



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:38 PM UTC

PDB ID : 4ZJT / pdb_00004zjt
Title : X-ray crystal structure of Lymnaea stagnalis acetylcholine binding protein (LsAChBP) in complex with 2-Thiophenylmethylene Anabaseine (2TAB)
Authors : Bobango, J.; Wu, J.; Talley, T.T.
Deposited on : 2015-04-29
Resolution : 1.85 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

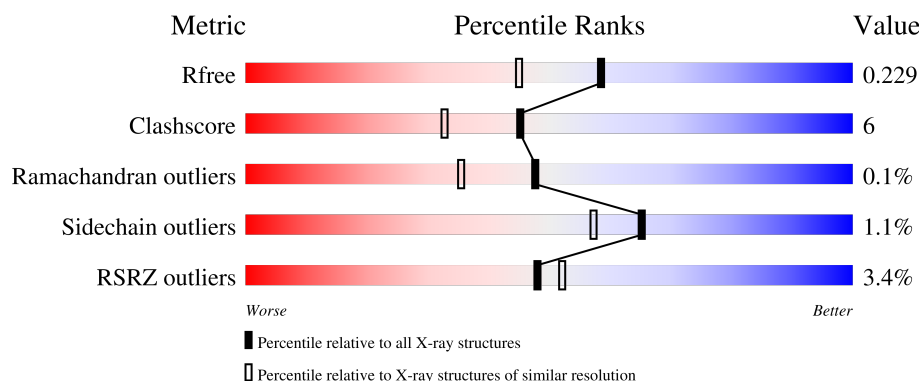
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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.



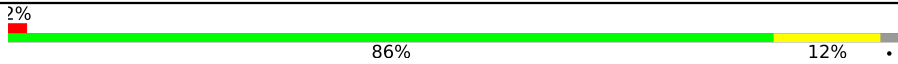
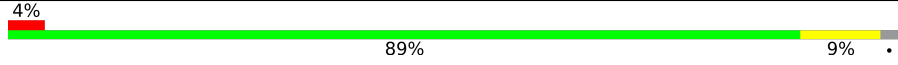
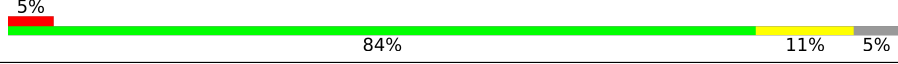
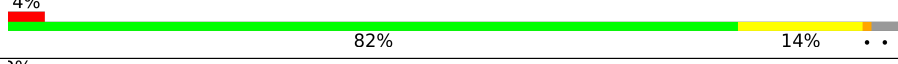
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	218	2% 87% 11% .
1	B	218	2% 82% 14% .
1	C	218	5% 90% 6% .
1	D	218	3% 87% 11% .
1	E	218	4% 87% 10% .

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Mol	Chain	Length	Quality of chain
1	F	218	
1	G	218	
1	H	218	
1	I	218	
1	J	218	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	4P6	E	301	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 17593 atoms, of which 140 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 Acetylcholine-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	212	Total	C	N	O	S	0	1	0
			1674	1048	282	339	5			
1	B	210	Total	C	N	O	S	0	1	0
			1655	1034	279	337	5			
1	C	209	Total	C	N	O	S	0	1	0
			1645	1032	276	332	5			
1	D	214	Total	C	N	O	S	0	0	0
			1674	1046	278	345	5			
1	E	212	Total	C	N	O	S	0	0	0
			1641	1028	275	333	5			
1	F	214	Total	C	N	O	S	0	0	0
			1660	1039	278	338	5			
1	G	214	Total	C	N	O	S	0	1	0
			1670	1045	278	342	5			
1	H	208	Total	C	N	O	S	0	1	0
			1651	1032	283	331	5			
1	I	211	Total	C	N	O	S	0	0	0
			1637	1025	270	337	5			
1	J	211	Total	C	N	O	S	0	0	0
			1660	1036	277	342	5			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	ASP	-	expression tag	UNP P58154
A	-6	TYR	-	expression tag	UNP P58154
A	-5	LYS	-	expression tag	UNP P58154
A	-4	ASP	-	expression tag	UNP P58154
A	-3	ASP	-	expression tag	UNP P58154
A	-2	ASP	-	expression tag	UNP P58154
A	-1	ASP	-	expression tag	UNP P58154
A	0	LYS	-	expression tag	UNP P58154
B	-7	ASP	-	expression tag	UNP P58154

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	TYR	-	expression tag	UNP P58154
B	-5	LYS	-	expression tag	UNP P58154
B	-4	ASP	-	expression tag	UNP P58154
B	-3	ASP	-	expression tag	UNP P58154
B	-2	ASP	-	expression tag	UNP P58154
B	-1	ASP	-	expression tag	UNP P58154
B	0	LYS	-	expression tag	UNP P58154
C	-7	ASP	-	expression tag	UNP P58154
C	-6	TYR	-	expression tag	UNP P58154
C	-5	LYS	-	expression tag	UNP P58154
C	-4	ASP	-	expression tag	UNP P58154
C	-3	ASP	-	expression tag	UNP P58154
C	-2	ASP	-	expression tag	UNP P58154
C	-1	ASP	-	expression tag	UNP P58154
C	0	LYS	-	expression tag	UNP P58154
D	-7	ASP	-	expression tag	UNP P58154
D	-6	TYR	-	expression tag	UNP P58154
D	-5	LYS	-	expression tag	UNP P58154
D	-4	ASP	-	expression tag	UNP P58154
D	-3	ASP	-	expression tag	UNP P58154
D	-2	ASP	-	expression tag	UNP P58154
D	-1	ASP	-	expression tag	UNP P58154
D	0	LYS	-	expression tag	UNP P58154
E	-7	ASP	-	expression tag	UNP P58154
E	-6	TYR	-	expression tag	UNP P58154
E	-5	LYS	-	expression tag	UNP P58154
E	-4	ASP	-	expression tag	UNP P58154
E	-3	ASP	-	expression tag	UNP P58154
E	-2	ASP	-	expression tag	UNP P58154
E	-1	ASP	-	expression tag	UNP P58154
E	0	LYS	-	expression tag	UNP P58154
F	-7	ASP	-	expression tag	UNP P58154
F	-6	TYR	-	expression tag	UNP P58154
F	-5	LYS	-	expression tag	UNP P58154
F	-4	ASP	-	expression tag	UNP P58154
F	-3	ASP	-	expression tag	UNP P58154
F	-2	ASP	-	expression tag	UNP P58154
F	-1	ASP	-	expression tag	UNP P58154
F	0	LYS	-	expression tag	UNP P58154
G	-7	ASP	-	expression tag	UNP P58154
G	-6	TYR	-	expression tag	UNP P58154
G	-5	LYS	-	expression tag	UNP P58154

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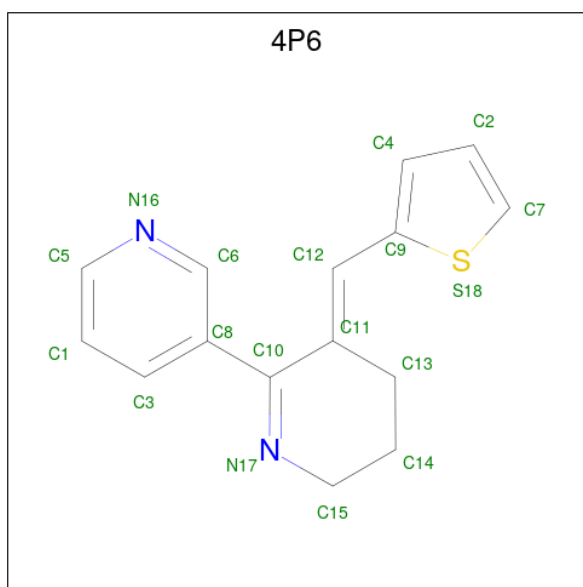
Chain	Residue	Modelled	Actual	Comment	Reference
G	-4	ASP	-	expression tag	UNP P58154
G	-3	ASP	-	expression tag	UNP P58154
G	-2	ASP	-	expression tag	UNP P58154
G	-1	ASP	-	expression tag	UNP P58154
G	0	LYS	-	expression tag	UNP P58154
H	-7	ASP	-	expression tag	UNP P58154
H	-6	TYR	-	expression tag	UNP P58154
H	-5	LYS	-	expression tag	UNP P58154
H	-4	ASP	-	expression tag	UNP P58154
H	-3	ASP	-	expression tag	UNP P58154
H	-2	ASP	-	expression tag	UNP P58154
H	-1	ASP	-	expression tag	UNP P58154
H	0	LYS	-	expression tag	UNP P58154
I	-7	ASP	-	expression tag	UNP P58154
I	-6	TYR	-	expression tag	UNP P58154
I	-5	LYS	-	expression tag	UNP P58154
I	-4	ASP	-	expression tag	UNP P58154
I	-3	ASP	-	expression tag	UNP P58154
I	-2	ASP	-	expression tag	UNP P58154
I	-1	ASP	-	expression tag	UNP P58154
I	0	LYS	-	expression tag	UNP P58154
J	-7	ASP	-	expression tag	UNP P58154
J	-6	TYR	-	expression tag	UNP P58154
J	-5	LYS	-	expression tag	UNP P58154
J	-4	ASP	-	expression tag	UNP P58154
J	-3	ASP	-	expression tag	UNP P58154
J	-2	ASP	-	expression tag	UNP P58154
J	-1	ASP	-	expression tag	UNP P58154
J	0	LYS	-	expression tag	UNP P58154

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	G	1	Total	O	P	0	0
			5	4	1		
2	G	1	Total	O	P	0	0
			5	4	1		
2	I	1	Total	O	P	0	0
			5	4	1		
2	I	1	Total	O	P	0	0
			5	4	1		
2	J	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is (3E)-3-(thiophen-2-ylmethyldiene)-3,4,5,6-tetrahydro-2,3'-bipyridine (CCD ID: 4P6) (formula: C₁₅H₁₄N₂S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	N	S	0	0
			32	15	14	2	1		
3	B	1	Total	C	H	N	S	0	0
			32	15	14	2	1		
3	C	1	Total	C	H	N	S	0	0
			32	15	14	2	1		
3	D	1	Total	C	H	N	S	0	0
			32	15	14	2	1		
3	E	1	Total	C	H	N	S	0	0
			32	15	14	2	1		
3	F	1	Total	C	H	N	S	0	0
			32	15	14	2	1		
3	G	1	Total	C	H	N	S	0	0
			32	15	14	2	1		
3	H	1	Total	C	H	N	S	0	0
			32	15	14	2	1		
3	I	1	Total	C	H	N	S	0	0
			32	15	14	2	1		
3	J	1	Total	C	H	N	S	0	0
			32	15	14	2	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	54	Total	O	0	0
			54	54		
4	B	67	Total	O	0	0
			67	67		

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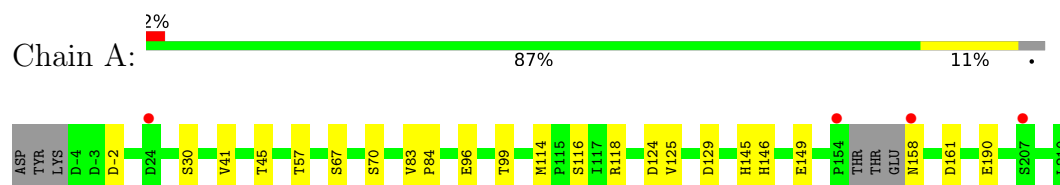
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	64	Total 64	O 64	0	0
4	D	75	Total 75	O 75	0	0
4	E	52	Total 52	O 52	0	0
4	F	67	Total 67	O 67	0	0
4	G	67	Total 67	O 67	0	0
4	H	62	Total 62	O 62	0	0
4	I	70	Total 70	O 70	0	0
4	J	78	Total 78	O 78	0	0

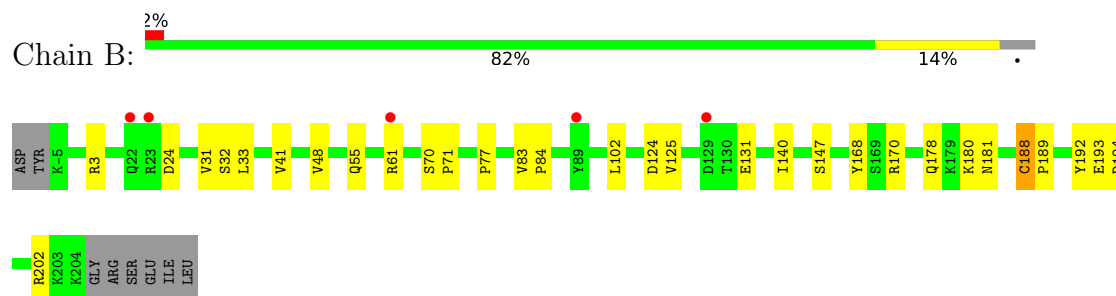
3 Residue-property plots [i](#)

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.

