



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 21, 2024 – 12:49 am GMT

PDB ID : 7AKR
Title : Human ADP-ribosylserine hydrolase ARH3 mutant E41A in complex with ADP-ribose dimer
Authors : Ariza, A.
Deposited on : 2020-10-02
Resolution : 1.95 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

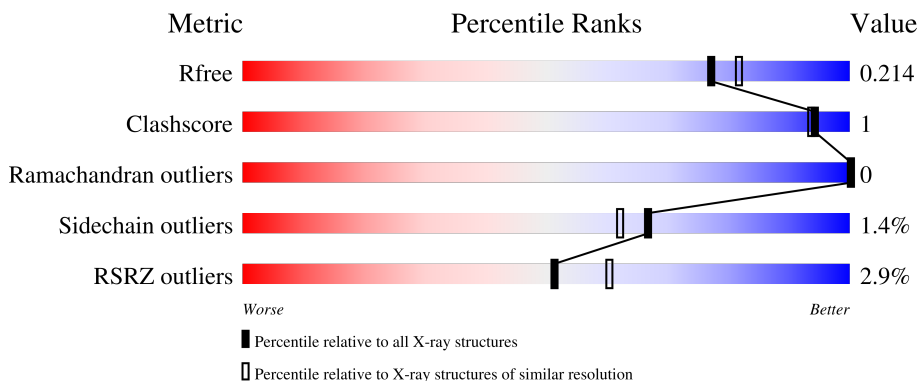
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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	349	 95%
1	BBB	349	 92%
1	CCC	349	 97%
1	DDD	349	 92%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11206 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 ADP-ribose glycohydrolase ARH3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	AAA	344	2647	1654	454	525	14	0	4	0
1	BBB	334	2568	1609	439	506	14	0	3	0
1	CCC	344	2652	1657	455	526	14	0	5	0
1	DDD	335	2574	1611	441	508	14	0	2	0

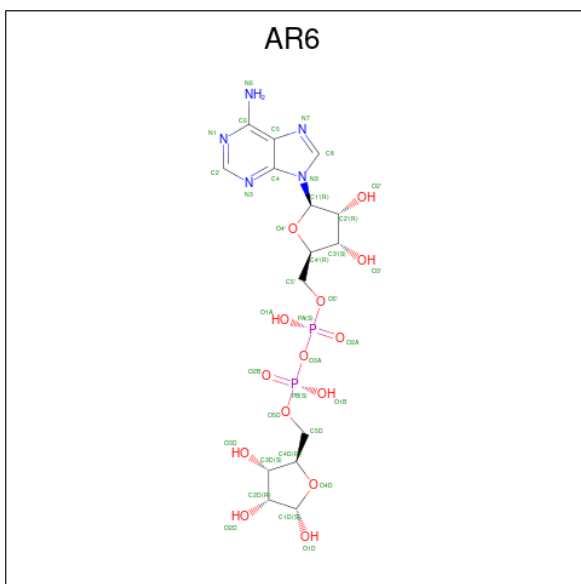
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AAA	15	GLY	-	expression tag	UNP Q9NX46
AAA	16	PRO	-	expression tag	UNP Q9NX46
AAA	17	GLY	-	expression tag	UNP Q9NX46
AAA	18	SER	-	expression tag	UNP Q9NX46
AAA	41	ALA	GLU	engineered mutation	UNP Q9NX46
BBB	15	GLY	-	expression tag	UNP Q9NX46
BBB	16	PRO	-	expression tag	UNP Q9NX46
BBB	17	GLY	-	expression tag	UNP Q9NX46
BBB	18	SER	-	expression tag	UNP Q9NX46
BBB	41	ALA	GLU	engineered mutation	UNP Q9NX46
CCC	15	GLY	-	expression tag	UNP Q9NX46
CCC	16	PRO	-	expression tag	UNP Q9NX46
CCC	17	GLY	-	expression tag	UNP Q9NX46
CCC	18	SER	-	expression tag	UNP Q9NX46
CCC	41	ALA	GLU	engineered mutation	UNP Q9NX46
DDD	15	GLY	-	expression tag	UNP Q9NX46
DDD	16	PRO	-	expression tag	UNP Q9NX46
DDD	17	GLY	-	expression tag	UNP Q9NX46
DDD	18	SER	-	expression tag	UNP Q9NX46
DDD	41	ALA	GLU	engineered mutation	UNP Q9NX46

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	AAA	2	Total	Mg	0	0
			2	2		
2	BBB	2	Total	Mg	0	0
			2	2		
2	CCC	2	Total	Mg	0	0
			2	2		
2	DDD	2	Total	Mg	0	0
			2	2		

- Molecule 3 is [(2R,3S,4R,5R)-5-(6-AMINOPURIN-9-YL)-3,4-DIHYDROXY-OXOLAN-2-YL]METHYL [HYDROXY-[[[(2R,3S,4R,5S)-3,4,5-TRIHYDROXYOXOLAN-2-YL]METHOXY]PHOSPHORYL] HYDROGEN PHOSPHATE (three-letter code: AR6) (formula: C₁₅H₂₃N₅O₁₄P₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	AAA	1	Total	C	N	O	P	0	0
			35	15	5	13	2		
3	AAA	1	Total	C	N	O	P	0	0
			24	10	5	7	2		
3	BBB	1	Total	C	N	O	P	0	0
			35	15	5	13	2		
3	BBB	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	CCC	1	Total	C	N	O	P	0	0
			35	15	5	13	2		

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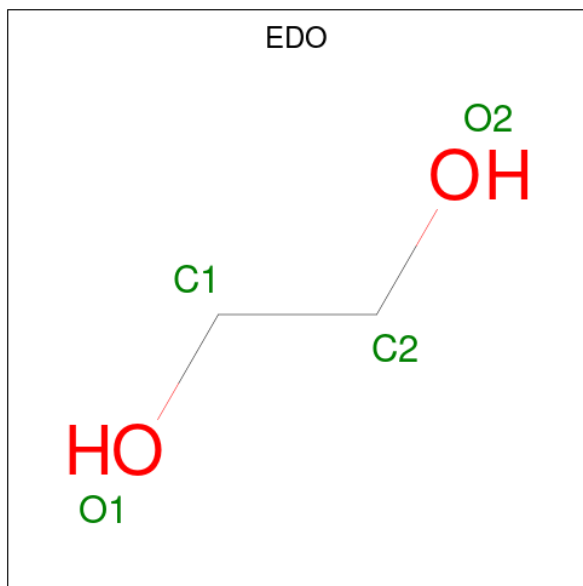
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	CCC	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	DDD	1	Total	C	N	O	P	0	0
			35	15	5	13	2		
3	DDD	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	AAA	1	Total	Cl	0	0
			1	1		
4	BBB	1	Total	Cl	0	0
			1	1		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
5	AAA	1	Total	C	O	0	0
			4	2	2		
5	AAA	1	Total	C	O	0	0
			4	2	2		
5	AAA	1	Total	C	O	0	0
			4	2	2		
5	AAA	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	AAA	1	Total C O 4 2 2	0	0
5	AAA	1	Total C O 4 2 2	0	0
5	AAA	1	Total C O 4 2 2	0	0
5	BBB	1	Total C O 4 2 2	0	0
5	BBB	1	Total C O 4 2 2	0	0
5	CCC	1	Total C O 4 2 2	0	0
5	CCC	1	Total C O 4 2 2	0	0
5	CCC	1	Total C O 4 2 2	0	0
5	CCC	1	Total C O 4 2 2	0	0
5	CCC	1	Total C O 4 2 2	0	0
5	CCC	1	Total C O 4 2 2	0	0
5	DDD	1	Total C O 4 2 2	0	0
5	DDD	1	Total C O 4 2 2	0	0
5	DDD	1	Total C O 4 2 2	0	0
5	DDD	1	Total C O 4 2 2	0	0

- Molecule 6 is water.

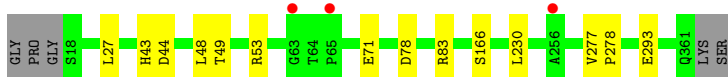
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	AAA	179	Total O 179 179	0	0
6	BBB	82	Total O 82 82	0	0
6	CCC	114	Total O 115 115	0	1
6	DDD	62	Total O 62 62	0	0

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.

