TAG	UNP O33245
2	539	HIS	-	EXPRESSION TAG	UNP O33245
2	540	HIS	-	EXPRESSION TAG	UNP O33245

- Molecule 3 is N 2 -[(3R)-3-hydroxydodecanoyl]-L-asparaginyl-N 1 -[(1S)-1-(hydroxymethyl)-3-methylbutyl]-L-glutamamide (three-letter code: FEB) (formula: C₂₇H₅₁N₅O₇).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	N	O	0	0
			32	20	5	7		
3	E	1	Total	C	N	O	0	0
			32	20	5	7		
3	G	1	Total	C	N	O	0	0
			32	20	5	7		
3	H	1	Total	C	N	O	0	0
			32	20	5	7		
3	J	1	Total	C	N	O	0	0
			32	20	5	7		
3	L	1	Total	C	N	O	0	0
			32	20	5	7		
3	N	1	Total	C	N	O	0	0
			32	20	5	7		
3	P	1	Total	C	N	O	0	0
			32	20	5	7		
3	R	1	Total	C	N	O	0	0
			32	20	5	7		
3	T	1	Total	C	N	O	0	0
			32	20	5	7		
3	V	1	Total	C	N	O	0	0
			32	20	5	7		
3	X	1	Total	C	N	O	0	0
			32	20	5	7		
3	Z	1	Total	C	N	O	0	0
			32	20	5	7		
3	2	1	Total	C	N	O	0	0
			32	20	5	7		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	32	Total O 32 32	0	0
4	B	42	Total O 42 42	0	0
4	C	131	Total O 131 131	0	0
4	D	25	Total O 25 25	0	0
4	E	130	Total O 130 130	0	0
4	F	41	Total O 41 41	0	0
4	G	113	Total O 113 113	0	0
4	H	126	Total O 126 126	0	0
4	I	38	Total O 38 38	0	0
4	J	122	Total O 122 122	0	0
4	K	33	Total O 33 33	0	0
4	L	132	Total O 132 132	0	0
4	M	42	Total O 42 42	0	0
4	N	121	Total O 121 121	0	0
4	O	39	Total O 39 39	0	0
4	P	111	Total O 111 111	0	0
4	Q	34	Total O 34 34	0	0
4	R	143	Total O 143 143	0	0
4	S	43	Total O 43 43	0	0
4	T	103	Total O 103 103	0	0
4	U	33	Total O 33 33	0	0

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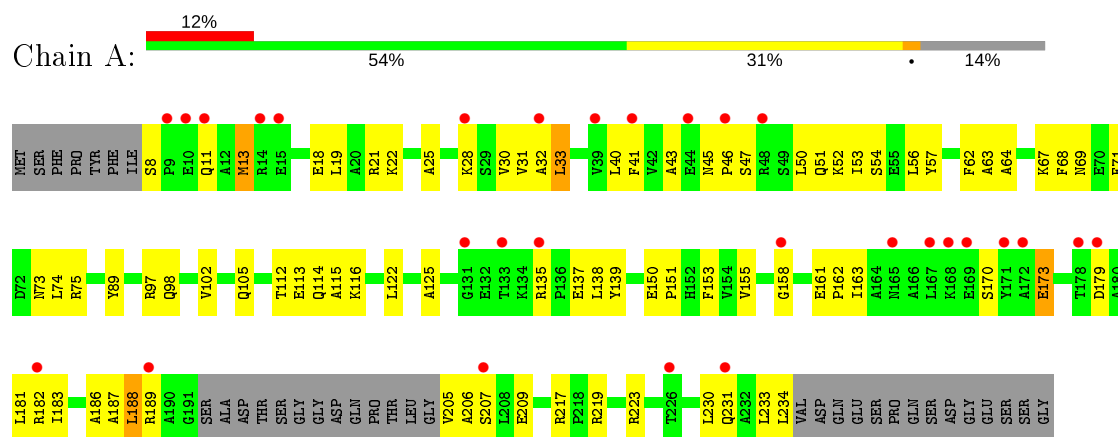
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	V	149	Total 149	O 149	0	0
4	W	26	Total 26	O 26	0	0
4	X	115	Total 115	O 115	0	0
4	Y	22	Total 22	O 22	0	0
4	Z	106	Total 106	O 106	0	0
4	1	55	Total 55	O 55	0	0
4	2	118	Total 118	O 118	0	0

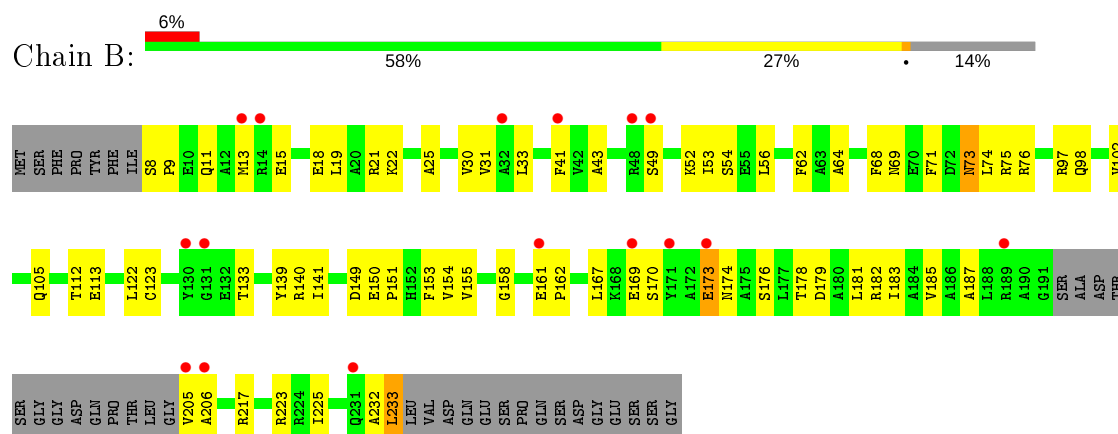
3 Residue-property plots [i](#)

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.

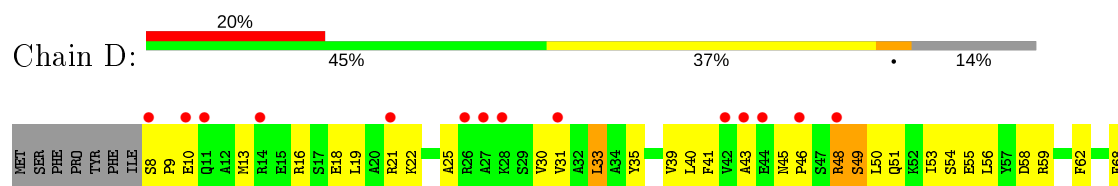
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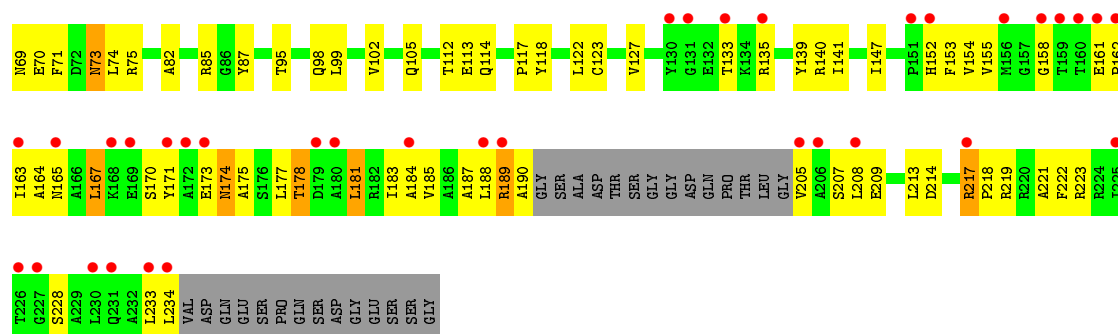


• Molecule 1: Proteasome subunit alpha

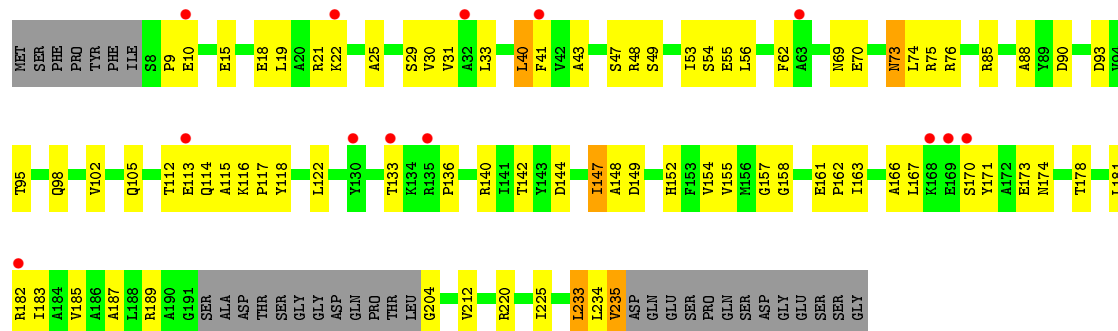


• Molecule 1: Proteasome subunit alpha

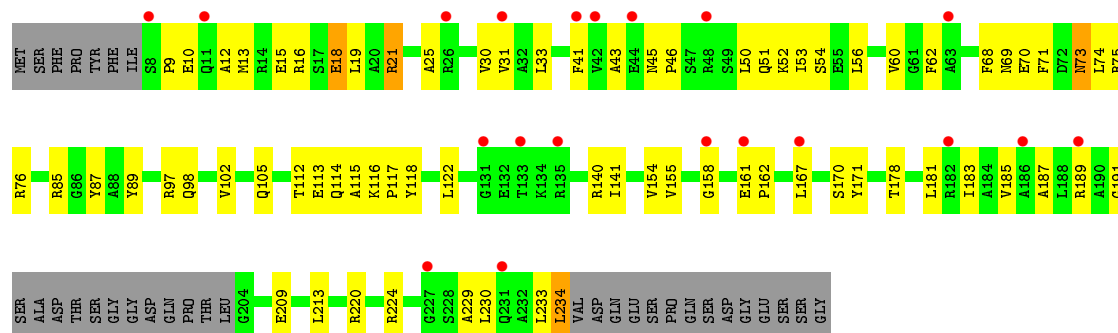




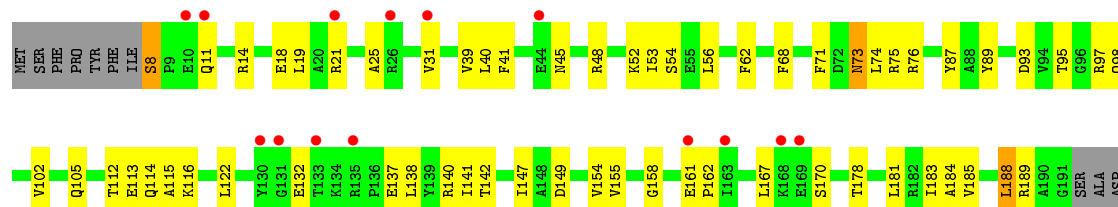
• Molecule 1: Proteasome subunit alpha



• Molecule 1: Proteasome subunit alpha

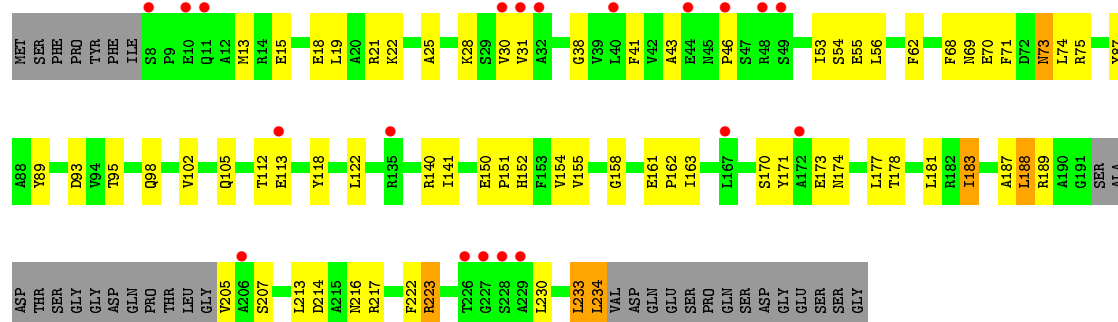


• Molecule 1: Proteasome subunit alpha

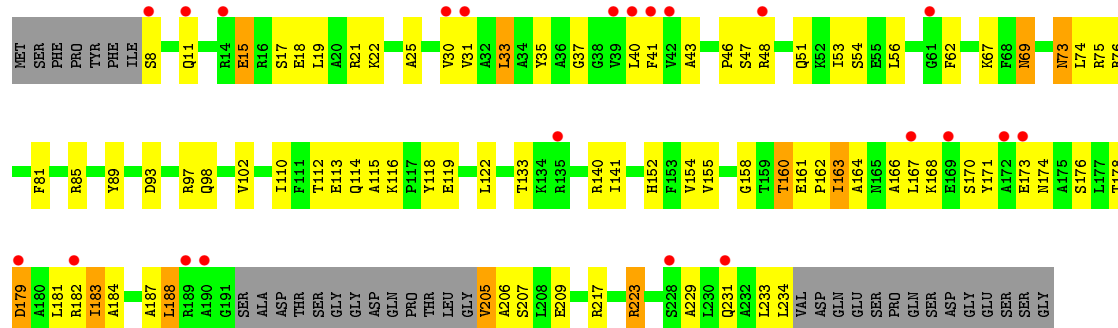




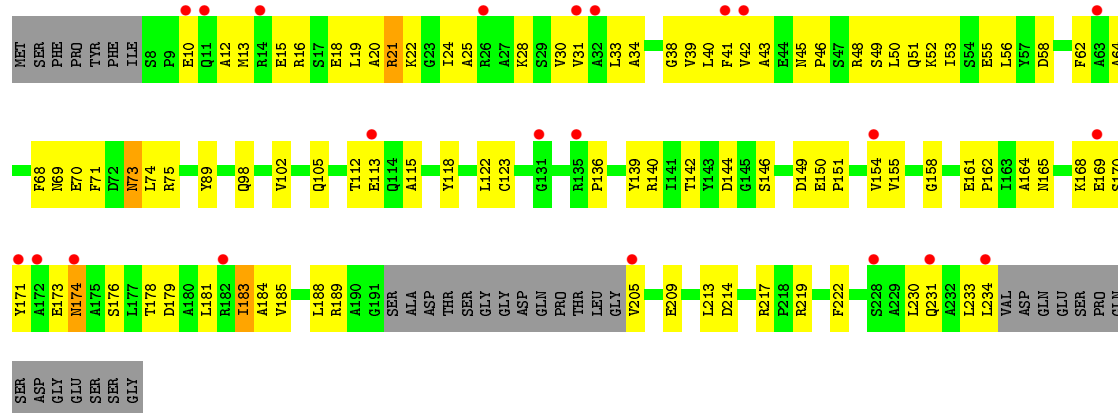
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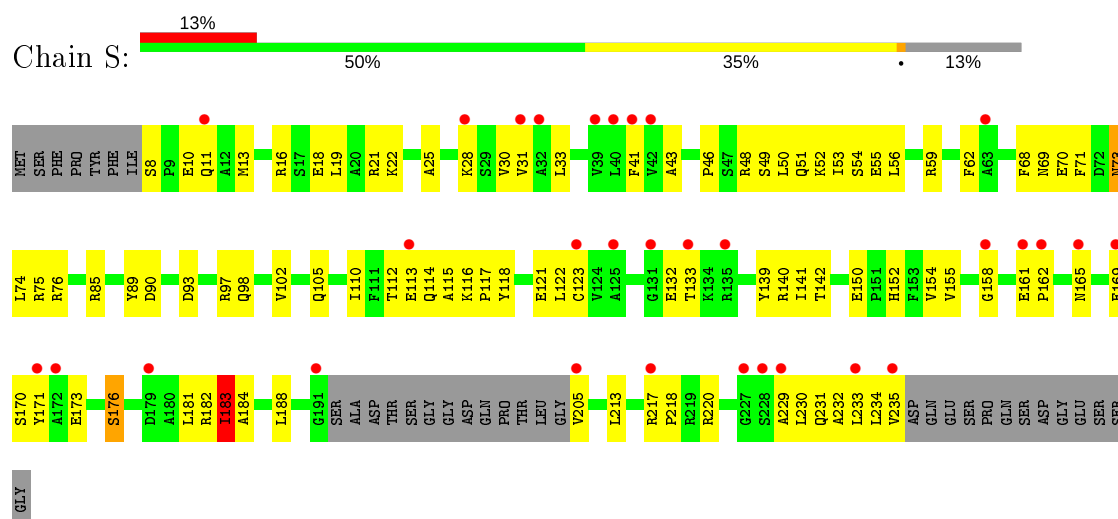
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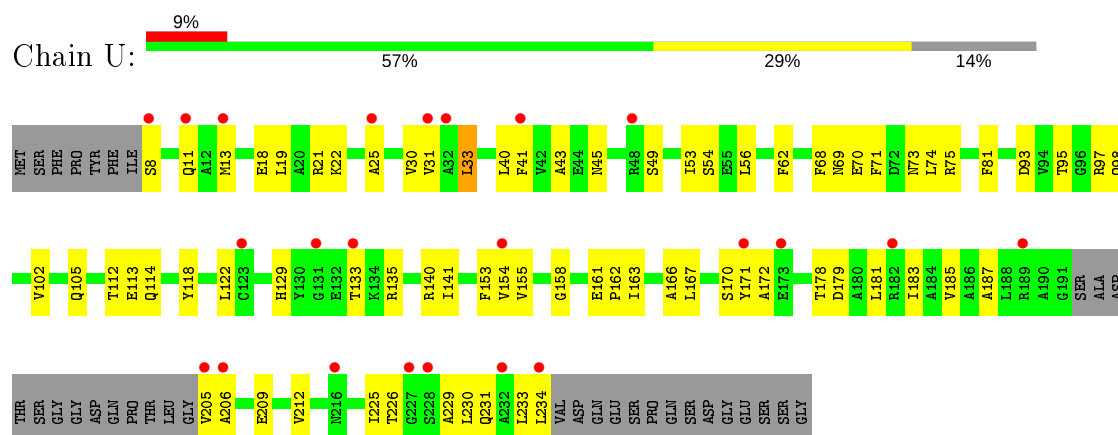
• Molecule 1: Proteasome subunit alpha