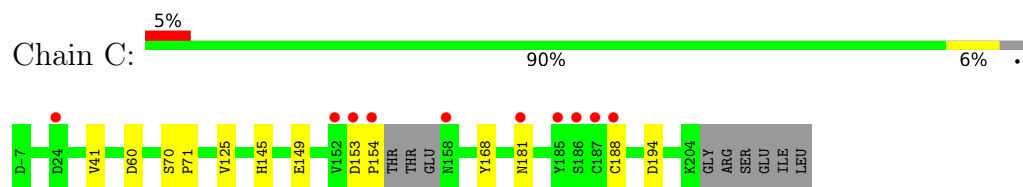
- Molecule 1: Acetylcholine-binding protein



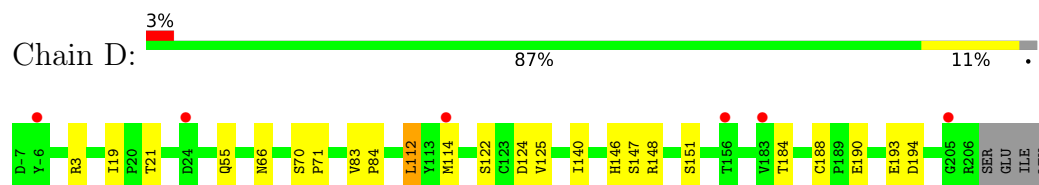
- Molecule 1: Acetylcholine-binding protein



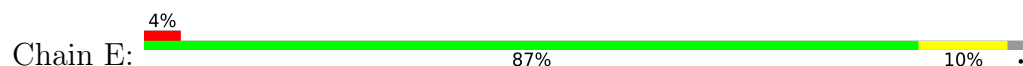
- Molecule 1: Acetylcholine-binding protein

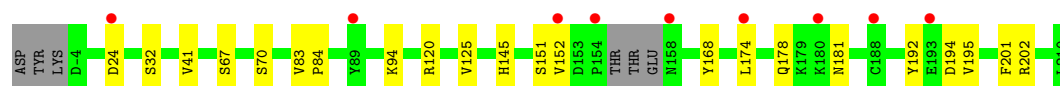


- Molecule 1: Acetylcholine-binding protein

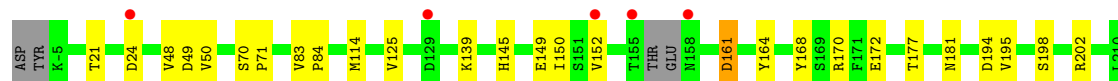
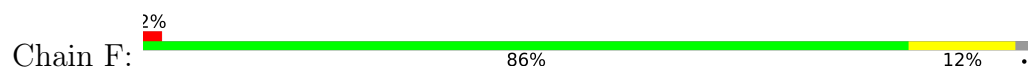


- Molecule 1: Acetylcholine-binding protein

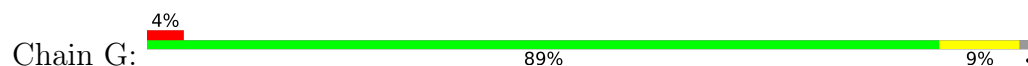




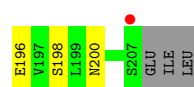
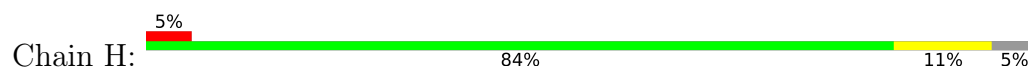
- Molecule 1: Acetylcholine-binding protein



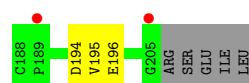
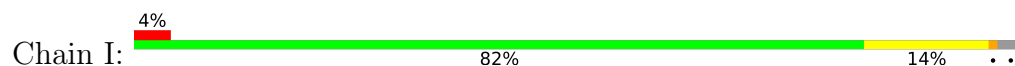
- Molecule 1: Acetylcholine-binding protein



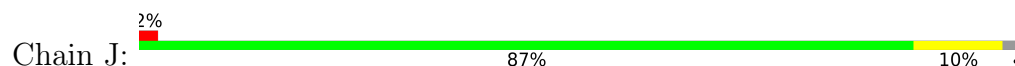
- Molecule 1: Acetylcholine-binding protein



- Molecule 1: Acetylcholine-binding protein



- Molecule 1: Acetylcholine-binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	85.84Å 127.60Å 122.47Å 90.00° 110.45° 90.00°	Depositor
Resolution (Å)	49.98 – 1.85 49.98 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.98-1.85) 93.8 (49.98-1.85)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.88 (at 1.74Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.189 , 0.225 0.198 , 0.229	Depositor DCC
R_{free} test set	2011 reflections (0.83%)	wwPDB-VP
Wilson B-factor (Å ²)	18.6	Xtriage
Anisotropy	0.098	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 33.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.011 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17593	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.29% 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: PO4, 4P6

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	0.42	0/1713	0.73	0/2339
1	B	0.46	0/1694	0.76	2/2317 (0.1%)
1	C	0.45	0/1685	0.71	0/2304
1	D	0.45	0/1711	0.74	0/2342
1	E	0.40	0/1676	0.71	0/2294
1	F	0.42	0/1695	0.74	0/2319
1	G	0.45	0/1708	0.74	0/2335
1	H	0.46	0/1689	0.75	0/2305
1	I	0.45	0/1673	0.78	0/2291
1	J	0.47	0/1696	0.73	0/2319
All	All	0.44	0/16940	0.74	2/23165 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	188	CYS	CA-C-N	5.28	126.44	119.84
1	B	188	CYS	C-N-CA	5.28	126.44	119.84

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	1674	0	1590	25	0
1	B	1655	0	1568	24	0
1	C	1645	0	1548	10	0
1	D	1674	0	1564	15	0
1	E	1641	0	1538	19	0
1	F	1660	0	1560	21	0
1	G	1670	0	1570	15	0
1	H	1651	0	1584	19	0
1	I	1637	0	1534	26	0
1	J	1660	0	1568	19	0
2	A	10	0	0	2	0
2	B	5	0	0	0	0
2	C	5	0	0	1	0
2	D	5	0	0	0	0
2	G	10	0	0	2	0
2	I	10	0	0	2	0
2	J	5	0	0	0	0
3	A	18	14	14	3	0
3	B	18	14	14	4	0
3	C	18	14	14	0	0
3	D	18	14	14	3	0
3	E	18	14	14	7	0
3	F	18	14	14	2	0
3	G	18	14	14	1	0
3	H	18	14	14	1	0
3	I	18	14	14	3	0
3	J	18	14	14	4	0
4	A	54	0	0	0	0
4	B	67	0	0	3	0
4	C	64	0	0	1	0
4	D	75	0	0	1	0
4	E	52	0	0	1	0
4	F	67	0	0	0	0
4	G	67	0	0	0	0
4	H	62	0	0	1	0
4	I	70	0	0	3	0
4	J	78	0	0	1	0
All	All	17453	140	15764	193	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 6.

