



wwPDB X-ray Structure Validation Summary Report ⓘ

May 13, 2020 – 11:39 am BST

PDB ID : 3SGV
Title : Crystal Structure of E. coli undecaprenyl pyrophosphate synthase in complex with BPH-1290
Authors : Cao, R.; Liu, Y.-L.; Oldfield, E.
Deposited on : 2011-06-15
Resolution : 1.61 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 : 2.11
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) : 2.11

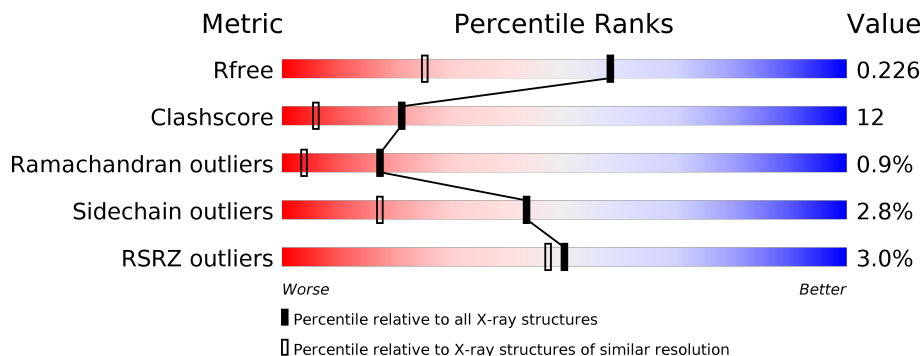
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.61 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	4693 (1.64-1.60)
Clashscore	141614	5002 (1.64-1.60)
Ramachandran outliers	138981	4888 (1.64-1.60)
Sidechain outliers	138945	4887 (1.64-1.60)
RSRZ outliers	127900	4609 (1.64-1.60)

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	A	253	 5% 75% 11% • 12%
1	B	253	 5% 67% 15% • 14%

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
2	2BJ	B	5001	-	-	X	-

2 Entry composition [i](#)

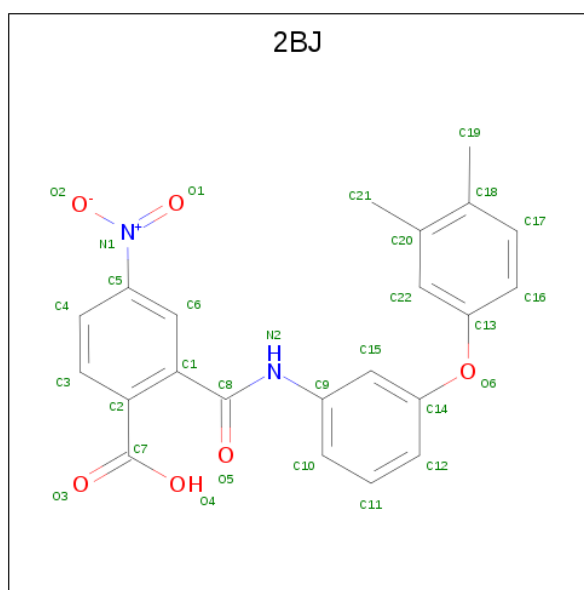
There are 3 unique types of molecules in this entry. The entry contains 4008 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Undecaprenyl pyrophosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	217	Total 1737	C 1099	N 322	O 309	S 7	0	4	0
1	A	222	Total 1786	C 1122	N 332	O 326	S 6	0	3	0

- Molecule 2 is 2-{[3-(3,4-dimethylphenoxy)phenyl]carbamoyl}-4-nitrobenzoic acid (three-letter code: 2BJ) (formula: C₂₂H₁₈N₂O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	B	1	Total 30	C 22	N 2	O 6	0	0
2	B	1	Total 30	C 22	N 2	O 6	0	0
2	A	1	Total 30	C 22	N 2	O 6	0	0
2	A	1	Total 30	C 22	N 2	O 6	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	15	14	1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	170	Total	O	0	0
			170	170		
3	A	180	Total	O	0	0
			180	180		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.14Å 69.00Å 112.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.61 – 1.61 43.61 – 1.61	Depositor EDS
% Data completeness (in resolution range)	97.3 (43.61-1.61) 97.2 (43.61-1.61)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.52 (at 1.61Å)	Xtrriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.174 , 0.214 0.193 , 0.226	Depositor DCC
R_{free} test set	3146 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	20.0	Xtrriage
Anisotropy	0.046	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4008	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.46% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 2BJ

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	1.20	5/1831 (0.3%)	1.13	5/2474 (0.2%)
1	B	1.29	7/1783 (0.4%)	1.27	13/2407 (0.5%)
All	All	1.25	12/3614 (0.3%)	1.20	18/4881 (0.4%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	174	GLU	CB-CG	-6.87	1.39	1.52
1	A	175	GLU	CG-CD	6.56	1.61	1.51
1	A	185	GLU	CD-OE2	6.32	1.32	1.25
1	B	18[A]	CYS	CB-SG	-6.32	1.71	1.82
1	B	18[B]	CYS	CB-SG	-6.32	1.71	1.82

