



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 10:09 AM UTC

PDB ID : 6IA9 / pdb_00006ia9
Title : urate oxidase under 2000 bar (220 MPa) of argon
Authors : Prange, T.; Colloc'h, N.; Carpentier, P.
Deposited on : 2018-11-26
Resolution : 1.80 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 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

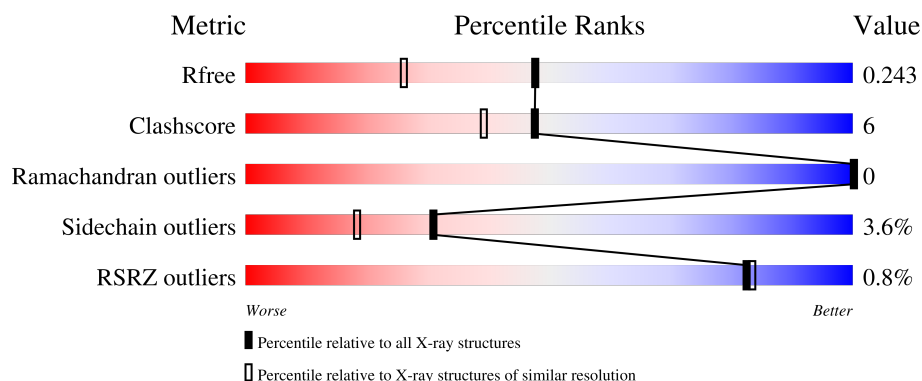
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	aaa	295	% . 87% 12%
1	aba	295	88% 12%

2 Entry composition [i](#)

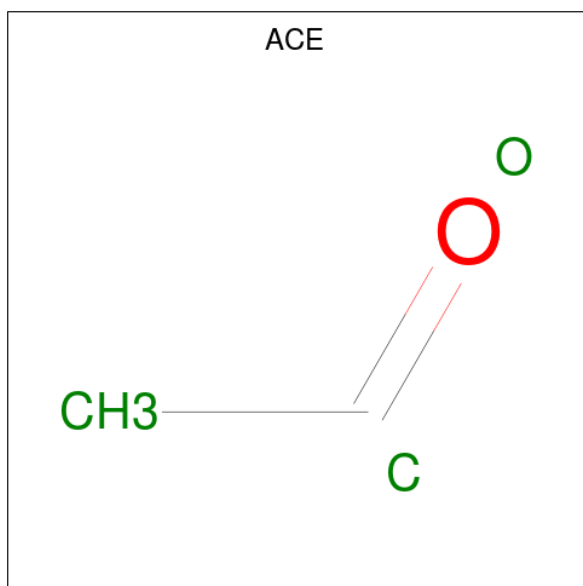
There are 6 unique types of molecules in this entry. The entry contains 5338 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Uricase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	aaa	295	Total	C	N	O	S	0	9	0
			2408	1525	418	457	8			
1	aba	295	Total	C	N	O	S	0	4	0
			2379	1502	412	456	9			

- Molecule 2 is ACETYL GROUP (CCD ID: ACE) (formula: C₂H₄O).

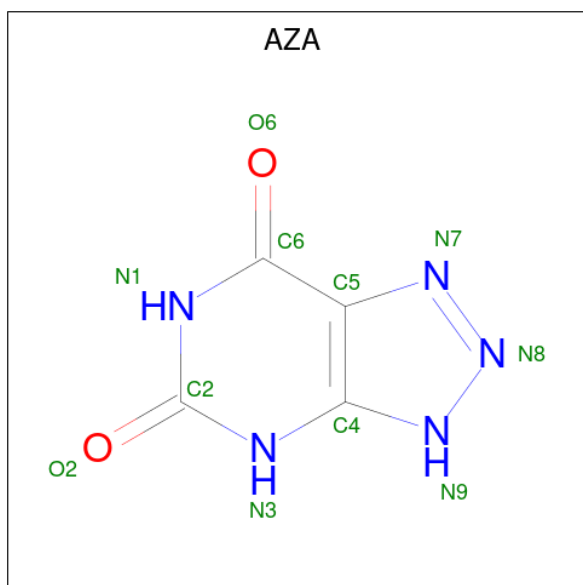


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	aaa	1	Total	C	O	0	0
			3	2	1		
2	aba	1	Total	C	O	0	0
			3	2	1		

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	aaa	1	Total Na 1 1	0	0
3	aba	1	Total Na 1 1	0	0

- Molecule 4 is 8-AZAXANTHINE (CCD ID: AZA) (formula: $C_4H_3N_5O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	aaa	1	Total C N O 11 4 5 2	0	0
4	aaa	1	Total C N O 11 4 5 2	0	0
4	aba	1	Total C N O 11 4 5 2	0	0
4	aba	1	Total C N O 11 4 5 2	0	0