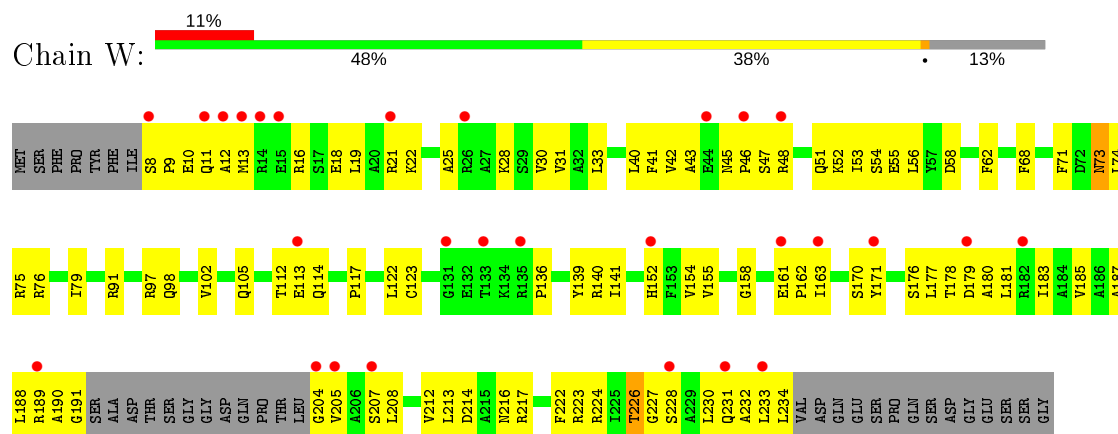
• Molecule 1: Proteasome subunit alpha



• Molecule 1: Proteasome subunit alpha

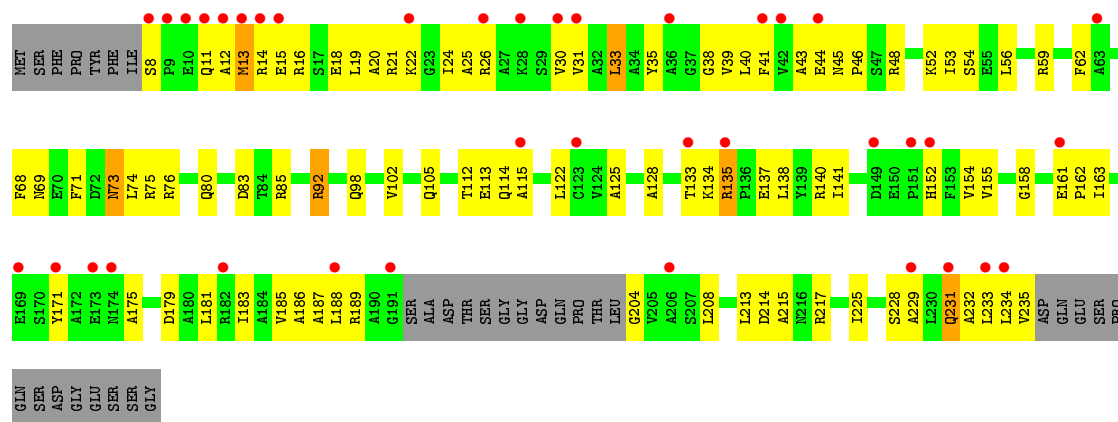


• Molecule 1: Proteasome subunit alpha

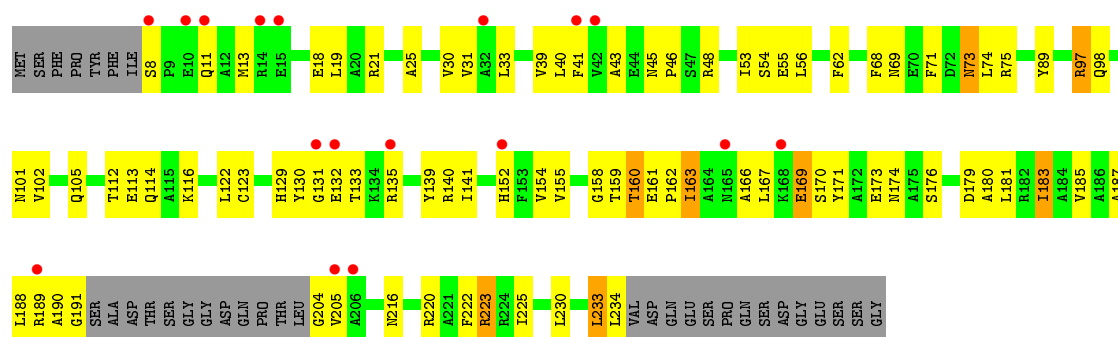


• Molecule 1: Proteasome subunit alpha

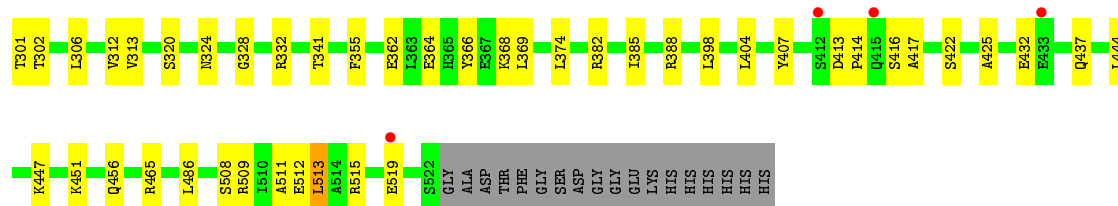
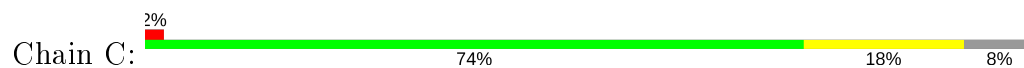




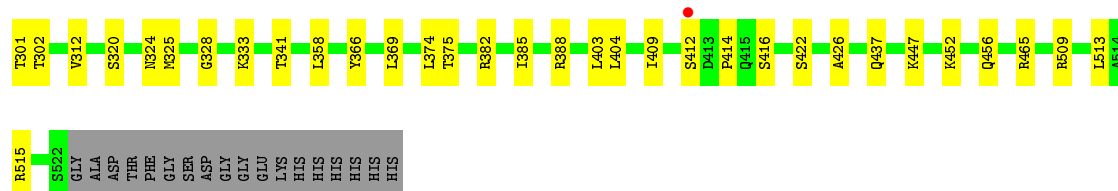
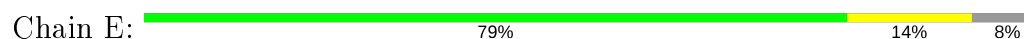
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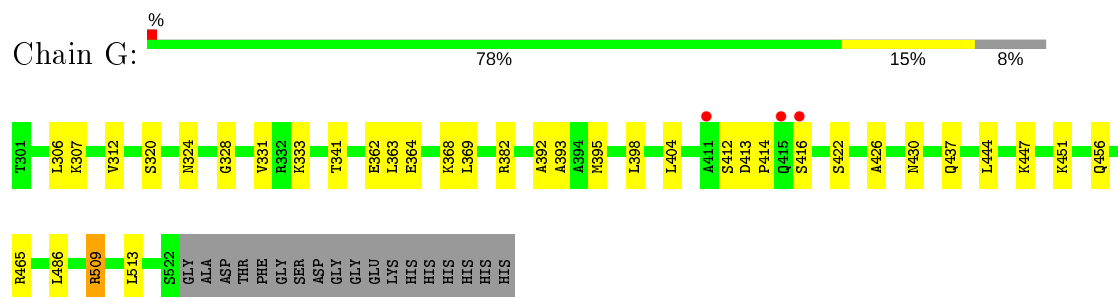
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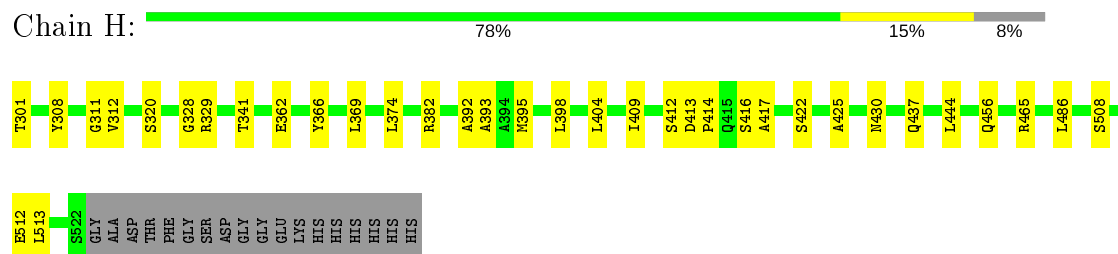
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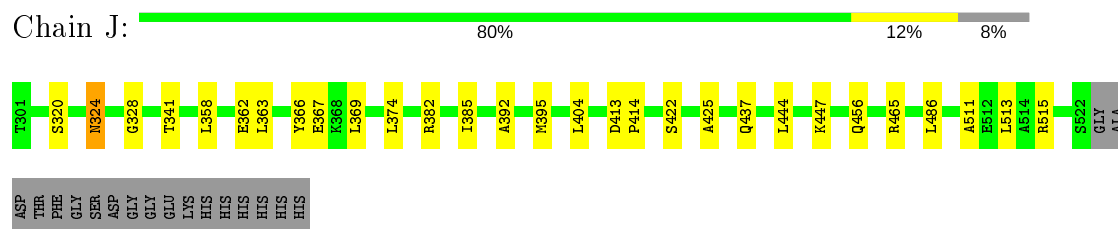
• Molecule 2: Proteasome subunit beta



