



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 02:02 PM UTC

PDB ID : 5VP1 / pdb_00005vp1
Title : Discovery of Clinical Candidate N-{(1S)-1-[3-Fluoro-4-(trifluoromethoxy)phenyl]-2-methoxyethyl}-7-methoxy-2-oxo-2,3-dihydropyrido[2,3-b]pyrazine-4(1H)-carboxamide (TAK-915), A Highly Potent, Selective, and Brain-Penetrating Phosphodiesterase 2A Inhibitor for the Treatment of Cognitive Disorders
Authors : Hoffman, I.D.
Deposited on : 2017-05-03
Resolution : 1.86 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](#) ①) 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

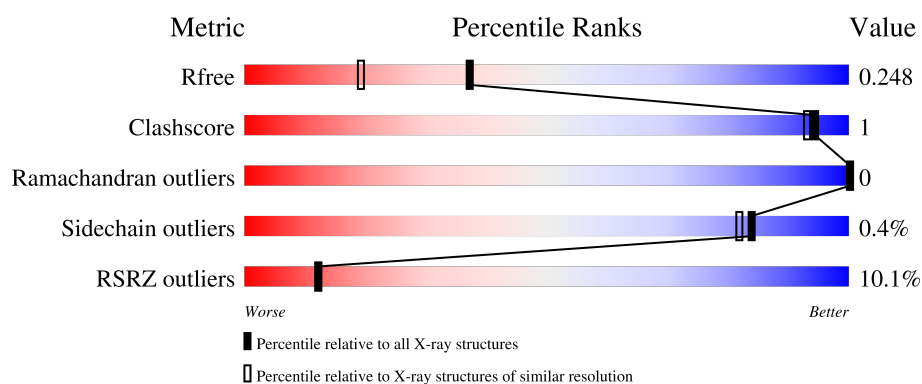
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.86 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	345	12% 94% ..
1	B	345	6% 95% ..
1	C	345	10% 89% 9% .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8522 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 cGMP-dependent 3',5'-cyclic phosphodiesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	338	Total	C	N	O	S	0	1	0
			2768	1763	474	506	25			
1	B	342	Total	C	N	O	S	0	1	0
			2793	1777	480	511	25			
1	C	313	Total	C	N	O	S	0	1	0
			2578	1644	445	464	25			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	575	GLY	-	expression tag	UNP O00408
A	576	HIS	-	expression tag	UNP O00408
A	577	ALA	-	expression tag	UNP O00408
B	575	GLY	-	expression tag	UNP O00408
B	576	HIS	-	expression tag	UNP O00408
B	577	ALA	-	expression tag	UNP O00408
C	575	GLY	-	expression tag	UNP O00408
C	576	HIS	-	expression tag	UNP O00408
C	577	ALA	-	expression tag	UNP O00408

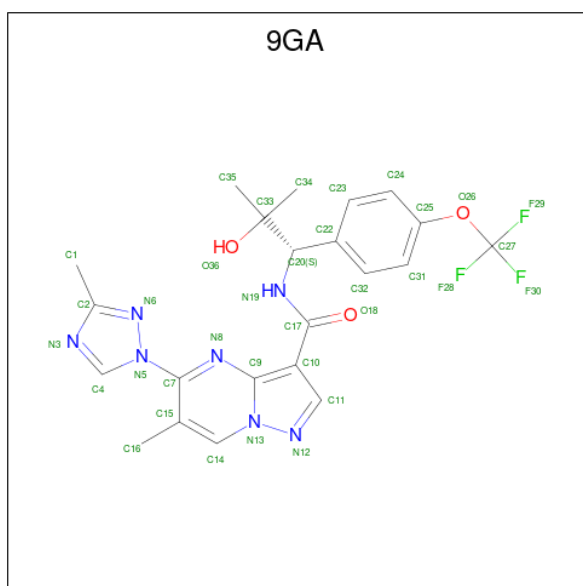
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Mg 2 2	0	0
3	B	1	Total Mg 1 1	0	0
3	C	1	Total Mg 1 1	0	0

- Molecule 4 is N-{(1S)-2-hydroxy-2-methyl-1-[4-(trifluoromethoxy)phenyl]propyl}-6-methyl-5-(3-methyl-1H-1,2,4-triazol-1-yl)pyrazolo[1,5-a]pyrimidine-3-carboxamide (CCD ID: 9GA) (formula: C₂₂H₂₂F₃N₇O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C F N O 35 22 3 7 3	0	0
4	B	1	Total C F N O 35 22 3 7 3	0	0
4	C	1	Total C F N O 35 22 3 7 3	0	0

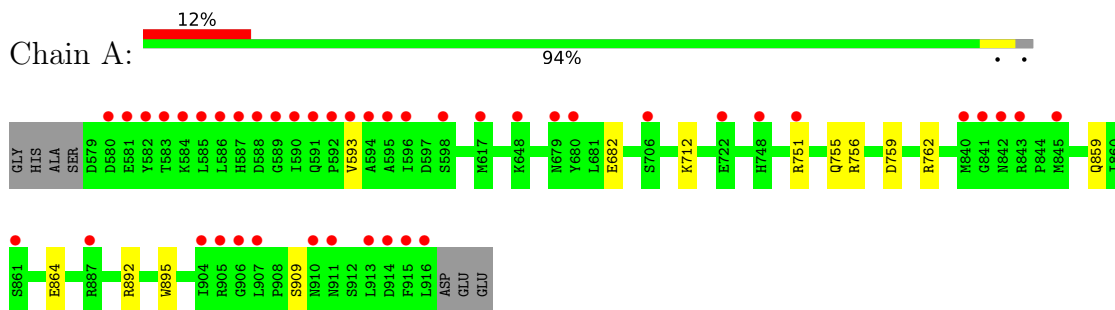
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	107	Total O 107 107	0	0
5	B	87	Total O 87 87	0	0
5	C	77	Total O 77 77	0	0