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	239	ARG	NE-CZ-NH1	-19.47	110.57	120.30
1	B	239	ARG	NE-CZ-NH2	13.27	126.94	120.30
1	B	239	ARG	CG-CD-NE	-7.62	95.81	111.80
1	A	204	PHE	CB-CG-CD2	-6.99	115.91	120.80
1	A	108	ARG	NE-CZ-NH1	6.93	123.76	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	68	TYR	Mainchain,Peptide

5.2 Too-close contacts [\(i\)](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1786	0	1753	26	0
1	B	1737	0	1726	60	0
2	A	75	0	47	3	0
2	B	60	0	34	13	0
3	A	180	0	0	8	1
3	B	170	0	0	6	1
All	All	4008	0	3560	87	1

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 12.

The worst 5 of 87 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:B:48[A]:LYS:HZ2	2:B:5001:2BJ:C7	1.33	1.40
1:B:48[A]:LYS:NZ	2:B:5001:2BJ:C7	1.99	1.26
1:B:48[A]:LYS:CD	1:B:51:ARG:HH12	1.55	1.19
1:B:48[A]:LYS:NZ	2:B:5001:2BJ:O4	1.76	1.16
1:B:48[A]:LYS:HD2	1:B:51:ARG:HH12	1.10	1.08

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:278:HOH:O	3:A:298:HOH:O[2_554]	1.46	0.74

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	223/253 (88%)	219 (98%)	3 (1%)	1 (0%)	34	15
1	B	217/253 (86%)	210 (97%)	4 (2%)	3 (1%)	11	1
All	All	440/506 (87%)	429 (98%)	7 (2%)	4 (1%)	17	4

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	91	TRP
1	B	92	ALA
1	A	204	PHE
1	B	204	PHE

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	184/206 (89%)	180 (98%)	4 (2%)	52	25
1	B	178/206 (86%)	169 (95%)	9 (5%)	24	5
All	All	362/412 (88%)	349 (96%)	13 (4%)	43	11

5 of 13 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	117	ASN
1	B	120[A]	LEU

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Mol	Chain	Res	Type
1	A	153[A]	GLN
1	B	94	ASP
1	A	115	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	166	ASN
1	A	168	GLN
1	A	43	HIS
1	B	117	ASN
1	A	144	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

5 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	2BJ	B	5001	-	29,32,32	3.10	6 (20%)	39,45,45	1.71	11 (28%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	2BJ	A	3001	-	29,32,32	2.25	8 (27%)	39,45,45	2.51	11 (28%)
2	2BJ	A	2001	-	29,32,32	2.01	9 (31%)	39,45,45	1.55	9 (23%)
2	2BJ	B	1001	-	29,32,32	3.63	8 (27%)	39,45,45	4.65	13 (33%)
2	2BJ	A	4001	-	16,16,32	1.08	0	21,21,45	1.07	2 (9%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	2BJ	B	5001	-	-	2/14/20/20	0/3/3/3
2	2BJ	A	3001	-	-	0/14/20/20	0/3/3/3
2	2BJ	A	2001	-	-	0/14/20/20	0/3/3/3
2	2BJ	B	1001	-	-	0/14/20/20	0/3/3/3
2	2BJ	A	4001	-	-	0/4/4/20	0/2/2/3

The worst 5 of 31 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1001	2BJ	O1-N1	16.54	1.50	1.22
2	B	5001	2BJ	O1-N1	14.66	1.47	1.22
2	B	1001	2BJ	C5-N1	-6.44	1.29	1.45
2	A	3001	2BJ	C2-C7	5.91	1.53	1.47
2	A	2001	2BJ	O1-N1	5.45	1.32	1.22

The worst 5 of 46 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1001	2BJ	C4-C5-N1	-19.03	105.05	119.38
2	B	1001	2BJ	C6-C5-N1	14.43	131.46	118.75
2	A	3001	2BJ	C4-C5-N1	-9.20	112.45	119.38
2	B	1001	2BJ	O1-N1-C5	-9.01	106.05	118.80
2	B	1001	2BJ	C1-C6-C5	-6.38	113.79	120.26