All (193) 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:E:152:VAL:HG12	1:E:195:VAL:HG23	1.59	0.84
1:A:146[B]:HIS:CD2	1:A:190:GLU:HG3	2.13	0.83
1:F:139:LYS:HE2	1:F:181:ASN:ND2	2.01	0.76
1:C:149:GLU:OE2	1:D:3:ARG:NH2	2.19	0.75
1:H:181:ASN:HB3	1:H:194:ASP:OD1	1.88	0.73
1:H:94:LYS:HE3	1:I:96:GLU:HG2	1.70	0.73
1:F:177:THR:HG22	1:F:198:SER:HB2	1.69	0.72
3:E:301:4P6:S18	3:E:301:4P6:H27	2.29	0.72
1:F:177:THR:CG2	1:F:198:SER:HB2	2.20	0.72
1:A:146[B]:HIS:CG	1:A:190:GLU:HG3	2.25	0.71
1:B:180:LYS:NZ	4:B:403:HOH:O	2.24	0.70
1:J:139:LYS:HE2	1:J:181:ASN:ND2	2.07	0.69
1:I:181:ASN:HB3	1:I:194:ASP:OD1	1.92	0.69
1:H:139:LYS:HG2	1:H:196:GLU:OE1	1.93	0.69
1:F:170:ARG:HD2	1:J:45:THR:HA	1.78	0.66
1:H:137:ARG:NH1	1:H:198:SER:OG	2.29	0.65
1:F:181:ASN:HB3	1:F:194:ASP:OD1	1.96	0.65
1:H:183:VAL:HG13	1:H:192:TYR:HB2	1.78	0.65
1:C:60:ASP:OD2	4:C:401:HOH:O	2.15	0.64
1:B:83:VAL:HG13	1:B:84:PRO:HD2	1.80	0.64
1:I:70:SER:HB2	1:I:71:PRO:HD2	1.79	0.64
1:A:149:GLU:OE2	1:B:3:ARG:NH2	2.24	0.64
1:H:70:SER:HB2	1:H:71:PRO:HD2	1.80	0.64
1:G:139:LYS:HE2	1:G:181:ASN:ND2	2.13	0.63
1:I:47:GLU:OE1	1:I:120:ARG:NH2	2.32	0.62
3:D:302:4P6:C13	3:D:302:4P6:S18	2.87	0.62
1:A:83:VAL:HG13	1:A:84:PRO:HD2	1.81	0.62
1:J:152:VAL:HG21	1:J:194:ASP:HA	1.82	0.61
1:G:127:GLY:O	1:G:130:THR:HB	2.01	0.60
1:J:83:VAL:HG13	1:J:84:PRO:HD2	1.83	0.60
1:A:45:THR:HG22	1:B:170:ARG:NH1	2.17	0.60
3:A:303:4P6:H26	3:A:303:4P6:C3	2.31	0.59
3:A:303:4P6:S18	3:A:303:4P6:H28	2.42	0.59
1:D:124:ASP:HB2	1:E:168:TYR:CE1	2.37	0.59
1:F:48:VAL:HG12	1:F:50:VAL:HG13	1.83	0.59
1:I:41:VAL:HG13	1:I:125:VAL:HG11	1.84	0.59
1:H:183:VAL:CG1	1:H:192:TYR:HB2	2.32	0.59
1:D:70:SER:HB2	1:D:71:PRO:HD2	1.84	0.59
1:H:183:VAL:HG12	1:H:192:TYR:O	2.03	0.59
3:A:303:4P6:S18	3:A:303:4P6:C13	2.91	0.59
1:F:83:VAL:HG13	1:F:84:PRO:HD2	1.83	0.58
1:E:174:LEU:HD11	1:E:202:ARG:HG2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:112:LEU:HB3	1:D:114:MET:HE3	1.84	0.58
1:I:152:VAL:HG12	1:I:195:VAL:HG23	1.84	0.58
1:D:66:ASN:ND2	4:D:401:HOH:O	2.32	0.57
1:A:41:VAL:CG1	1:A:125:VAL:HG11	2.35	0.57
3:B:302:4P6:S18	3:B:302:4P6:C13	2.93	0.56
3:I:303:4P6:C12	1:J:114:MET:HE1	2.35	0.56
3:J:302:4P6:S18	3:J:302:4P6:C13	2.95	0.55
1:G:70:SER:HB2	1:G:71:PRO:HD2	1.90	0.54
1:A:114:MET:HE1	3:E:301:4P6:C12	2.37	0.54
1:I:7:LEU:O	1:I:11:ARG:HG3	2.07	0.54
3:F:301:4P6:C13	3:F:301:4P6:S18	2.96	0.54
1:F:21:THR:OG1	1:G:-2:ASP:HB3	2.06	0.54
1:H:122:SER:HB2	1:I:37:ASN:ND2	2.22	0.54
1:H:94:LYS:HE3	1:I:96:GLU:CG	2.37	0.54
1:F:152:VAL:HG12	1:F:195:VAL:HG23	1.89	0.53
1:A:-2:ASP:HB2	1:E:24:ASP:HA	1.91	0.53
1:C:181:ASN:HB3	1:C:194:ASP:OD2	2.08	0.53
3:D:302:4P6:S18	3:D:302:4P6:H28	2.50	0.52
1:E:32:SER:HA	1:E:178:GLN:HE22	1.74	0.52
3:E:301:4P6:S18	3:E:301:4P6:C13	2.97	0.52
1:A:114:MET:CE	3:E:301:4P6:C12	2.88	0.51
1:A:146[B]:HIS:CE1	1:A:190:GLU:CG	2.94	0.51
3:B:302:4P6:S18	3:B:302:4P6:H28	2.51	0.51
1:D:125:VAL:O	1:D:125:VAL:HG12	2.09	0.51
1:J:110:GLU:OE2	4:J:401:HOH:O	2.19	0.51
1:A:146[B]:HIS:CE1	1:A:190:GLU:HG3	2.46	0.51
1:I:157:GLU:HG2	1:I:158:ASN:N	2.26	0.51
1:B:181:ASN:HB3	1:B:194:ASP:OD1	2.12	0.50
1:D:146:HIS:CE1	1:D:148:ARG:HB2	2.46	0.50
1:F:70:SER:HB2	1:F:71:PRO:HD2	1.93	0.50
1:I:60:ASP:OD2	4:I:401:HOH:O	2.19	0.50
1:I:70:SER:HB2	1:I:71:PRO:CD	2.41	0.50
1:A:146[B]:HIS:NE2	1:A:190:GLU:HG3	2.27	0.50
1:B:83:VAL:CG1	1:B:84:PRO:HD2	2.41	0.50
3:I:303:4P6:C12	1:J:114:MET:CE	2.90	0.50
1:D:83:VAL:HG13	1:D:84:PRO:HD2	1.93	0.50
1:I:154:PRO:HG3	1:I:180:LYS:HB2	1.94	0.50
1:B:125:VAL:HG12	1:B:125:VAL:O	2.11	0.50
3:B:302:4P6:H26	3:B:302:4P6:C3	2.40	0.50
1:H:70:SER:HB2	1:H:71:PRO:CD	2.41	0.50
1:B:170:ARG:NH2	4:B:405:HOH:O	2.43	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:70:SER:HB2	1:J:71:PRO:HD2	1.94	0.49
1:E:174:LEU:HD13	1:E:201:PHE:HA	1.94	0.49
1:F:114:MET:HE1	3:J:302:4P6:C12	2.42	0.49
1:B:70:SER:HB2	1:B:71:PRO:HD2	1.94	0.49
1:E:120:ARG:NH1	4:E:402:HOH:O	2.44	0.49
1:D:55:GLN:HG3	1:D:114:MET:SD	2.53	0.49
1:G:130:THR:HG22	1:G:132:SER:H	1.78	0.49
3:F:301:4P6:C3	3:F:301:4P6:H26	2.42	0.48
1:I:94:LYS:HE3	1:J:96:GLU:HG3	1.95	0.48
1:E:174:LEU:HD12	1:E:174:LEU:N	2.28	0.48
3:D:302:4P6:C3	3:D:302:4P6:H26	2.42	0.48
1:J:70:SER:HB2	1:J:71:PRO:CD	2.43	0.48
1:G:158:ASN:OD1	1:G:177:THR:HG22	2.14	0.48
1:H:151:SER:HB2	1:H:193:GLU:OE1	2.13	0.48
1:A:41:VAL:CG1	1:A:125:VAL:CG1	2.91	0.48
1:A:96:GLU:HG3	1:E:94:LYS:HD2	1.95	0.48
1:J:181:ASN:ND2	1:J:194:ASP:OD1	2.40	0.48
1:I:125:VAL:HG12	1:I:125:VAL:O	2.13	0.47
1:D:70:SER:HB2	1:D:71:PRO:CD	2.44	0.47
1:E:181:ASN:HB3	1:E:194:ASP:OD1	2.13	0.47
1:A:124:ASP:HB2	1:B:168:TYR:CE1	2.49	0.47
1:A:125:VAL:HG12	1:A:125:VAL:O	2.13	0.47
1:A:83:VAL:CG1	1:A:84:PRO:HD2	2.44	0.47
1:B:31:VAL:HA	1:B:55:GLN:O	2.14	0.47
1:H:30:SER:HB3	1:H:155:THR:HG22	1.96	0.47
1:H:83:VAL:HG13	1:H:84:PRO:HD2	1.95	0.47
1:G:139:LYS:HE3	1:G:196:GLU:OE1	2.15	0.47
1:E:192:TYR:OH	3:E:301:4P6:H21	2.14	0.47
3:H:301:4P6:C13	3:H:301:4P6:S18	3.03	0.47
1:E:125:VAL:HG12	1:E:125:VAL:O	2.14	0.46
1:I:163:GLU:HG2	4:I:464:HOH:O	2.14	0.46
1:J:152:VAL:CG2	1:J:194:ASP:HA	2.44	0.46
2:A:301:PO4:O2	1:E:145:HIS:HE1	1.98	0.46
3:I:303:4P6:S18	3:I:303:4P6:C13	3.04	0.46
1:J:188:CYS:HB3	1:J:189:PRO:HD2	1.97	0.46
1:B:147:SER:HB3	1:B:193:GLU:HG3	1.97	0.46
3:G:303:4P6:C13	3:G:303:4P6:S18	3.04	0.46
1:I:145:HIS:HE1	2:I:302:PO4:O2	1.99	0.46
1:A:114:MET:HE1	3:E:301:4P6:C11	2.46	0.45
1:B:193:GLU:O	4:B:402:HOH:O	2.20	0.45
1:F:83:VAL:CG1	1:F:84:PRO:HD2	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:124:ASP:HB2	1:H:168:TYR:CE1	2.51	0.45
1:C:125:VAL:O	1:C:125:VAL:HG12	2.16	0.45
1:A:41:VAL:HG12	1:A:125:VAL:HG11	1.98	0.45
1:E:67:SER:HA	1:E:70:SER:OG	2.17	0.44
1:A:30:SER:HB2	1:A:57:THR:OG1	2.17	0.44
1:D:140:ILE:O	1:D:194:ASP:HB2	2.17	0.44
1:A:145:HIS:HE1	2:A:302:PO4:O4	2.00	0.44
1:I:179:LYS:HE2	1:I:179:LYS:HB2	1.73	0.44
1:B:41:VAL:HG22	1:B:48:VAL:HG12	2.00	0.44
1:F:168:TYR:CE1	1:J:124:ASP:HB2	2.53	0.44
1:G:152:VAL:HG12	1:G:195:VAL:HG23	2.00	0.44
1:J:139:LYS:HE2	1:J:181:ASN:HD21	1.81	0.44
1:A:-2:ASP:CB	1:E:24:ASP:HA	2.48	0.43
1:D:19:ILE:HD12	1:D:21:THR:HG23	2.00	0.43
1:G:145:HIS:HE1	2:G:302:PO4:O4	2.00	0.43
1:J:181:ASN:HD22	1:J:194:ASP:CG	2.25	0.43
1:I:157:GLU:HG2	1:I:158:ASN:H	1.83	0.43
1:I:180:LYS:HG2	1:I:181:ASN:N	2.33	0.43
1:H:125:VAL:O	1:H:125:VAL:HG12	2.18	0.43
1:G:70:SER:HB2	1:G:71:PRO:CD	2.48	0.43
3:J:302:4P6:C3	3:J:302:4P6:H26	2.49	0.43
1:E:41:VAL:CG1	1:E:125:VAL:HG11	2.48	0.43
1:G:130:THR:HG22	1:G:133:GLY:H	1.84	0.43
1:H:145:HIS:HE1	2:I:301:PO4:O2	2.02	0.43
1:I:152:VAL:HG12	1:I:195:VAL:CG2	2.49	0.43
1:F:145:HIS:HE1	2:G:301:PO4:O1	2.00	0.43
1:I:55:GLN:HA	1:I:114:MET:HG3	2.01	0.43
1:F:161:ASP:HB2	1:F:164:TYR:HD2	1.84	0.42
1:C:41:VAL:CG1	1:C:125:VAL:HG11	2.50	0.42
1:F:149:GLU:OE2	1:G:3:ARG:NH2	2.38	0.42
1:G:139:LYS:HG3	1:G:196:GLU:CD	2.44	0.42
3:E:301:4P6:C3	3:E:301:4P6:H26	2.47	0.42
1:B:124:ASP:HB2	1:C:168:TYR:CE1	2.55	0.42
1:F:145:HIS:HB2	1:F:150:ILE:HD12	2.01	0.42
1:C:153:ASP:HA	1:C:154:PRO:HD3	1.88	0.42
1:I:41:VAL:CG1	1:I:125:VAL:HG11	2.49	0.42
1:B:131:GLU:O	1:B:202:ARG:NH1	2.50	0.42
1:B:188:CYS:HA	1:B:189:PRO:HD3	1.86	0.42
1:C:41:VAL:CG1	1:C:125:VAL:CG1	2.98	0.42
1:F:125:VAL:O	1:F:125:VAL:HG12	2.18	0.42
1:B:32:SER:HA	1:B:178:GLN:OE1	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:152:VAL:O	1:J:152:VAL:HG23	2.20	0.42
1:C:145:HIS:HE1	2:C:301:PO4:O2	2.02	0.42
1:I:32:SER:HA	1:I:178:GLN:HE22	1.85	0.42
1:C:70:SER:HB2	1:C:71:PRO:CD	2.50	0.41
1:B:140:ILE:O	1:B:194:ASP:HB2	2.19	0.41
1:B:192:TYR:OH	3:B:302:4P6:H21	2.20	0.41
1:F:172:GLU:OE2	1:F:202:ARG:NE	2.50	0.41
1:H:175:ASP:HB3	1:H:200[B]:ASN:HD22	1.85	0.41
1:D:151:SER:HB2	1:D:193:GLU:OE1	2.20	0.41
1:F:114:MET:CE	3:J:302:4P6:C12	2.98	0.41
1:E:83:VAL:HG13	1:E:84:PRO:HD2	2.01	0.41
1:J:99:THR:HG23	1:J:116:SER:HB3	2.01	0.41
1:A:118:ARG:HH12	1:E:120:ARG:HH21	1.69	0.41
1:A:99:THR:HG23	1:A:116:SER:HB3	2.03	0.41
1:B:147:SER:CB	1:B:193:GLU:HG3	2.51	0.41
1:H:15:ARG:HD3	4:H:458:HOH:O	2.20	0.41
1:E:41:VAL:CG1	1:E:125:VAL:CG1	2.99	0.41
1:G:153:ASP:HA	1:G:154:PRO:HD3	1.88	0.41
1:I:167:GLN:HG2	4:I:457:HOH:O	2.21	0.41
1:B:70:SER:HB2	1:B:71:PRO:CD	2.50	0.40
1:B:77:PRO:HA	1:B:102:LEU:HD23	2.02	0.40
1:A:67:SER:HA	1:A:70:SER:OG	2.20	0.40
1:D:147:SER:HB3	1:D:193:GLU:HG3	2.04	0.40
1:D:184:THR:HG23	1:D:190:GLU:O	2.22	0.40
1:I:139:LYS:HG2	1:I:196:GLU:HG2	2.02	0.40
1:B:33:LEU:H	1:B:178:GLN:HE22	1.69	0.40
1:F:48:VAL:HG12	1:F:49:ASP:N	2.35	0.40
1:J:155:THR:OG1	1:J:156:THR:N	2.54	0.40

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	209/218 (96%)	208 (100%)	1 (0%)	0	100	100
1	B	209/218 (96%)	208 (100%)	1 (0%)	0	100	100
1	C	206/218 (94%)	206 (100%)	0	0	100	100
1	D	212/218 (97%)	211 (100%)	1 (0%)	0	100	100
1	E	208/218 (95%)	206 (99%)	2 (1%)	0	100	100
1	F	210/218 (96%)	208 (99%)	1 (0%)	1 (0%)	24	13
1	G	211/218 (97%)	209 (99%)	1 (0%)	1 (0%)	24	13
1	H	205/218 (94%)	205 (100%)	0	0	100	100
1	I	209/218 (96%)	207 (99%)	2 (1%)	0	100	100
1	J	209/218 (96%)	207 (99%)	2 (1%)	0	100	100
All	All	2088/2180 (96%)	2075 (99%)	11 (0%)	2 (0%)	48	35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	161	ASP
1	G	189	PRO

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	193/204 (95%)	190 (98%)	3 (2%)	55	44
1	B	191/204 (94%)	189 (99%)	2 (1%)	68	60
1	C	188/204 (92%)	187 (100%)	1 (0%)	81	77
1	D	191/204 (94%)	188 (98%)	3 (2%)	55	44
1	E	186/204 (91%)	185 (100%)	1 (0%)	81	77
1	F	189/204 (93%)	188 (100%)	1 (0%)	81	77
1	G	191/204 (94%)	189 (99%)	2 (1%)	68	60
1	H	191/204 (94%)	191 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	I	187/204 (92%)	183 (98%)	4 (2%)	47	33
1	J	192/204 (94%)	188 (98%)	4 (2%)	47	33
All	All	1899/2040 (93%)	1878 (99%)	21 (1%)	65	57

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	129	ASP
1	A	158	ASN
1	A	161	ASP
1	B	24	ASP
1	B	61	ARG
1	C	188	CYS
1	D	112	LEU
1	D	122	SER
1	D	188	CYS
1	E	151	SER
1	F	24	ASP
1	G	130	THR
1	G	188	CYS
1	I	11	ARG
1	I	163	GLU
1	I	183	VAL
1	I	186	SER
1	J	24	ASP
1	J	180	LYS
1	J	188	CYS
1	J	196	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	54	GLN
1	A	55	GLN
1	A	69	HIS
1	A	158	ASN
1	B	101	GLN
1	B	200	ASN
1	C	54	GLN

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Mol	Chain	Res	Type
1	C	119	GLN
1	D	54	GLN
1	D	158	ASN
1	E	12	GLN
1	E	119	GLN
1	E	200	ASN
1	F	12	GLN
1	F	54	GLN
1	F	69	HIS
1	F	101	GLN
1	F	145	HIS
1	F	200	ASN
1	G	9	ASN
1	G	119	GLN
1	G	200	ASN
1	H	9	ASN
1	H	22	GLN
1	H	55	GLN
1	H	69	HIS
1	H	101	GLN
1	H	181	ASN
1	I	12	GLN
1	I	101	GLN
1	I	200	ASN
1	J	69	HIS
1	J	200	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

20 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PO4	D	301	-	4,4,4	0.90	0	6,6,6	0.16	0
3	4P6	C	302	-	20,20,20	0.69	0	24,26,26	1.78	6 (25%)
3	4P6	B	302	-	20,20,20	0.65	0	24,26,26	1.63	5 (20%)
2	PO4	G	301	-	4,4,4	0.80	0	6,6,6	0.41	0
3	4P6	H	301	-	20,20,20	0.70	1 (5%)	24,26,26	1.58	7 (29%)
3	4P6	G	303	-	20,20,20	0.85	1 (5%)	24,26,26	1.59	4 (16%)
3	4P6	J	302	-	20,20,20	0.72	0	24,26,26	1.88	8 (33%)
3	4P6	E	301	-	20,20,20	0.63	0	24,26,26	1.71	4 (16%)
2	PO4	A	302	-	4,4,4	0.87	0	6,6,6	0.67	0
3	4P6	F	301	-	20,20,20	0.67	0	24,26,26	1.78	7 (29%)
3	4P6	A	303	-	20,20,20	0.75	0	24,26,26	1.82	6 (25%)
3	4P6	I	303	-	20,20,20	0.70	0	24,26,26	1.66	6 (25%)
3	4P6	D	302	-	20,20,20	0.64	0	24,26,26	2.02	9 (37%)
2	PO4	C	301	-	4,4,4	0.81	0	6,6,6	0.47	0
2	PO4	G	302	-	4,4,4	0.83	0	6,6,6	0.61	0
2	PO4	I	301	-	4,4,4	1.06	0	6,6,6	0.76	0
2	PO4	J	301	-	4,4,4	0.70	0	6,6,6	0.57	0
2	PO4	A	301	-	4,4,4	0.95	0	6,6,6	0.58	0
2	PO4	I	302	-	4,4,4	0.98	0	6,6,6	0.71	0
2	PO4	B	301	-	4,4,4	0.81	0	6,6,6	0.28	0