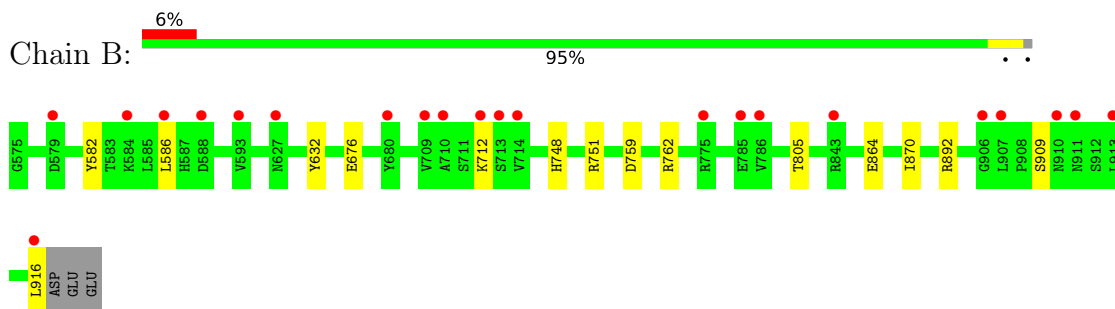
3 Residue-property plots [i](#)

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.

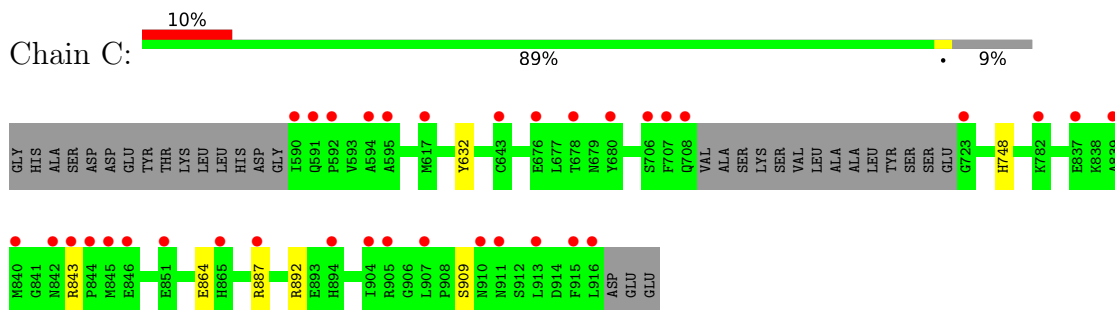
- Molecule 1: cGMP-dependent 3',5'-cyclic phosphodiesterase



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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	165.13Å 72.71Å 90.06Å 90.00° 109.15° 90.00°	Depositor
Resolution (Å)	30.00 – 1.86 30.00 – 1.86	Depositor EDS
% Data completeness (in resolution range)	98.6 (30.00-1.86) 98.7 (30.00-1.86)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 1.85Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.225 , 0.243 0.230 , 0.248	Depositor DCC
R_{free} test set	4114 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å ²)	26.7	Xtriage
Anisotropy	0.524	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 29.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8522	wwPDB-VP
Average B, all atoms (Å ²)	36.0	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 28.37 % 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 1.8661e-03. 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: MG, ZN, 9GA

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	0.59	0/2840	0.84	1/3832 (0.0%)
1	B	0.63	1/2866 (0.0%)	0.84	1/3867 (0.0%)
1	C	0.58	0/2646	0.82	0/3567
All	All	0.60	1/8352 (0.0%)	0.83	2/11266 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	916	LEU	C-O	13.15	1.49	1.23

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	593	VAL	N-CA-CB	5.86	118.51	110.54
1	B	916	LEU	CA-C-O	-5.52	111.42	120.80

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	A	2768	0	2710	9	0
1	B	2793	0	2730	13	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2578	0	2527	3	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	35	0	0	0	0
4	B	35	0	0	1	0
4	C	35	0	0	0	0
5	A	107	0	0	0	0
5	B	87	0	0	0	0
5	C	77	0	0	0	0
All	All	8522	0	7967	21	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 1.