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	5001	2BJ	C4-C5-N1-O1

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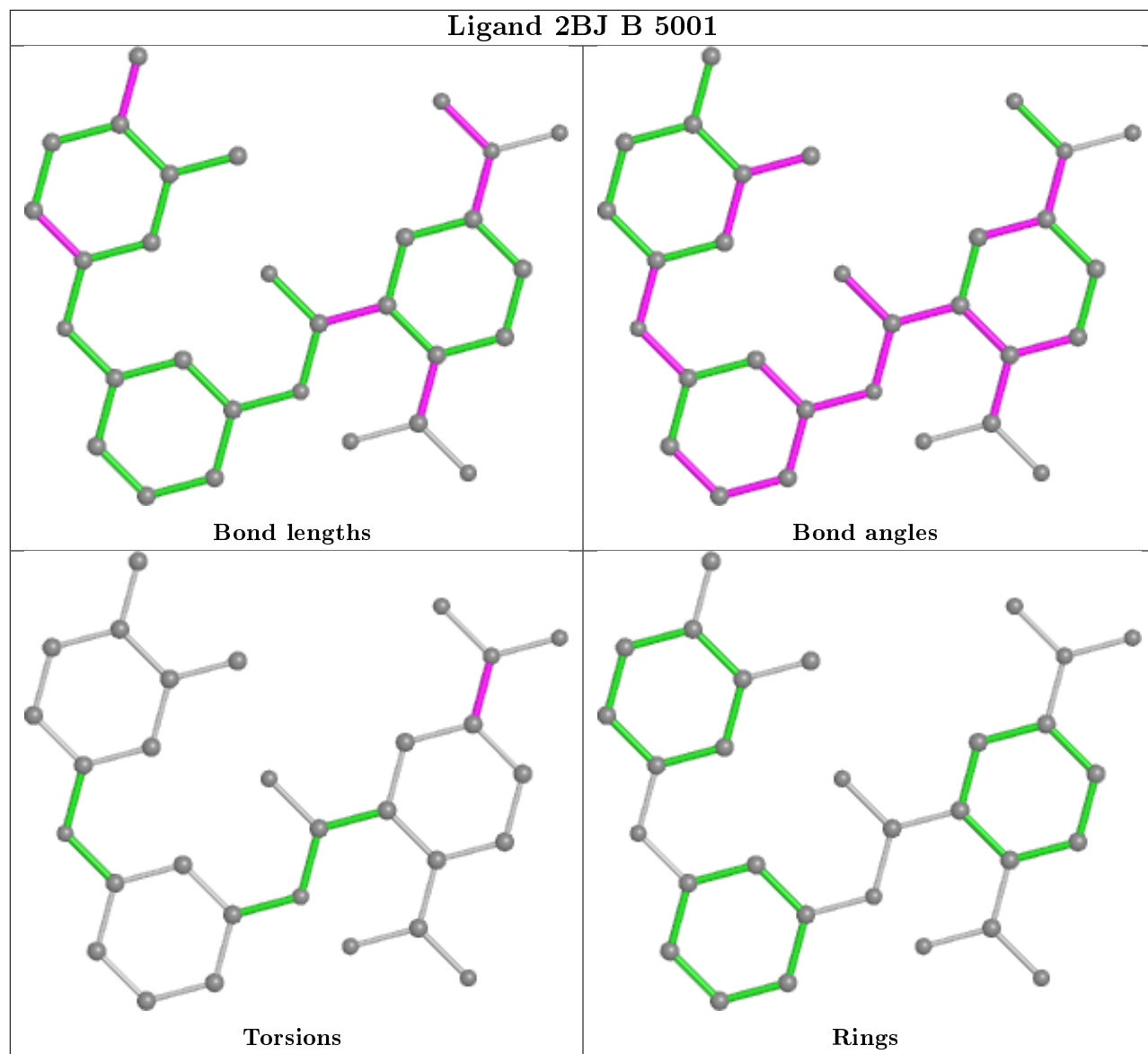
Mol	Chain	Res	Type	Atoms
2	B	5001	2BJ	C6-C5-N1-O1