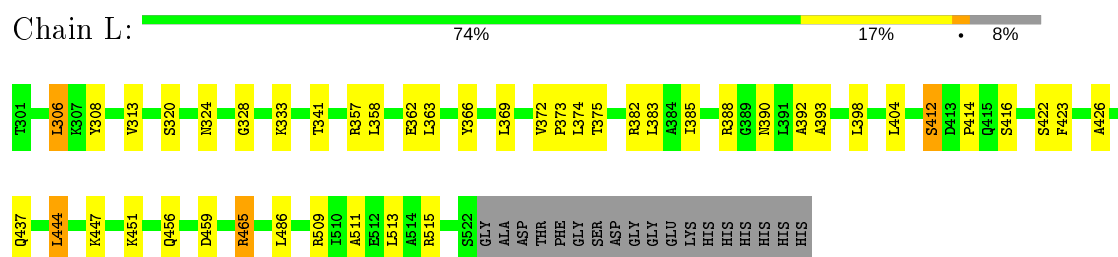
- Molecule 2: Proteasome subunit beta

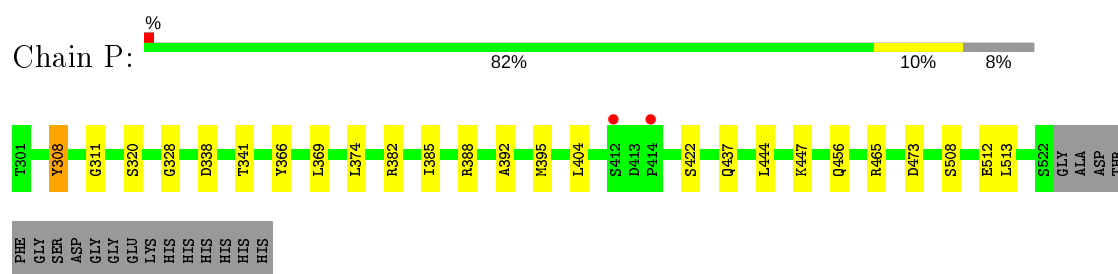


- Molecule 2: Proteasome subunit beta

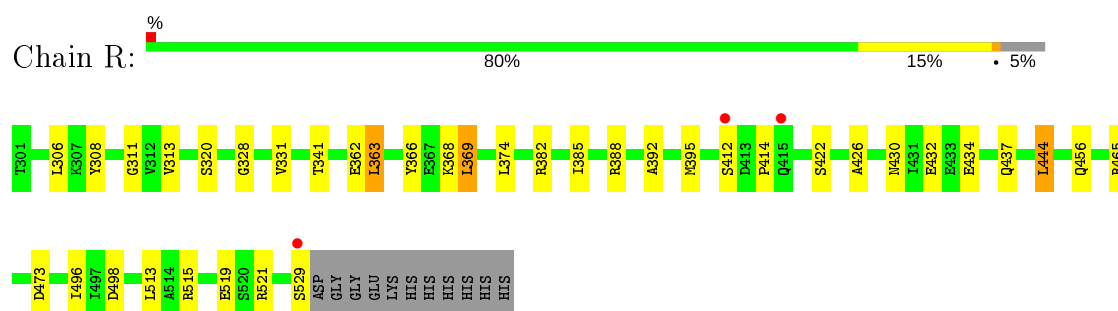


- Molecule 2: Proteasome subunit beta

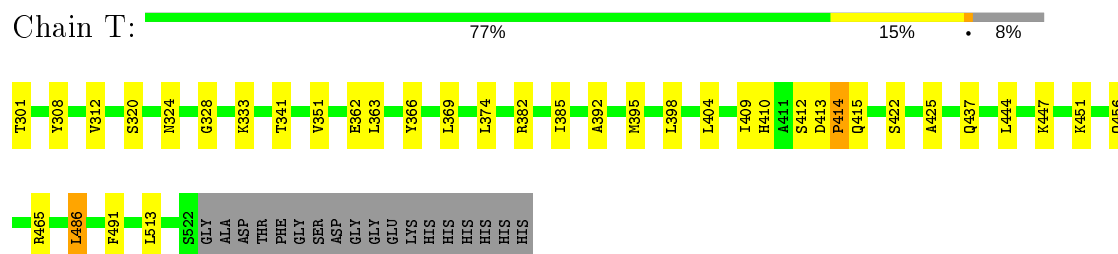




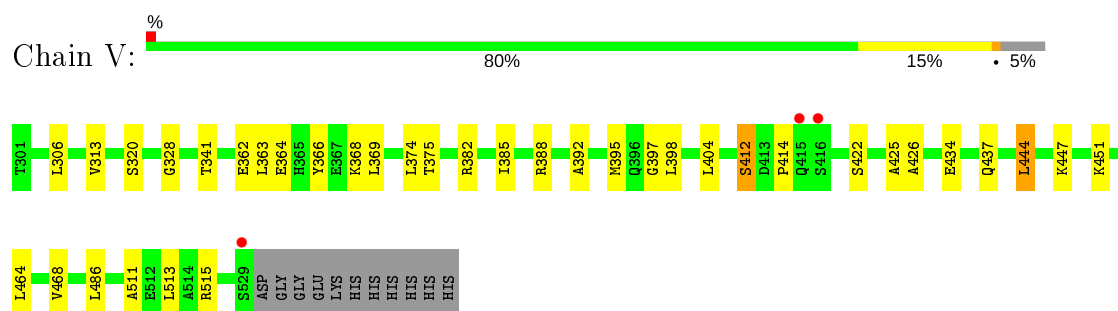
- Molecule 2: Proteasome subunit beta



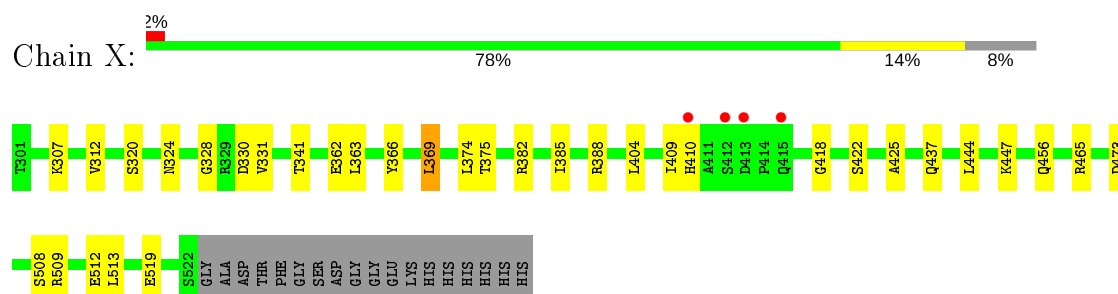
- Molecule 2: Proteasome subunit beta



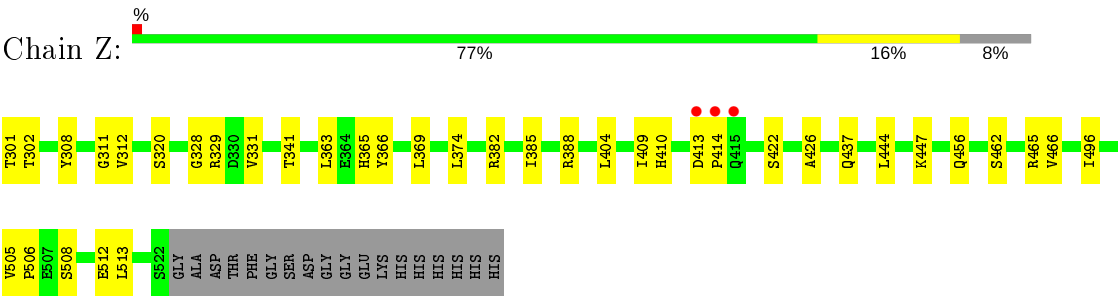
- Molecule 2: Proteasome subunit beta



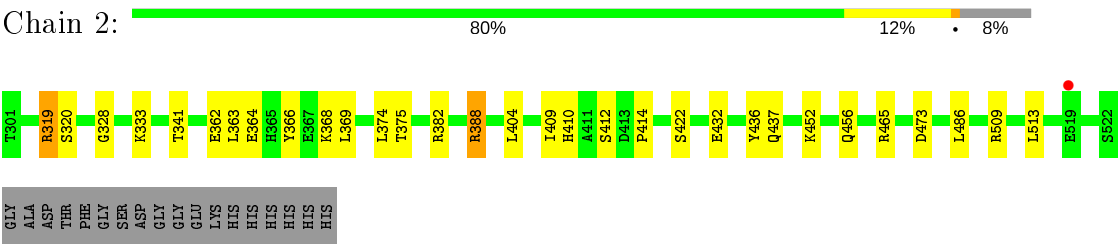
- Molecule 2: Proteasome subunit beta



● Molecule 2: Proteasome subunit beta



● Molecule 2: Proteasome subunit beta



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	170.19Å 118.10Å 194.35Å 90.00° 112.62° 90.00°	Depositor
Resolution (Å)	25.00 – 2.50 25.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	96.5 (25.00-2.50) 96.6 (25.00-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 2.50Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.208 , 0.229 0.198 , 0.217	Depositor DCC
R_{free} test set	12081 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	30.7	Xtriage
Anisotropy	0.279	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 53.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	48828	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.82% 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: FEB

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	1	0.81	0/1679	0.76	0/2268
1	A	0.73	1/1675 (0.1%)	0.70	0/2263
1	B	0.62	0/1667	0.69	1/2252 (0.0%)
1	D	0.62	0/1671	0.79	1/2258 (0.0%)
1	F	0.70	0/1686	0.74	0/2278
1	I	0.67	0/1679	0.75	1/2268 (0.0%)
1	K	0.76	0/1679	0.77	0/2268
1	M	0.65	0/1675	0.71	0/2263
1	O	0.70	0/1675	0.76	0/2263
1	Q	0.69	0/1675	0.73	1/2263 (0.0%)
1	S	0.67	0/1682	0.73	2/2273 (0.1%)
1	U	0.66	0/1675	0.71	0/2263
1	W	0.56	0/1679	0.74	1/2268 (0.0%)
1	Y	0.73	0/1686	0.75	0/2278
2	2	0.77	0/1662	0.77	1/2254 (0.0%)
2	C	0.70	0/1662	0.78	2/2254 (0.1%)
2	E	0.76	0/1662	0.79	0/2254
2	G	0.70	0/1662	0.75	0/2254
2	H	0.73	0/1662	0.76	0/2254
2	J	0.75	0/1662	0.78	0/2254
2	L	0.74	1/1662 (0.1%)	0.77	1/2254 (0.0%)
2	N	0.66	0/1662	0.74	0/2254
2	P	0.76	1/1662 (0.1%)	0.76	2/2254 (0.1%)
2	R	0.82	0/1708	0.79	1/2316 (0.0%)
2	T	0.76	0/1662	0.82	0/2254
2	V	0.82	0/1708	0.75	0/2316
2	X	0.70	0/1662	0.78	1/2254 (0.0%)
2	Z	0.63	0/1662	0.76	1/2254 (0.0%)
All	All	0.71	3/46843 (0.0%)	0.75	16/63406 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	308	TYR	CD1-CE1	-5.57	1.30	1.39
2	P	308	TYR	CD1-CE1	-5.11	1.31	1.39
1	A	57	TYR	CD1-CE1	-5.10	1.31	1.39

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	S	183	ILE	CB-CA-C	-5.96	99.67	111.60
1	Q	183	ILE	CB-CA-C	-5.70	100.20	111.60
2	C	413	ASP	C-N-CD	-5.67	108.12	120.60
2	Z	388	ARG	NE-CZ-NH2	-5.42	117.59	120.30
2	2	388	ARG	NE-CZ-NH2	-5.37	117.61	120.30

There are no chirality outliers.

There are no planarity outliers.

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	1	1654	0	1651	93	0
1	A	1650	0	1648	74	0
1	B	1642	0	1637	64	0
1	D	1646	0	1645	144	0
1	F	1661	0	1660	68	0
1	I	1654	0	1651	62	0
1	K	1654	0	1651	82	0
1	M	1650	0	1648	69	0
1	O	1650	0	1648	96	0
1	Q	1650	0	1648	89	0
1	S	1657	0	1657	98	0
1	U	1650	0	1648	64	0
1	W	1654	0	1651	90	0
1	Y	1661	0	1660	120	0
2	2	1638	0	1629	26	0
2	C	1638	0	1629	36	0
2	E	1638	0	1629	30	0
2	G	1638	0	1629	34	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	H	1638	0	1629	22	0
2	J	1638	0	1629	30	0
2	L	1638	0	1629	49	0
2	N	1638	0	1629	24	0
2	P	1638	0	1629	18	0
2	R	1683	0	1665	37	0
2	T	1638	0	1629	37	0
2	V	1683	0	1665	31	0
2	X	1638	0	1629	27	0
2	Z	1638	0	1629	24	0
3	2	32	0	33	2	0
3	C	32	0	33	3	0
3	E	32	0	33	2	0
3	G	32	0	33	3	0
3	H	32	0	33	1	0
3	J	32	0	33	1	0
3	L	32	0	33	1	0
3	N	32	0	33	3	0
3	P	32	0	33	2	0
3	R	32	0	33	3	0
3	T	32	0	33	3	0
3	V	32	0	33	0	0
3	X	32	0	33	2	0
3	Z	32	0	33	2	0
4	1	55	0	0	2	0
4	2	118	0	0	6	0
4	A	32	0	0	3	0
4	B	42	0	0	3	0
4	C	131	0	0	11	0
4	D	25	0	0	3	0
4	E	130	0	0	7	0
4	F	41	0	0	6	0
4	G	113	0	0	8	0
4	H	126	0	0	6	0
4	I	38	0	0	6	0
4	J	122	0	0	3	0
4	K	33	0	0	6	0
4	L	132	0	0	9	0
4	M	42	0	0	0	0
4	N	121	0	0	3	0
4	O	39	0	0	6	0
4	P	111	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	Q	34	0	0	0	0
4	R	143	0	0	9	0
4	S	43	0	0	19	0
4	T	103	0	0	3	0
4	U	33	0	0	2	0
4	V	149	0	0	5	0
4	W	26	0	0	3	0
4	X	115	0	0	4	0
4	Y	22	0	0	3	0
4	Z	106	0	0	1	0
All	All	48828	0	46443	1514	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 16.

The worst 5 of 1514 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:D:217:ARG:NH1	1:D:217:ARG:HB2	1.41	1.30
1:A:13:MET:CE	1:O:116:LYS:HD3	1.62	1.27
1:M:217:ARG:HD3	1:M:223:ARG:NH2	1.53	1.24
1:B:170:SER:OG	1:B:183:ILE:HG23	1.39	1.23
1:Y:229:ALA:O	1:Y:233:LEU:HD13	1.41	1.20

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	1	211/248 (85%)	199 (94%)	11 (5%)	1 (0%)	29	48
1	A	210/248 (85%)	203 (97%)	7 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	209/248 (84%)	203 (97%)	6 (3%)	0	100	100
1	D	209/248 (84%)	203 (97%)	6 (3%)	0	100	100
1	F	212/248 (86%)	208 (98%)	4 (2%)	0	100	100
1	I	211/248 (85%)	208 (99%)	3 (1%)	0	100	100
1	K	211/248 (85%)	204 (97%)	7 (3%)	0	100	100
1	M	210/248 (85%)	204 (97%)	6 (3%)	0	100	100
1	O	210/248 (85%)	203 (97%)	7 (3%)	0	100	100
1	Q	210/248 (85%)	202 (96%)	8 (4%)	0	100	100
1	S	211/248 (85%)	204 (97%)	7 (3%)	0	100	100
1	U	210/248 (85%)	206 (98%)	4 (2%)	0	100	100
1	W	211/248 (85%)	205 (97%)	6 (3%)	0	100	100
1	Y	212/248 (86%)	206 (97%)	6 (3%)	0	100	100
2	2	220/240 (92%)	220 (100%)	0	0	100	100
2	C	220/240 (92%)	220 (100%)	0	0	100	100
2	E	220/240 (92%)	218 (99%)	2 (1%)	0	100	100
2	G	220/240 (92%)	218 (99%)	2 (1%)	0	100	100
2	H	220/240 (92%)	217 (99%)	3 (1%)	0	100	100
2	J	220/240 (92%)	218 (99%)	2 (1%)	0	100	100
2	L	220/240 (92%)	218 (99%)	2 (1%)	0	100	100
2	N	220/240 (92%)	218 (99%)	2 (1%)	0	100	100
2	P	220/240 (92%)	218 (99%)	2 (1%)	0	100	100
2	R	227/240 (95%)	224 (99%)	3 (1%)	0	100	100
2	T	220/240 (92%)	219 (100%)	1 (0%)	0	100	100
2	V	227/240 (95%)	226 (100%)	1 (0%)	0	100	100
2	X	220/240 (92%)	219 (100%)	1 (0%)	0	100	100
2	Z	220/240 (92%)	217 (99%)	3 (1%)	0	100	100
All	All	6041/6832 (88%)	5928 (98%)	112 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1	131	GLY

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	1	164/192 (85%)	151 (92%)	13 (8%)	12	24
1	A	164/192 (85%)	160 (98%)	4 (2%)	49	74
1	B	163/192 (85%)	158 (97%)	5 (3%)	40	67
1	D	164/192 (85%)	152 (93%)	12 (7%)	14	27
1	F	165/192 (86%)	153 (93%)	12 (7%)	14	27
1	I	164/192 (85%)	157 (96%)	7 (4%)	29	53
1	K	164/192 (85%)	156 (95%)	8 (5%)	25	47
1	M	164/192 (85%)	154 (94%)	10 (6%)	18	36
1	O	164/192 (85%)	151 (92%)	13 (8%)	12	24
1	Q	164/192 (85%)	154 (94%)	10 (6%)	18	36
1	S	165/192 (86%)	158 (96%)	7 (4%)	30	54
1	U	164/192 (85%)	159 (97%)	5 (3%)	41	68
1	W	164/192 (85%)	159 (97%)	5 (3%)	41	68
1	Y	165/192 (86%)	156 (94%)	9 (6%)	21	41
2	2	165/178 (93%)	160 (97%)	5 (3%)	41	68
2	C	165/178 (93%)	161 (98%)	4 (2%)	49	74
2	E	165/178 (93%)	161 (98%)	4 (2%)	49	74
2	G	165/178 (93%)	160 (97%)	5 (3%)	41	68
2	H	165/178 (93%)	158 (96%)	7 (4%)	30	54
2	J	165/178 (93%)	160 (97%)	5 (3%)	41	68
2	L	165/178 (93%)	155 (94%)	10 (6%)	18	36
2	N	165/178 (93%)	159 (96%)	6 (4%)	35	61
2	P	165/178 (93%)	163 (99%)	2 (1%)	71	88
2	R	169/178 (95%)	163 (96%)	6 (4%)	35	61
2	T	165/178 (93%)	157 (95%)	8 (5%)	25	48
2	V	169/178 (95%)	163 (96%)	6 (4%)	35	61

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	X	165/178 (93%)	159 (96%)	6 (4%)	35	61
2	Z	165/178 (93%)	159 (96%)	6 (4%)	35	61
All	All	4616/5180 (89%)	4416 (96%)	200 (4%)	29	53

5 of 200 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	M	223	ARG
1	O	223	ARG
1	1	97	ARG
2	N	363	LEU
1	O	48	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 114 such sidechains are listed below:

Mol	Chain	Res	Type
1	O	69	ASN
1	Q	174	ASN
1	1	69	ASN
1	O	105	GLN
2	P	324	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

14 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	$\# Z > 2$
3	FEB	J	300	2	31,31,38	1.32	4 (12%)	39,40,47	1.29	6 (15%)
3	FEB	L	300	2	31,31,38	1.29	6 (19%)	39,40,47	1.22	3 (7%)
3	FEB	N	300	2	31,31,38	1.33	4 (12%)	39,40,47	1.26	5 (12%)
3	FEB	G	300	2	31,31,38	1.27	4 (12%)	39,40,47	1.30	5 (12%)
3	FEB	E	300	2	31,31,38	1.14	2 (6%)	39,40,47	1.15	3 (7%)
3	FEB	V	300	2	31,31,38	1.34	6 (19%)	39,40,47	1.21	4 (10%)
3	FEB	X	300	2	31,31,38	1.22	3 (9%)	39,40,47	1.27	6 (15%)
3	FEB	Z	300	2	31,31,38	1.16	3 (9%)	39,40,47	1.30	7 (17%)
3	FEB	C	300	2	31,31,38	1.12	3 (9%)	39,40,47	1.27	5 (12%)
3	FEB	P	300	2	31,31,38	1.33	7 (22%)	39,40,47	1.34	3 (7%)
3	FEB	R	300	2	31,31,38	1.23	5 (16%)	39,40,47	1.24	5 (12%)
3	FEB	T	300	2	31,31,38	1.30	3 (9%)	39,40,47	1.35	5 (12%)
3	FEB	2	300	2	31,31,38	1.19	3 (9%)	39,40,47	1.28	4 (10%)
3	FEB	H	300	2	31,31,38	1.35	6 (19%)	39,40,47	1.23	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
3	FEB	J	300	2	-	2/41/41/48	-
3	FEB	L	300	2	-	3/41/41/48	-
3	FEB	N	300	2	-	2/41/41/48	-
3	FEB	G	300	2	-	2/41/41/48	-
3	FEB	E	300	2	-	2/41/41/48	-
3	FEB	V	300	2	-	3/41/41/48	-
3	FEB	X	300	2	-	3/41/41/48	-
3	FEB	Z	300	2	-	2/41/41/48	-
3	FEB	C	300	2	-	2/41/41/48	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	FEB	P	300	2	-	3/41/41/48	-
3	FEB	R	300	2	-	3/41/41/48	-
3	FEB	T	300	2	-	4/41/41/48	-
3	FEB	2	300	2	-	2/41/41/48	-
3	FEB	H	300	2	-	3/41/41/48	-

The worst 5 of 59 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	J	300	FEB	C19-N18	3.50	1.53	1.45
3	R	300	FEB	O49-C38	3.18	1.29	1.23
3	N	300	FEB	C3-C2	3.08	1.57	1.52
3	V	300	FEB	O21-C20	3.06	1.29	1.23
3	T	300	FEB	C19-N18	2.98	1.52	1.45

The worst 5 of 66 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	T	300	FEB	C10-N9-C20	3.29	128.73	121.67
3	2	300	FEB	C19-C20-N9	-3.16	109.77	116.70
3	X	300	FEB	C19-C20-N9	-3.15	109.79	116.70
3	T	300	FEB	C19-C20-N9	-3.15	109.79	116.70
3	J	300	FEB	C19-C20-N9	-3.12	109.87	116.70

There are no chirality outliers.