All (21) 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:A:755:GLN:OE1	1:B:712:LYS:NZ	2.10	0.84
1:B:805:THR:CG2	1:B:870:ILE:HD13	2.27	0.64
1:B:805:THR:HG22	1:B:870:ILE:HD13	1.85	0.59
1:A:712:LYS:HE3	1:B:751:ARG:HG2	1.89	0.54
1:B:805:THR:HG23	4:B:1003:9GA:F30	1.99	0.53
1:A:864:GLU:HG3	1:A:892:ARG:HD2	1.93	0.51
1:C:864:GLU:CG	1:C:892:ARG:HD2	2.41	0.51
1:B:805:THR:CG2	1:B:870:ILE:CD1	2.89	0.50
1:B:864:GLU:CG	1:B:892:ARG:HD2	2.42	0.50
1:C:864:GLU:HG3	1:C:892:ARG:HD2	1.93	0.50
1:A:755:GLN:HB2	1:B:712:LYS:HD3	1.94	0.50
1:B:864:GLU:HG3	1:B:892:ARG:HD2	1.93	0.50
1:A:864:GLU:CG	1:A:892:ARG:HD2	2.41	0.49
1:A:751:ARG:HG2	1:B:712:LYS:HD2	1.97	0.47
1:B:632:TYR:O	1:B:748:HIS:NE2	2.41	0.46
1:B:759:ASP:OD1	1:B:762:ARG:NH2	2.49	0.45
1:A:759:ASP:OD1	1:A:762:ARG:NH2	2.50	0.43
1:C:632:TYR:O	1:C:748:HIS:NE2	2.41	0.42
1:B:582:TYR:O	1:B:586:LEU:HB2	2.20	0.41
1:A:859:GLN:HG2	1:A:895:TRP:CE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:682:GLU:OE1	1:A:756:ARG:NH2	2.55	0.40

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)
1:B:676:GLU:OE2	1:C:843:ARG:NH1[3_555]	2.12	0.08

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	337/345 (98%)	335 (99%)	2 (1%)	0	100	100
1	B	341/345 (99%)	340 (100%)	1 (0%)	0	100	100
1	C	310/345 (90%)	308 (99%)	2 (1%)	0	100	100
All	All	988/1035 (96%)	983 (100%)	5 (0%)	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	A	306/310 (99%)	305 (100%)	1 (0%)	86	85

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	308/310 (99%)	307 (100%)	1 (0%)	86	85
1	C	285/310 (92%)	283 (99%)	2 (1%)	76	70
All	All	899/930 (97%)	895 (100%)	4 (0%)	84	82

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	909	SER
1	B	909	SER
1	C	887	ARG
1	C	909	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	591	GLN
1	A	708	GLN
1	A	741	HIS
1	A	781	GLN
1	A	794	GLN
1	B	576	HIS
1	B	591	GLN
1	B	627	ASN
1	B	708	GLN
1	B	794	GLN
1	C	591	GLN
1	C	708	GLN
1	C	781	GLN
1	C	794	GLN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 7 are monoatomic - leaving 3 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
4	9GA	C	1003	-	37,38,38	1.20	4 (10%)	46,58,58	2.69	16 (34%)
4	9GA	B	1003	-	37,38,38	1.24	1 (2%)	46,58,58	2.85	16 (34%)
4	9GA	A	1004	-	37,38,38	1.11	2 (5%)	46,58,58	2.69	14 (30%)

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	9GA	C	1003	-	-	0/24/27/27	0/4/4/4
4	9GA	B	1003	-	-	1/24/27/27	0/4/4/4
4	9GA	A	1004	-	-	1/24/27/27	0/4/4/4

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1003	9GA	N5-N6	3.98	1.41	1.37
4	C	1003	9GA	N5-N6	3.01	1.40	1.37
4	A	1004	9GA	N5-N6	2.98	1.40	1.37
4	C	1003	9GA	C7-N5	-2.36	1.39	1.41
4	A	1004	9GA	O26-C27	2.14	1.43	1.31
4	C	1003	9GA	C2-N3	-2.06	1.33	1.37
4	C	1003	9GA	O26-C27	2.02	1.42	1.31

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1003	9GA	C4-N3-C2	8.61	111.02	102.88
4	A	1004	9GA	C4-N3-C2	8.44	110.86	102.88
4	B	1003	9GA	C4-N3-C2	8.37	110.80	102.88
4	C	1003	9GA	C11-N12-N13	6.70	110.36	103.81
4	B	1003	9GA	C11-N12-N13	6.67	110.33	103.81
4	B	1003	9GA	C2-N6-N5	6.60	107.59	102.40
4	A	1004	9GA	C11-N12-N13	6.54	110.20	103.81
4	C	1003	9GA	C2-N6-N5	6.05	107.16	102.40
4	A	1004	9GA	C2-N6-N5	5.92	107.06	102.40
4	C	1003	9GA	N6-C2-N3	-5.52	108.11	115.24
4	B	1003	9GA	N6-C2-N3	-5.47	108.18	115.24
4	A	1004	9GA	N6-C2-N3	-5.36	108.32	115.24
4	B	1003	9GA	C7-N5-N6	5.23	126.63	120.04
4	A	1004	9GA	C7-N5-N6	5.07	126.42	120.04
4	A	1004	9GA	C16-C15-C14	-4.47	116.80	122.85
4	C	1003	9GA	C7-N5-N6	4.41	125.59	120.04
4	B	1003	9GA	C4-N5-N6	-4.33	105.70	109.45
4	B	1003	9GA	C14-N13-N12	4.27	132.62	126.49
4	C	1003	9GA	C14-N13-N12	4.18	132.50	126.49
4	B	1003	9GA	C11-C10-C9	4.06	107.60	104.57
4	A	1004	9GA	C14-N13-N12	4.04	132.30	126.49
4	B	1003	9GA	C16-C15-C14	-3.71	117.83	122.85
4	B	1003	9GA	C10-C11-N12	-3.68	107.89	112.45
4	A	1004	9GA	C4-N5-N6	-3.64	106.30	109.45
4	C	1003	9GA	C4-N5-N6	-3.47	106.45	109.45
4	B	1003	9GA	C22-C20-N19	3.35	117.24	112.71
4	A	1004	9GA	C10-C9-N8	3.34	132.61	123.54
4	C	1003	9GA	C16-C15-C14	-3.31	118.36	122.85
4	C	1003	9GA	C10-C11-N12	-3.24	108.42	112.45
4	C	1003	9GA	C10-C9-N8	3.18	132.19	123.54
4	C	1003	9GA	N5-C4-N3	-3.16	107.21	110.74
4	B	1003	9GA	C10-C9-N8	3.16	132.12	123.54
4	C	1003	9GA	C35-C33-C34	-3.06	106.04	110.52
4	A	1004	9GA	C16-C15-C7	3.05	124.76	118.86
4	A	1004	9GA	N5-C4-N3	-3.01	107.37	110.74
4	A	1004	9GA	C10-C11-N12	-2.98	108.75	112.45
4	B	1003	9GA	C35-C33-C34	-2.97	106.18	110.52
4	C	1003	9GA	C11-C10-C9	2.91	106.74	104.57
4	B	1003	9GA	N5-C4-N3	-2.75	107.66	110.74
4	A	1004	9GA	C11-C10-C9	2.58	106.50	104.57
4	A	1004	9GA	C22-C20-N19	2.42	115.97	112.71
4	C	1003	9GA	C16-C15-C7	2.32	123.35	118.86
4	B	1003	9GA	C32-C22-C23	2.28	121.13	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1003	9GA	C22-C20-N19	2.28	115.78	112.71
4	B	1003	9GA	C16-C15-C7	2.21	123.14	118.86
4	C	1003	9GA	C35-C33-C20	2.04	114.76	111.23