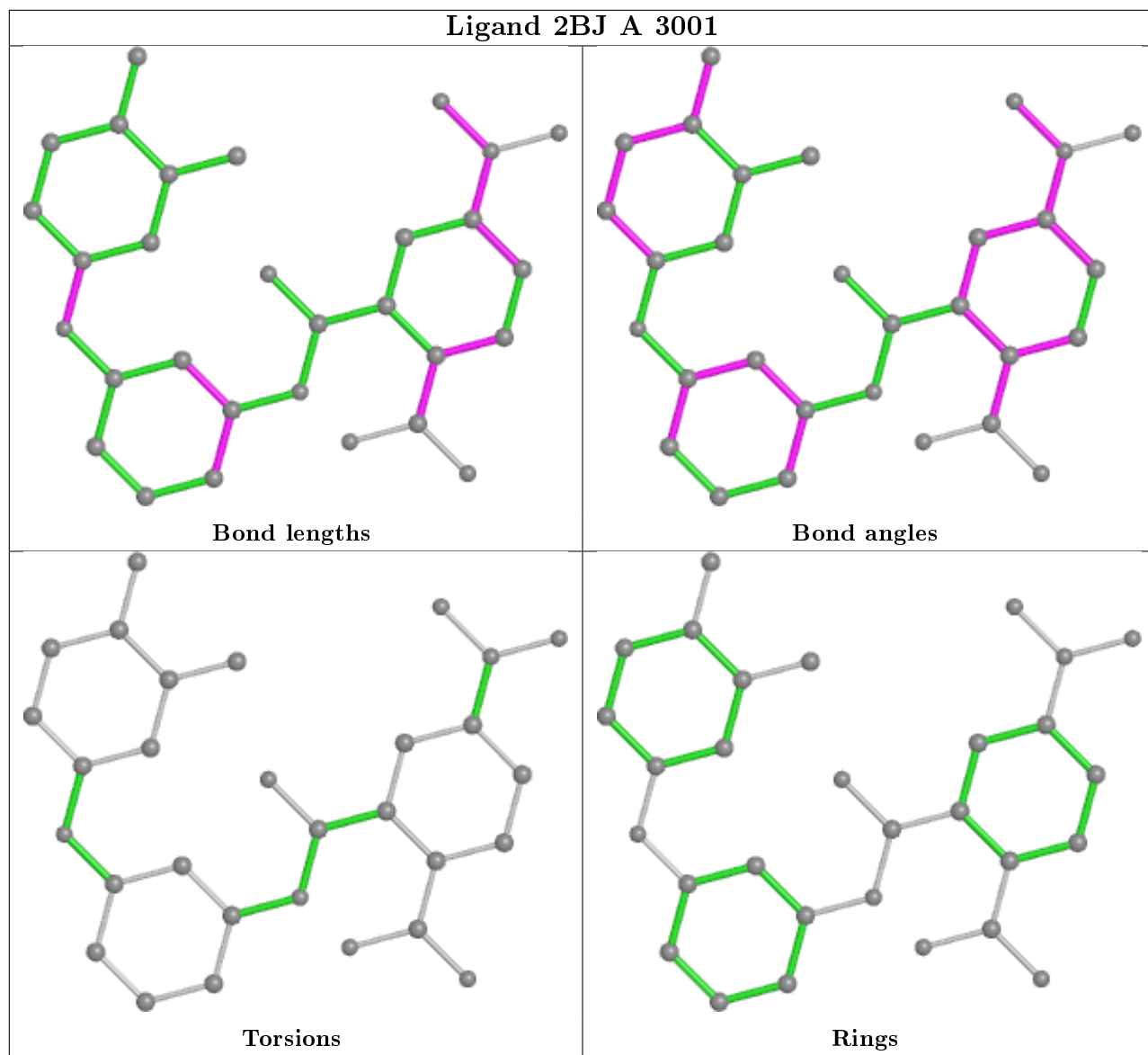
There are no ring outliers.

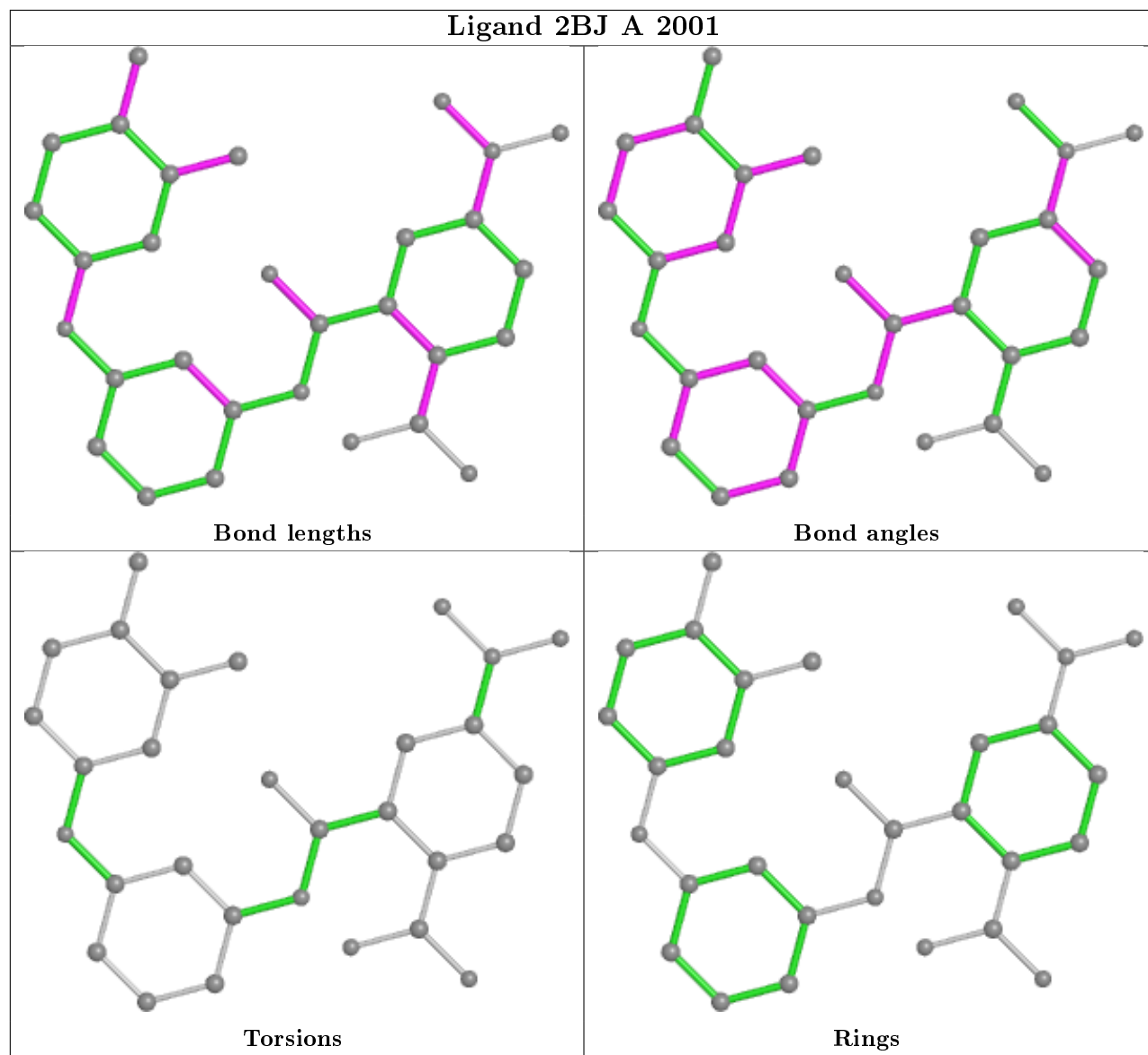
3 monomers are involved in 16 short contacts:

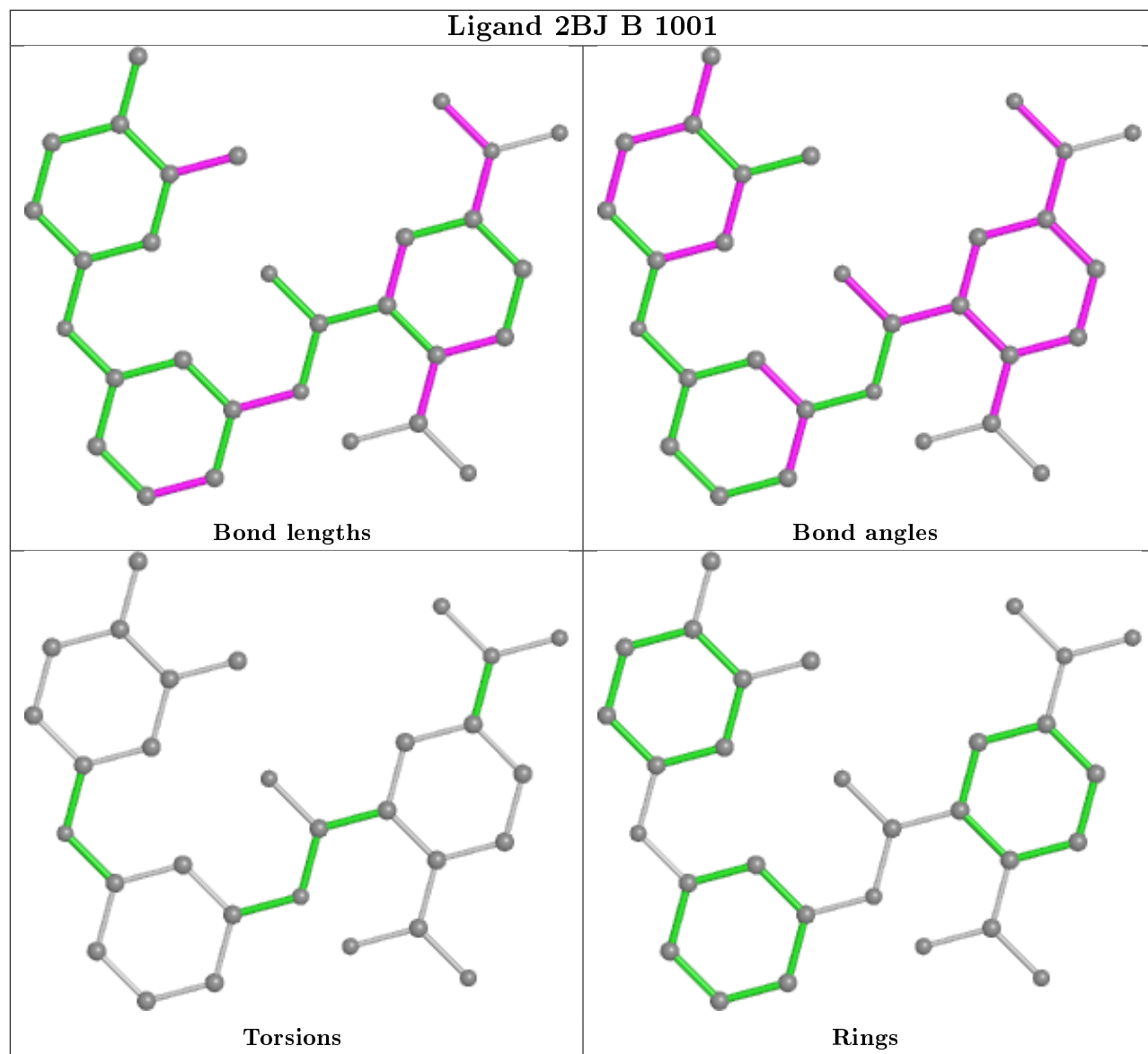
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	5001	2BJ	10	0
2	A	3001	2BJ	3	0
2	B	1001	2BJ	3	0