5 of 36 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	J	300	FEB	C27-C36-C38-O49
3	L	300	FEB	C27-C36-C38-O49
3	E	300	FEB	C27-C36-C38-O49
3	X	300	FEB	C27-C36-C38-O49
3	C	300	FEB	C27-C36-C38-N18

There are no ring outliers.

13 monomers are involved in 28 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	J	300	FEB	1	0
3	L	300	FEB	1	0

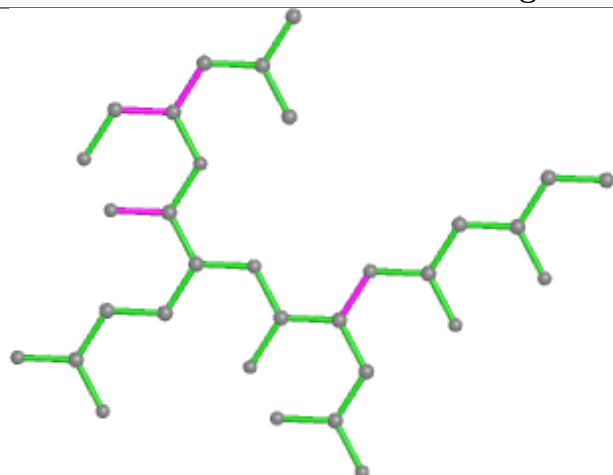
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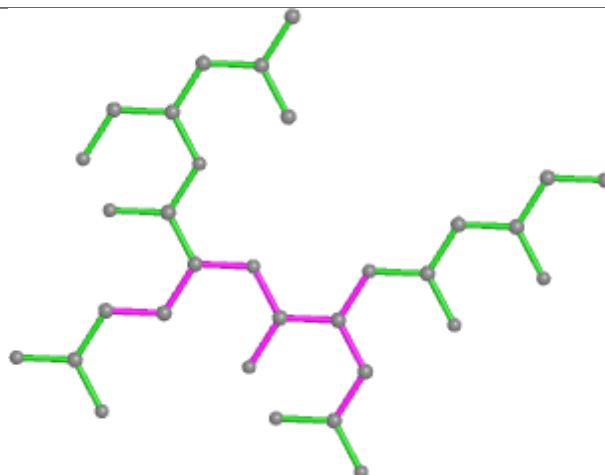
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	N	300	FEB	3	0
3	G	300	FEB	3	0
3	E	300	FEB	2	0
3	X	300	FEB	2	0
3	Z	300	FEB	2	0
3	C	300	FEB	3	0
3	P	300	FEB	2	0
3	R	300	FEB	3	0
3	T	300	FEB	3	0
3	2	300	FEB	2	0
3	H	300	FEB	1	0

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.