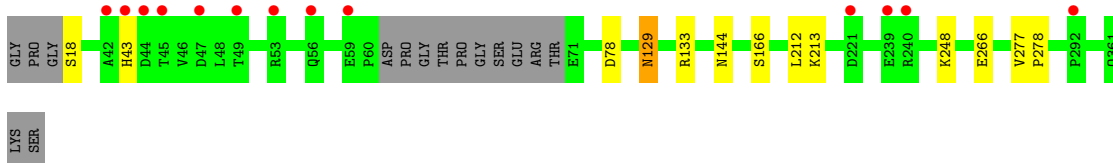
- Molecule 1: ADP-ribose glycohydrolase ARH3

Chain AAA:  95%



- Molecule 1: ADP-ribose glycohydrolase ARH3

Chain BBB:  92%



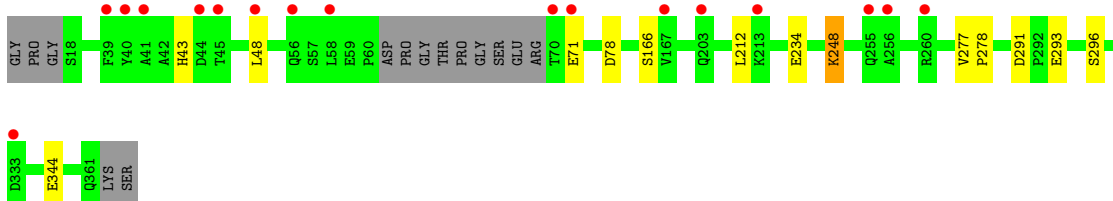
- Molecule 1: ADP-ribose glycohydrolase ARH3

Chain CCC:  97%



- Molecule 1: ADP-ribose glycohydrolase ARH3

Chain DDD:  92%



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	86.62Å 91.60Å 91.09Å 90.00° 105.47° 90.00°	Depositor
Resolution (Å)	70.63 – 1.95 70.63 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.9 (70.63-1.95) 92.4 (70.63-1.95)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.22 (at 1.95Å)	Xtrriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.187 , 0.211 0.192 , 0.214	Depositor DCC
R_{free} test set	4757 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	30.5	Xtrriage
Anisotropy	0.324	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.019 for l,-k,h	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11206	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.35% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: AR6, CL, MG, EDO

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	0.69	0/2705	0.76	0/3659
1	BBB	0.69	0/2622	0.75	1/3542 (0.0%)
1	CCC	0.68	0/2713	0.74	1/3670 (0.0%)
1	DDD	0.66	0/2623	0.72	0/3545
All	All	0.68	0/10663	0.74	2/14416 (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	BBB	129	ASN	CB-CA-C	5.03	120.45	110.40
1	CCC	129	ASN	CB-CA-C	5.03	120.45	110.40

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	AAA	2647	0	2567	10	0
1	BBB	2568	0	2502	8	0
1	CCC	2652	0	2573	2	0
1	DDD	2574	0	2496	7	0
2	AAA	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	BBB	2	0	0	0	0
2	CCC	2	0	0	0	0
2	DDD	2	0	0	0	0
3	AAA	59	0	28	2	0
3	BBB	62	0	28	0	0
3	CCC	62	0	28	0	0
3	DDD	62	0	28	1	0
4	AAA	1	0	0	0	0
4	BBB	1	0	0	1	0
5	AAA	28	0	42	3	0
5	BBB	8	0	12	0	0
5	CCC	20	0	30	0	0
5	DDD	16	0	24	0	0
6	AAA	179	0	0	1	0
6	BBB	82	0	0	1	0
6	CCC	115	0	0	1	0
6	DDD	62	0	0	0	0
All	All	11206	0	10358	27	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 1.

The worst 5 of 27 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:BBB:213:LYS:HG2	1:DDD:296:SER:HB3	1.59	0.85
1:AAA:71:GLU:OE2	1:BBB:18:SER:HA	1.86	0.76
1:CCC:129:ASN:ND2	6:CCC:501:HOH:O	2.25	0.69
1:DDD:71:GLU:HB3	1:DDD:344:GLU:HA	1.76	0.68
1:BBB:248[A]:LYS:HE3	1:BBB:266:GLU:OE1	1.97	0.63

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	AAA	346/349 (99%)	338 (98%)	8 (2%)	0	100	100
1	BBB	333/349 (95%)	327 (98%)	6 (2%)	0	100	100
1	CCC	347/349 (99%)	340 (98%)	7 (2%)	0	100	100
1	DDD	333/349 (95%)	328 (98%)	5 (2%)	0	100	100
All	All	1359/1396 (97%)	1333 (98%)	26 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	AAA	280/279 (100%)	278 (99%)	2 (1%)	84	82
1	BBB	271/279 (97%)	267 (98%)	4 (2%)	65	60
1	CCC	281/279 (101%)	277 (99%)	4 (1%)	67	62
1	DDD	271/279 (97%)	266 (98%)	5 (2%)	59	53
All	All	1103/1116 (99%)	1088 (99%)	15 (1%)	67	62

5 of 15 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	CCC	78	ASP
1	DDD	234	GLU
1	CCC	129	ASN
1	DDD	248	LYS
1	DDD	78	ASP

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

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 36 ligands modelled in this entry, 10 are monoatomic - leaving 26 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	$\# Z > 2$
3	AR6	BBB	403	3,2	33,38,39	0.66	0	37,58,60	0.93	2 (5%)
5	EDO	CCC	407	-	3,3,3	0.12	0	2,2,2	0.42	0
5	EDO	AAA	409	-	3,3,3	0.15	0	2,2,2	0.24	0
5	EDO	CCC	405	-	3,3,3	0.15	0	2,2,2	0.43	0
5	EDO	AAA	408	-	3,3,3	0.04	0	2,2,2	0.18	0
5	EDO	BBB	406	-	3,3,3	0.27	0	2,2,2	0.47	0
5	EDO	DDD	407	-	3,3,3	0.04	0	2,2,2	0.07	0
5	EDO	AAA	406	-	3,3,3	0.31	0	2,2,2	0.70	0
3	AR6	CCC	403	3,2	33,38,39	0.59	0	37,58,60	0.86	2 (5%)
3	AR6	DDD	403	3,2	33,38,39	0.62	0	37,58,60	0.92	1 (2%)
5	EDO	AAA	407	-	3,3,3	0.22	0	2,2,2	0.19	0
5	EDO	DDD	406	-	3,3,3	0.17	0	2,2,2	0.27	0
5	EDO	AAA	412	-	3,3,3	0.03	0	2,2,2	0.05	0
3	AR6	BBB	404	3	24,29,39	0.69	0	29,45,60	0.91	2 (6%)
5	EDO	CCC	406	-	3,3,3	0.24	0	2,2,2	0.65	0
5	EDO	DDD	408	-	3,3,3	0.07	0	2,2,2	0.17	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	AR6	DDD	404	3	24,29,39	1.00	1 (4%)	29,45,60	1.11	2 (6%)
5	EDO	CCC	408	-	3,3,3	0.12	0	2,2,2	0.19	0
5	EDO	CCC	409	-	3,3,3	0.08	0	2,2,2	0.36	0
5	EDO	BBB	407	-	3,3,3	0.16	0	2,2,2	0.17	0
5	EDO	AAA	410	-	3,3,3	0.06	0	2,2,2	0.31	0
5	EDO	AAA	411	-	3,3,3	0.25	0	2,2,2	0.34	0
3	AR6	CCC	404	3	24,29,39	0.74	0	29,45,60	0.94	2 (6%)
5	EDO	DDD	405	-	3,3,3	0.17	0	2,2,2	0.77	0
3	AR6	AAA	403	3,2	33,38,39	0.70	0	37,58,60	0.78	0
3	AR6	AAA	404	3	21,26,39	0.72	0	24,39,60	0.78	1 (4%)

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	AR6	BBB	403	3,2	-	2/18/51/54	0/4/4/4
5	EDO	CCC	407	-	-	0/1/1/1	-
5	EDO	AAA	409	-	-	1/1/1/1	-
5	EDO	CCC	405	-	-	1/1/1/1	-
5	EDO	AAA	408	-	-	1/1/1/1	-
5	EDO	BBB	406	-	-	0/1/1/1	-
5	EDO	DDD	407	-	-	1/1/1/1	-
5	EDO	AAA	406	-	-	0/1/1/1	-
3	AR6	CCC	403	3,2	-	4/18/51/54	0/4/4/4
3	AR6	DDD	403	3,2	-	4/18/51/54	0/4/4/4
5	EDO	AAA	407	-	-	1/1/1/1	-
5	EDO	DDD	406	-	-	1/1/1/1	-
5	EDO	AAA	412	-	-	1/1/1/1	-
3	AR6	BBB	404	3	-	3/12/32/54	0/3/3/4
5	EDO	CCC	406	-	-	0/1/1/1	-
5	EDO	DDD	408	-	-	1/1/1/1	-
3	AR6	DDD	404	3	-	5/12/32/54	0/3/3/4
5	EDO	CCC	408	-	-	1/1/1/1	-
5	EDO	CCC	409	-	-	0/1/1/1	-
5	EDO	BBB	407	-	-	1/1/1/1	-
5	EDO	AAA	410	-	-	1/1/1/1	-
5	EDO	AAA	411	-	-	1/1/1/1	-
3	AR6	CCC	404	3	-	6/12/32/54	0/3/3/4
5	EDO	DDD	405	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	AR6	AAA	403	3,2	-	3/18/51/54	0/4/4/4
3	AR6	AAA	404	3	-	0/6/29/54	0/3/3/4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	DDD	404	AR6	PB-O2B	3.41	1.61	1.50

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	DDD	403	AR6	O4D-C4D-C3D	2.74	107.13	104.70
3	DDD	404	AR6	PA-O3A-PB	2.63	141.84	132.83
3	CCC	404	AR6	PA-O3A-PB	2.44	141.22	132.83
3	DDD	404	AR6	O5D-PB-O1B	2.39	116.78	107.64
3	CCC	403	AR6	O4D-C4D-C3D	2.32	106.76	104.70

There are no chirality outliers.