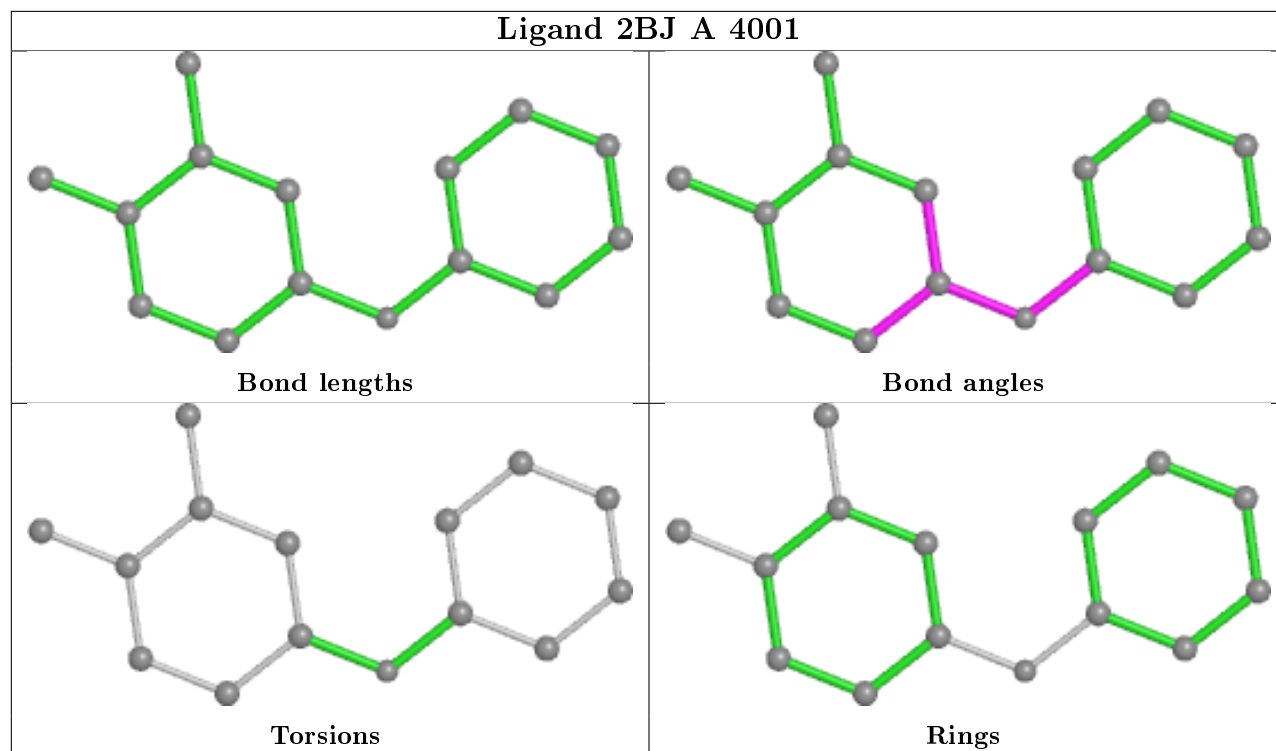
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.











5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled '#RSRZ > 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	222/253 (87%)	-0.06	0 100 100	12, 22, 43, 56	0
1	B	217/253 (85%)	0.21	13 (5%) 21 19	11, 21, 50, 73	0
All	All	439/506 (86%)	0.08	13 (2%) 50 47	11, 21, 47, 73	0

The worst 5 of 13 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	91	TRP	8.5
1	B	92	ALA	6.5
1	B	69	ALA	4.9
1	B	70	PHE	4.8
1	B	116	PHE	4.8

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](#)