- Molecule 5 is ARGON (CCD ID: AR) (formula: Ar).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	aaa	2	Total Ar 2 2	0	0
5	aba	2	Total Ar 2 2	0	0

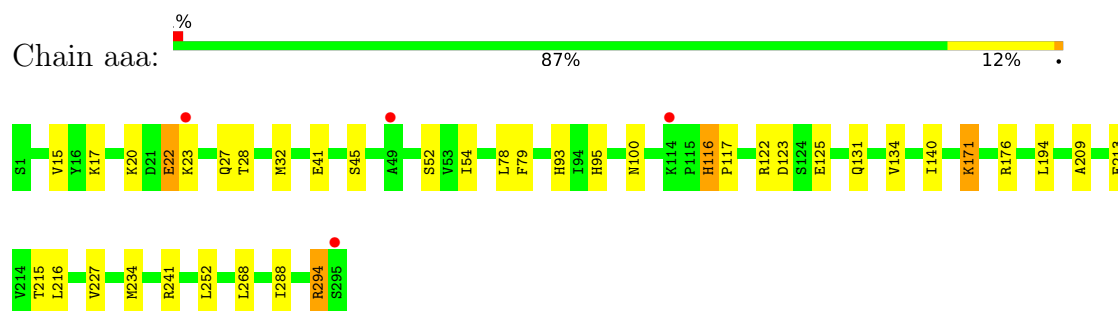
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	aaa	217	Total 219	O 219	0	2
6	aba	274	Total 276	O 276	0	2

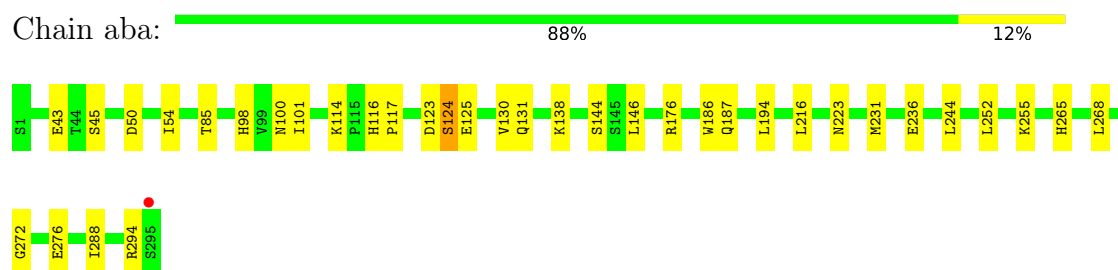
3 Residue-property plots

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: Uricase



- Molecule 1: Uricase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	78.55Å 96.68Å 105.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.33 – 1.80 46.33 – 1.80	Depositor EDS
% Data completeness (in resolution range)	96.2 (46.33-1.80) 96.2 (46.33-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.25 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.209 , 0.243 0.209 , 0.243	Depositor DCC
R_{free} test set	3583 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	29.6	Xtriage
Anisotropy	0.103	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 32.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5338	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 58.27 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.0947e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: AZA, AR, NA, ACE

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	aaa	0.58	0/2489	1.04	2/3373 (0.1%)
1	aba	0.59	0/2443	1.02	2/3314 (0.1%)
All	All	0.58	0/4932	1.03	4/6687 (0.1%)

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	aba	0	1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	aba	236	GLU	CB-CA-C	5.96	121.65	110.63
1	aba	85	THR	CA-CB-OG1	-5.27	101.69	109.60
1	aaa	28	THR	CA-CB-OG1	-5.18	101.83	109.60
1	aaa	116	HIS	CB-CA-C	5.04	116.67	109.26

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	aba	223	ASN	Peptide

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	aaa	2408	0	2393	31	0
1	aba	2379	0	2334	26	0
2	aaa	3	0	3	0	0
2	aba	3	0	3	0	0
3	aaa	1	0	0	0	0
3	aba	1	0	0	0	0
4	aaa	22	0	6	1	0
4	aba	22	0	6	1	0
5	aaa	2	0	0	1	0
5	aba	2	0	0	0	0
6	aaa	219	0	0	3	0
6	aba	276	0	0	7	0
All	All	5338	0	4745	54	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.