5 of 39 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	AAA	403	AR6	PB-O3A-PA-O5'
3	BBB	404	AR6	C5'-O5'-PA-O1A
3	BBB	404	AR6	C5'-O5'-PA-O3A
3	CCC	403	AR6	PB-O3A-PA-O5'
3	CCC	404	AR6	PB-O3A-PA-O5'

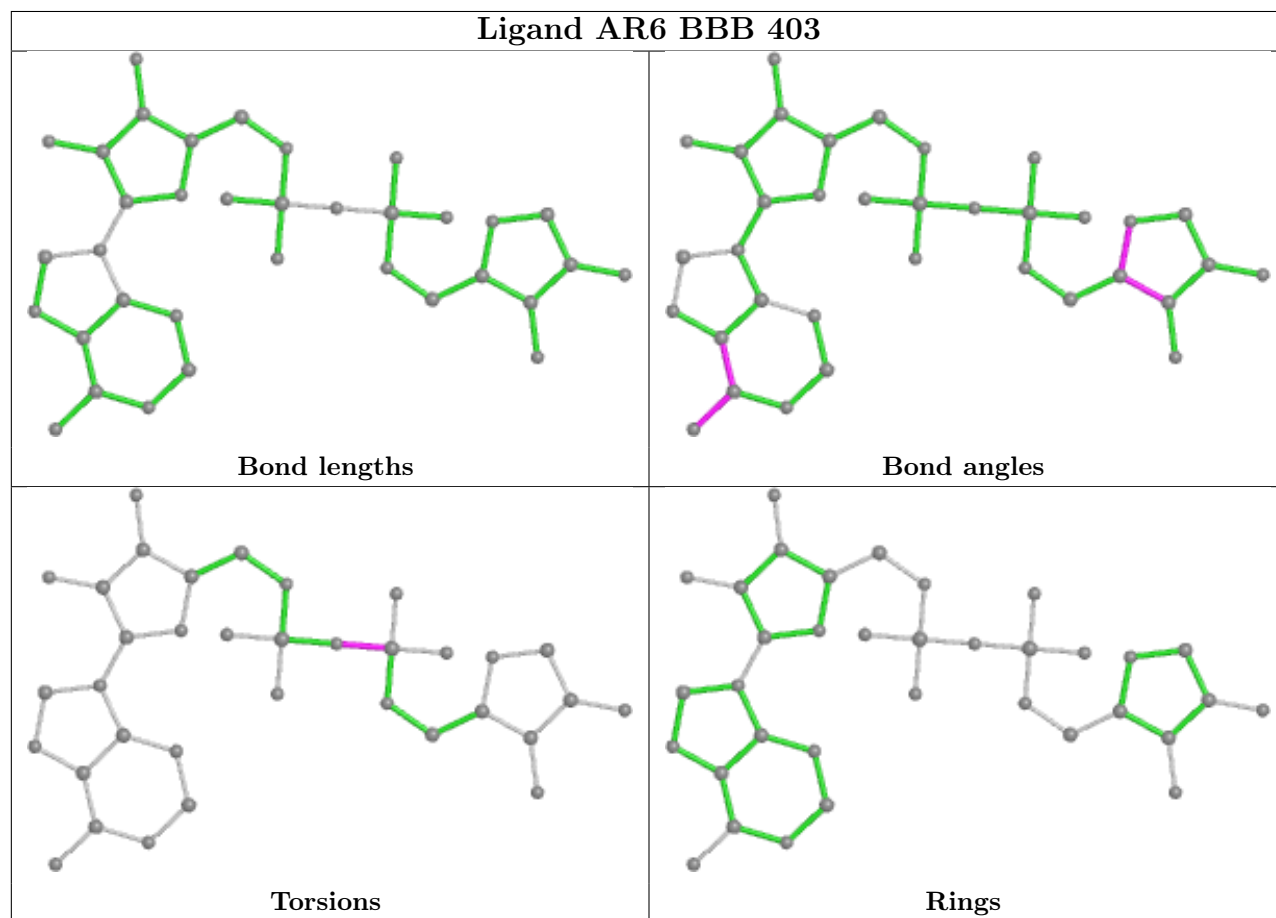
There are no ring outliers.

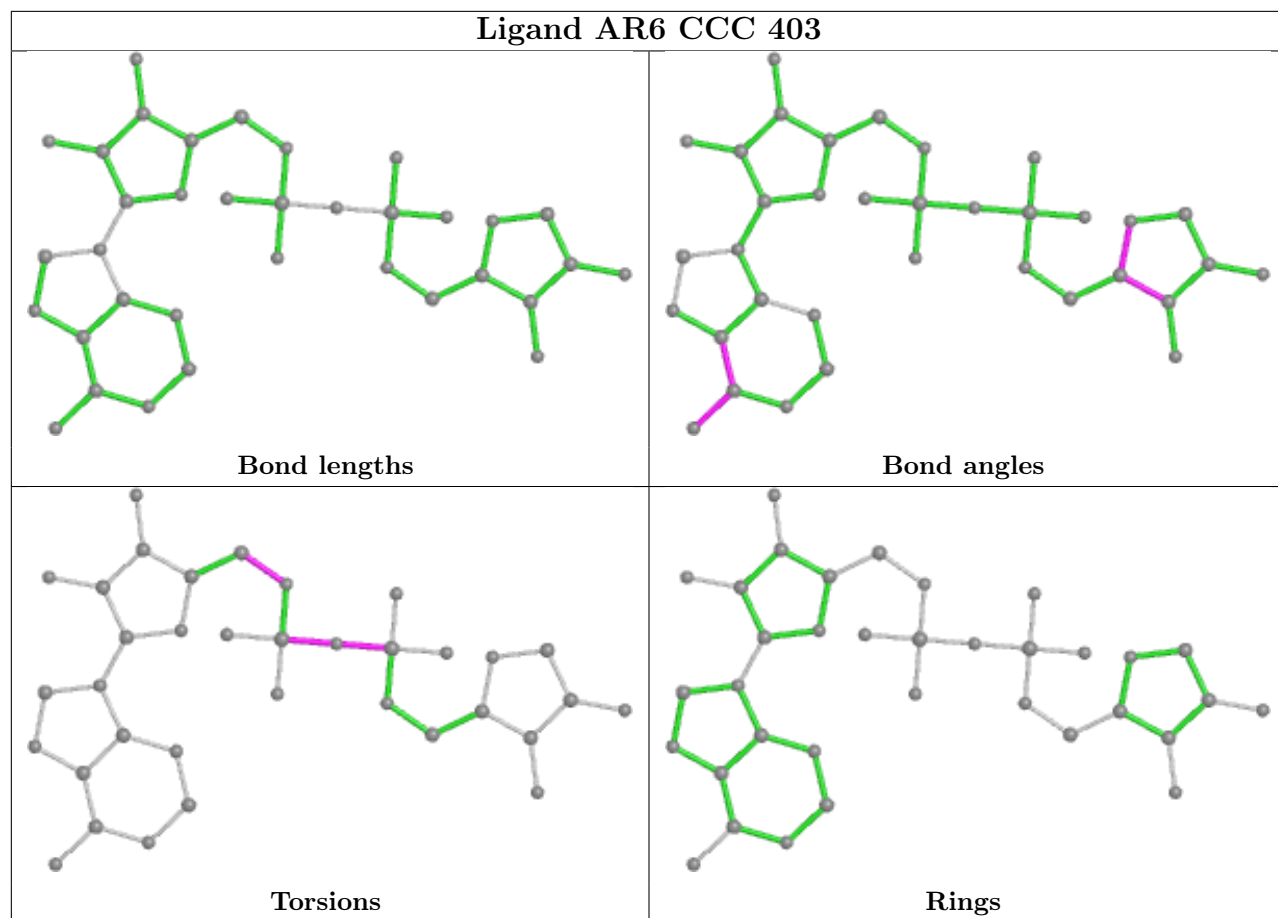
4 monomers are involved in 6 short contacts:

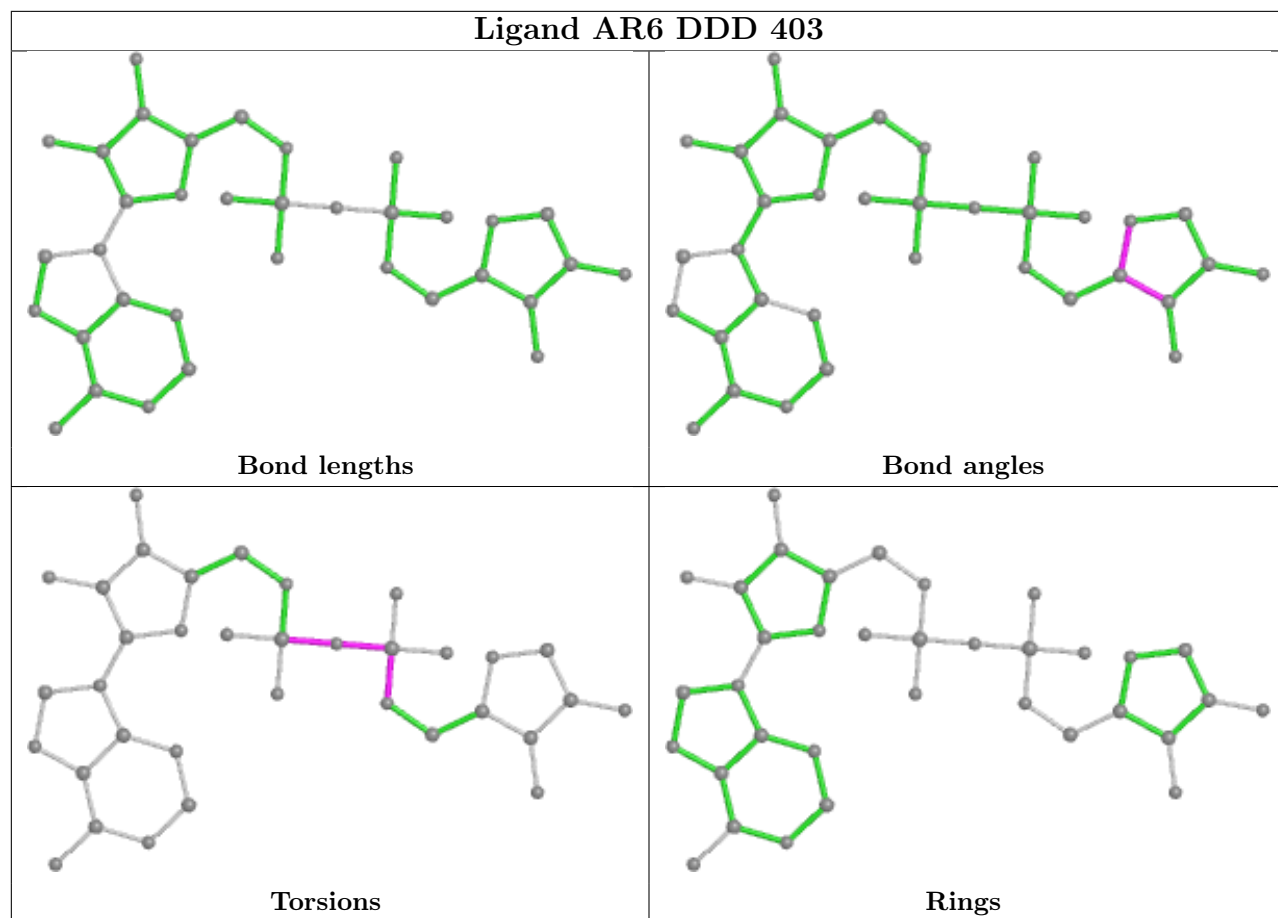
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	AAA	409	EDO	1	0
5	AAA	406	EDO	2	0
3	DDD	404	AR6	1	0
3	AAA	404	AR6	2	0

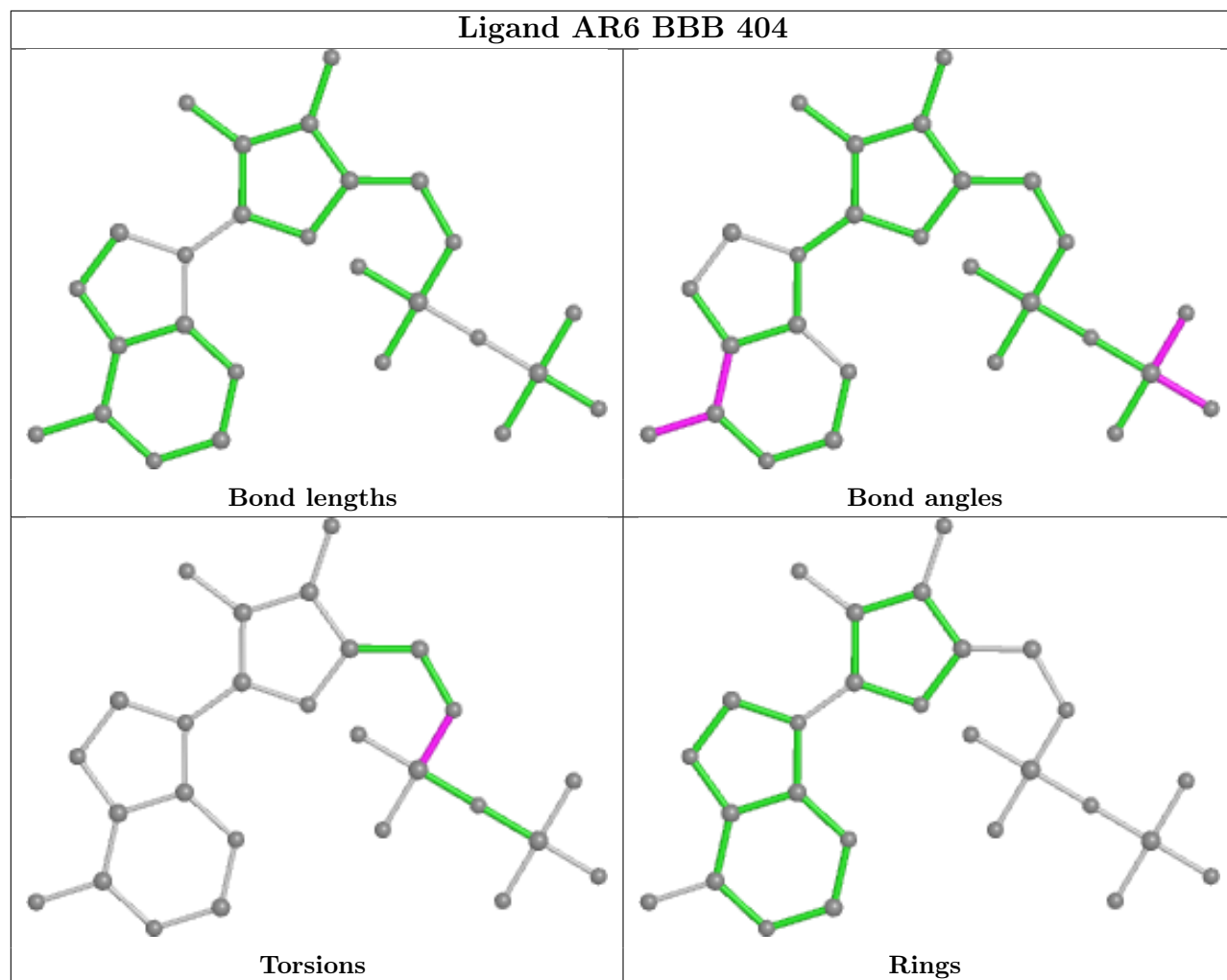
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