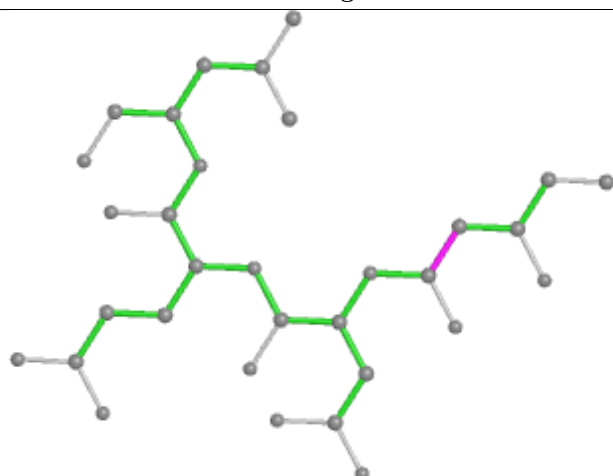
Ligand FEB J 300



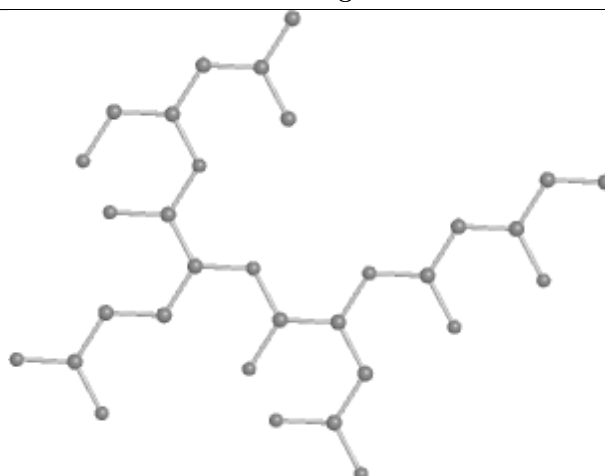
Bond lengths



Bond angles

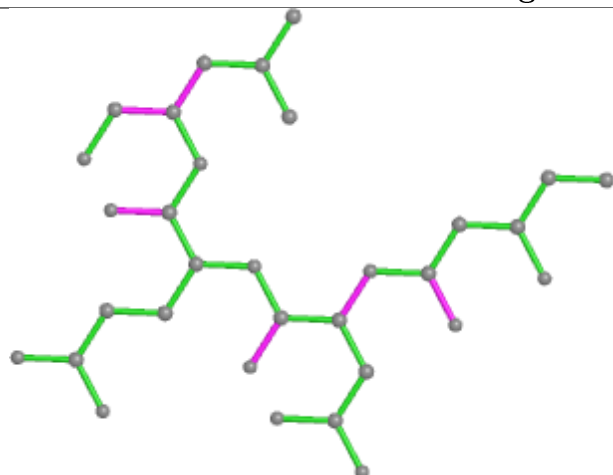


Torsions

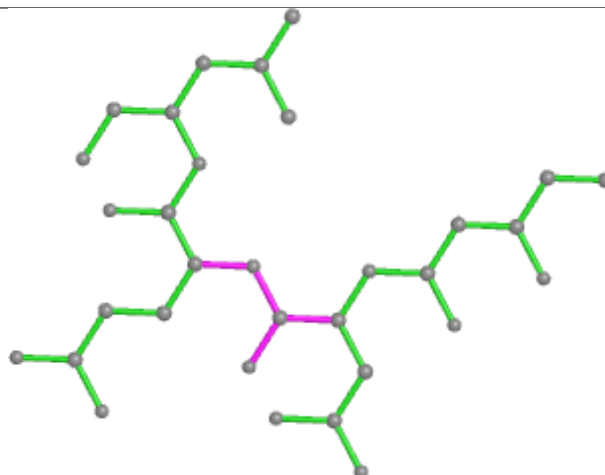


Rings

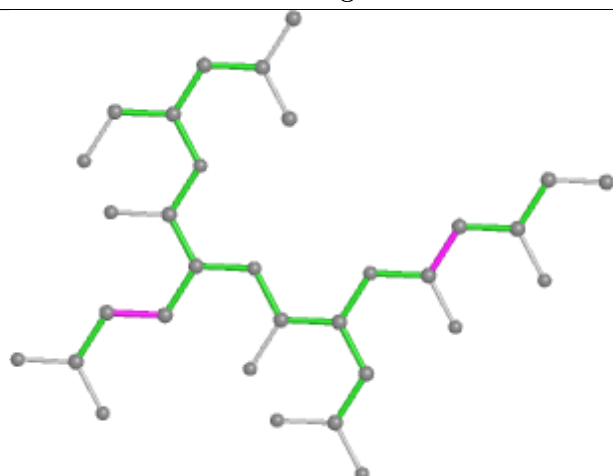
Ligand FEB L 300



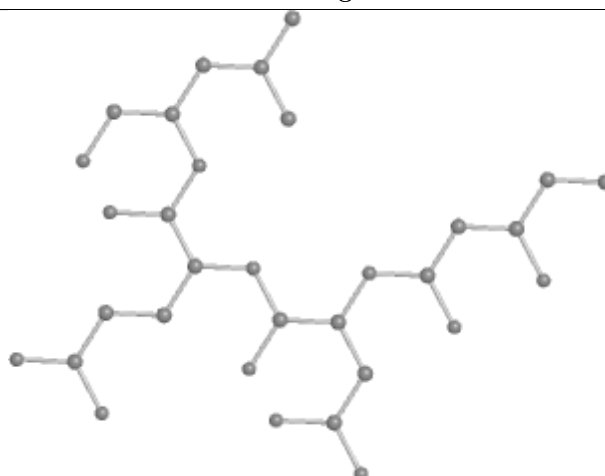
Bond lengths



Bond angles

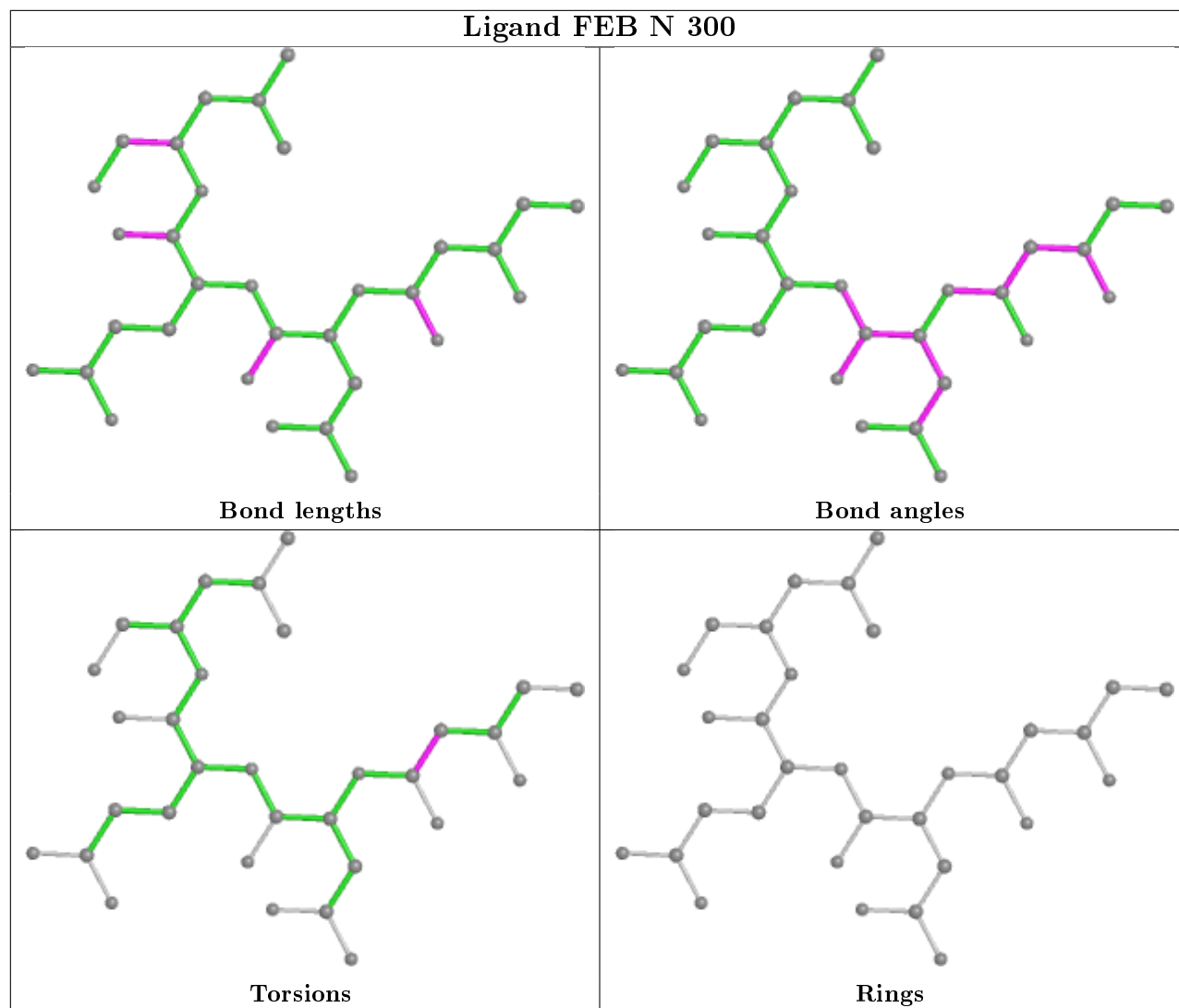


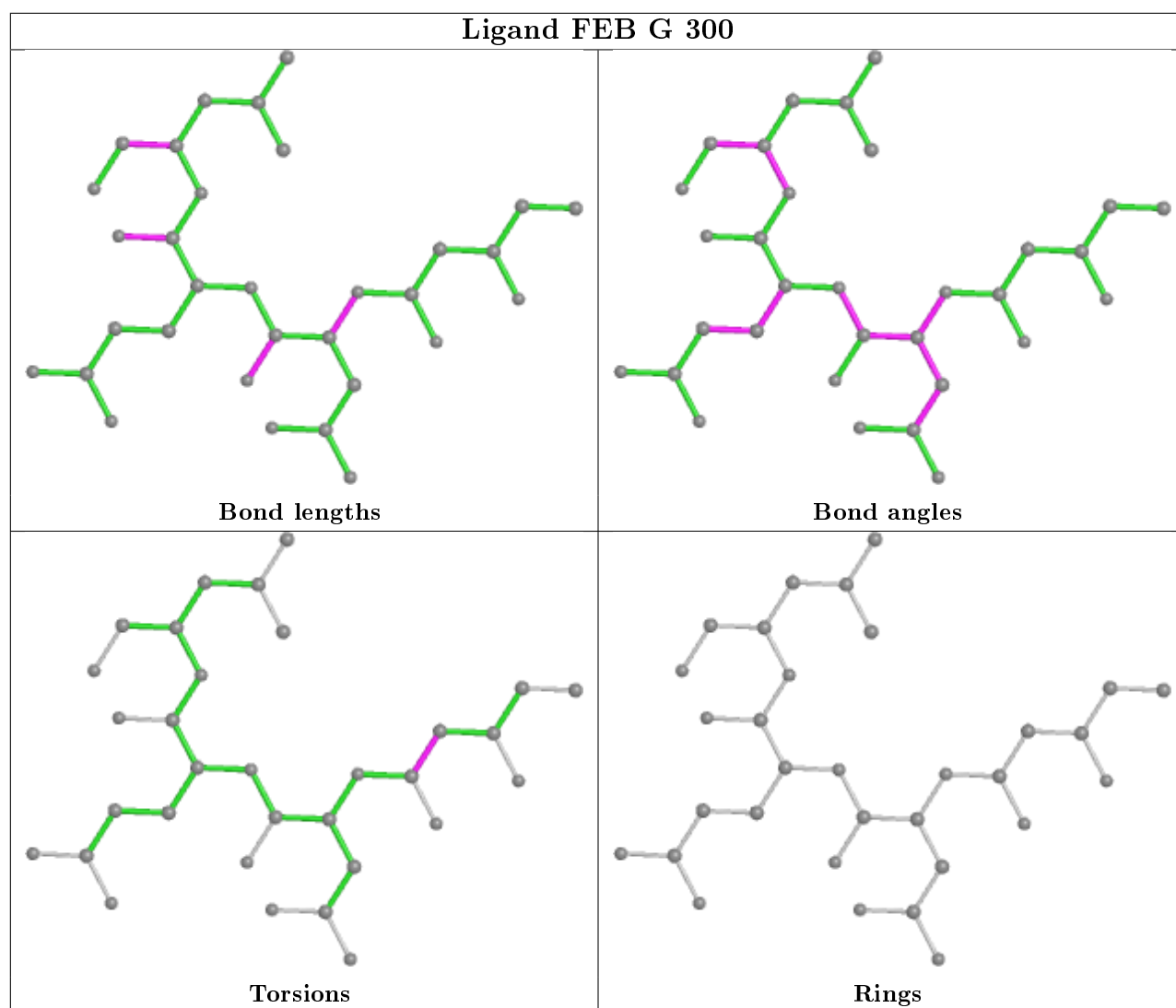
Torsions



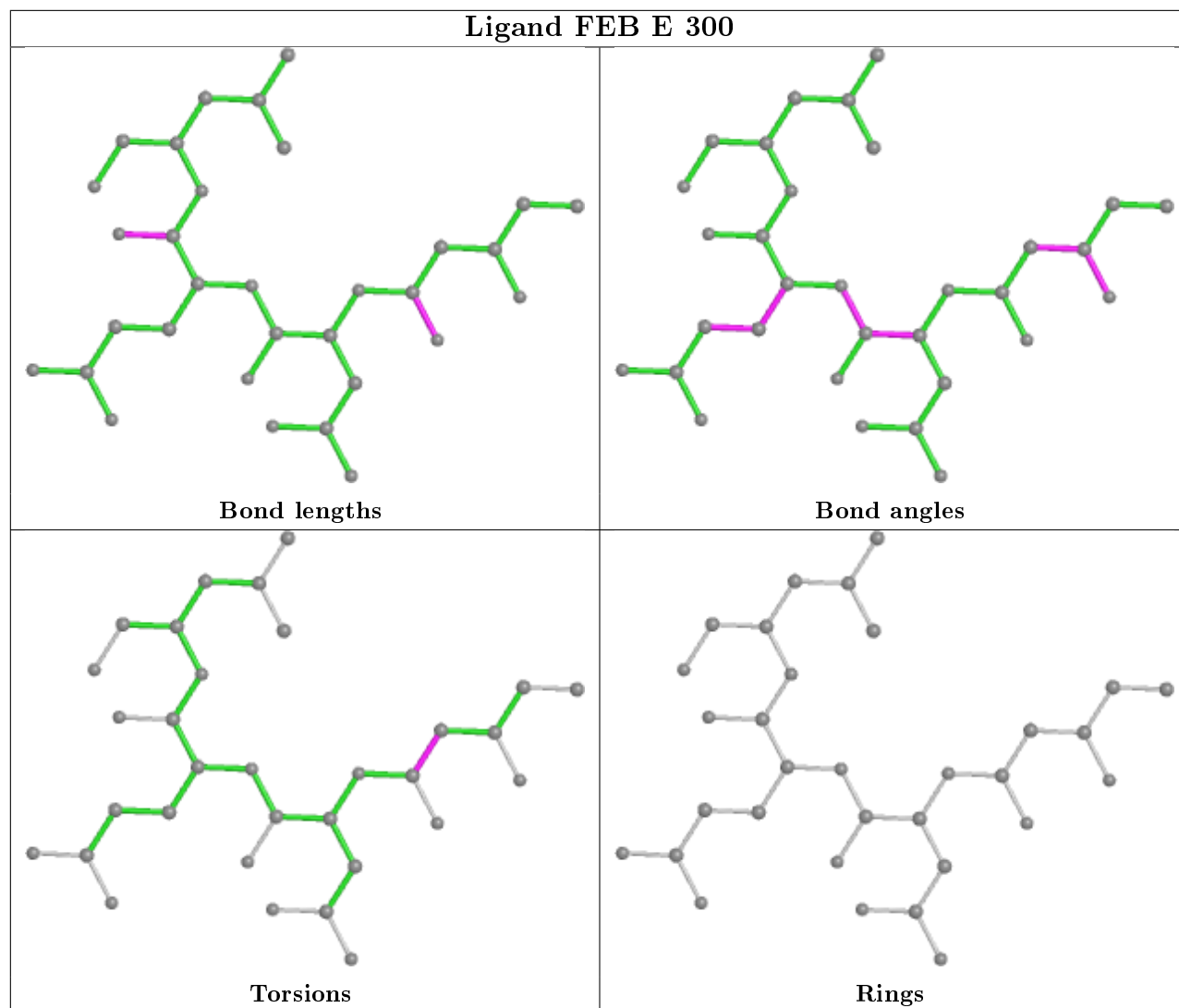
Rings

Ligand FEB N 300

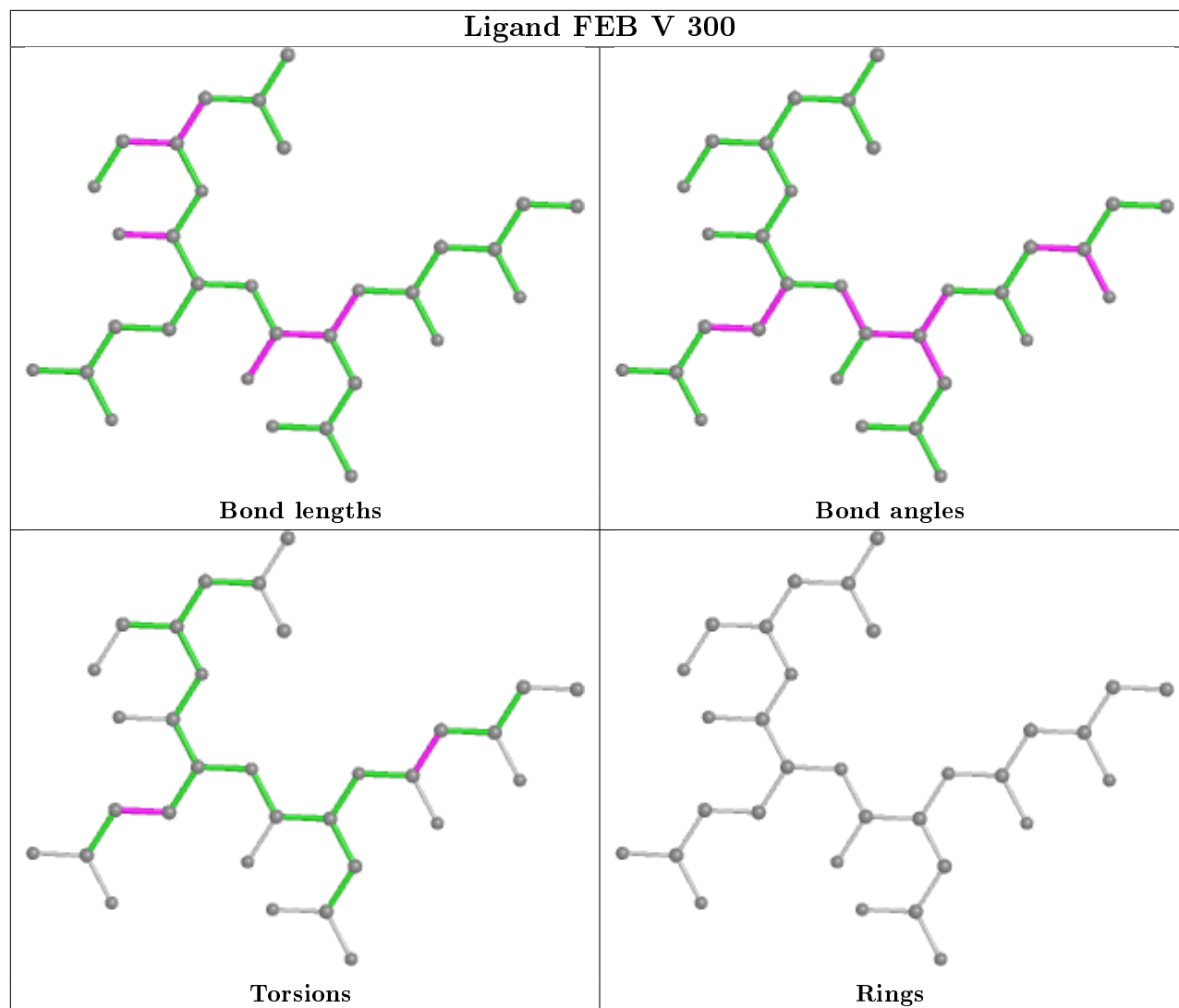




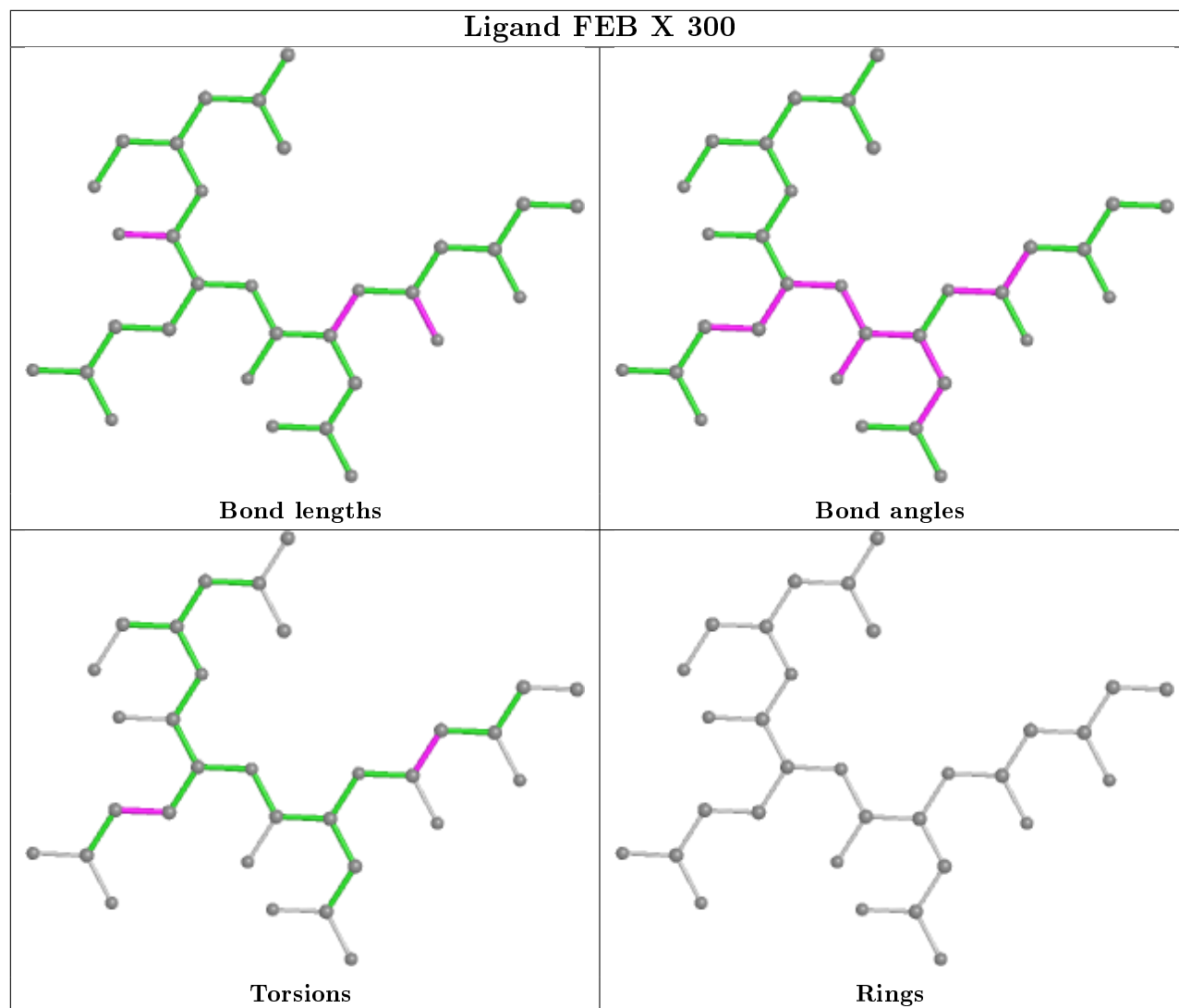
Ligand FEB E 300



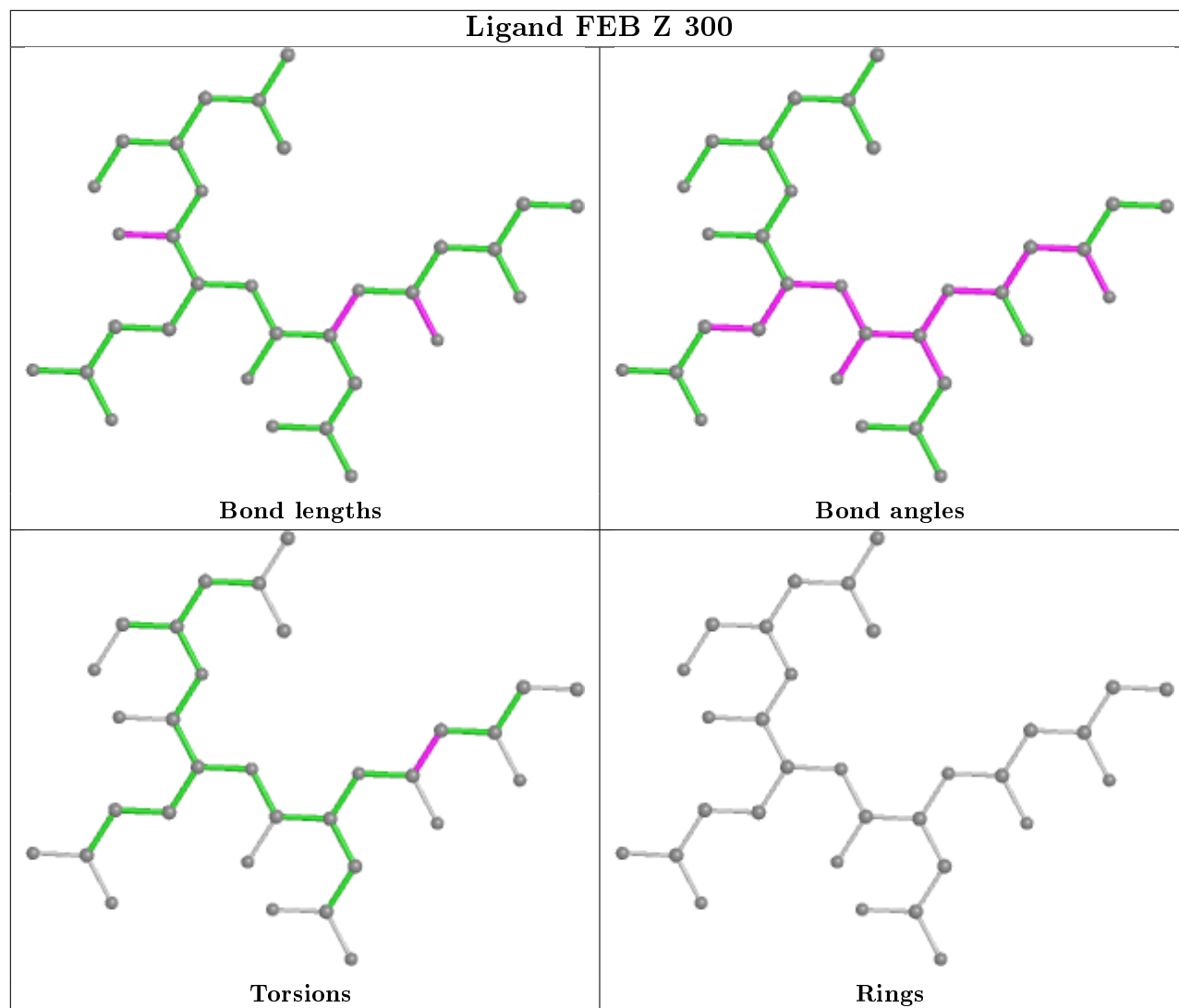
Ligand FEB V 300

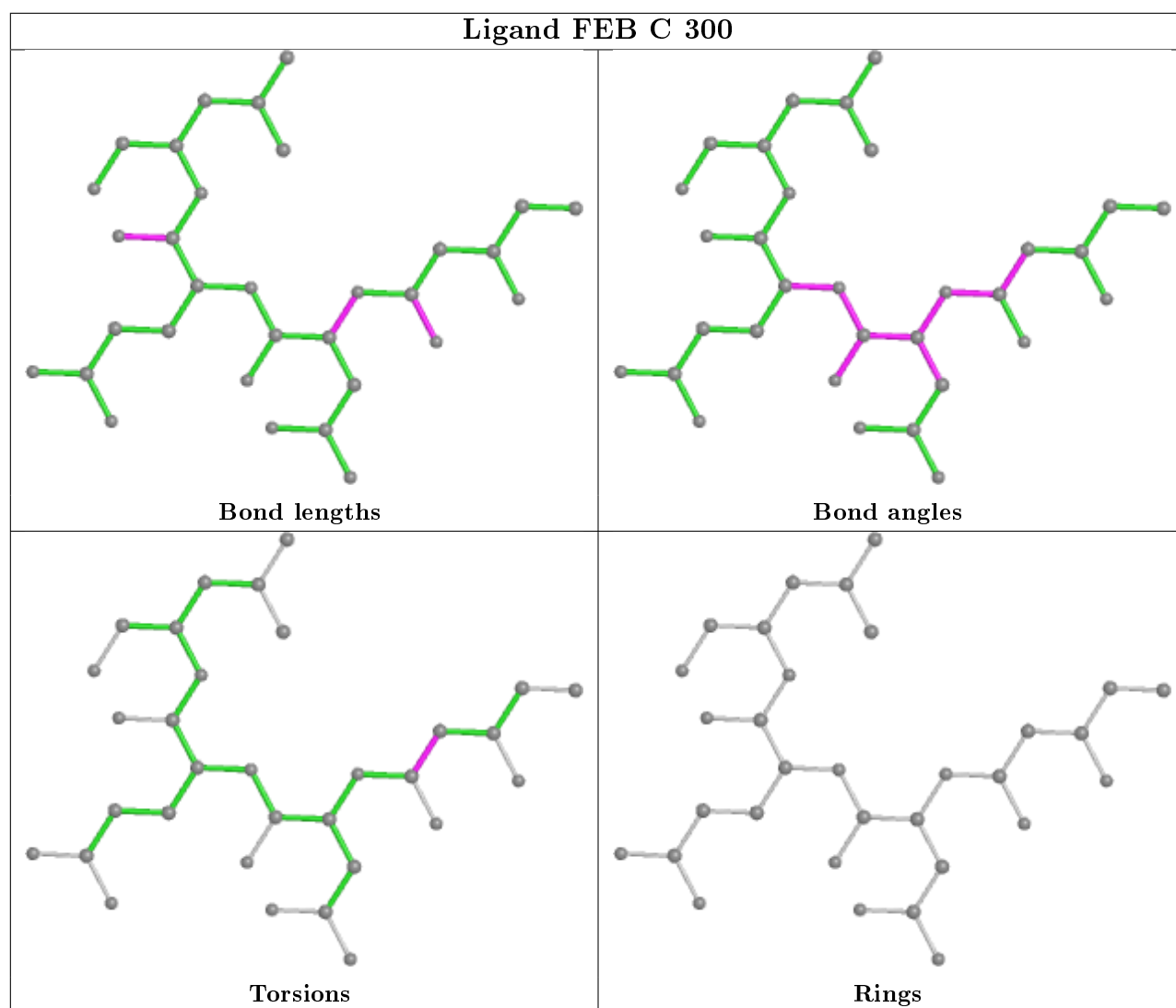


Ligand FEB X 300

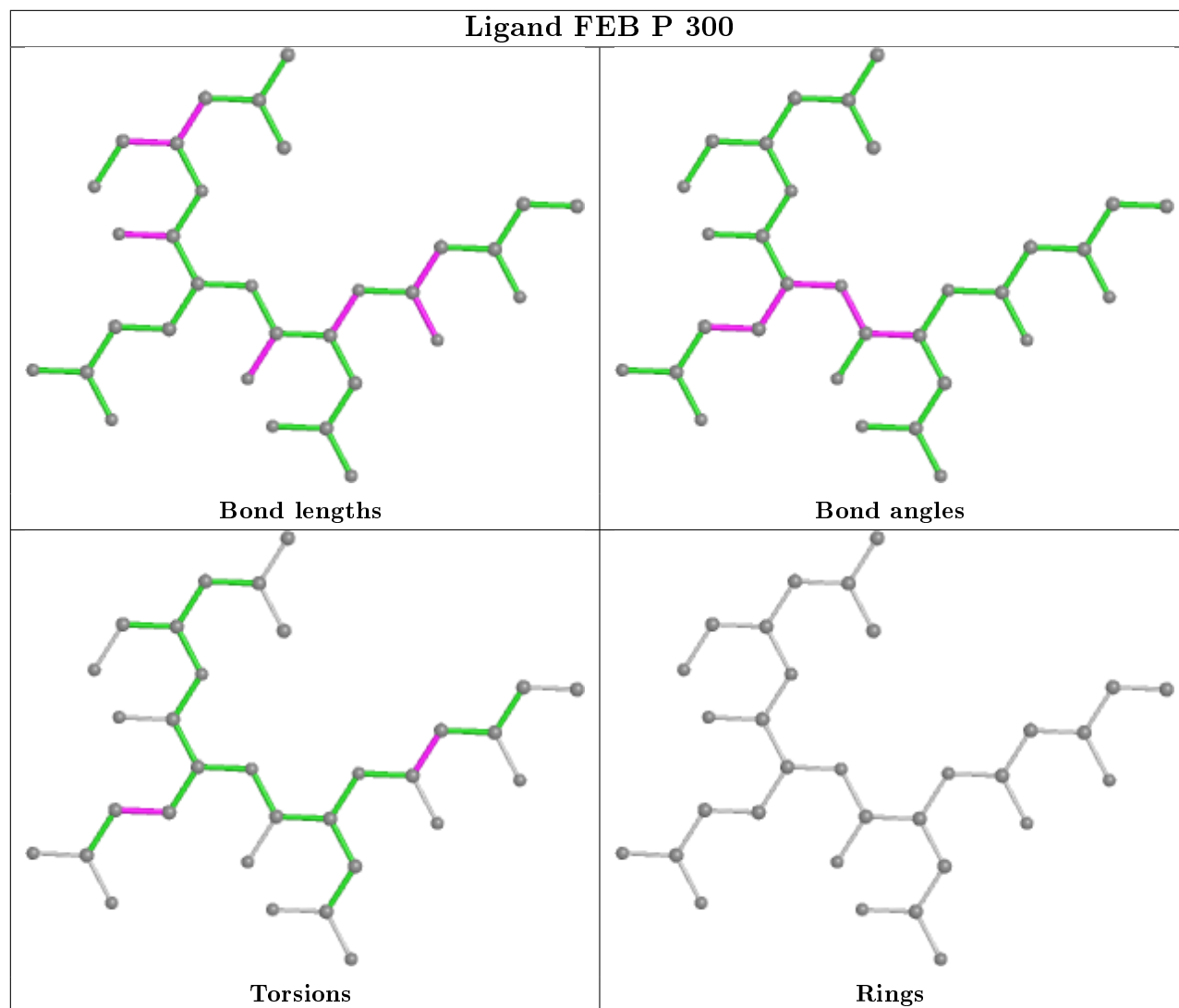


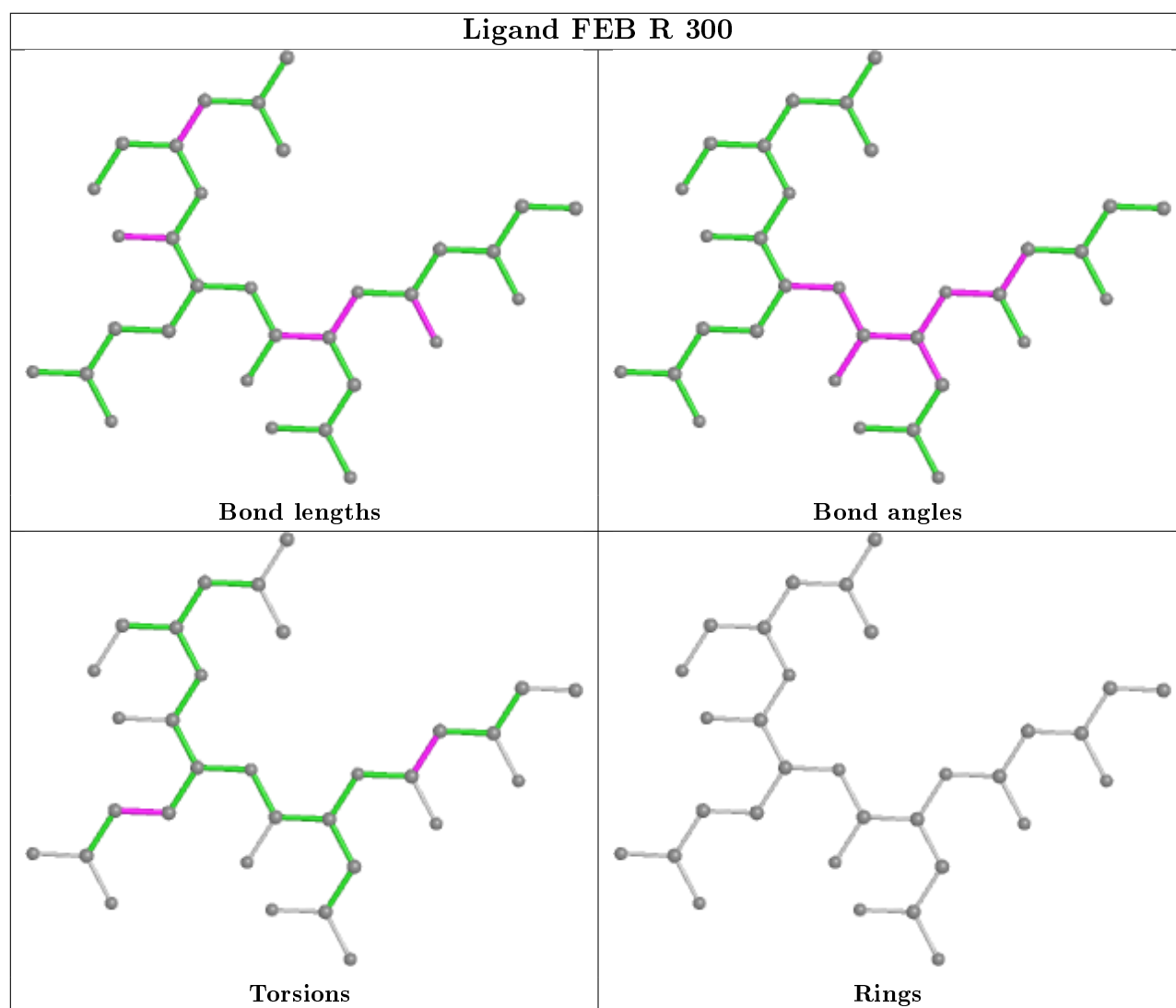
Ligand FEB Z 300

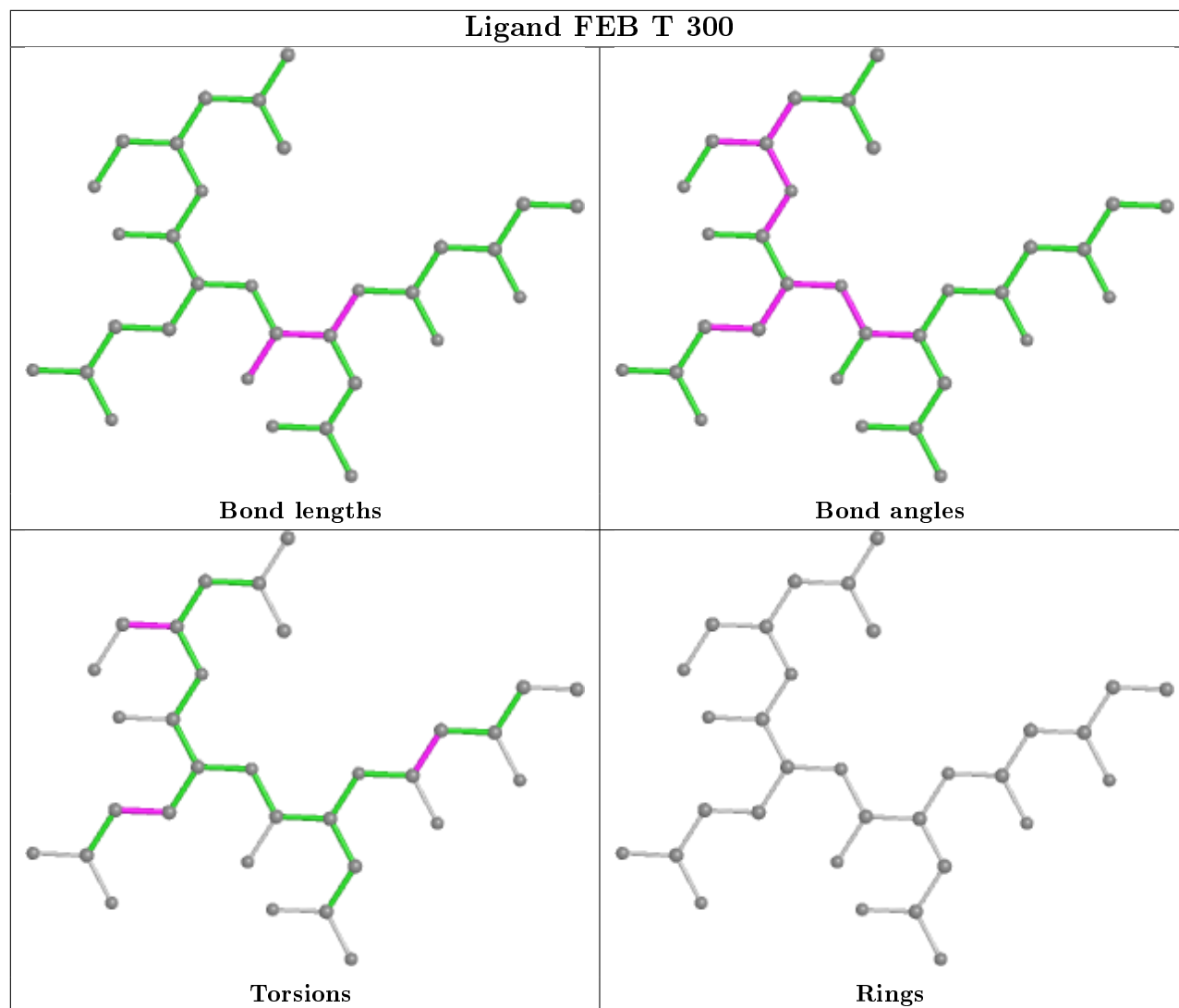




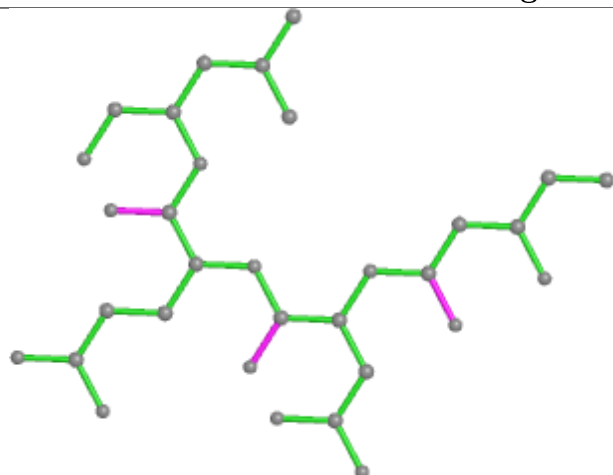
Ligand FEB P 300



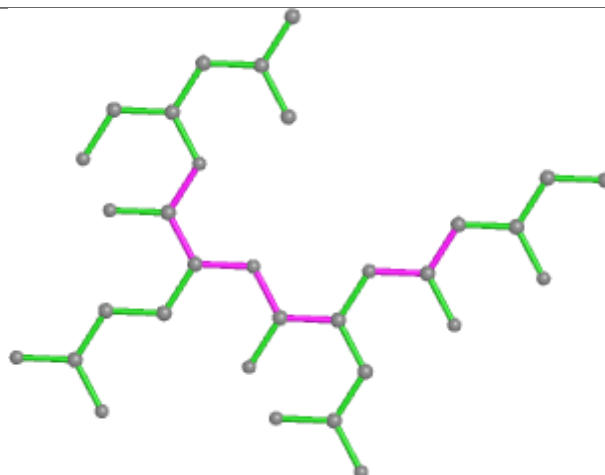




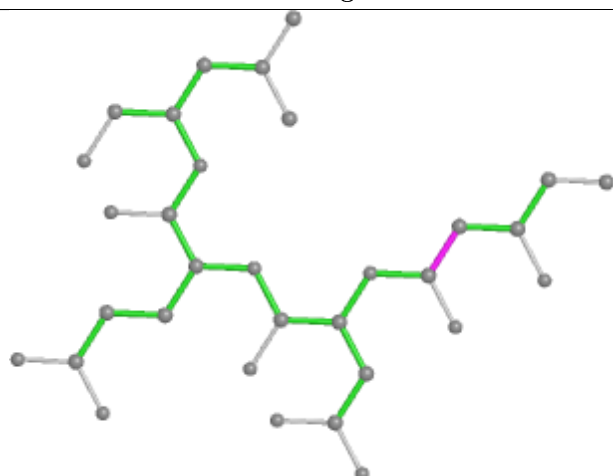
Ligand FEB 2 300



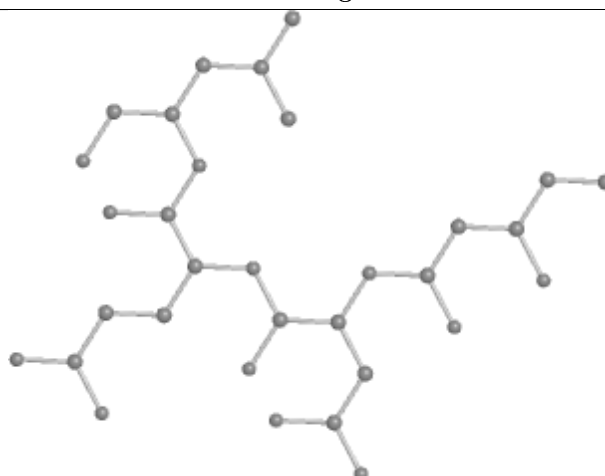
Bond lengths



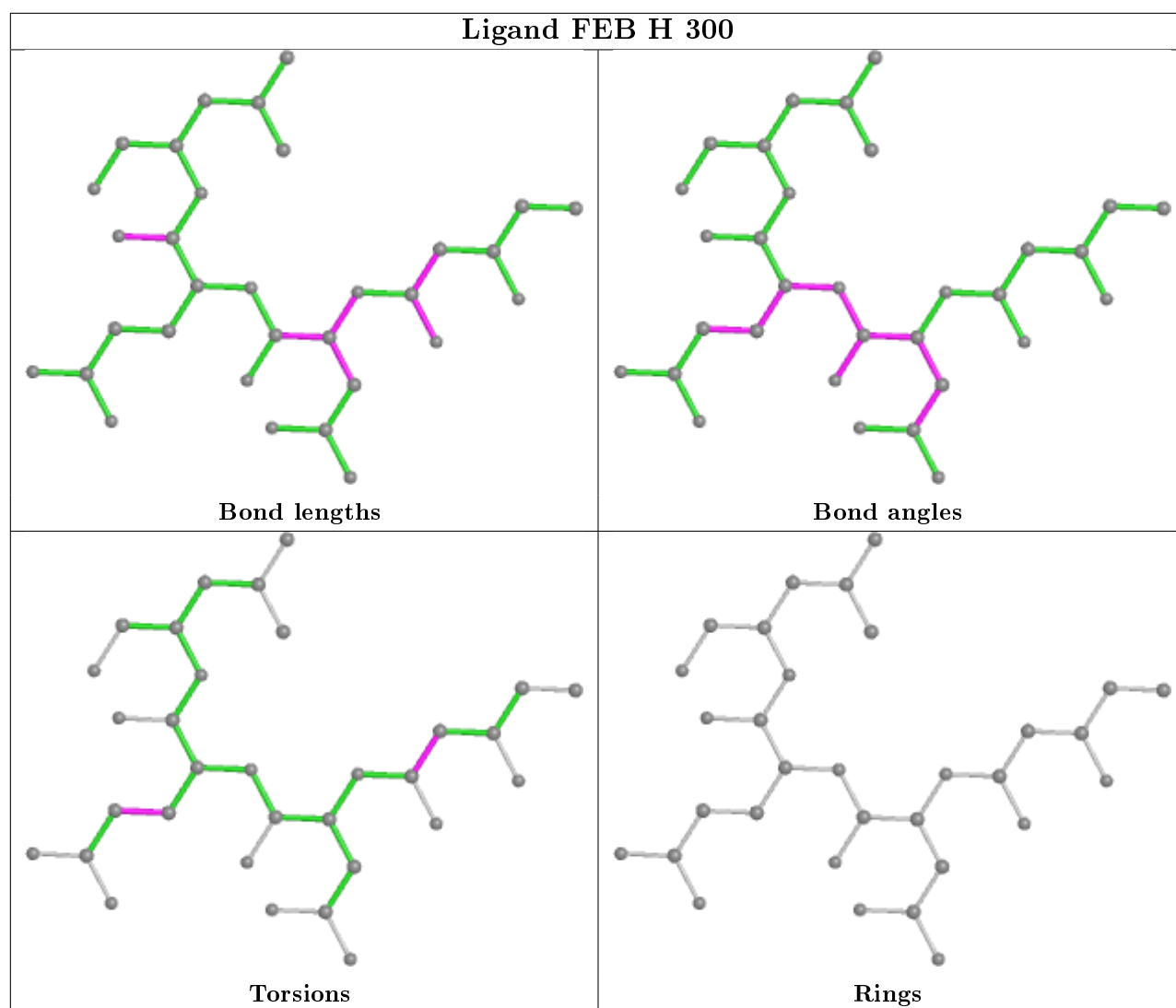
Bond angles



Torsions



Rings



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	1	215/248 (86%)	0.47	17 (7%) 12 12	23, 51, 86, 94	0
1	A	214/248 (86%)	0.62	29 (13%) 3 2	24, 53, 88, 100	0
1	B	213/248 (85%)	0.54	16 (7%) 14 14	24, 52, 87, 94	0
1	D	213/248 (85%)	0.97	49 (23%) 0 0	25, 55, 93, 112	0
1	F	216/248 (87%)	0.46	13 (6%) 21 22	23, 53, 87, 93	0
1	I	215/248 (86%)	0.53	20 (9%) 8 8	24, 52, 87, 95	0