The worst 5 of 54 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:aaa:294:ARG:HH11	1:aaa:294:ARG:HG2	1.29	0.98
1:aaa:22:GLU:HA	1:aaa:22:GLU:OE1	1.68	0.89
1:aaa:171[A]:LYS:HE3	1:aaa:171[A]:LYS:H	1.39	0.86
1:aba:176:ARG:HH12	4:aba:305:AZA:HN3	1.29	0.78
1:aba:43:GLU:CB	6:aba:601:HOH:O	2.35	0.73

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	aaa	302/295 (102%)	292 (97%)	10 (3%)	0	100	100
1	aba	297/295 (101%)	292 (98%)	5 (2%)	0	100	100
All	All	599/590 (102%)	584 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	aaa	270/261 (103%)	258 (96%)	12 (4%)	25	13
1	aba	265/261 (102%)	255 (96%)	10 (4%)	29	17
All	All	535/522 (102%)	513 (96%)	22 (4%)	31	15

5 of 22 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	aba	125	GLU
1	aba	187[B]	GLN
1	aba	187[A]	GLN
1	aba	194	LEU
1	aaa	171[A]	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 6 are monoatomic - leaving 6 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	ACE	aaa	301	1	1,2,2	0.14	0	0,1,1	-	-
4	AZA	aba	305	-	12,12,12	0.77	0	12,17,17	1.80	2 (16%)
4	AZA	aaa	304	-	12,12,12	0.71	0	12,17,17	1.78	1 (8%)
2	ACE	aba	301	1	1,2,2	0.19	0	0,1,1	-	-
4	AZA	aba	306	-	12,12,12	0.70	0	12,17,17	1.87	2 (16%)
4	AZA	aaa	303	-	12,12,12	0.81	0	12,17,17	1.86	2 (16%)

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
4	AZA	aaa	304	-	-	-	0/2/2/2
4	AZA	aba	306	-	-	-	0/2/2/2
4	AZA	aba	305	-	-	-	0/2/2/2
4	AZA	aaa	303	-	-	-	0/2/2/2

There are no bond length outliers.

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	aba	306	AZA	N3-C4-N9	5.61	133.72	124.57
4	aaa	304	AZA	N3-C4-N9	5.56	133.65	124.57
4	aba	305	AZA	N3-C4-N9	5.55	133.62	124.57
4	aaa	303	AZA	N3-C4-N9	5.37	133.33	124.57
4	aaa	303	AZA	N3-C2-N1	2.30	119.36	115.74

There are no chirality outliers.

There are no torsion outliers.

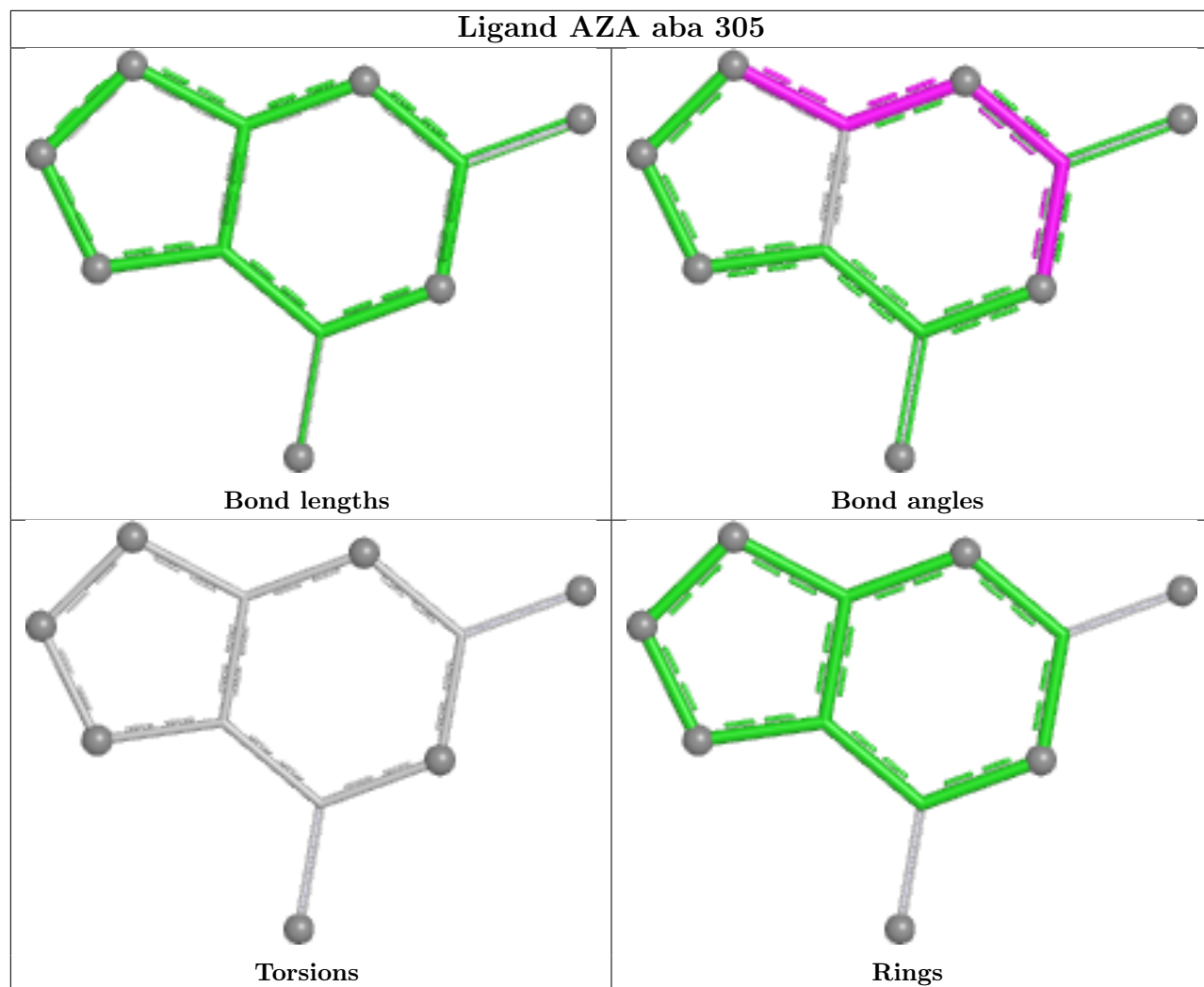
There are no ring outliers.