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	4P6	H	301	-	-	0/8/19/19	0/3/3/3
3	4P6	A	303	-	-	0/8/19/19	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	4P6	C	302	-	-	2/8/19/19	0/3/3/3
3	4P6	G	303	-	-	0/8/19/19	0/3/3/3
3	4P6	J	302	-	-	0/8/19/19	0/3/3/3
3	4P6	B	302	-	-	0/8/19/19	0/3/3/3
3	4P6	I	303	-	-	0/8/19/19	0/3/3/3
3	4P6	E	301	-	-	4/8/19/19	0/3/3/3
3	4P6	D	302	-	-	0/8/19/19	0/3/3/3
3	4P6	F	301	-	-	0/8/19/19	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	301	4P6	C10-N17	2.31	1.30	1.28
3	G	303	4P6	C10-N17	2.30	1.30	1.28

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	302	4P6	C9-C12-C11	-5.09	124.00	130.40
3	A	303	4P6	C9-C12-C11	-4.87	124.27	130.40
3	E	301	4P6	C9-C12-C11	-4.61	124.60	130.40
3	F	301	4P6	C9-C12-C11	-4.59	124.63	130.40
3	B	302	4P6	C9-C12-C11	-4.50	124.74	130.40
3	J	302	4P6	C9-C12-C11	-4.40	124.86	130.40
3	C	302	4P6	C3-C8-C10	3.94	127.19	120.74
3	J	302	4P6	C8-C10-N17	-3.91	114.25	117.97
3	E	301	4P6	C8-C10-N17	-3.89	114.26	117.97
3	G	303	4P6	C9-C12-C11	-3.81	125.61	130.40
3	A	303	4P6	C8-C10-N17	-3.78	114.37	117.97
3	I	303	4P6	C8-C10-N17	-3.75	114.40	117.97
3	D	302	4P6	C8-C10-N17	-3.66	114.48	117.97
3	I	303	4P6	C9-C12-C11	-3.58	125.89	130.40
3	A	303	4P6	C1-C3-C8	-3.52	116.90	120.36
3	C	302	4P6	C9-C12-C11	-3.50	126.00	130.40
3	D	302	4P6	C1-C3-C8	-3.48	116.94	120.36
3	G	303	4P6	C8-C10-N17	-3.33	114.80	117.97
3	H	301	4P6	C9-C12-C11	-3.28	126.28	130.40
3	E	301	4P6	C13-C11-C12	3.20	128.12	124.19
3	C	302	4P6	C6-C8-C10	-3.13	114.98	121.14
3	B	302	4P6	C12-C11-C10	-3.11	113.58	120.17
3	H	301	4P6	C1-C3-C8	-3.10	117.31	120.36
3	F	301	4P6	C8-C10-N17	-3.06	115.06	117.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	301	4P6	C12-C11-C10	-3.00	113.81	120.17
3	E	301	4P6	C12-C11-C10	-2.97	113.87	120.17
3	J	302	4P6	C1-C3-C8	-2.95	117.47	120.36
3	D	302	4P6	C12-C11-C10	-2.92	113.99	120.17
3	J	302	4P6	C15-N17-C10	-2.91	112.05	115.99
3	C	302	4P6	C12-C11-C10	-2.85	114.13	120.17
3	J	302	4P6	C12-C11-C10	-2.80	114.24	120.17
3	H	301	4P6	C12-C11-C10	-2.80	114.25	120.17
3	C	302	4P6	C1-C3-C8	-2.77	117.64	120.36
3	D	302	4P6	C3-C8-C6	2.68	120.60	117.61
3	F	301	4P6	C1-C3-C8	-2.64	117.77	120.36
3	A	303	4P6	C3-C8-C6	2.63	120.54	117.61
3	B	302	4P6	C8-C10-N17	-2.59	115.51	117.97
3	D	302	4P6	C6-C8-C10	-2.55	116.12	121.14
3	H	301	4P6	C8-C10-N17	-2.55	115.54	117.97
3	G	303	4P6	C1-C3-C8	-2.51	117.90	120.36
3	B	302	4P6	C13-C11-C12	2.44	127.20	124.19
3	I	303	4P6	C1-C3-C8	-2.44	117.96	120.36
3	I	303	4P6	C12-C11-C10	-2.37	115.14	120.17
3	D	302	4P6	C13-C11-C12	2.37	127.11	124.19
3	D	302	4P6	C14-C13-C11	-2.29	107.47	112.14
3	J	302	4P6	C3-C8-C6	2.29	120.16	117.61
3	F	301	4P6	C6-C8-C10	-2.28	116.66	121.14
3	G	303	4P6	C14-C13-C11	-2.26	107.53	112.14
3	H	301	4P6	C6-C8-C10	-2.25	116.71	121.14
3	F	301	4P6	C13-C11-C12	2.25	126.96	124.19
3	C	302	4P6	C8-C10-N17	-2.25	115.83	117.97
3	J	302	4P6	C14-C13-C11	-2.24	107.58	112.14
3	H	301	4P6	C3-C8-C10	2.18	124.31	120.74
3	A	303	4P6	C6-C8-C10	-2.17	116.87	121.14
3	J	302	4P6	C6-C8-C10	-2.15	116.91	121.14
3	D	302	4P6	C15-N17-C10	-2.12	113.12	115.99
3	H	301	4P6	C13-C11-C12	2.12	126.80	124.19
3	I	303	4P6	C7-S18-C9	2.09	95.02	91.86
3	B	302	4P6	C1-C3-C8	-2.08	118.32	120.36
3	I	303	4P6	C3-C8-C6	2.04	119.89	117.61
3	F	301	4P6	C14-C13-C11	-2.04	107.98	112.14
3	A	303	4P6	C12-C11-C10	-2.02	115.90	120.17

There are no chirality outliers.

All (6) torsion outliers are listed below:

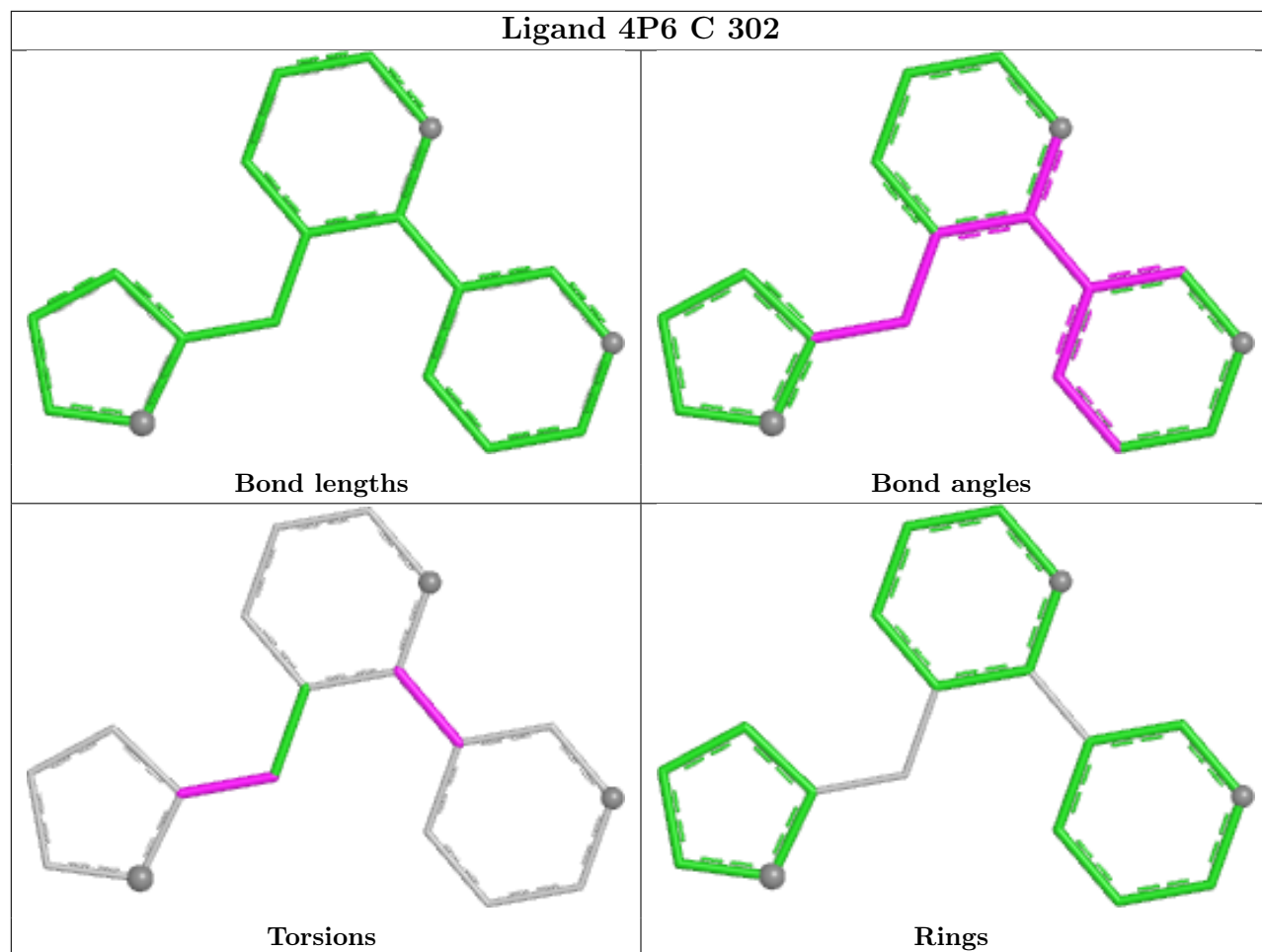
Mol	Chain	Res	Type	Atoms
3	E	301	4P6	N17-C10-C8-C3
3	E	301	4P6	C11-C10-C8-C6
3	E	301	4P6	N17-C10-C8-C6
3	E	301	4P6	C11-C10-C8-C3
3	C	302	4P6	C11-C10-C8-C6
3	C	302	4P6	C11-C12-C9-C4

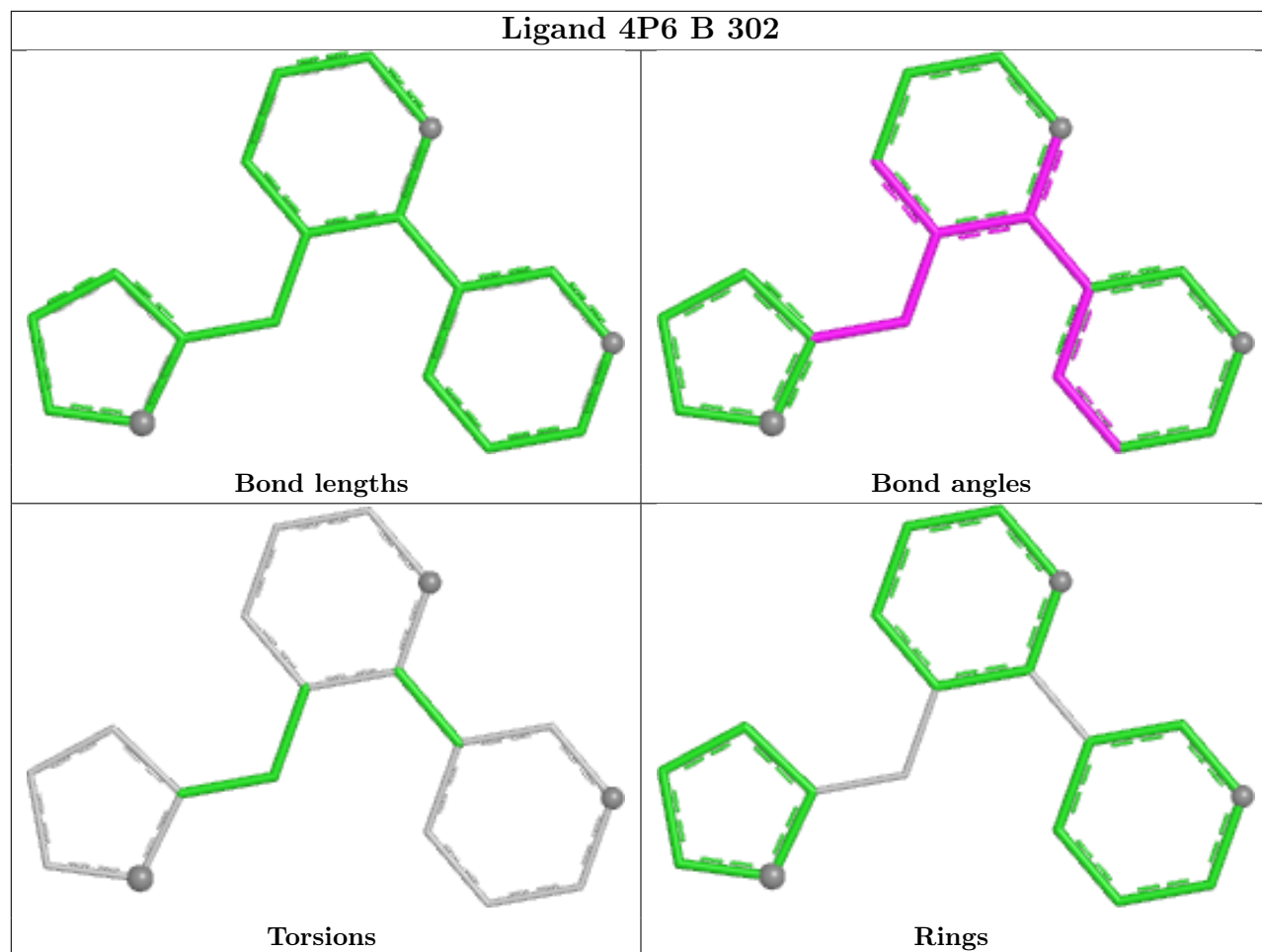
There are no ring outliers.