1	K	215/248 (86%)	0.43	17 (7%) 12 12	24, 53, 87, 95	0
1	M	214/248 (86%)	0.57	20 (9%) 8 8	23, 54, 88, 96	0
1	O	214/248 (86%)	0.56	22 (10%) 6 6	22, 53, 88, 95	0
1	Q	214/248 (86%)	0.63	22 (10%) 6 6	26, 52, 87, 95	0
1	S	215/248 (86%)	0.69	31 (14%) 2 2	24, 53, 89, 99	0
1	U	214/248 (86%)	0.59	23 (10%) 6 5	23, 53, 87, 95	0
1	W	215/248 (86%)	0.68	28 (13%) 3 3	24, 54, 93, 98	0
1	Y	216/248 (87%)	0.76	38 (17%) 1 1	26, 55, 90, 98	0
2	2	222/240 (92%)	-0.36	1 (0%) 91 91	12, 23, 41, 66	0
2	C	222/240 (92%)	-0.35	4 (1%) 68 71	13, 24, 41, 68	0
2	E	222/240 (92%)	-0.38	1 (0%) 91 91	13, 24, 41, 65	0
2	G	222/240 (92%)	-0.33	3 (1%) 75 77	14, 24, 43, 68	0
2	H	222/240 (92%)	-0.37	0 100 100	11, 24, 42, 66	0
2	J	222/240 (92%)	-0.35	0 100 100	12, 24, 43, 65	0
2	L	222/240 (92%)	-0.38	0 100 100	13, 24, 42, 65	0
2	N	222/240 (92%)	-0.29	5 (2%) 60 63	13, 24, 43, 67	0
2	P	222/240 (92%)	-0.37	2 (0%) 84 86	14, 24, 42, 67	0
2	R	229/240 (95%)	-0.39	3 (1%) 77 79	13, 24, 46, 67	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
2	T	222/240 (92%)	-0.35	0 100 100	15, 25, 43, 67	0
2	V	229/240 (95%)	-0.39	3 (1%) 77 79	12, 23, 44, 67	0
2	X	222/240 (92%)	-0.38	4 (1%) 68 71	13, 25, 43, 67	0
2	Z	222/240 (92%)	-0.33	3 (1%) 75 77	13, 25, 43, 67	0
All	All	6125/6832 (89%)	0.12	374 (6%) 21 22	11, 35, 84, 112	0

The worst 5 of 374 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	233	LEU	7.0
1	B	205	VAL	6.5
1	Y	12	ALA	6.4
1	Y	10	GLU	6.3
1	S	235	VAL	5.0

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, 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
3	FEB	X	300	32/39	0.92	0.14	21,26,30,33	0
3	FEB	E	300	32/39	0.93	0.14	10,21,35,37	0