There are no chirality outliers.

All (2) torsion outliers are listed below:

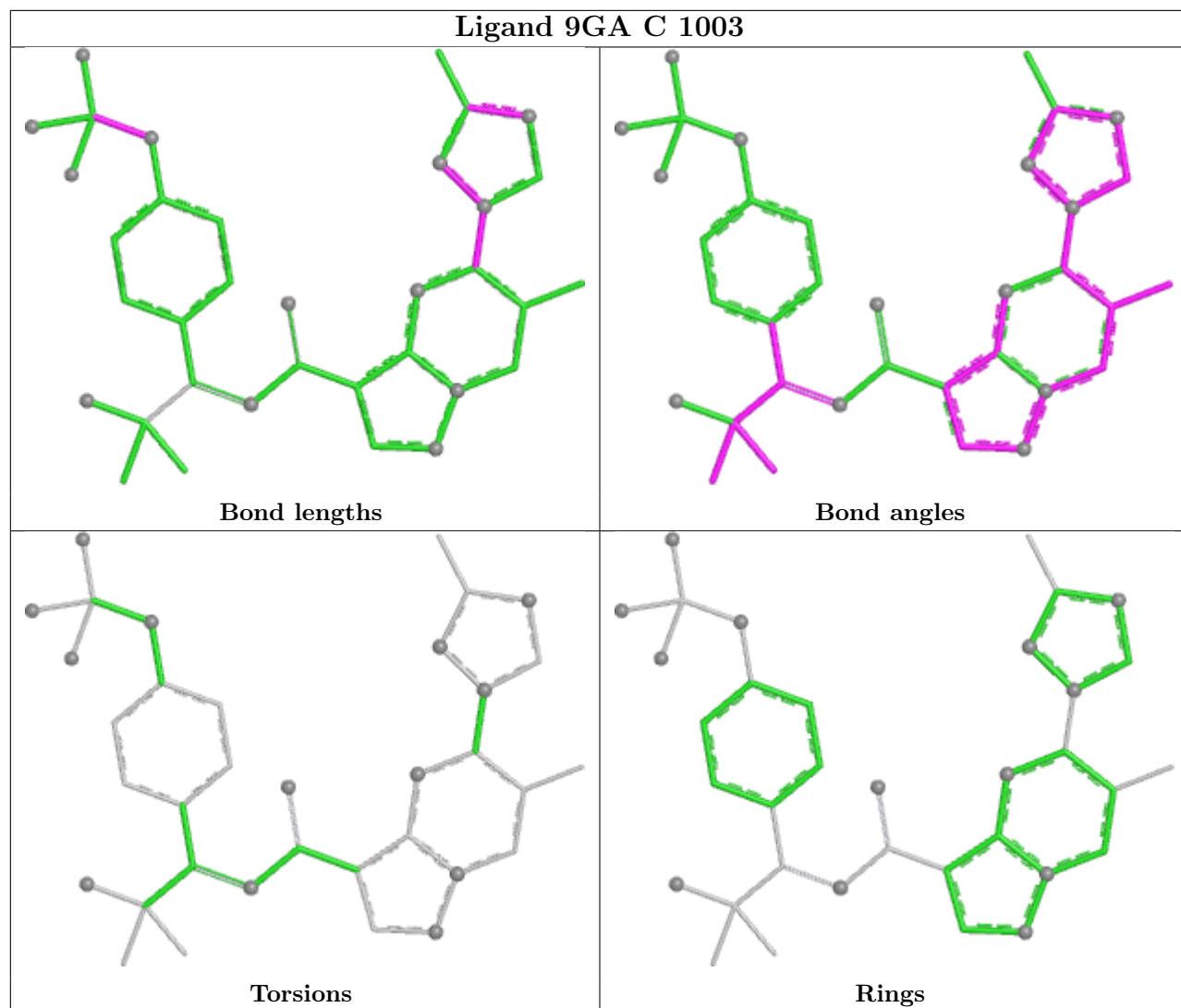
Mol	Chain	Res	Type	Atoms
4	A	1004	9GA	N8-C7-N5-N6
4	B	1003	9GA	N8-C7-N5-N6