16 monomers are involved in 35 short contacts:

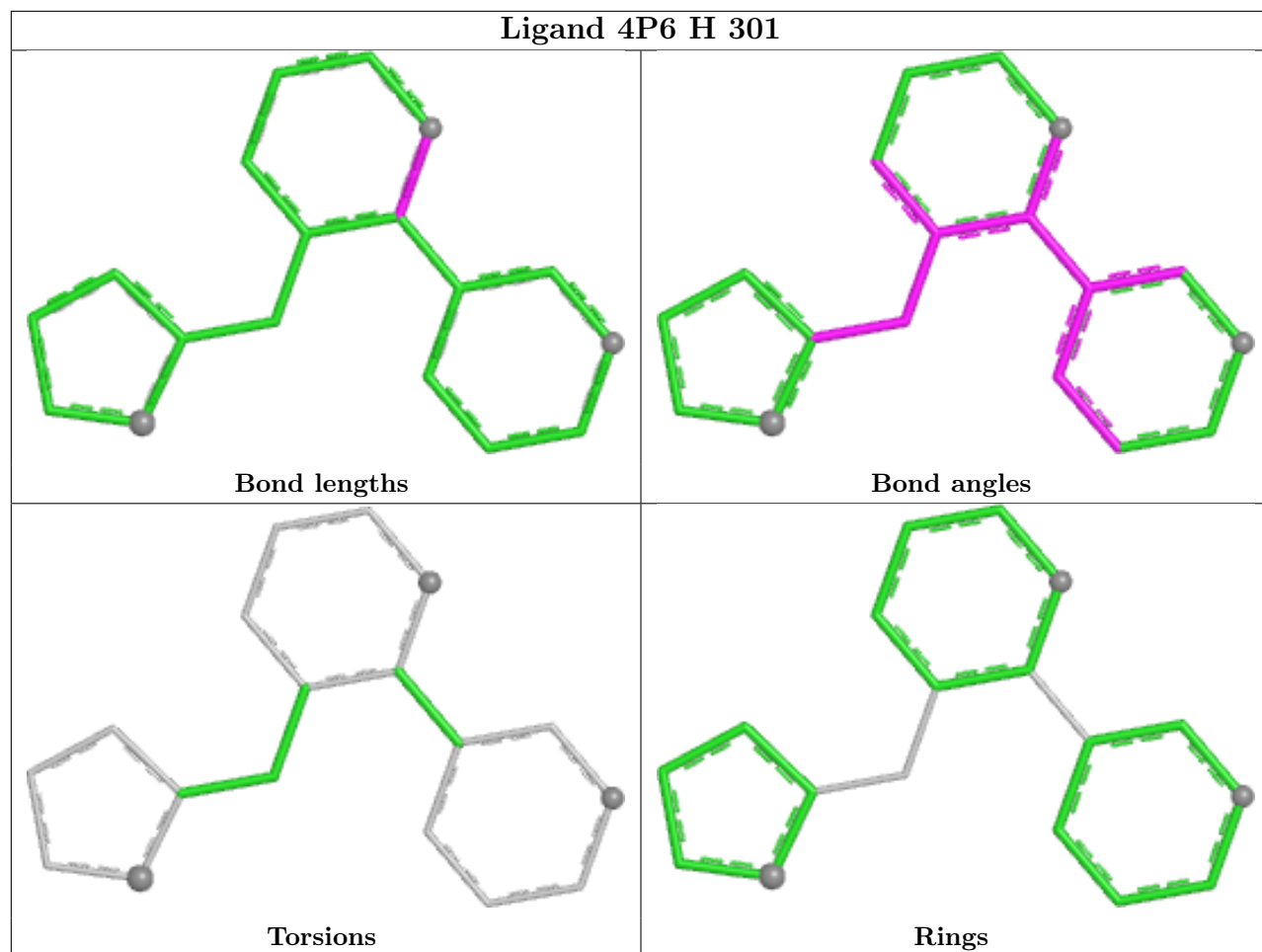
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	302	4P6	4	0
2	G	301	PO4	1	0
3	H	301	4P6	1	0
3	G	303	4P6	1	0
3	J	302	4P6	4	0
3	E	301	4P6	7	0
2	A	302	PO4	1	0
3	F	301	4P6	2	0
3	A	303	4P6	3	0
3	I	303	4P6	3	0
3	D	302	4P6	3	0
2	C	301	PO4	1	0
2	G	302	PO4	1	0
2	I	301	PO4	1	0
2	A	301	PO4	1	0
2	I	302	PO4	1	0

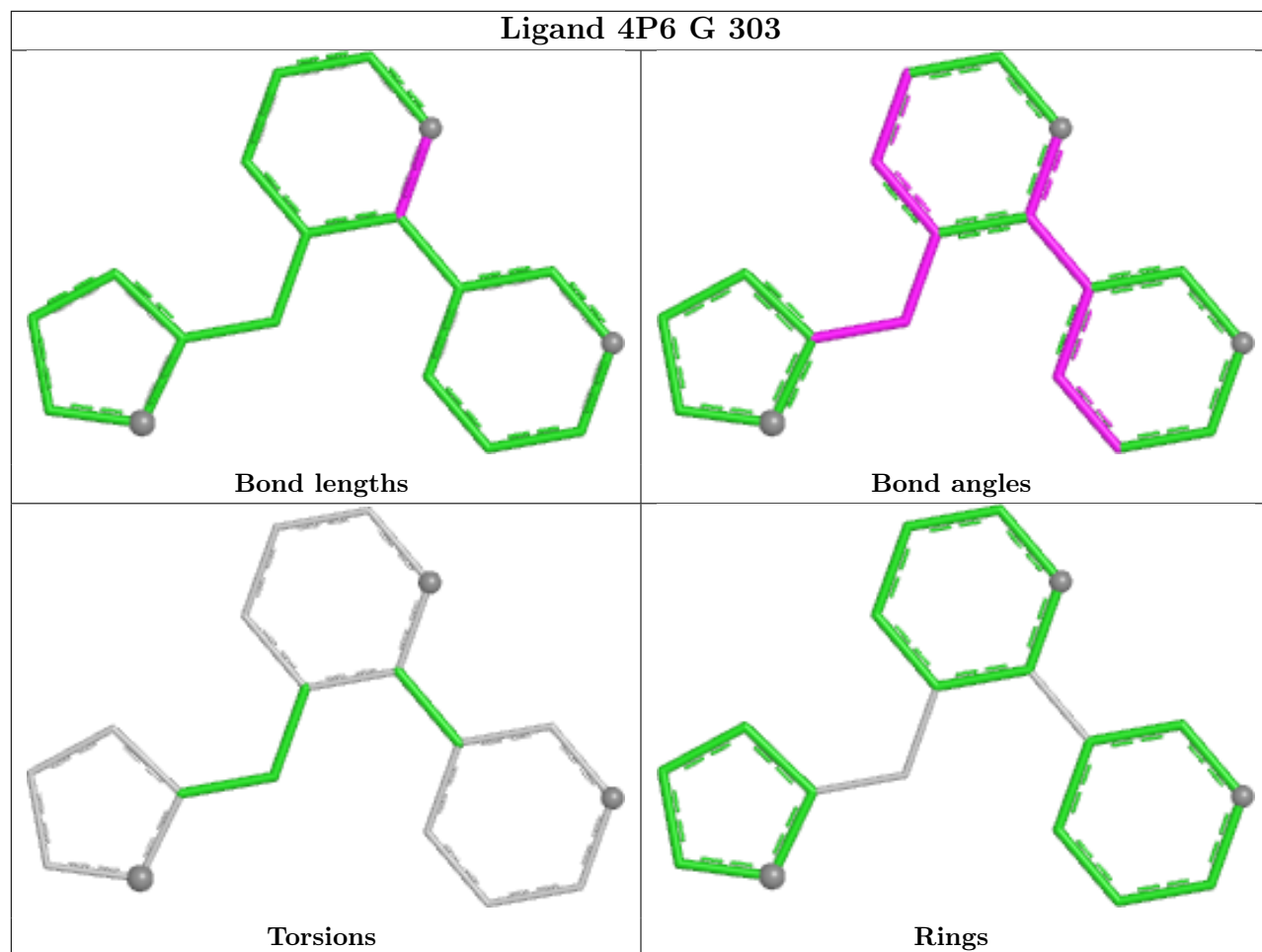
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