2 monomers are involved in 2 short contacts:

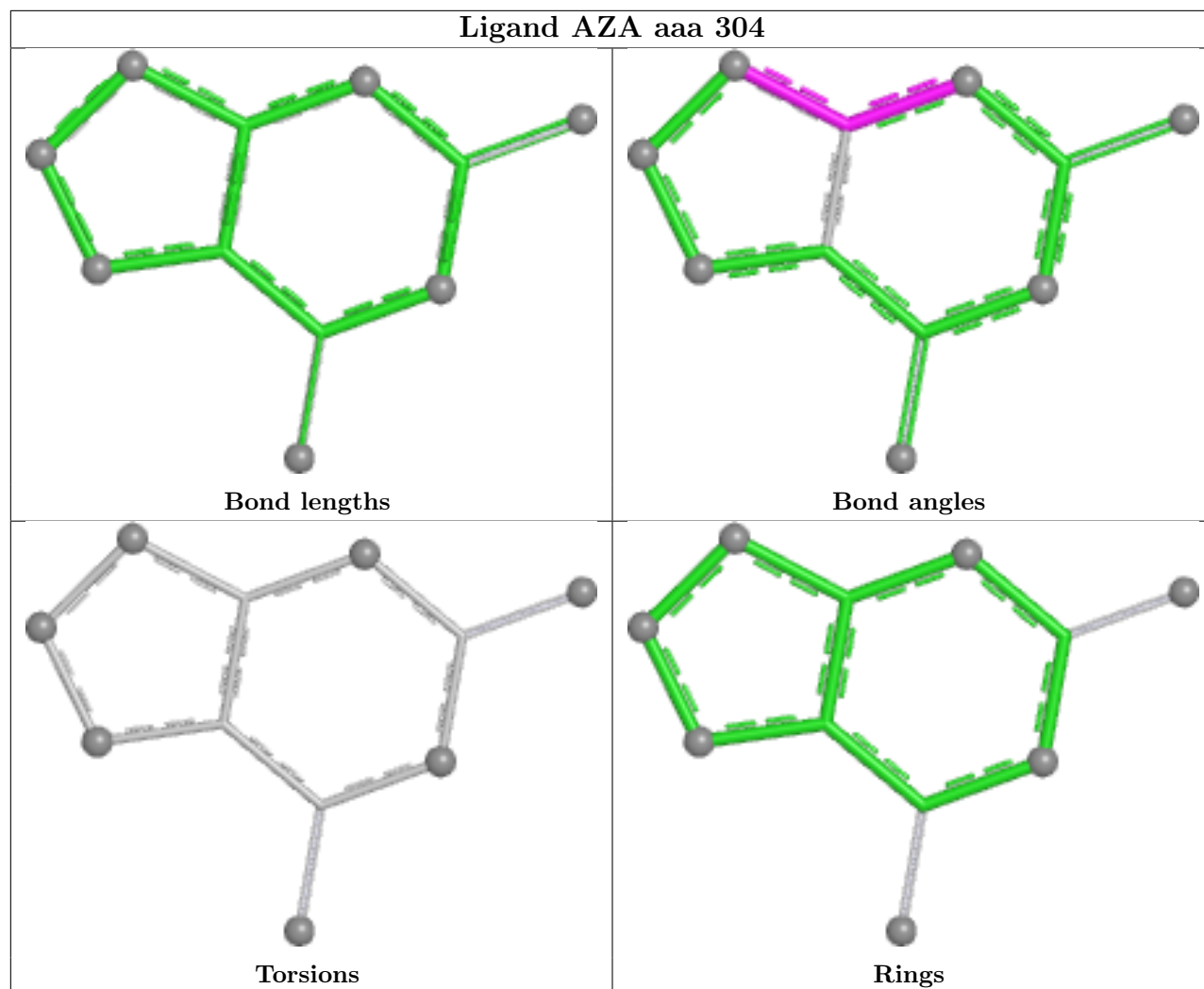
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	aba	305	AZA	1	0
4	aaa	303	AZA	1	0