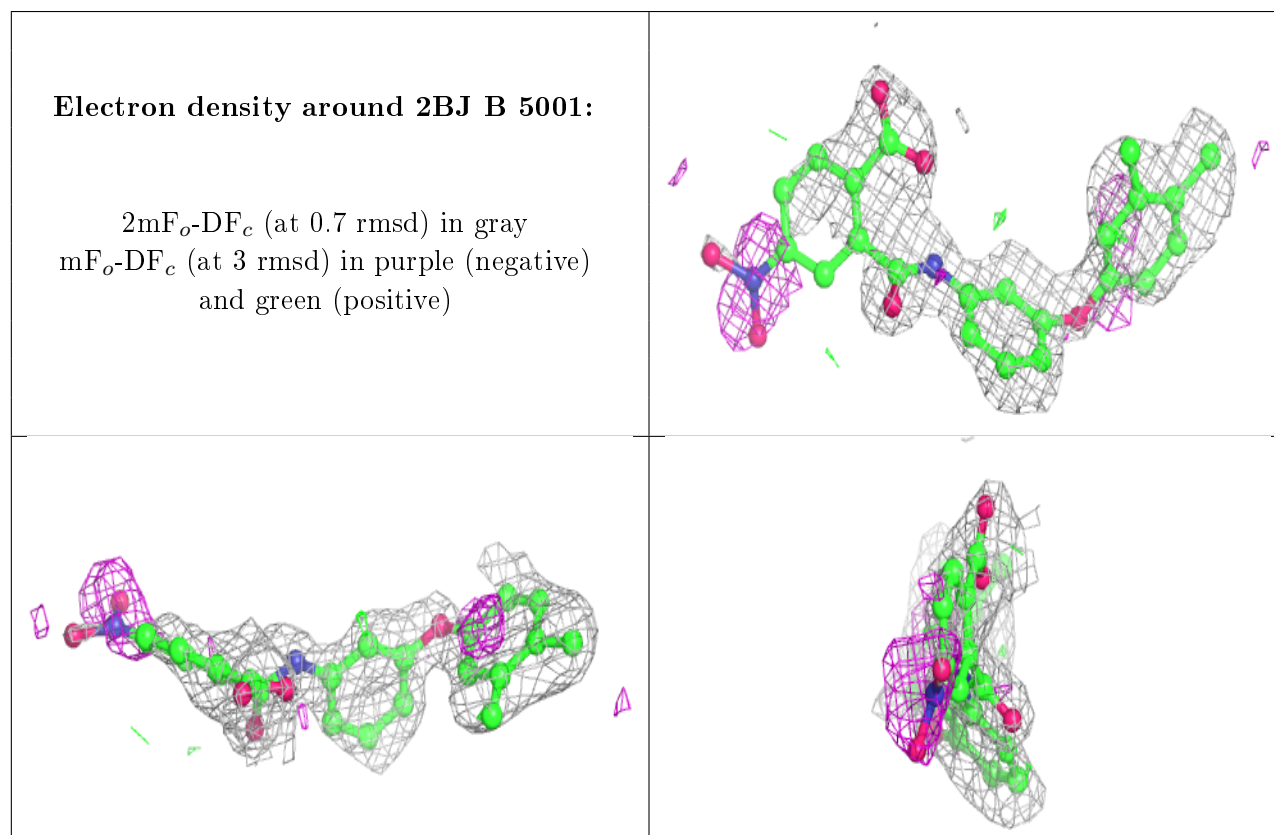
There are no carbohydrates in this entry.

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.

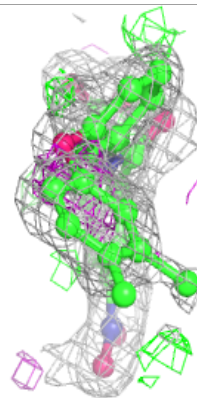
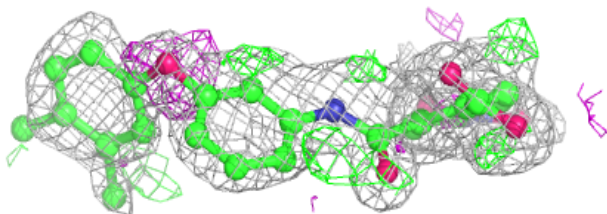
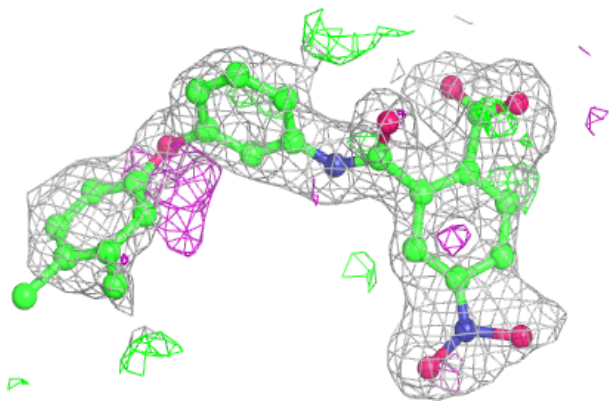
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	2BJ	B	5001	30/30	0.64	0.30	38,47,59,63	0
2	2BJ	B	1001	30/30	0.85	0.18	24,33,45,47	0
2	2BJ	A	4001	15/30	0.85	0.24	27,37,59,60	0
2	2BJ	A	3001	30/30	0.92	0.11	14,21,37,43	0
2	2BJ	A	2001	30/30	0.93	0.11	13,21,30,40	0

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.