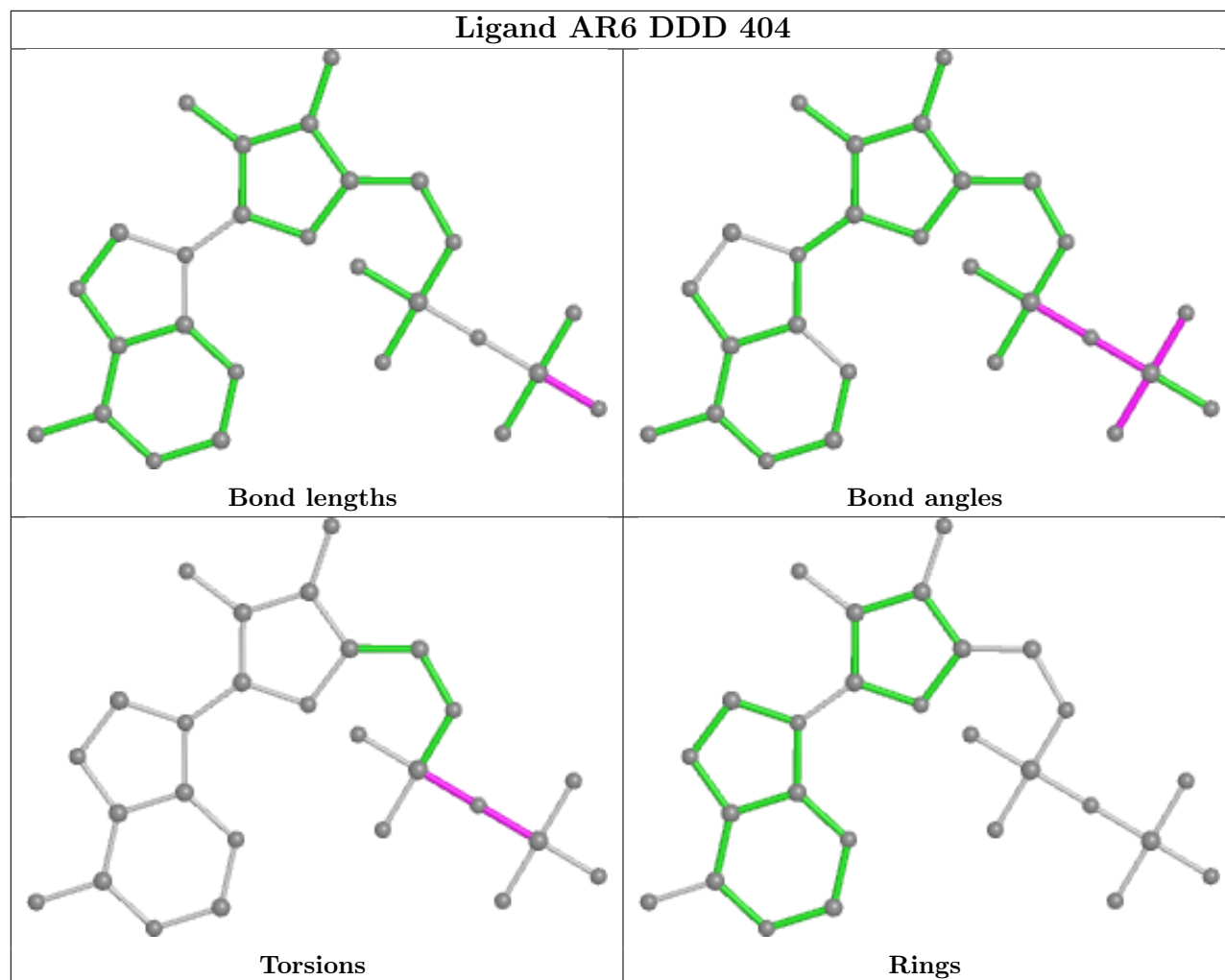
Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD 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.