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.

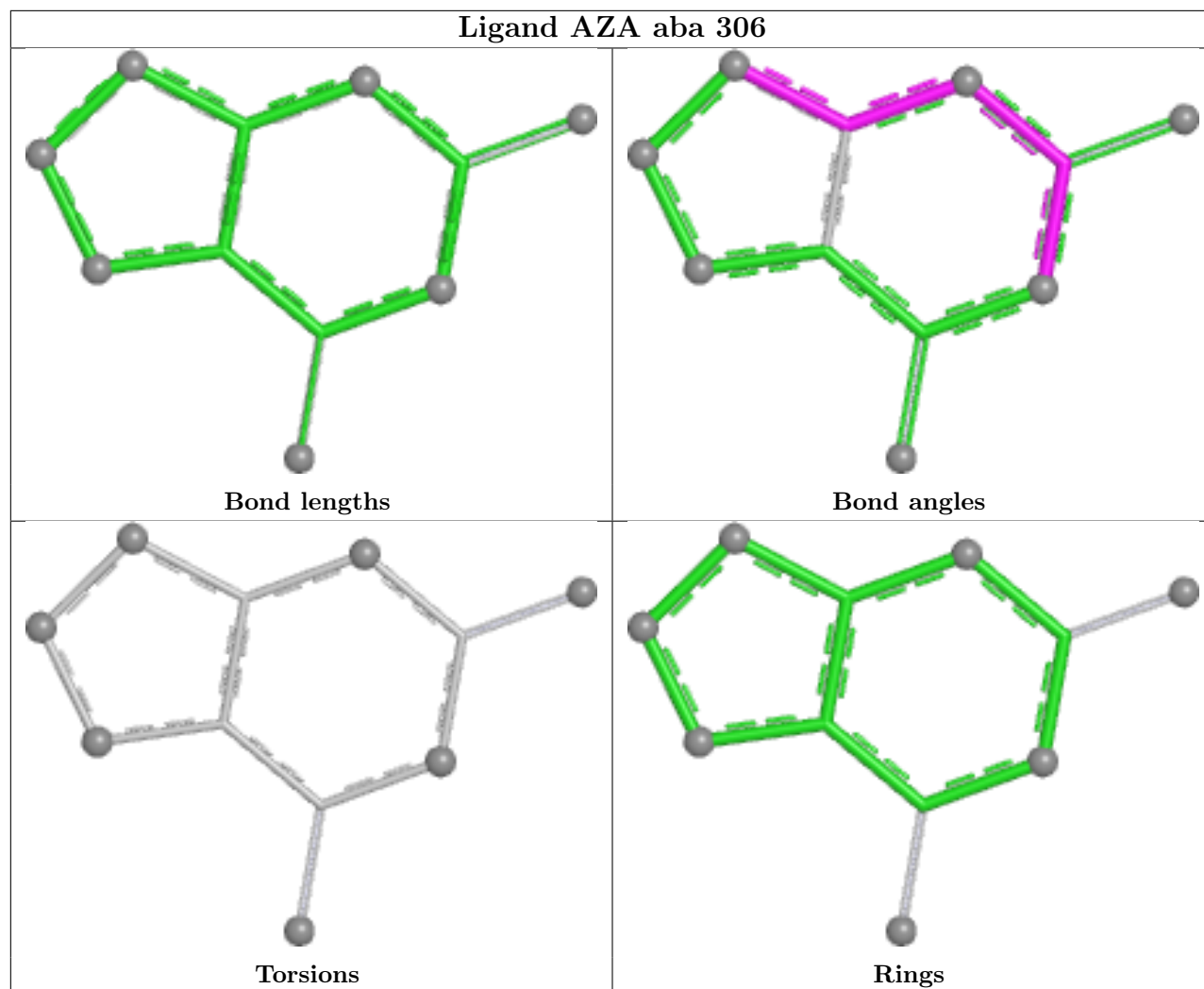
Ligand AZA aba 305

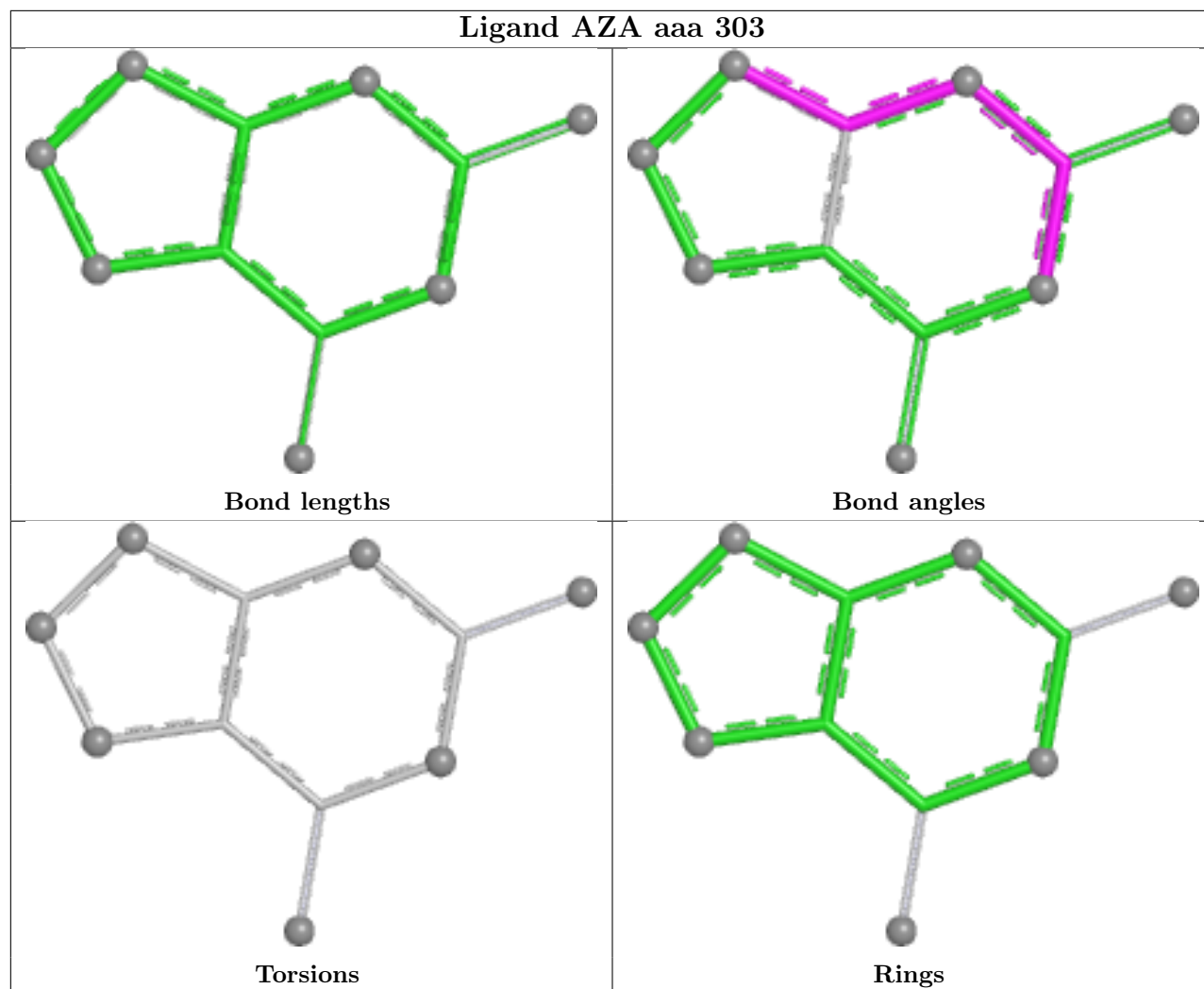


Ligand AZA aaa 304



Ligand AZA aba 306





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	aaa	295/295 (100%)	0.13	4 (1%) 73 73	16, 34, 54, 80	9 (3%)
1	aba	295/295 (100%)	-0.24	1 (0%) 90 90	12, 29, 48, 79	4 (1%)
All	All	590/590 (100%)	-0.05	5 (0%) 82 83	12, 32, 52, 80	13 (2%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	aaa	23[A]	LYS	4.2
1	aaa	49	ALA	2.6
1	aaa	295	SER	2.1
1	aaa	114[A]	LYS	2.1
1	aba	295	SER	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](#)