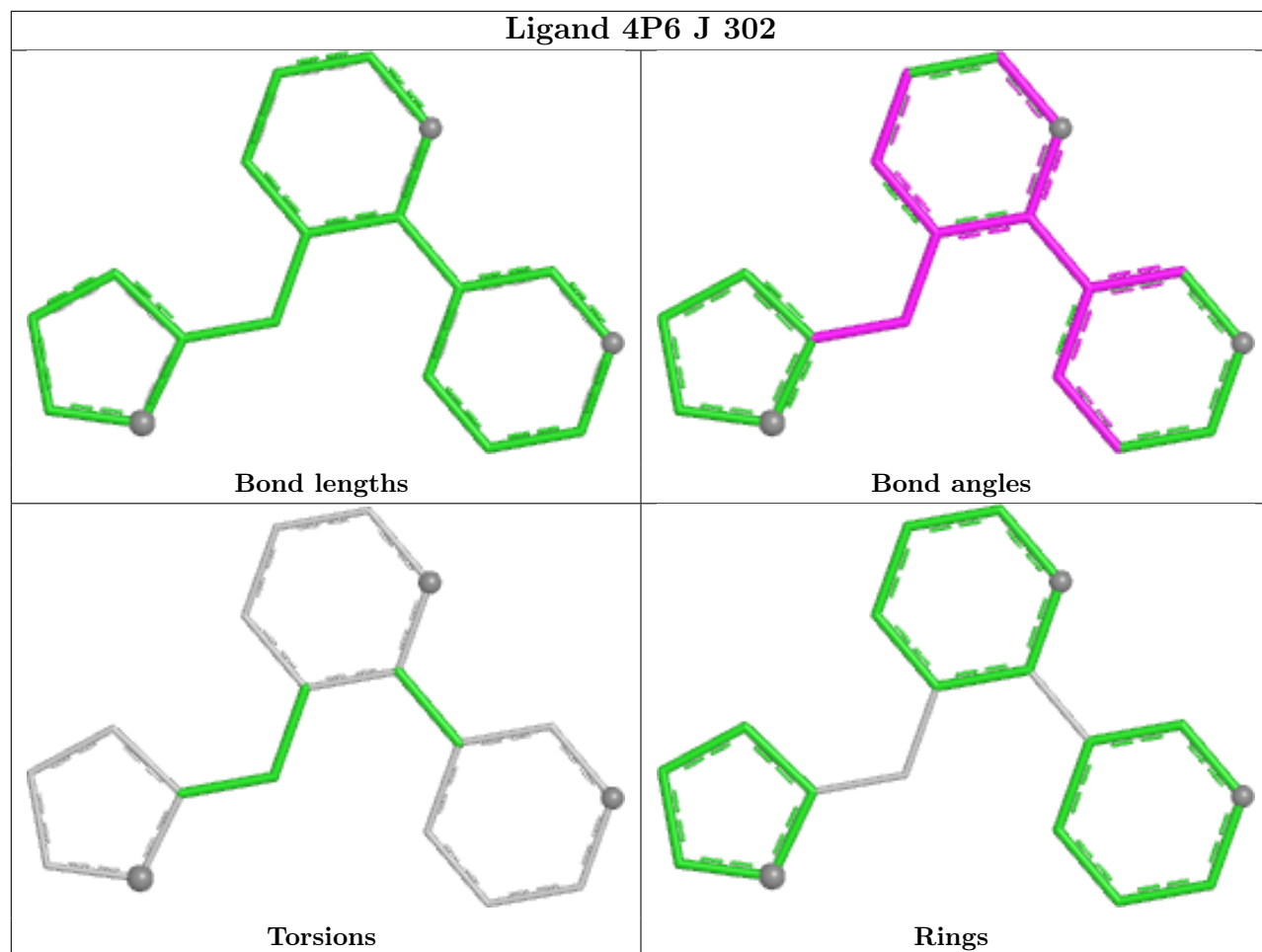


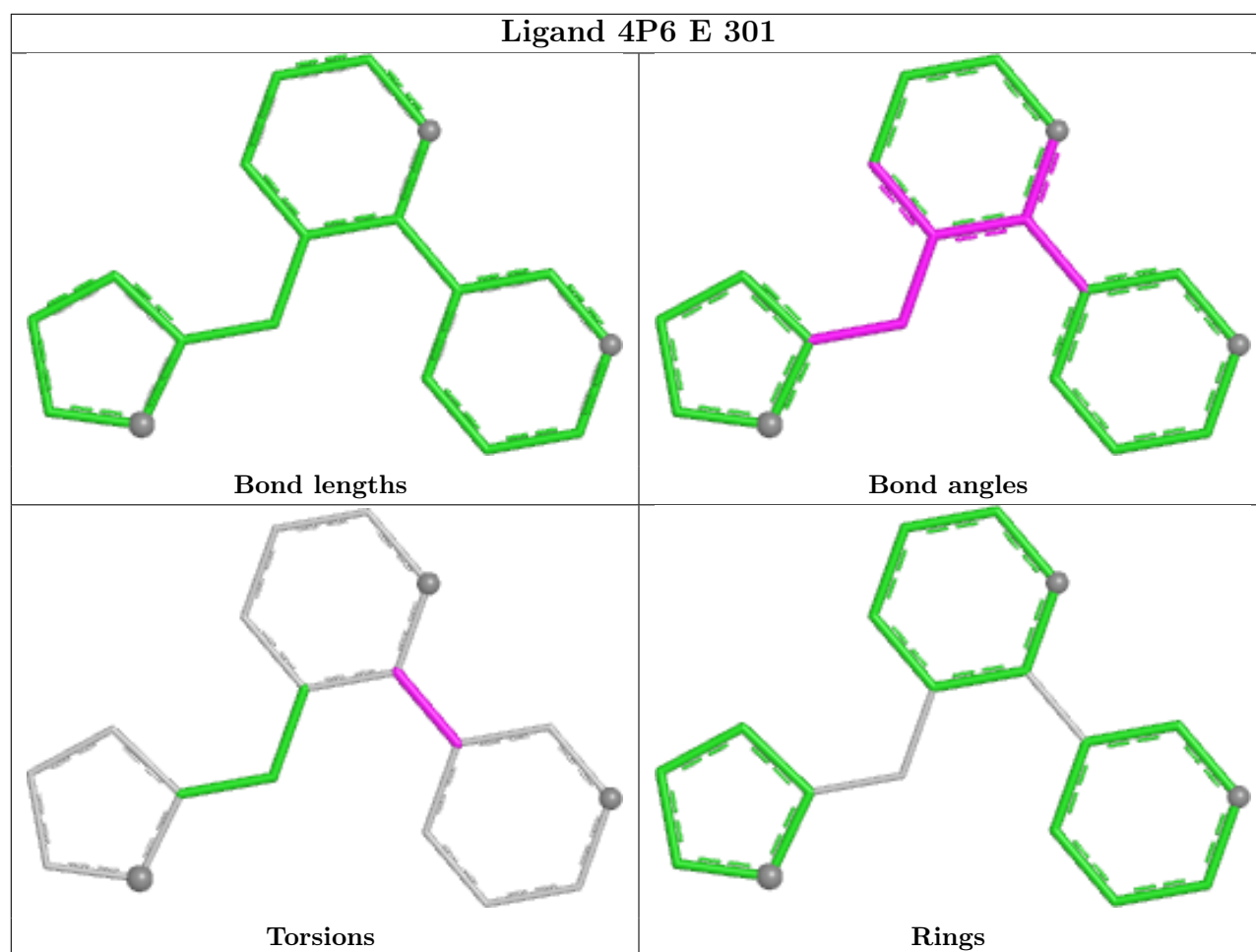
Ligand 4P6 H 301

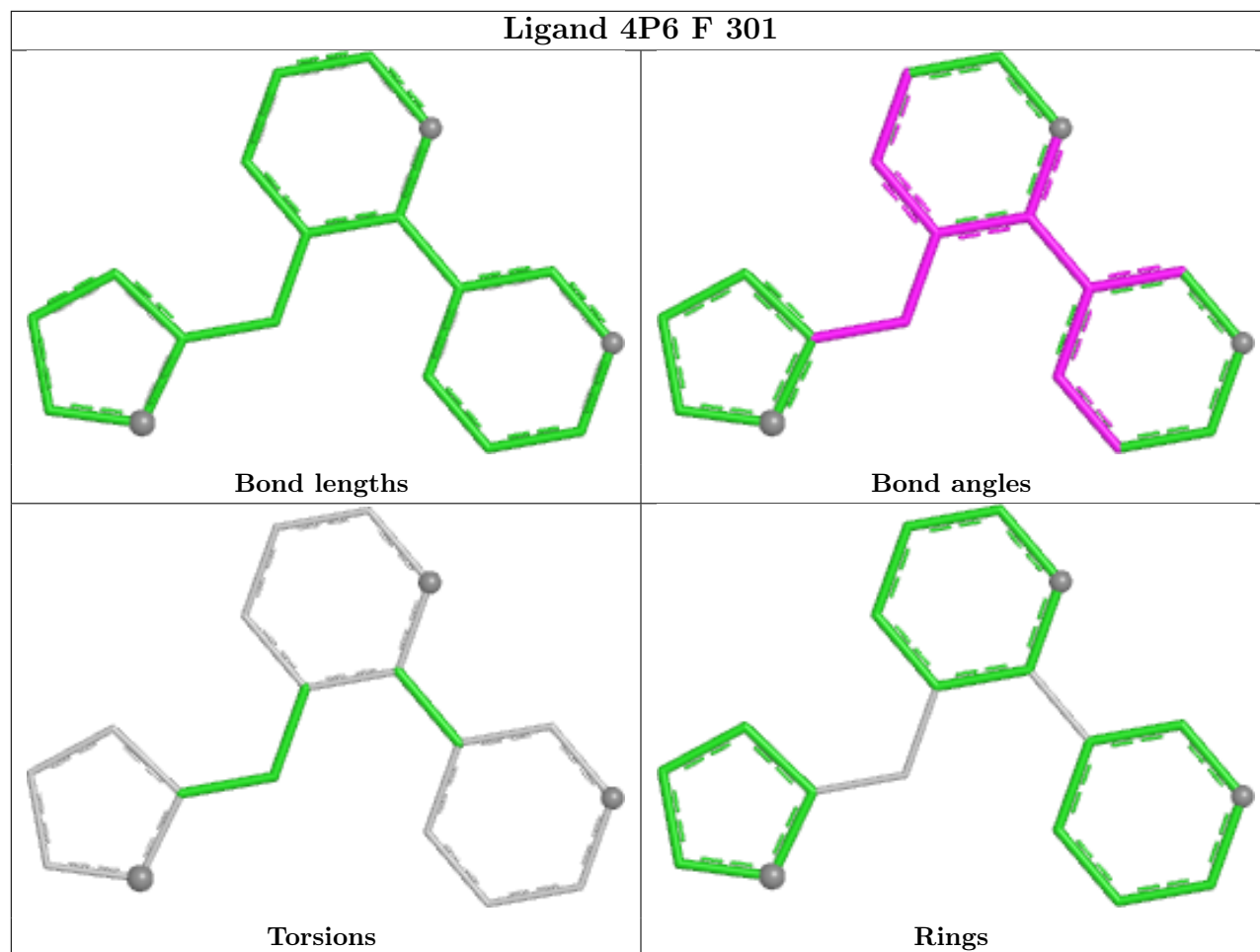




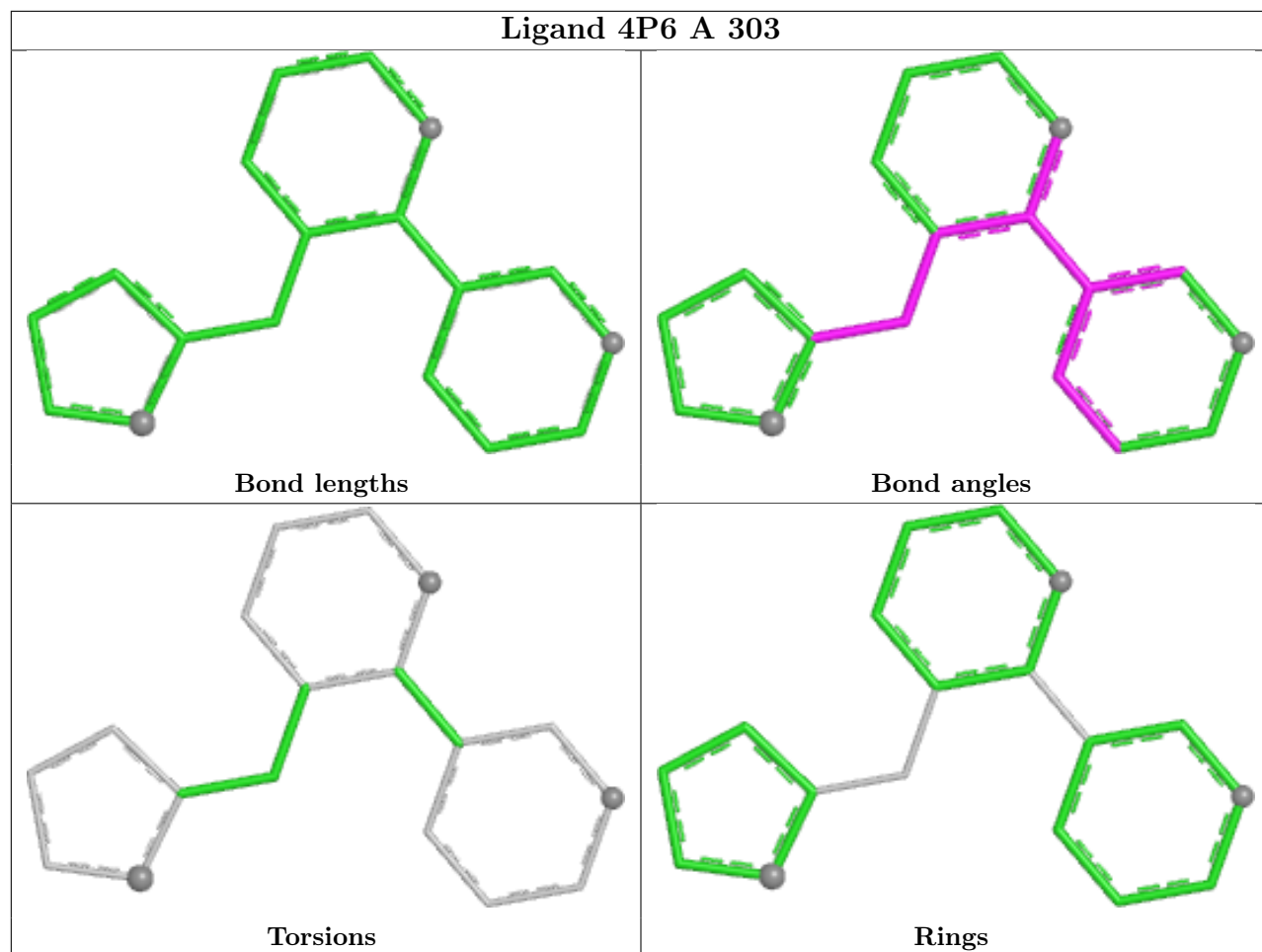
Ligand 4P6 J 302



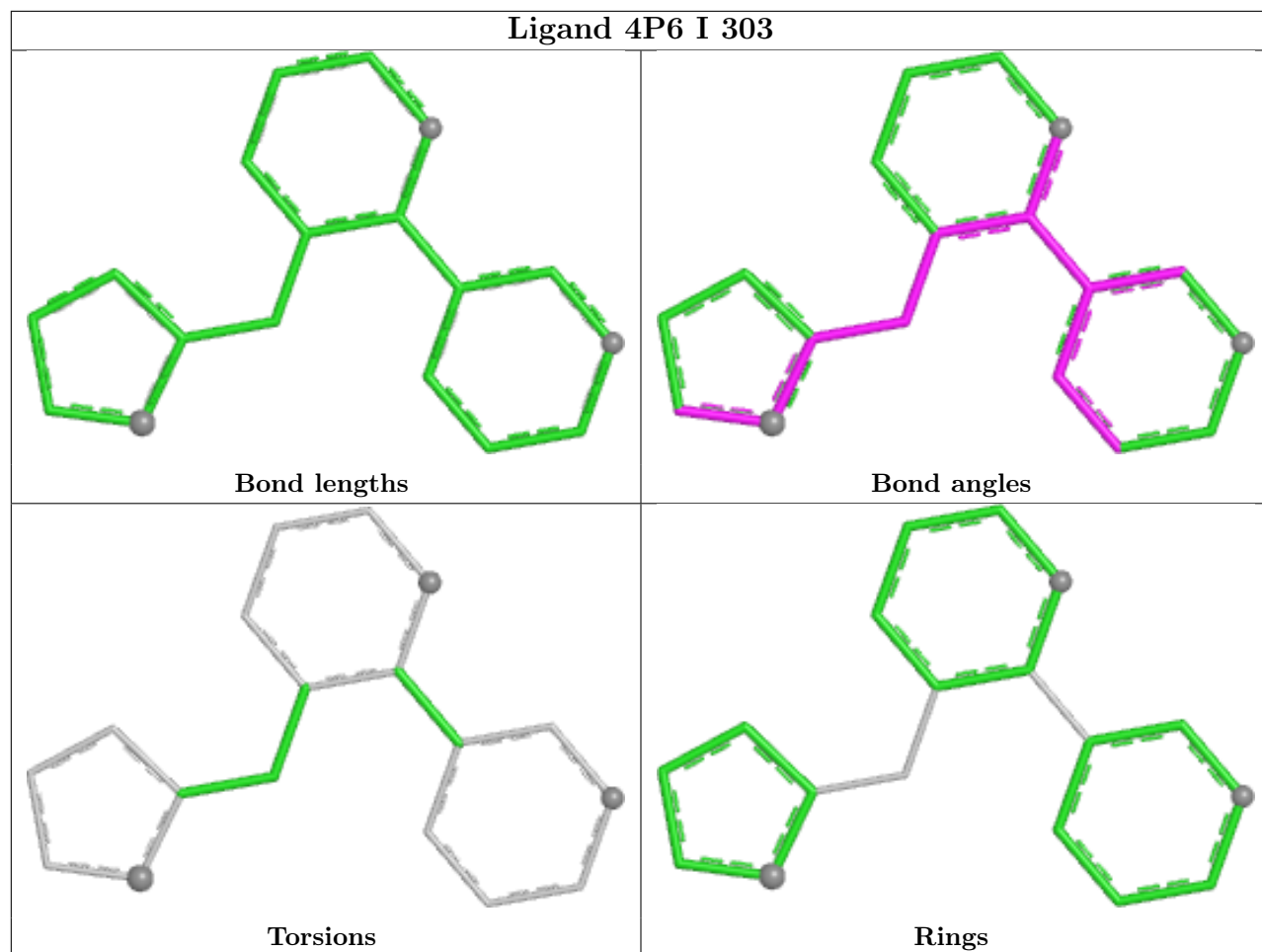


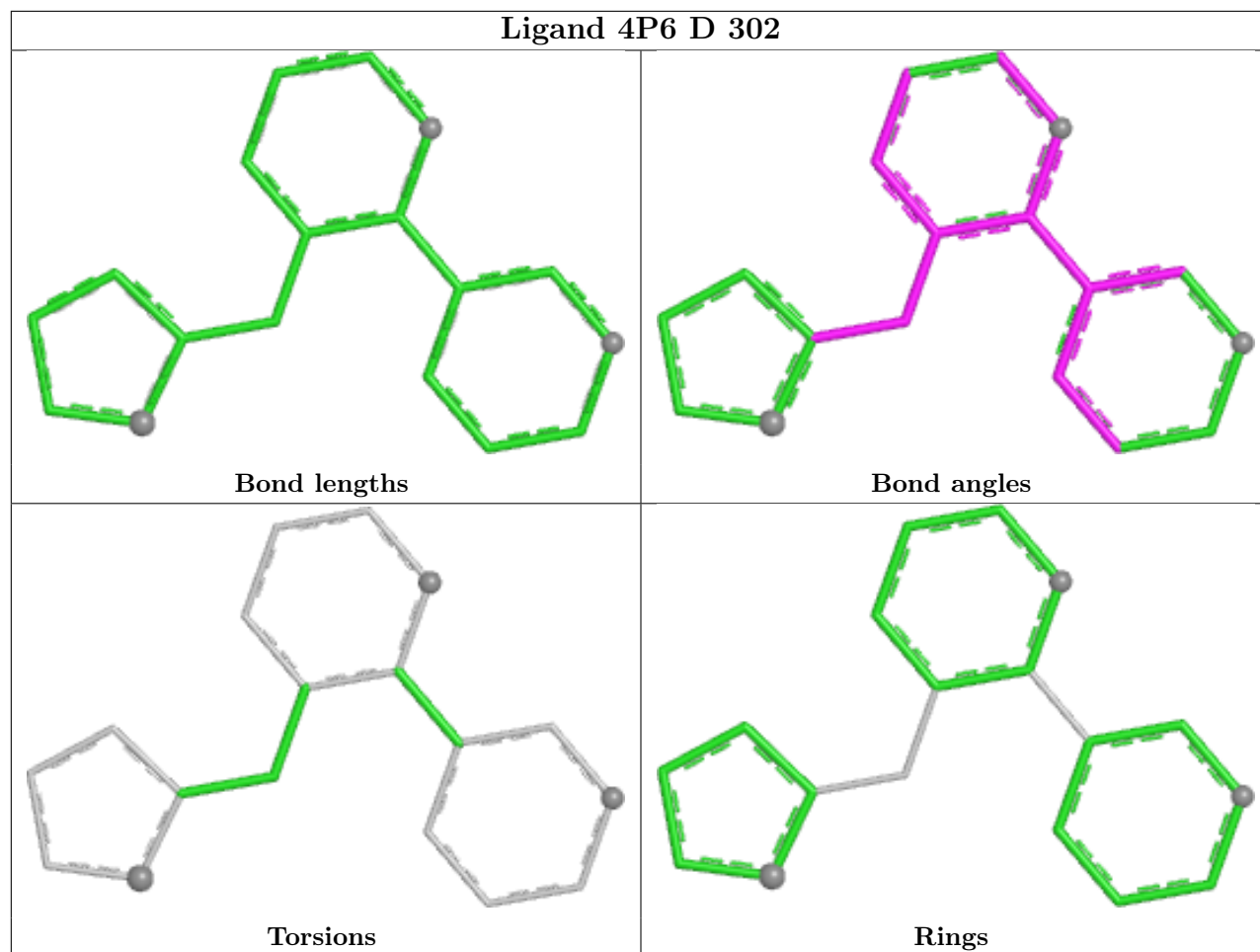


Ligand 4P6 A 303



Ligand 4P6 I 303





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	212/218 (97%)	0.19	4 (1%) 66 70	16, 27, 40, 49	1 (0%)
1	B	210/218 (96%)	0.04	5 (2%) 59 63	14, 24, 37, 51	1 (0%)
1	C	209/218 (95%)	0.17	10 (4%) 35 39	15, 24, 43, 62	1 (0%)
1	D	214/218 (98%)	0.10	6 (2%) 55 58	15, 24, 40, 51	0
1	E	212/218 (97%)	0.27	9 (4%) 40 44	19, 26, 49, 63	0
1	F	214/218 (98%)	0.15	5 (2%) 61 64	16, 25, 43, 51	0
1	G	214/218 (98%)	0.17	8 (3%) 45 49	15, 25, 44, 65	1 (0%)
1	H	208/218 (95%)	0.11	11 (5%) 32 35	15, 22, 41, 61	1 (0%)
1	I	211/218 (96%)	0.06	8 (3%) 44 48	15, 24, 42, 57	0
1	J	211/218 (96%)	-0.01	5 (2%) 59 63	14, 23, 38, 53	0
All	All	2115/2180 (97%)	0.12	71 (3%) 48 52	14, 24, 42, 65	5 (0%)

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	154	PRO	5.5
1	C	154	PRO	4.3
1	C	187	CYS	3.9
1	J	152	VAL	3.6
1	E	188	CYS	3.4
1	B	23	ARG	3.3
1	I	24	ASP	3.3
1	C	188	CYS	3.3
1	H	162	SER	3.1
1	D	24	ASP	3.0
1	H	24	ASP	3.0
1	G	189	PRO	3.0
1	E	152	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
1	E	158	ASN	3.0
1	C	158	ASN	2.9
1	I	183	VAL	2.9
1	F	155	THR	2.9
1	C	186	SER	2.9
1	G	24	ASP	2.8
1	G	186	SER	2.8
1	G	158	ASN	2.8
1	H	207	SER	2.7
1	G	-6	TYR	2.7
1	G	188	CYS	2.7
1	I	187	CYS	2.7
1	H	23	ARG	2.7
1	C	24	ASP	2.6
1	H	61	ARG	2.6
1	J	23	ARG	2.6
1	D	-6	TYR	2.6
1	D	156	THR	2.6
1	E	89	TYR	2.5
1	H	25	ARG	2.5
1	H	-3	ASP	2.4
1	F	152	VAL	2.4
1	B	22	GLN	2.4
1	H	156	THR	2.4
1	I	205	GLY	2.4
1	E	180	LYS	2.4
1	C	185	TYR	2.4
1	C	152	VAL	2.4
1	G	154	PRO	2.3
1	E	24	ASP	2.3
1	A	154	PRO	2.3
1	G	194	ASP	2.3
1	H	68	SER	2.2
1	C	181	ASN	2.2
1	I	23	ARG	2.2
1	I	189	PRO	2.2
1	D	114	MET	2.2
1	E	193	GLU	2.2
1	I	186	SER	2.2
1	A	24	ASP	2.2