3	FEB	2	300	32/39	0.93	0.14	16,27,34,34	0
3	FEB	Z	300	32/39	0.93	0.15	23,29,36,37	0
3	FEB	J	300	32/39	0.94	0.12	12,23,36,38	0
3	FEB	N	300	32/39	0.94	0.13	15,25,33,36	0
3	FEB	R	300	32/39	0.94	0.15	21,25,37,40	0

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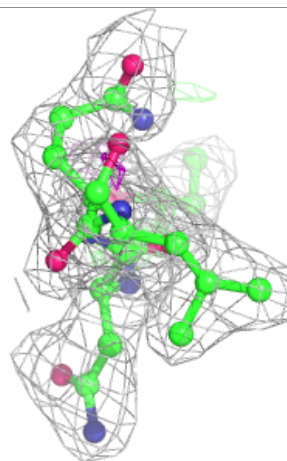
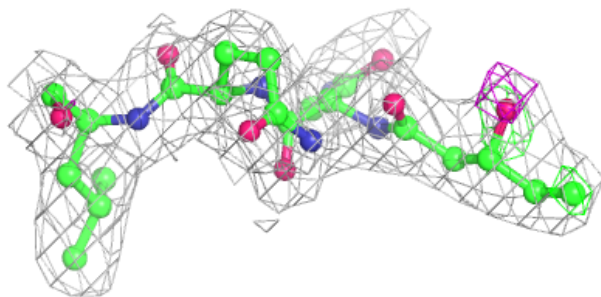
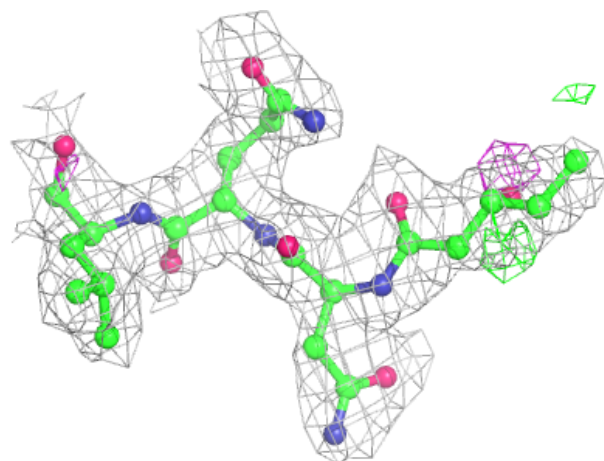
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	FEB	T	300	32/39	0.94	0.14	18,27,36,37	0
3	FEB	H	300	32/39	0.94	0.13	17,25,34,38	0
3	FEB	P	300	32/39	0.95	0.13	20,26,35,37	0
3	FEB	L	300	32/39	0.95	0.14	18,24,34,37	0
3	FEB	G	300	32/39	0.95	0.14	17,25,32,34	0
3	FEB	V	300	32/39	0.95	0.12	17,27,36,37	0
3	FEB	C	300	32/39	0.95	0.14	19,25,32,35	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.

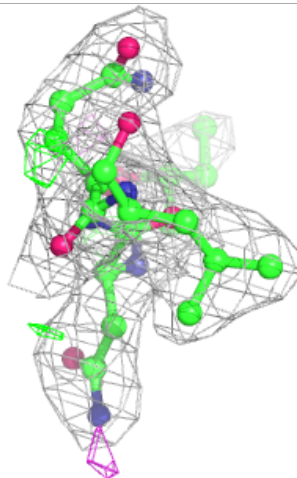
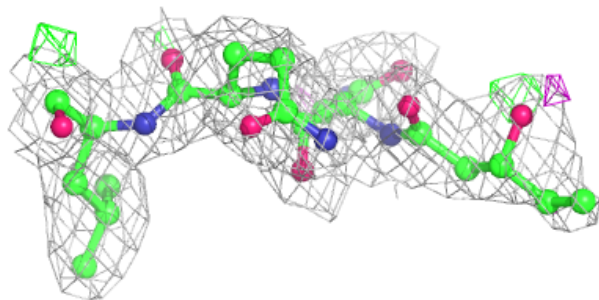
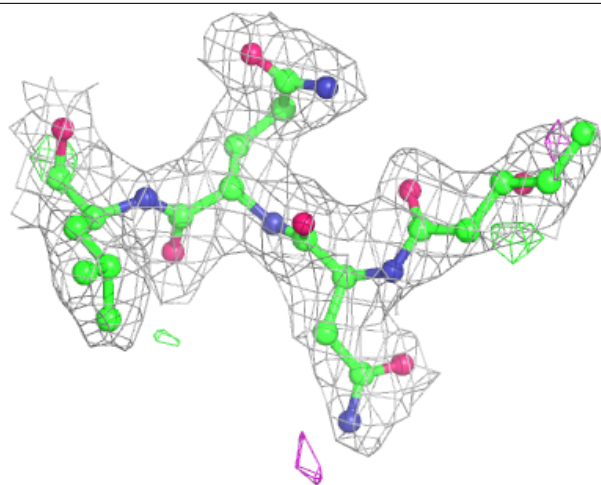
Electron density around FEB X 300:

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



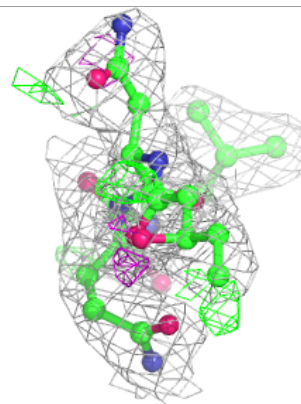
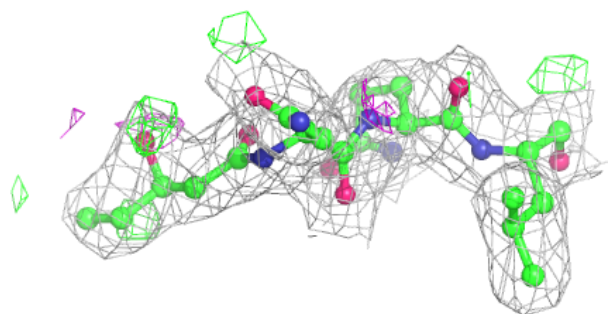
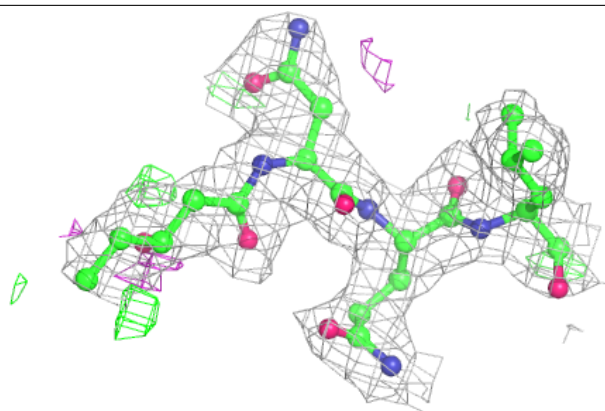
Electron density around FEB E 300:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

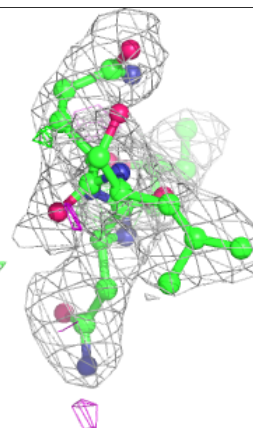
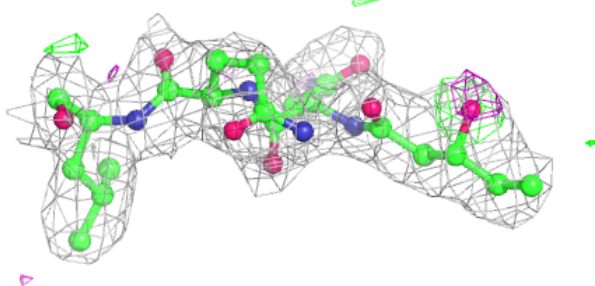
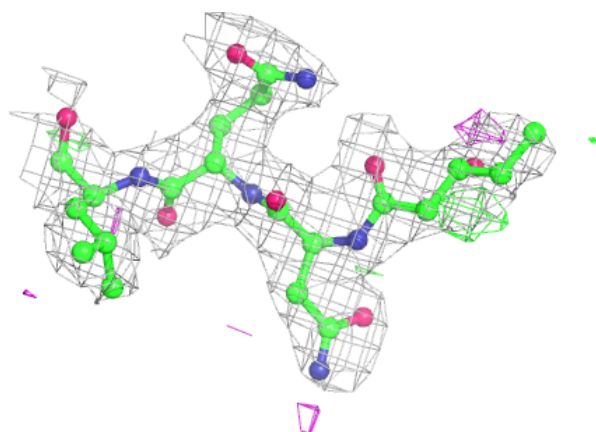


Electron density around FEB 2 300:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

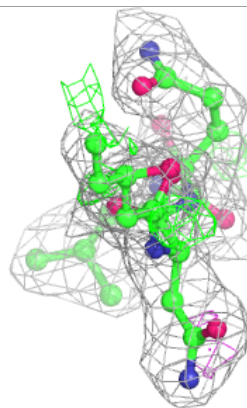
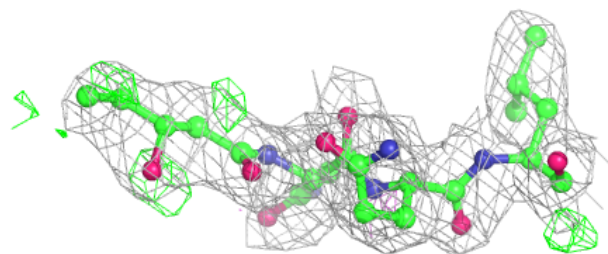
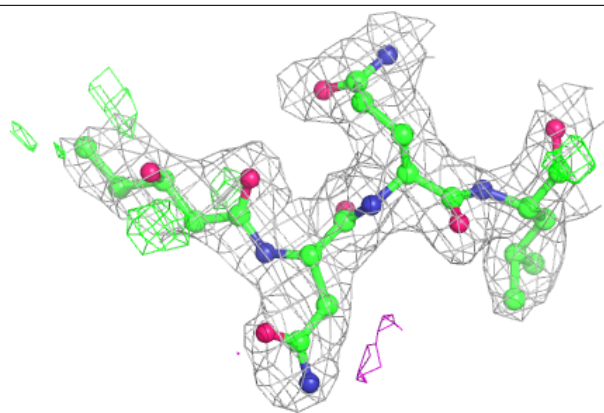
**Electron density around FEB Z 300:**