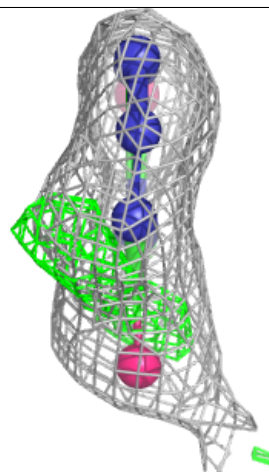
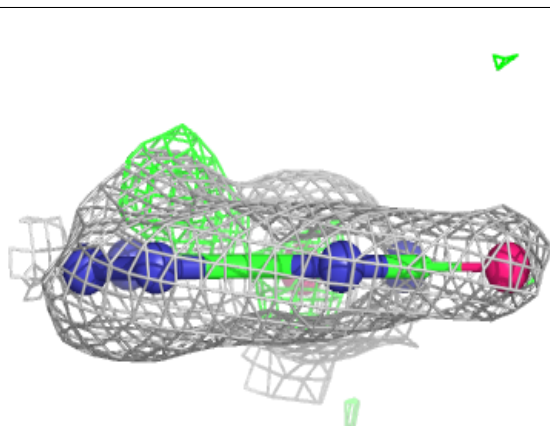
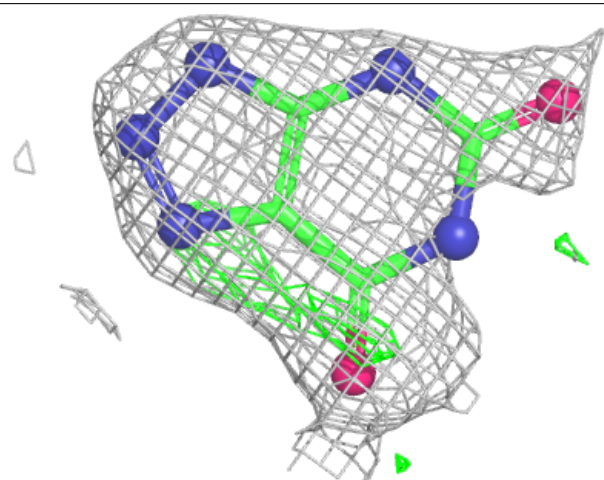
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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
2	ACE	aba	301	3/3	0.82	0.15	55,55,61,64	0
2	ACE	aaa	301	3/3	0.86	0.13	58,58,61,65	0
4	AZA	aaa	304	11/11	0.87	0.18	34,38,45,50	11
3	NA	aaa	302	1/1	0.89	0.11	46,46,46,46	0
4	AZA	aba	306	11/11	0.90	0.16	26,37,41,59	11
5	AR	aaa	306	1/1	0.90	0.38	44,44,44,44	1
5	AR	aaa	305	1/1	0.93	0.31	56,56,56,56	1
3	NA	aba	302	1/1	0.95	0.07	38,38,38,38	0
5	AR	aba	304	1/1	0.95	0.25	42,42,42,42	1
4	AZA	aaa	303	11/11	0.96	0.06	25,28,29,30	0
5	AR	aba	303	1/1	0.97	0.29	57,57,57,57	0
4	AZA	aba	305	11/11	0.98	0.04	20,23,26,26	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.

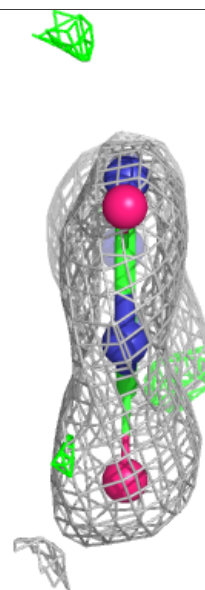
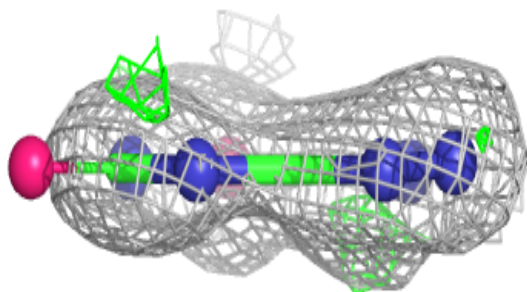
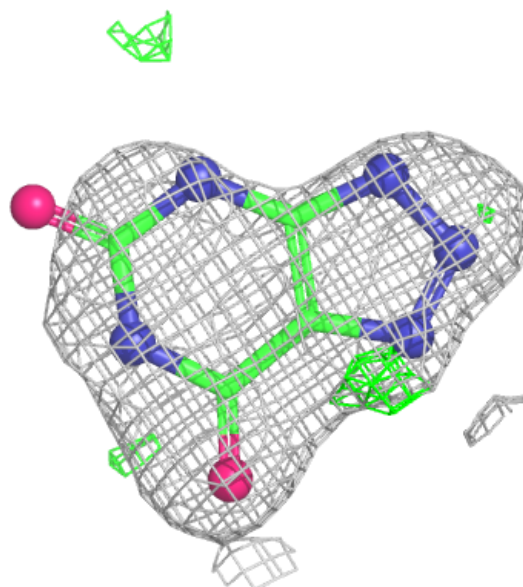
Electron density around AZA aaa 304:

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



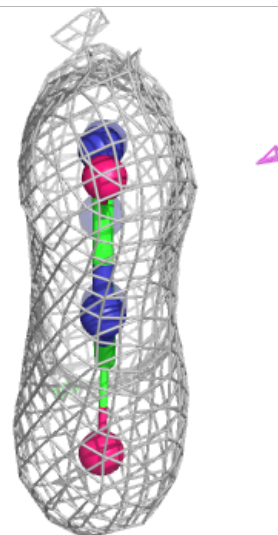
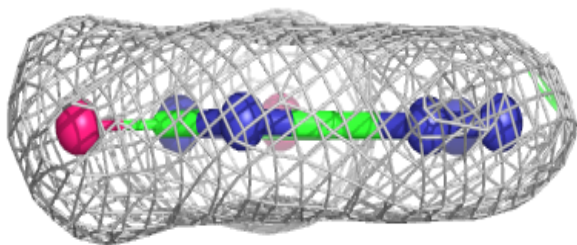
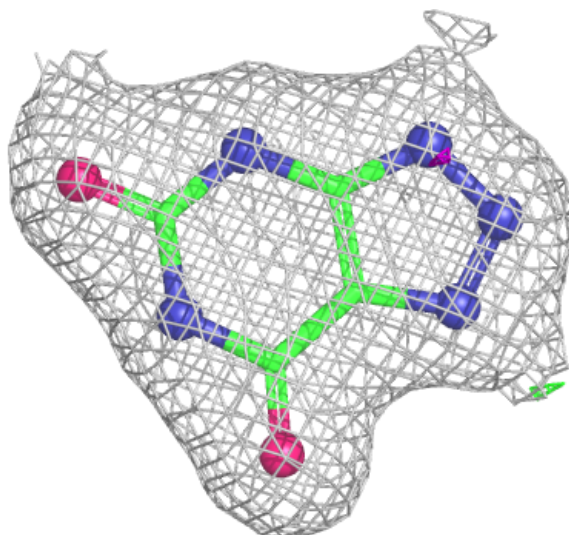
Electron density around AZA aba 306:

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



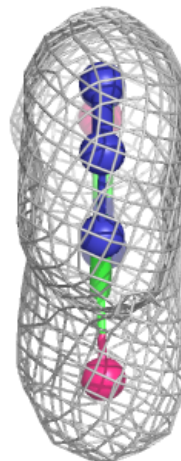
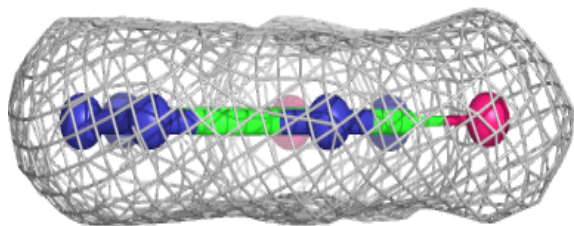
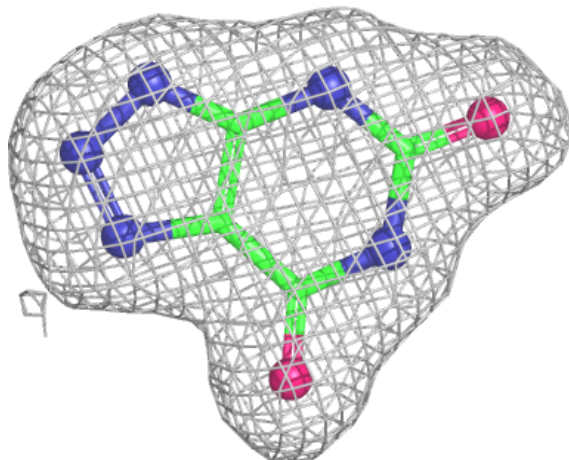
Electron density around AZA aaa 303:

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



Electron density around AZA aba 305:

$2mF_o-DF_c$ (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.