1	J	-4	ASP	2.2
1	D	205	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	158	ASN	2.1
1	B	129	ASP	2.1
1	E	174	LEU	2.1
1	J	193	GLU	2.1
1	D	183	VAL	2.1
1	F	158	ASN	2.1
1	F	24	ASP	2.1
1	F	129	ASP	2.1
1	I	-5	LYS	2.1
1	C	153	ASP	2.1
1	A	207	SER	2.0
1	B	89	TYR	2.0
1	H	-2	ASP	2.0
1	H	155	THR	2.0
1	J	205	GLY	2.0
1	B	61	ARG	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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($\text{\AA}^2$)	Q<0.9
3	4P6	C	302	18/18	0.87	0.11	19,32,40,46	0
3	4P6	A	303	18/18	0.93	0.08	20,30,37,41	0
3	4P6	E	301	18/18	0.93	0.08	24,32,39,40	0
3	4P6	G	303	18/18	0.94	0.07	19,28,37,43	0
3	4P6	B	302	18/18	0.95	0.07	17,26,35,39	0
3	4P6	D	302	18/18	0.95	0.07	22,30,39,41	0
3	4P6	I	303	18/18	0.95	0.07	18,27,36,39	0

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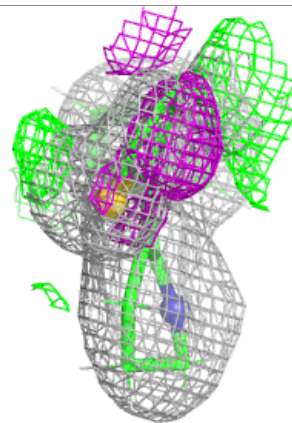
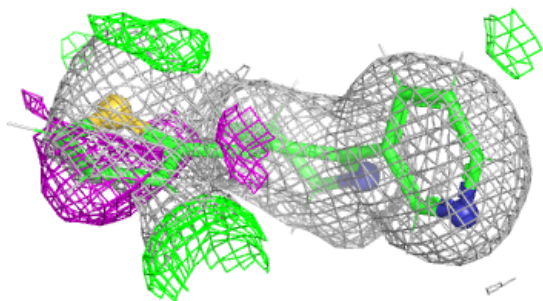
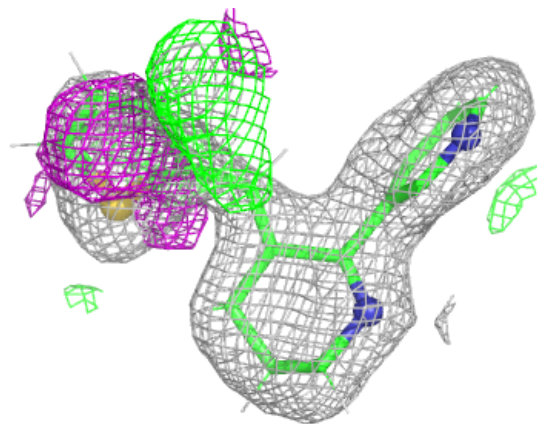
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	4P6	J	302	18/18	0.95	0.07	17,29,37,43	0
3	4P6	F	301	18/18	0.96	0.07	19,31,38,44	0
2	PO4	A	302	5/5	0.97	0.05	18,20,22,23	0
3	4P6	H	301	18/18	0.97	0.05	16,24,33,36	0
2	PO4	I	301	5/5	0.98	0.04	17,18,18,19	0
2	PO4	I	302	5/5	0.98	0.06	20,20,21,24	0
2	PO4	C	301	5/5	0.98	0.05	17,17,19,20	0
2	PO4	A	301	5/5	0.99	0.05	21,21,23,25	0
2	PO4	J	301	5/5	0.99	0.05	20,20,20,23	0
2	PO4	D	301	5/5	0.99	0.04	18,19,20,21	0
2	PO4	G	301	5/5	0.99	0.04	17,18,19,22	0
2	PO4	G	302	5/5	0.99	0.04	18,18,21,21	0
2	PO4	B	301	5/5	0.99	0.04	16,16,17,17	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.

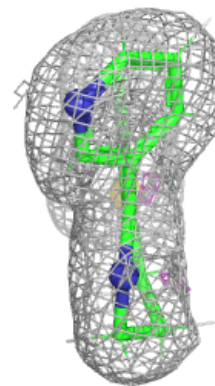
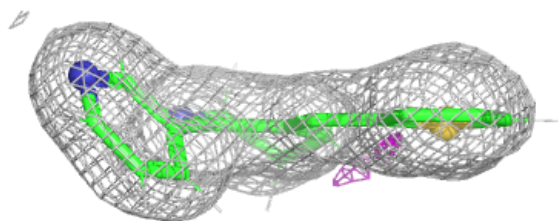
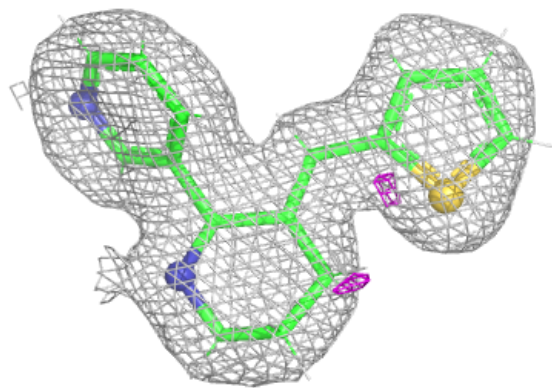
Electron density around 4P6 C 302:

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



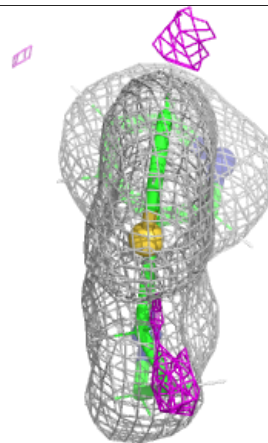
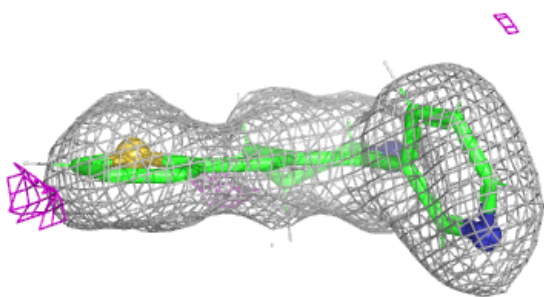
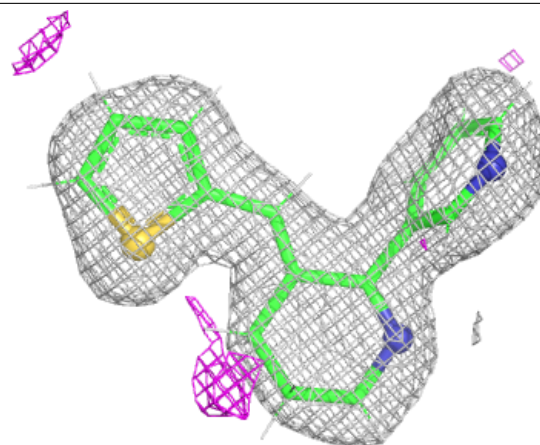
Electron density around 4P6 A 303:

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



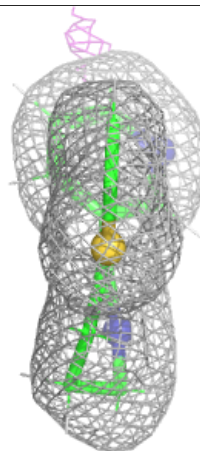
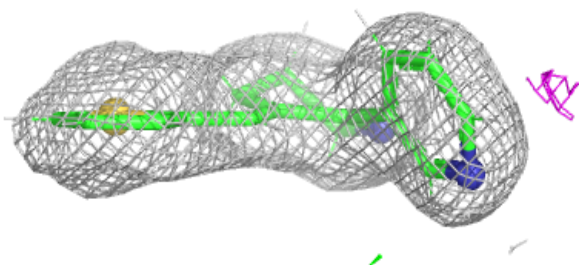
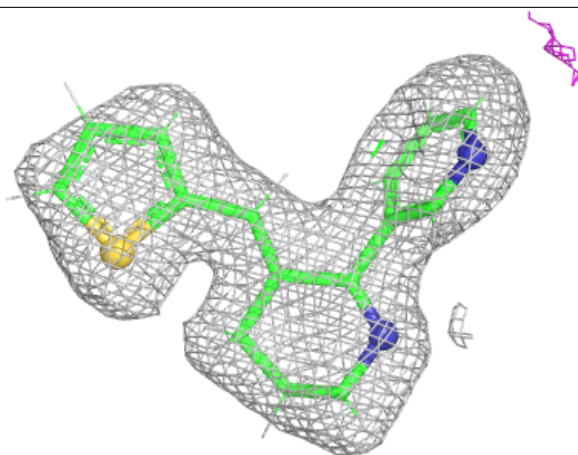
Electron density around 4P6 E 301:

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



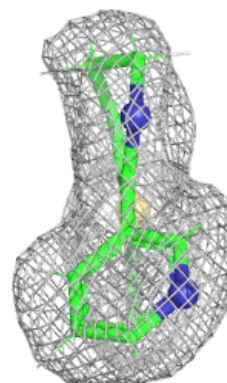
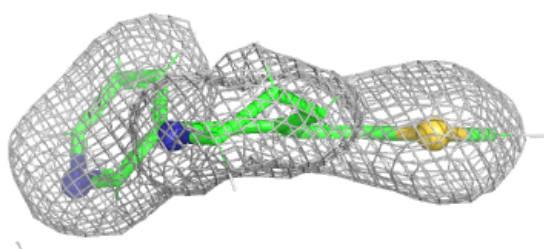
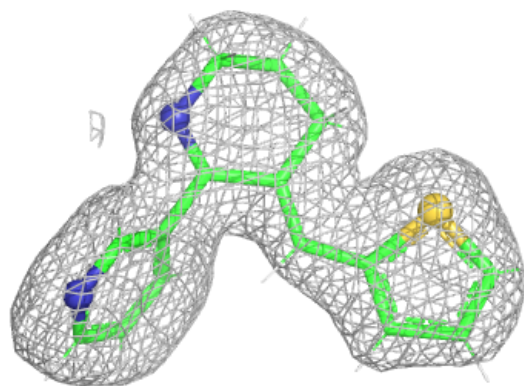
Electron density around 4P6 G 303:

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

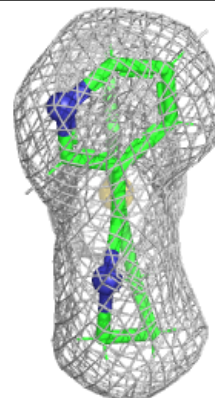
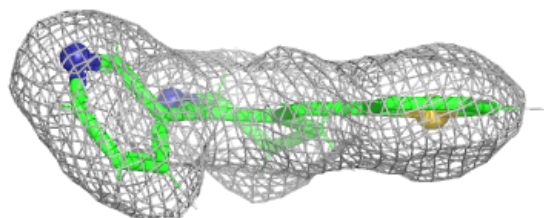
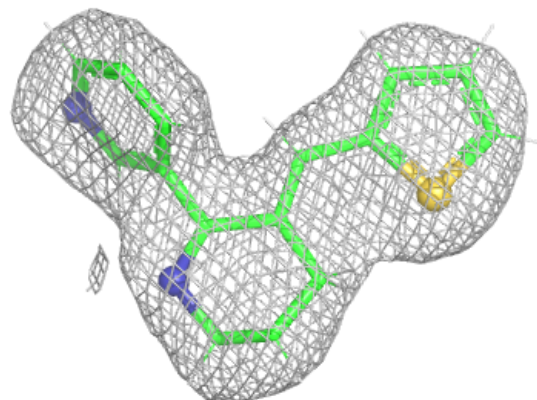


Electron density around 4P6 B 302:

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

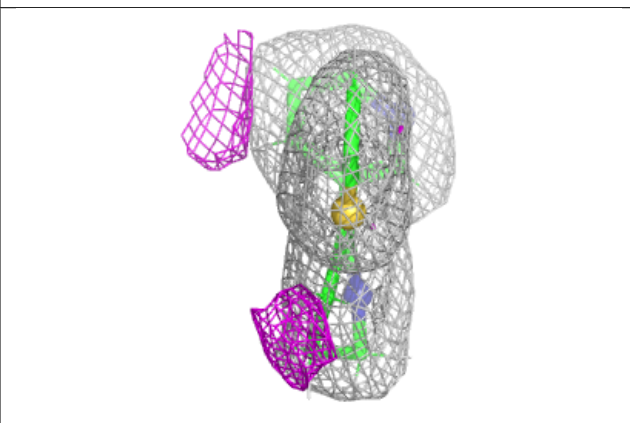
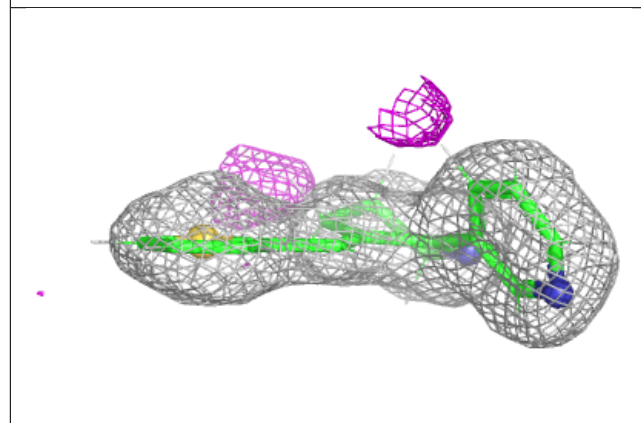
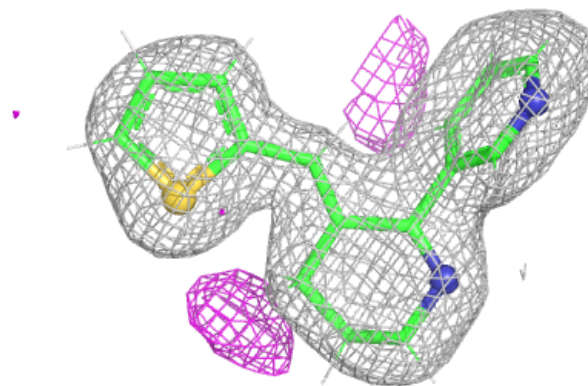
**Electron density around 4P6 D 302:**

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

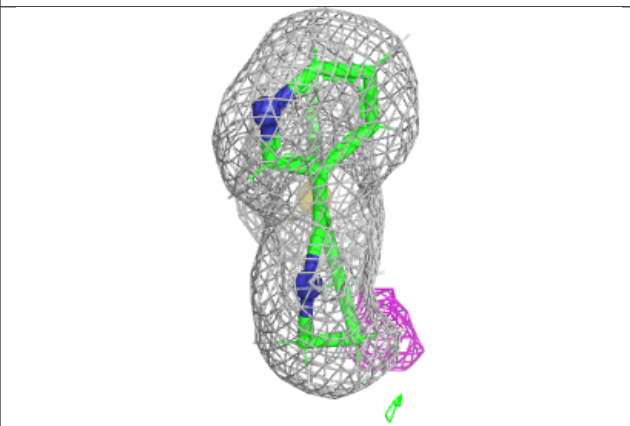
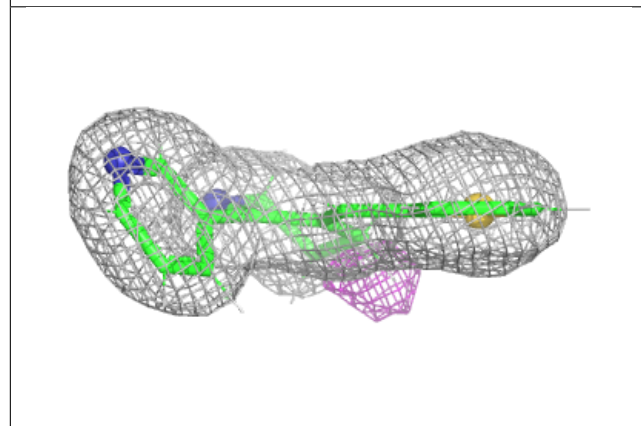
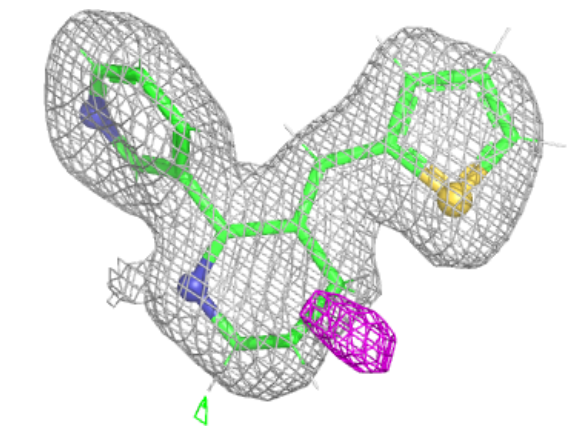


Electron density around 4P6 I 303:

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

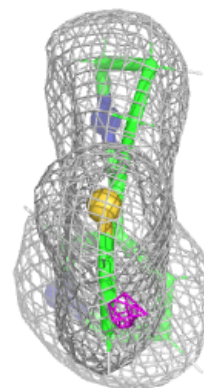
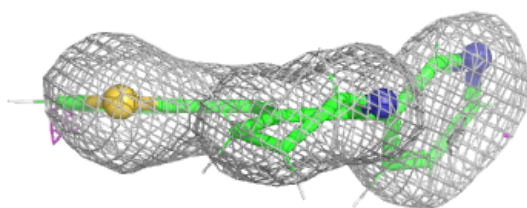
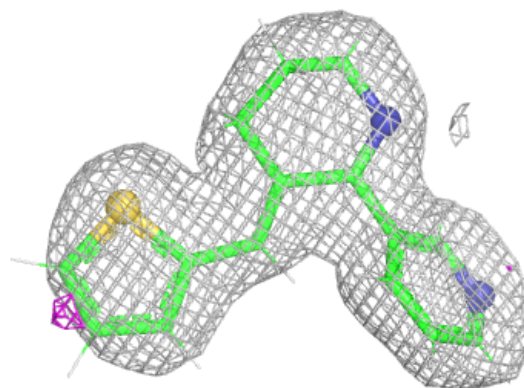
**Electron density around 4P6 J 302:**

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

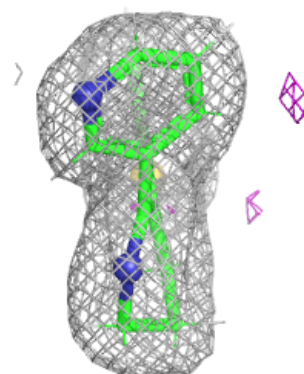
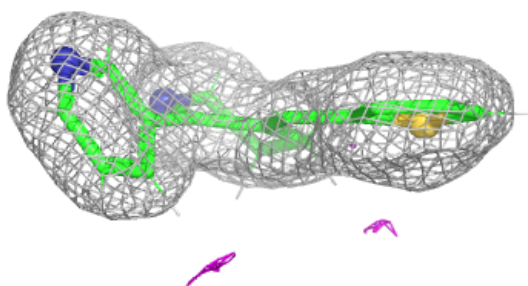
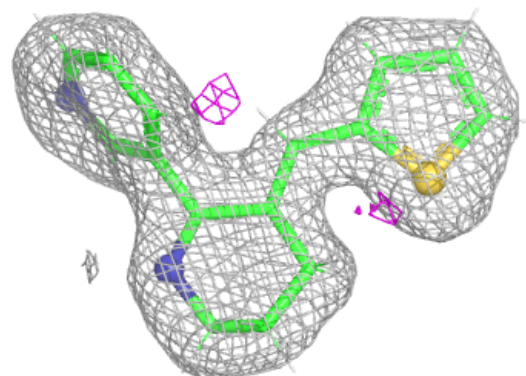


Electron density around 4P6 F 301:

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

**Electron density around 4P6 H 301:**

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



6.5 Other polymers ⓘ

There are no such residues in this entry.