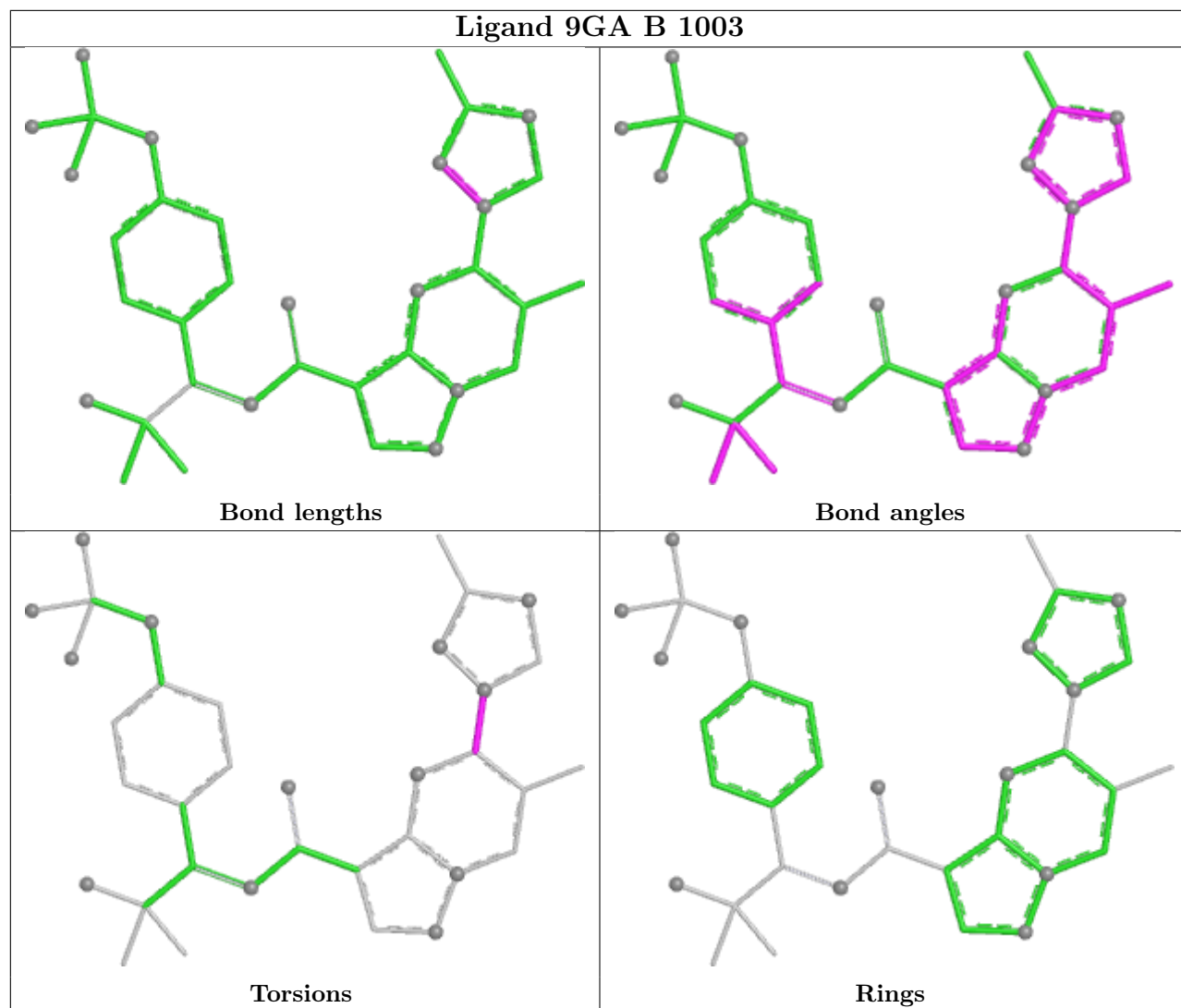
There are no ring outliers.