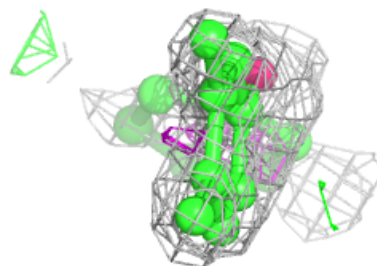
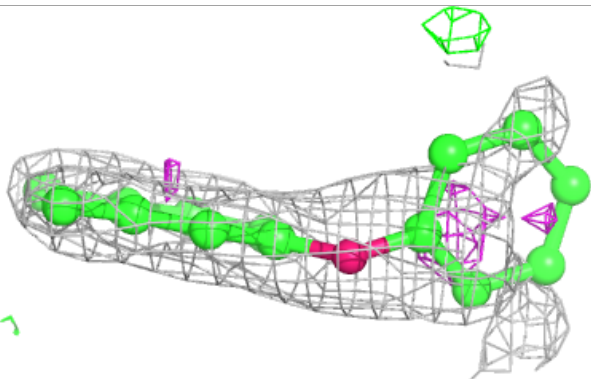
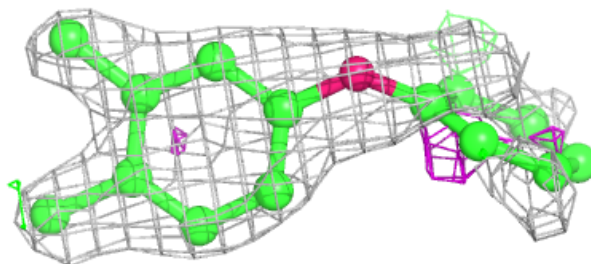


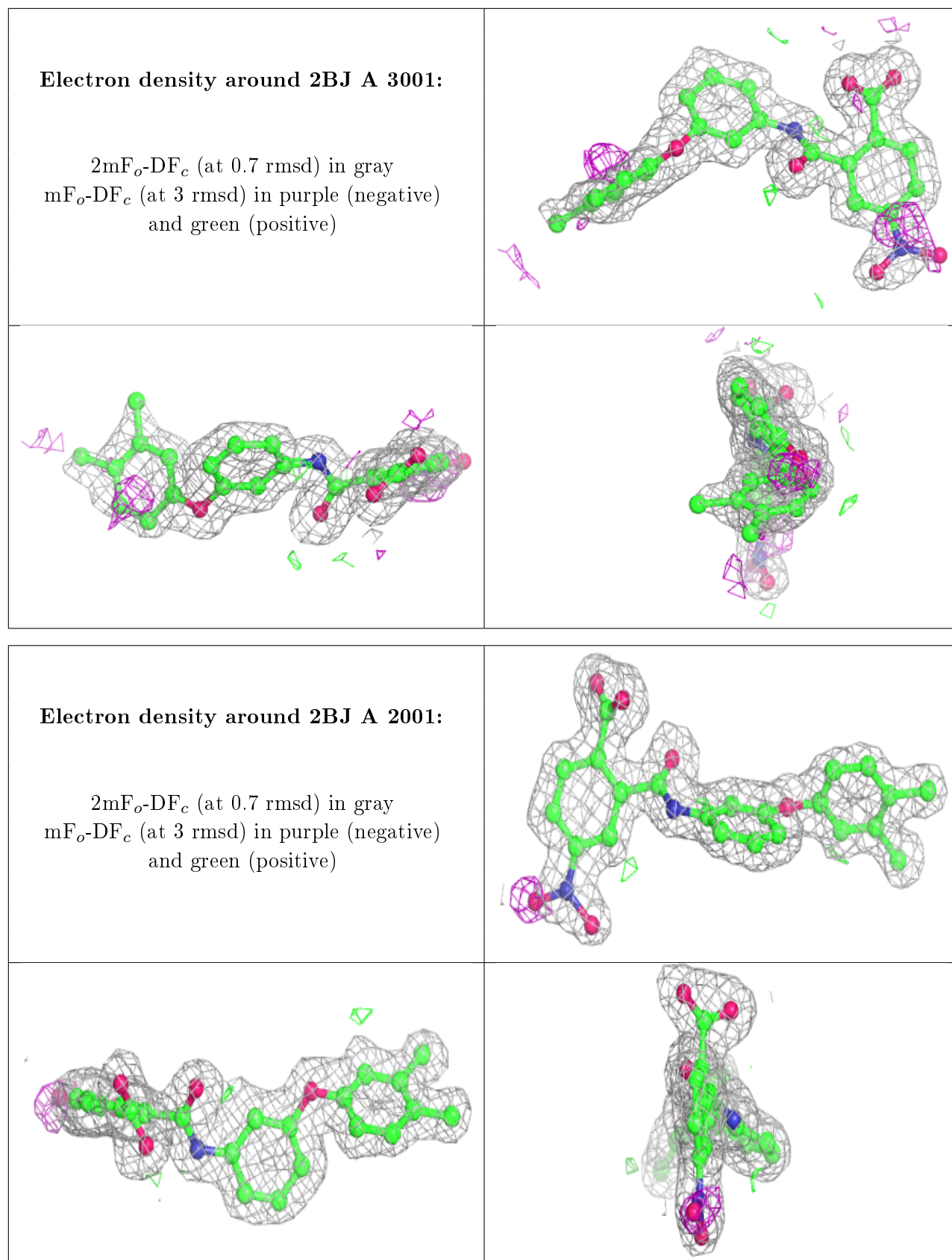
Electron density around 2BJ B 1001:

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

**Electron density around 2BJ A 4001:**

$2mF_o-DF_c$ (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.