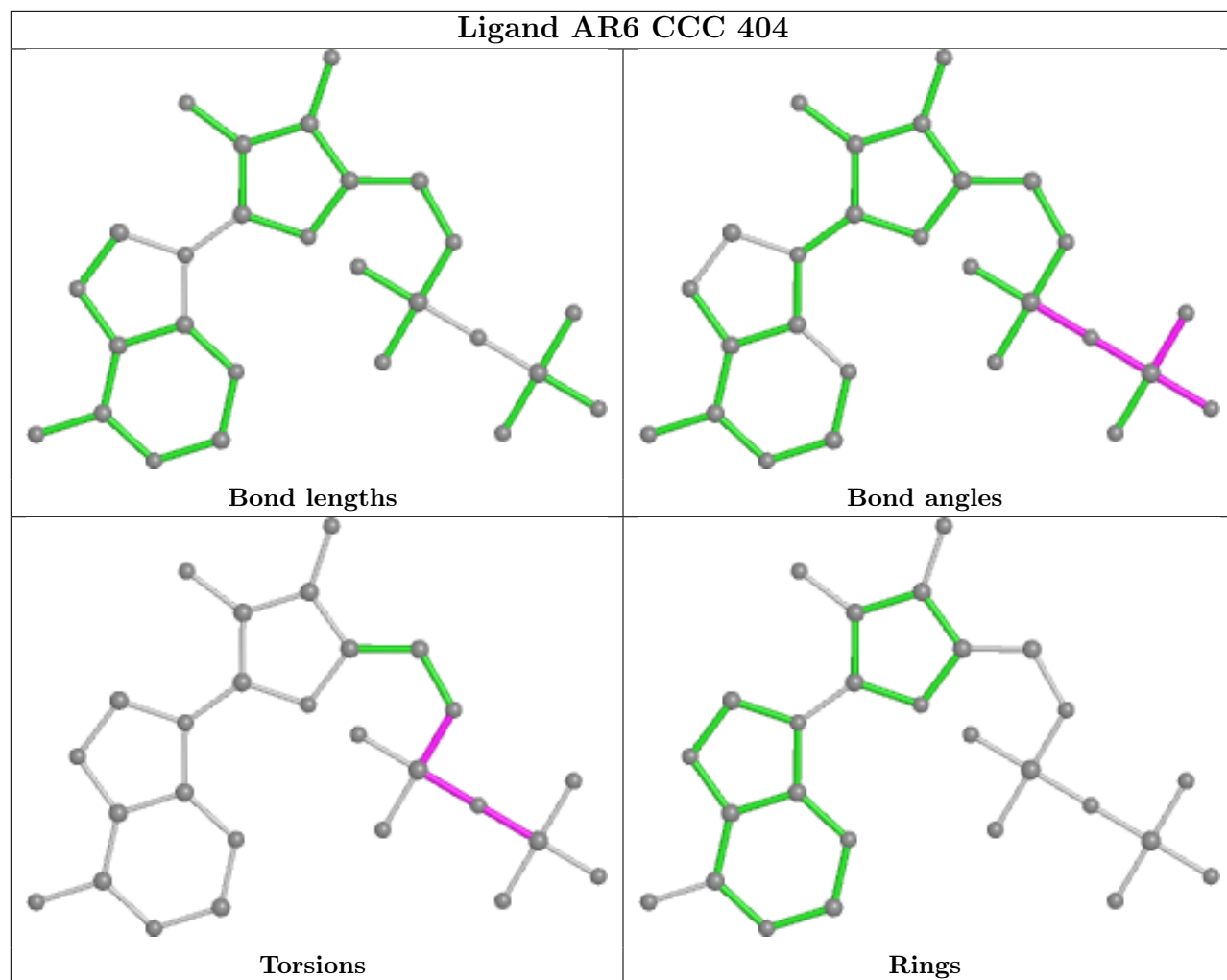


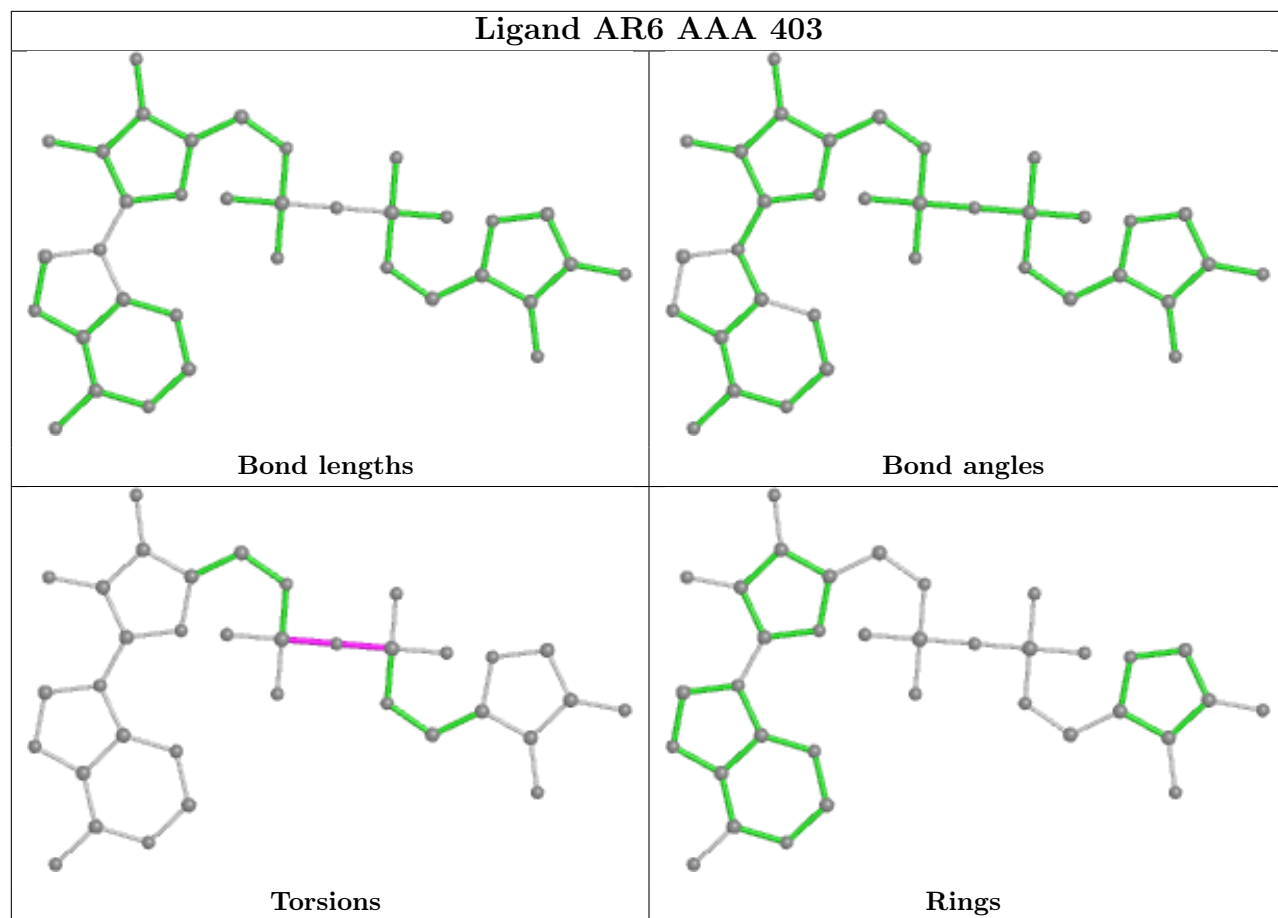


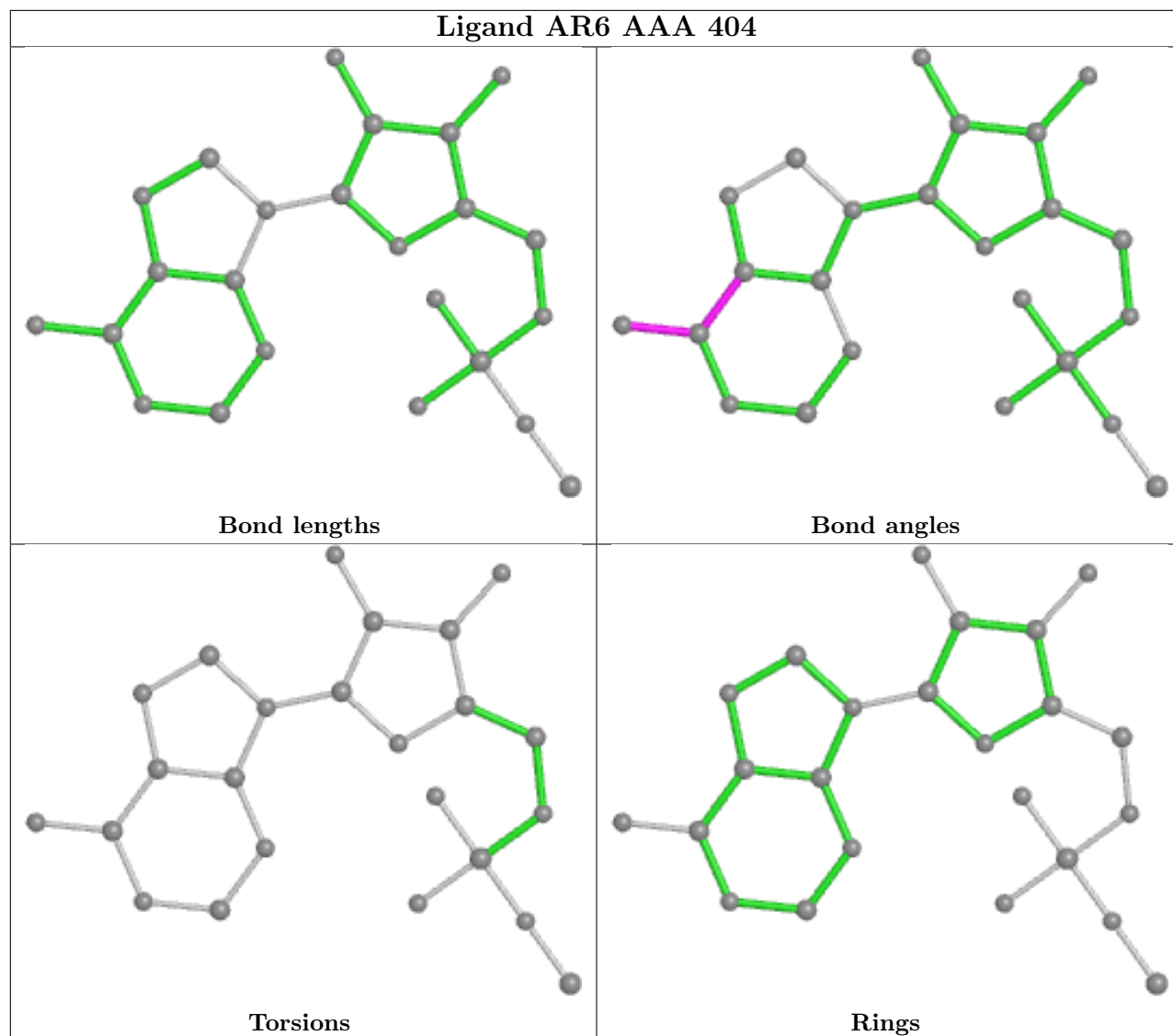












5.7 Other polymers [i](#)

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	AAA	344/349 (98%)	0.09	3 (0%) 84 89	21, 32, 52, 65	0
1	BBB	334/349 (95%)	0.09	13 (3%) 39 49	29, 43, 80, 114	0
1	CCC	344/349 (98%)	0.01	6 (1%) 70 77	29, 41, 72, 91	0
1	DDD	335/349 (95%)	0.24	17 (5%) 28 37	37, 52, 85, 103	0
All	All	1357/1396 (97%)	0.11	39 (2%) 51 60	21, 42, 77, 114	0

The worst 5 of 39 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BBB	44	ASP	8.8
1	DDD	44	ASP	6.5
1	BBB	42	ALA	5.5
1	DDD	70	THR	4.8
1	BBB	45	THR	4.7