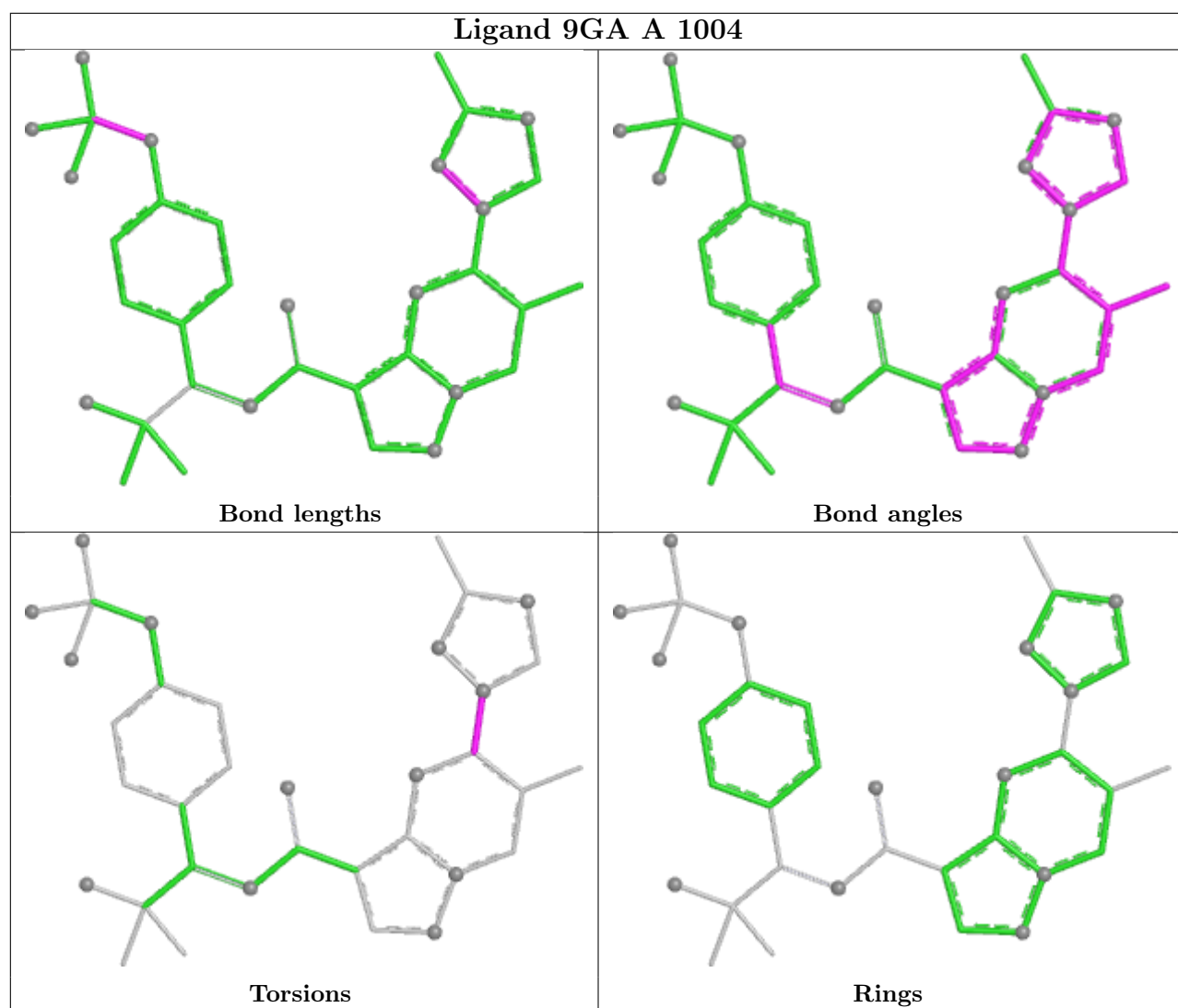
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1003	9GA	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.







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	338/345 (97%)	0.73	43 (12%) 8 7	19, 34, 66, 89	1 (0%)
1	B	342/345 (99%)	0.61	22 (6%) 25 27	20, 35, 52, 70	1 (0%)
1	C	313/345 (90%)	0.74	35 (11%) 10 9	21, 37, 62, 77	1 (0%)
All	All	993/1035 (95%)	0.69	100 (10%) 12 12	19, 35, 61, 89	3 (0%)

All (100) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	589	GLY	6.2
1	A	586	LEU	5.9
1	C	723	GLY	5.6
1	A	590	ILE	4.7
1	A	910	ASN	4.7
1	A	587	HIS	4.6
1	A	592	PRO	4.3
1	C	913	LEU	4.3
1	C	840	MET	4.2
1	A	916	LEU	4.2
1	A	591	GLN	4.1
1	B	916	LEU	4.1
1	B	714	VAL	4.0
1	A	840	MET	4.0
1	C	845	MET	4.0
1	A	907	LEU	3.8
1	C	916	LEU	3.8
1	A	680	TYR	3.7
1	A	748	HIS	3.7
1	A	588	ASP	3.6
1	A	580	ASP	3.5
1	B	584	LYS	3.5
1	C	617	MET	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	907	LEU	3.4
1	B	627	ASN	3.4
1	A	582	TYR	3.3
1	C	904	ILE	3.3
1	A	583	THR	3.3
1	A	845	MET	3.2
1	C	595	ALA	3.2
1	A	913	LEU	3.2
1	A	593	VAL	3.2
1	C	708	GLN	3.2
1	B	843	ARG	3.1
1	C	676	GLU	3.1
1	A	585	LEU	3.1
1	B	712	LYS	3.1
1	A	617	MET	3.1
1	C	707	PHE	2.9
1	C	594	ALA	2.9
1	C	591	GLN	2.9
1	C	592	PRO	2.8
1	C	905	ARG	2.8
1	A	595	ALA	2.8
1	C	839	ALA	2.8
1	B	588	ASP	2.8
1	A	843	ARG	2.8
1	C	643	CYS	2.8
1	A	722	GLU	2.8
1	A	904	ILE	2.8
1	B	775	ARG	2.7
1	C	911	ASN	2.7
1	C	843	ARG	2.7
1	A	648	LYS	2.6
1	B	911	ASN	2.6
1	C	590	ILE	2.6
1	B	709	VAL	2.6
1	A	584	LYS	2.6
1	B	710	ALA	2.6
1	B	910	ASN	2.5
1	C	910	ASN	2.5
1	A	905	ARG	2.5
1	B	785	GLU	2.5
1	A	842	ASN	2.5
1	C	680	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	842	ASN	2.4
1	A	915	PHE	2.4
1	A	598	SER	2.4
1	B	593	VAL	2.4
1	B	680	TYR	2.4
1	C	915	PHE	2.4
1	A	679	ASN	2.4
1	B	907	LEU	2.3
1	C	846	GLU	2.3
1	B	786	VAL	2.3
1	A	911	ASN	2.3
1	C	851	GLU	2.3
1	A	594	ALA	2.3
1	A	861	SER	2.3
1	A	914	ASP	2.3
1	B	906	GLY	2.2
1	A	581	GLU	2.2
1	A	706	SER	2.2
1	A	596	ILE	2.2
1	B	586	LEU	2.2
1	A	906	GLY	2.2
1	C	844	PRO	2.2
1	B	713	SER	2.2
1	C	706	SER	2.1
1	B	913	LEU	2.1
1	B	579	ASP	2.1
1	A	887	ARG	2.1
1	A	751	ARG	2.1
1	C	887	ARG	2.1
1	C	782	LYS	2.0
1	C	678	THR	2.0
1	A	841	GLY	2.0
1	C	865	HIS	2.0
1	C	837	GLU	2.0
1	C	894	HIS	2.0