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and green (positive)



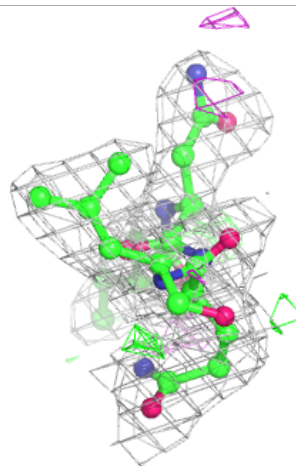
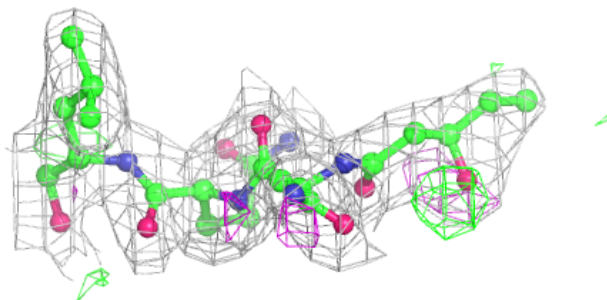
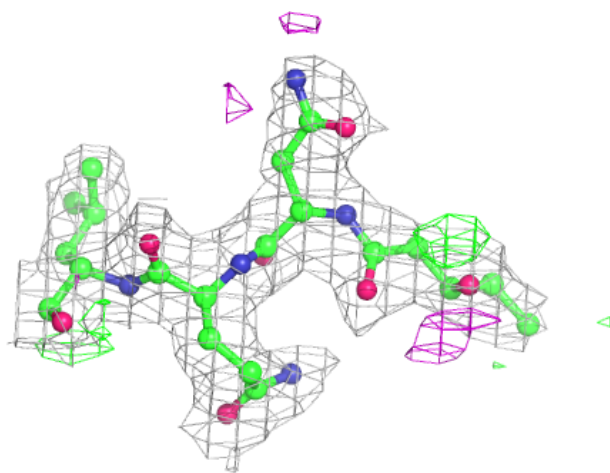
Electron density around FEB J 300:

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



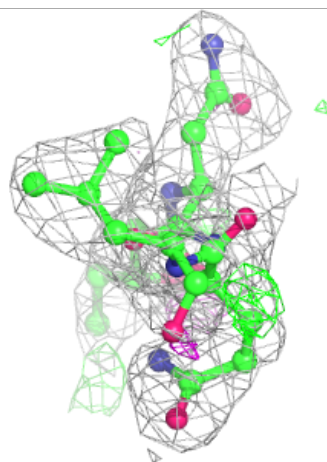
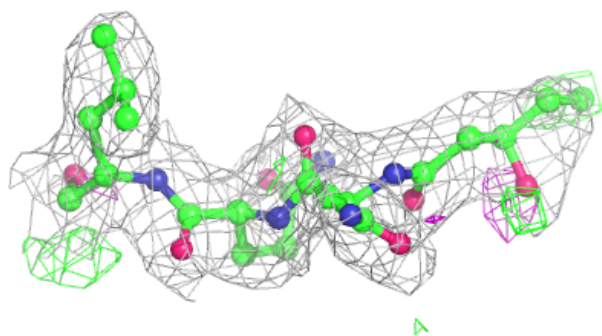
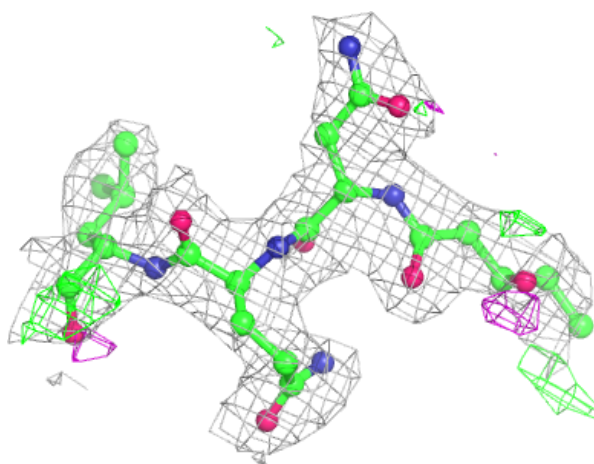
Electron density around FEB N 300:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



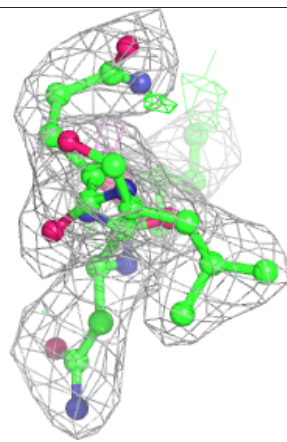
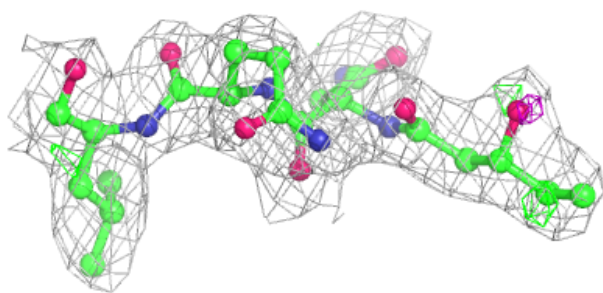
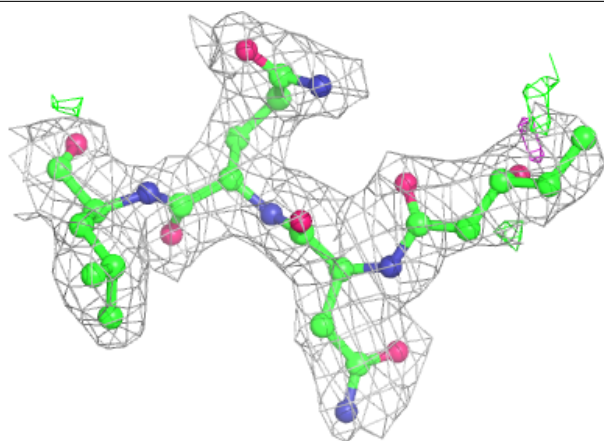
Electron density around FEB R 300:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



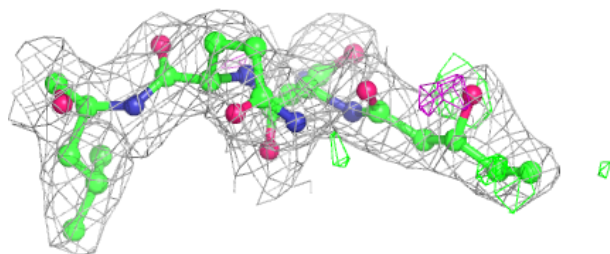
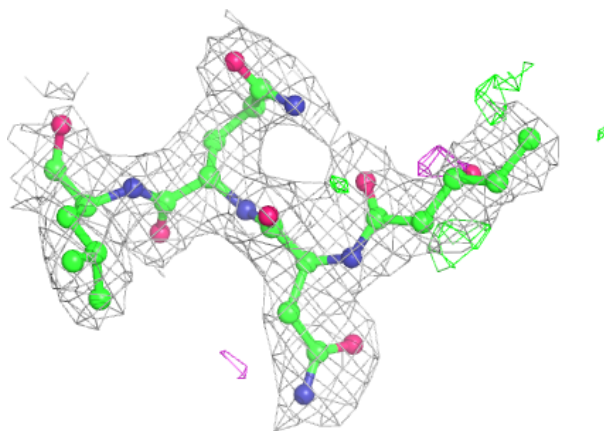
Electron density around FEB T 300:

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and green (positive)

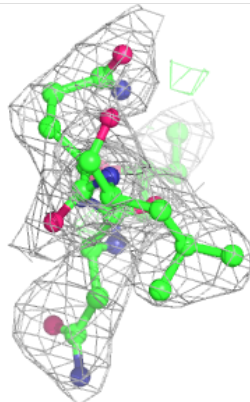
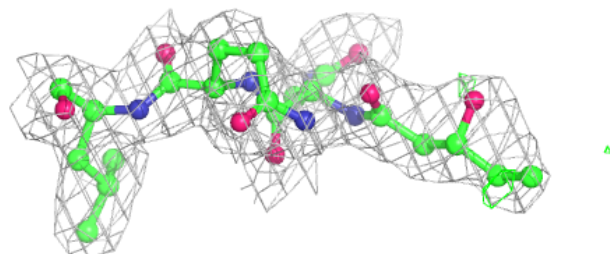
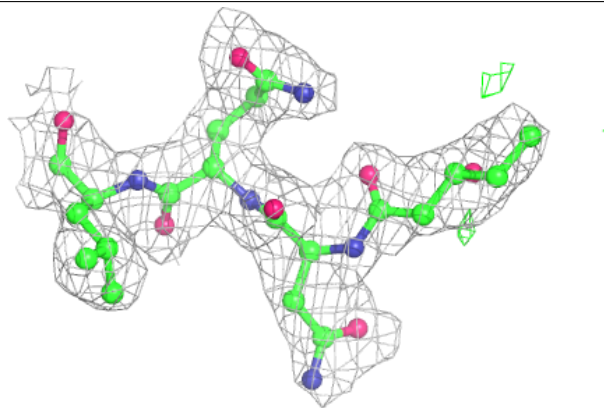


Electron density around FEB H 300:

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and green (positive)

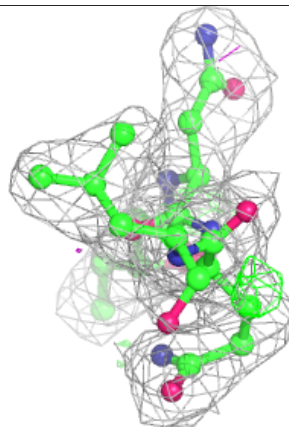
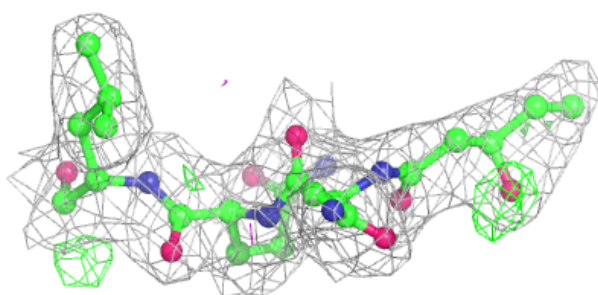
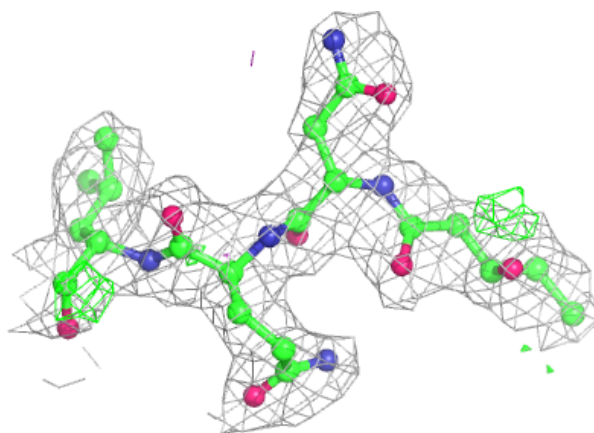
**Electron density around FEB P 300:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

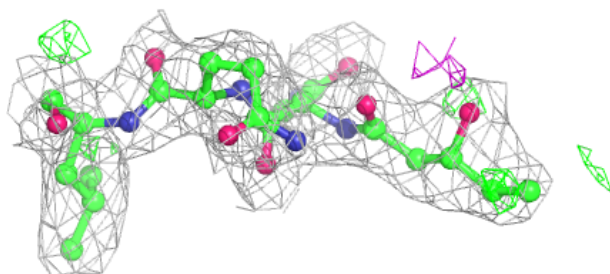
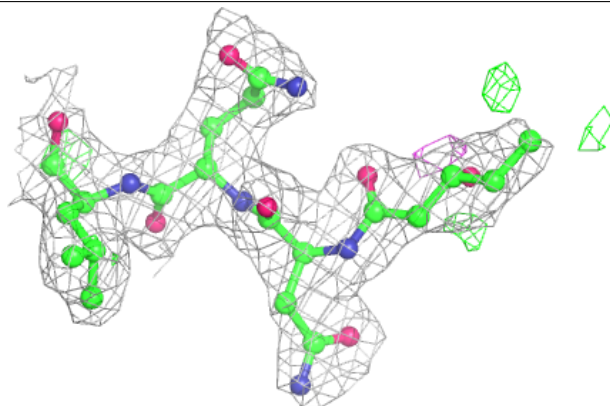


Electron density around FEB L 300:

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and green (positive)

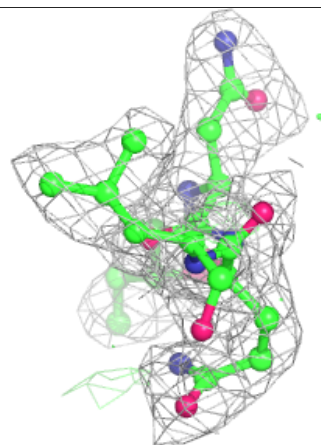
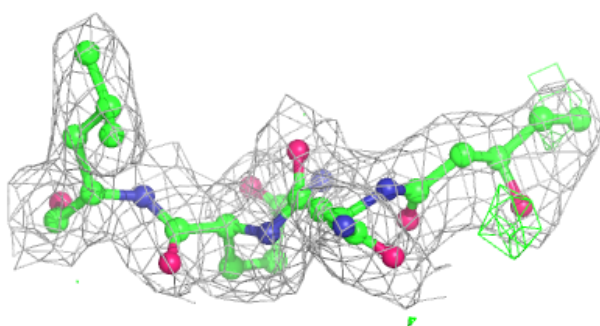
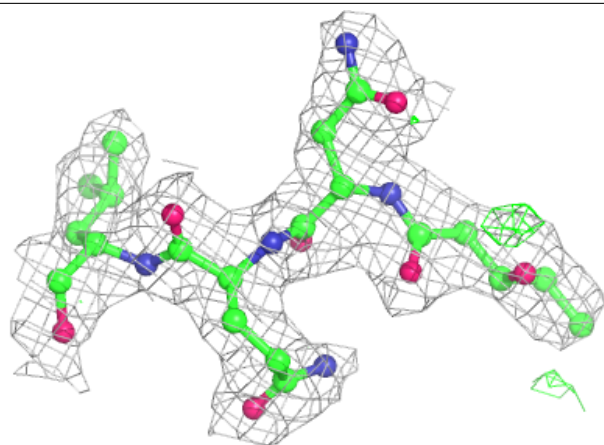
**Electron density around FEB G 300:**

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and green (positive)



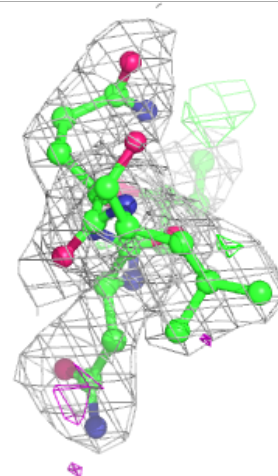
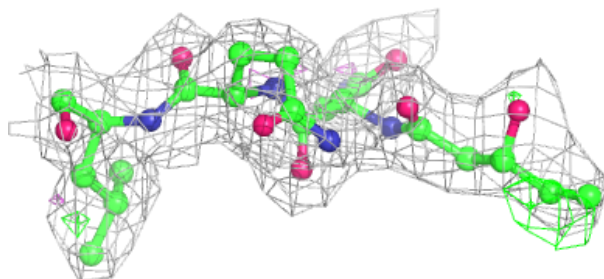
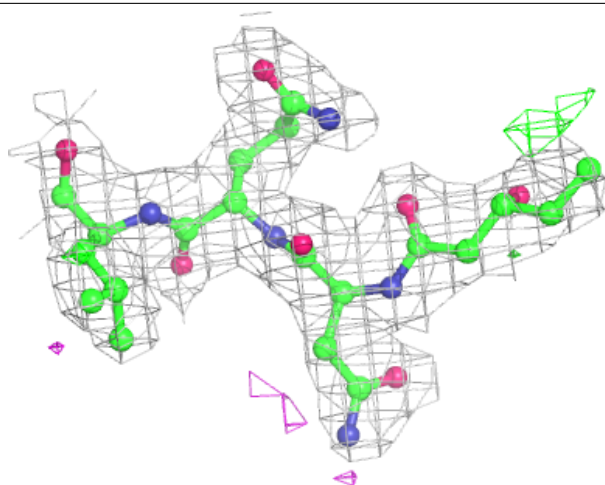
Electron density around FEB V 300:

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



Electron density around FEB C 300:

$2mF_o-DF_c$ (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.