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

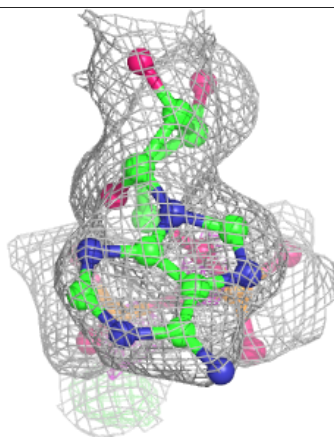
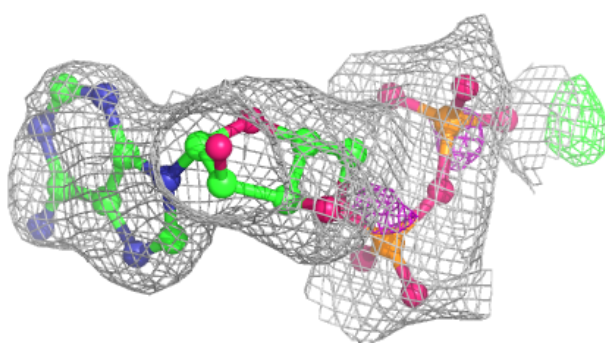
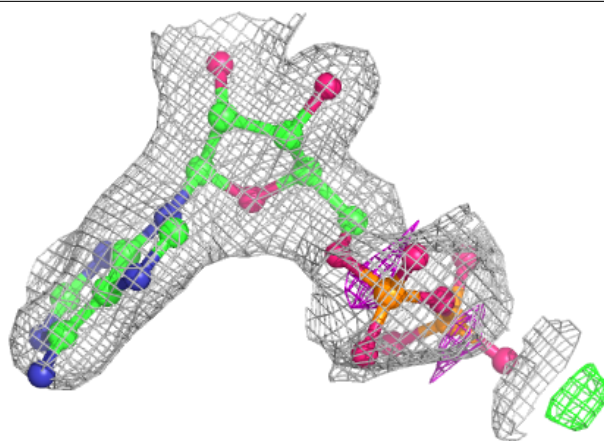
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	EDO	AAA	411	4/4	0.76	0.20	60,65,65,69	0
5	EDO	AAA	409	4/4	0.80	0.26	50,55,58,62	0
4	CL	AAA	405	1/1	0.81	0.07	61,61,61,61	0
5	EDO	AAA	408	4/4	0.84	0.13	58,58,59,59	0
5	EDO	DDD	408	4/4	0.85	0.13	68,69,69,70	0
5	EDO	CCC	408	4/4	0.86	0.17	62,66,66,67	0
3	AR6	CCC	404	27/36	0.87	0.15	39,56,90,95	0
5	EDO	AAA	412	4/4	0.87	0.28	55,59,61,61	0
5	EDO	BBB	407	4/4	0.87	0.21	59,60,60,60	0
3	AR6	BBB	404	27/36	0.87	0.14	39,57,96,100	0
5	EDO	AAA	410	4/4	0.87	0.18	68,70,72,73	0
3	AR6	DDD	404	27/36	0.88	0.14	44,57,94,100	0
5	EDO	DDD	405	4/4	0.89	0.13	48,48,49,52	0
4	CL	BBB	405	1/1	0.92	0.13	67,67,67,67	0
5	EDO	CCC	409	4/4	0.92	0.14	51,52,53,54	0
5	EDO	DDD	406	4/4	0.93	0.25	52,53,54,56	0
5	EDO	AAA	406	4/4	0.93	0.23	39,42,44,44	0
5	EDO	CCC	405	4/4	0.94	0.16	54,55,57,57	0
5	EDO	BBB	406	4/4	0.94	0.12	46,46,46,47	0
3	AR6	AAA	404	24/36	0.95	0.11	27,34,45,70	0
5	EDO	AAA	407	4/4	0.95	0.15	49,51,52,54	0
2	MG	DDD	401	1/1	0.95	0.09	39,39,39,39	0
5	EDO	CCC	407	4/4	0.96	0.11	43,44,44,47	0
5	EDO	DDD	407	4/4	0.96	0.10	62,63,63,64	0
2	MG	CCC	402	1/1	0.96	0.03	29,29,29,29	0
2	MG	BBB	402	1/1	0.97	0.03	34,34,34,34	0
2	MG	DDD	402	1/1	0.97	0.06	42,42,42,42	0
2	MG	CCC	401	1/1	0.97	0.08	29,29,29,29	0
2	MG	AAA	402	1/1	0.97	0.13	23,23,23,23	0
5	EDO	CCC	406	4/4	0.97	0.08	42,47,47,48	0
3	AR6	CCC	403	35/36	0.97	0.08	29,38,46,47	0
3	AR6	AAA	403	35/36	0.98	0.11	23,26,30,30	0
2	MG	BBB	401	1/1	0.98	0.14	32,32,32,32	0
3	AR6	DDD	403	35/36	0.98	0.07	38,41,45,45	0
2	MG	AAA	401	1/1	0.98	0.11	23,23,23,23	0
3	AR6	BBB	403	35/36	0.99	0.08	29,32,35,37	0

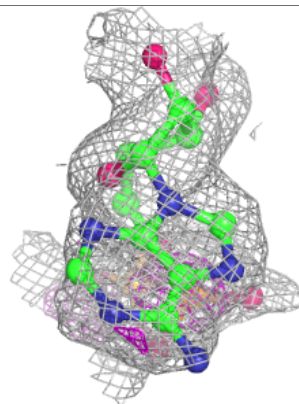
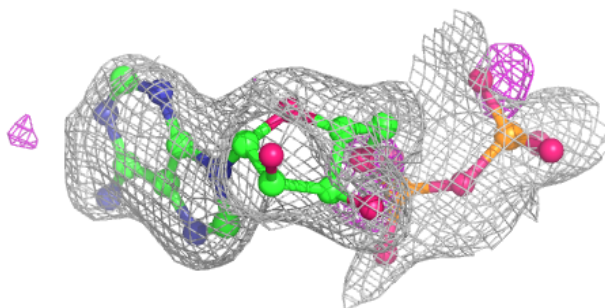
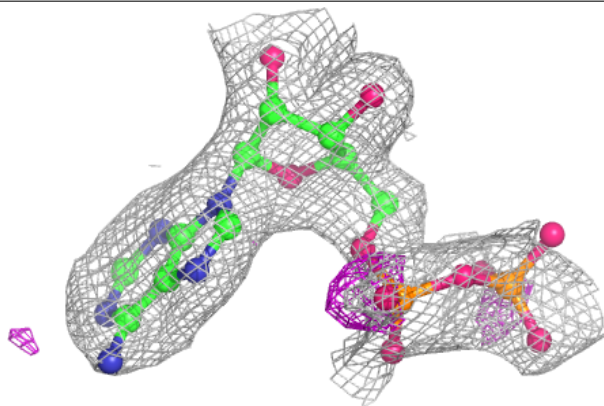
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around AR6 CCC 404:

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

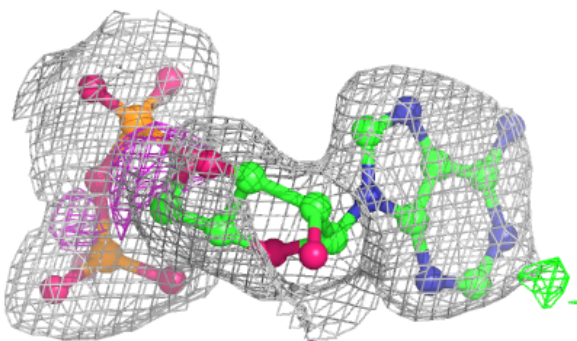
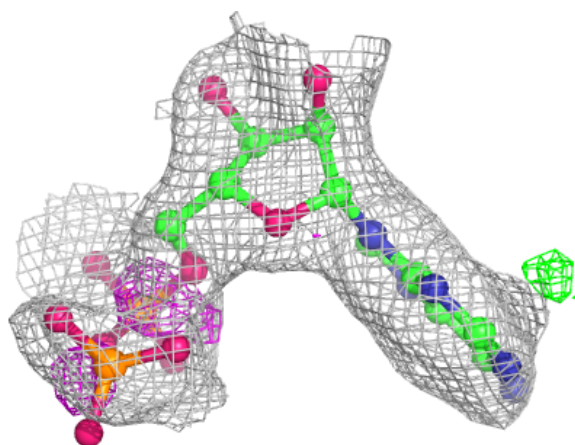
**Electron density around AR6 BBB 404:**

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



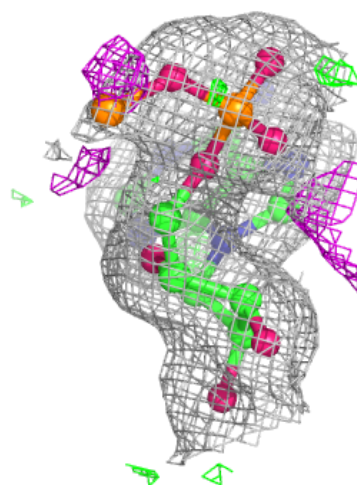
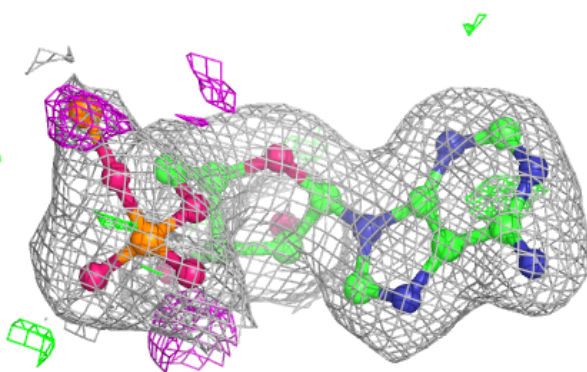
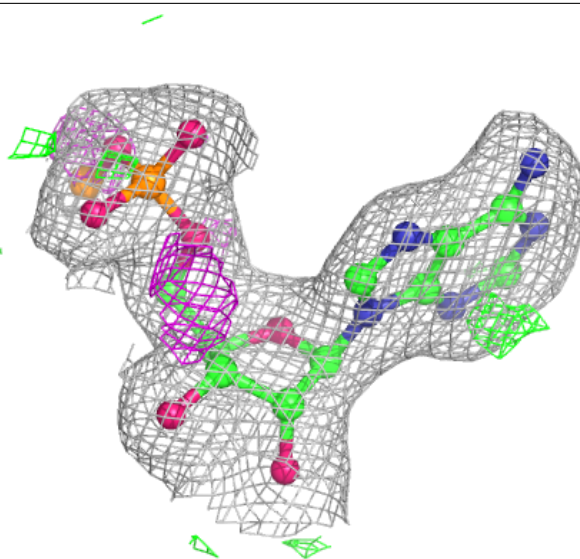
Electron density around AR6 DDD 404:

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



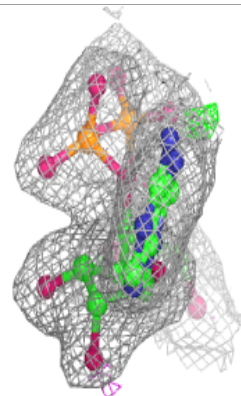
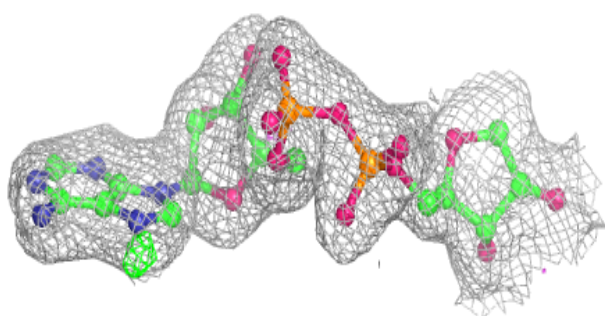
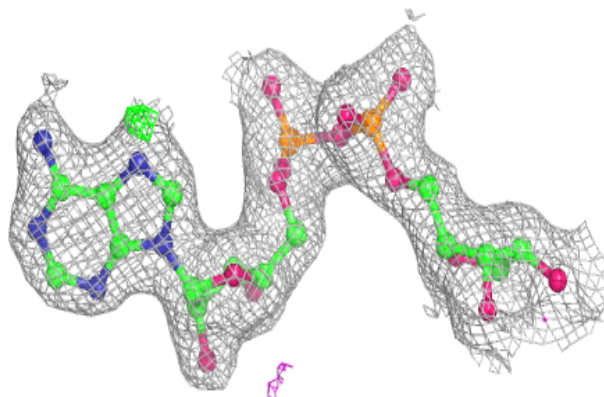
Electron density around AR6 AAA 404:

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

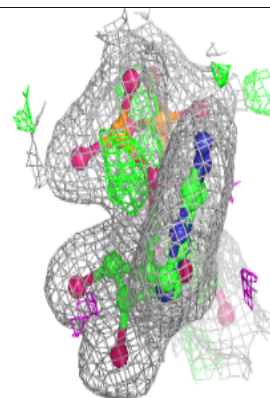
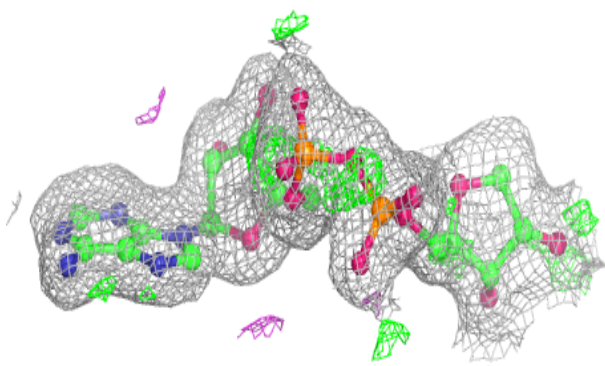
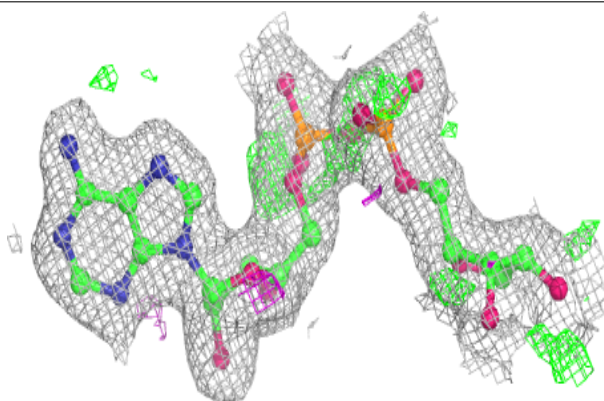


Electron density around AR6 CCC 403:

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

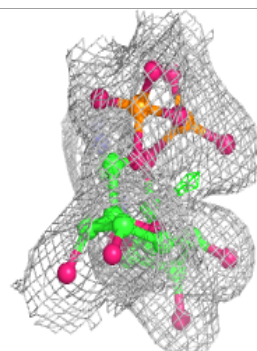
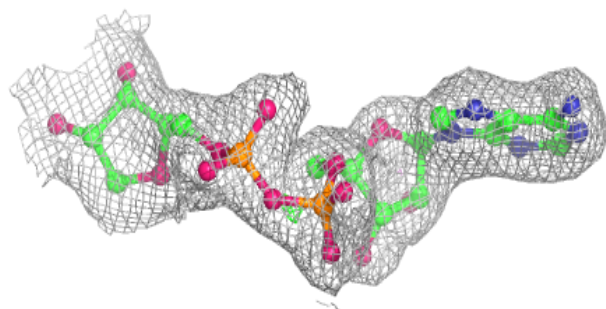
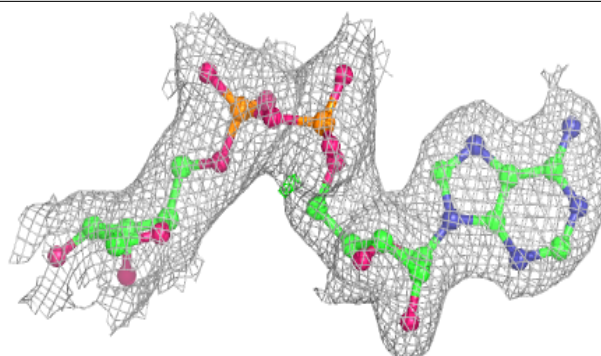
**Electron density around AR6 AAA 403:**

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and green (positive)

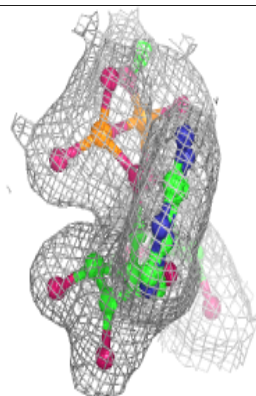
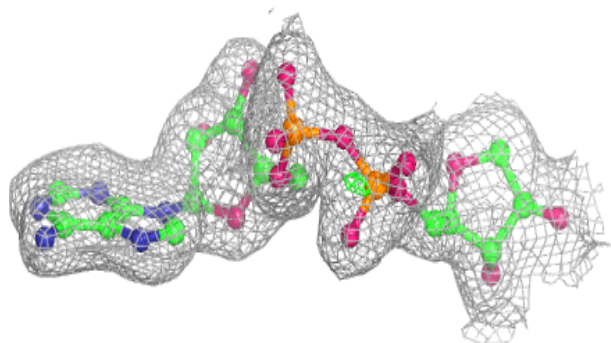
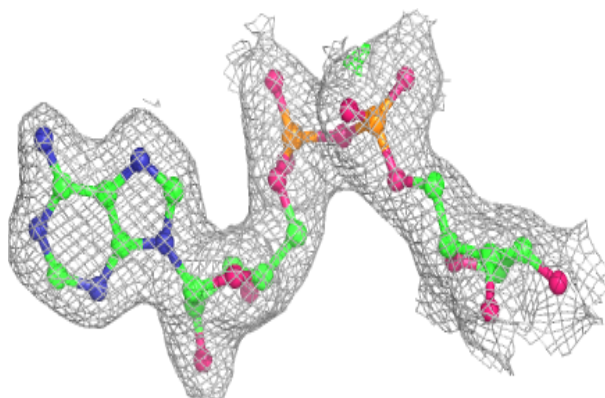


Electron density around AR6 DDD 403:

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

**Electron density around AR6 BBB 403:**

$2mF_o-DF_c$ (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.