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

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

6.3 Carbohydrates

There are no oligosaccharides in this entry.

6.4 Ligands

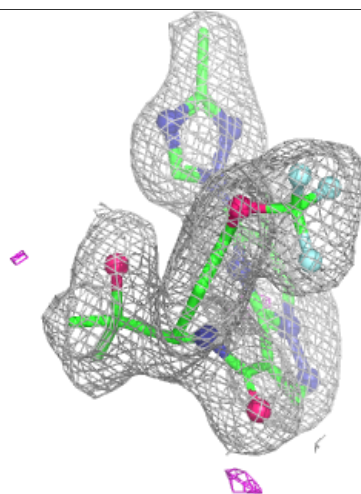
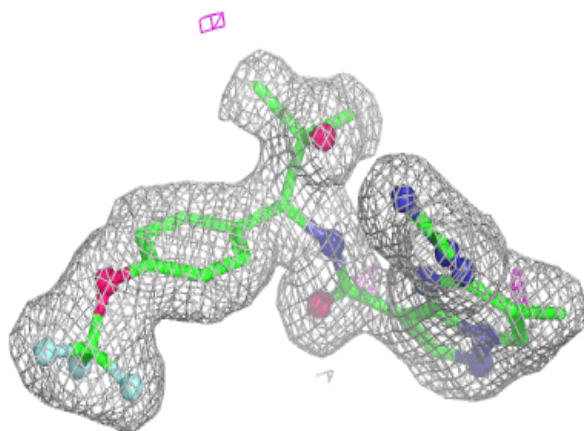
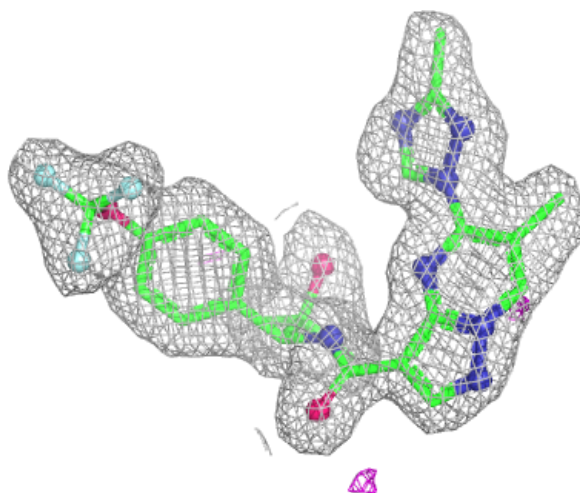
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	MG	A	1003	1/1	0.77	0.13	49,49,49,49	0
4	9GA	A	1004	35/35	0.94	0.07	22,25,31,33	0
4	9GA	B	1003	35/35	0.94	0.07	22,25,30,31	0
4	9GA	C	1003	35/35	0.95	0.07	22,26,31,31	0
3	MG	C	1002	1/1	0.97	0.05	21,21,21,21	0
3	MG	A	1002	1/1	0.98	0.03	22,22,22,22	0
3	MG	B	1002	1/1	0.99	0.02	22,22,22,22	0
2	ZN	B	1001	1/1	1.00	0.01	25,25,25,25	0
2	ZN	C	1001	1/1	1.00	0.01	24,24,24,24	0
2	ZN	A	1001	1/1	1.00	0.01	23,23,23,23	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.

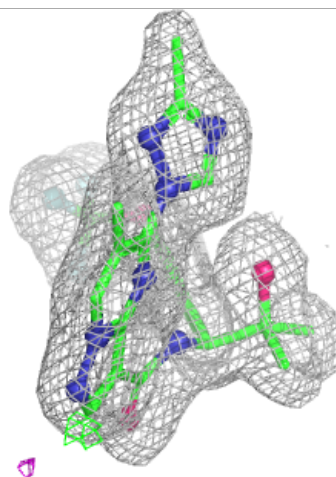
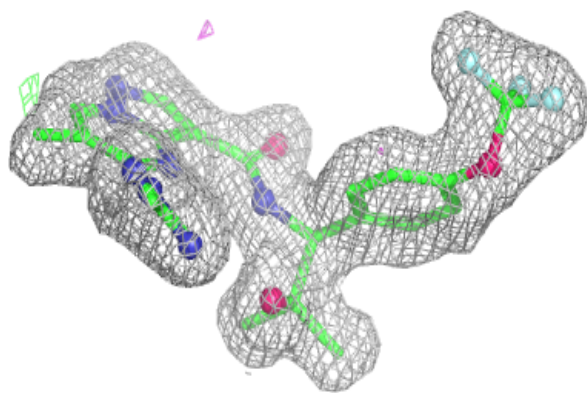
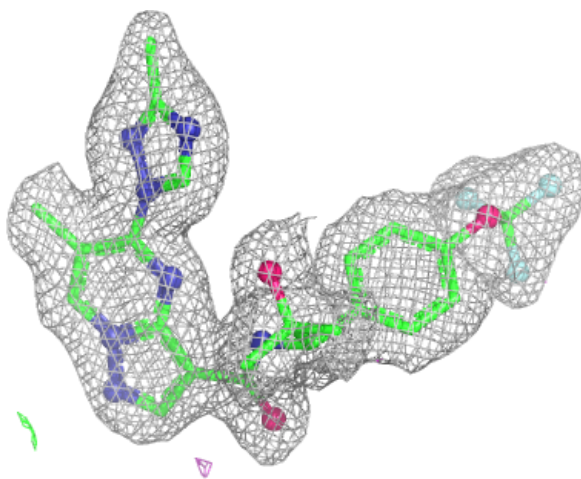
Electron density around 9GA A 1004:

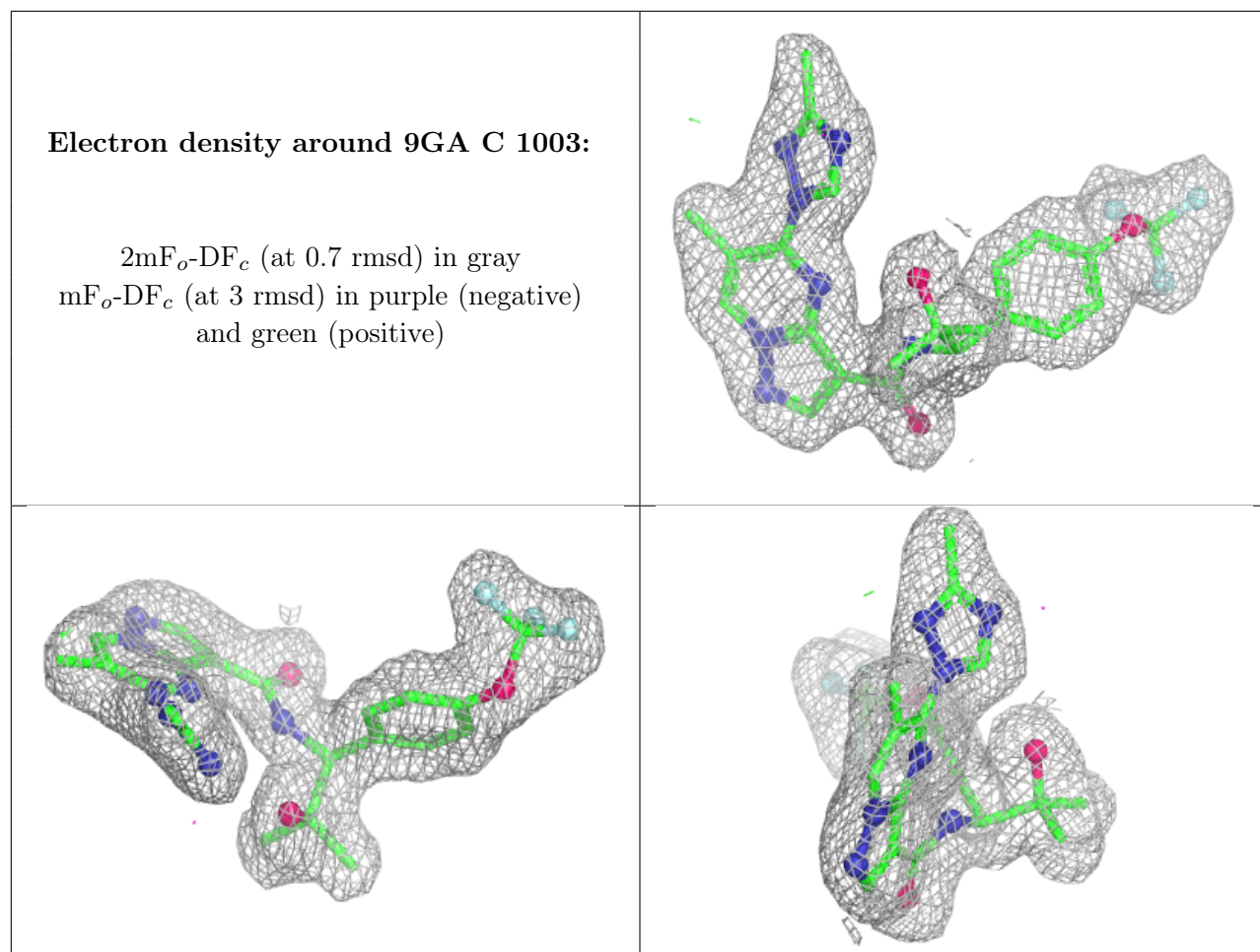
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 9GA B 1003:

$2mF_o-DF_c$ (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.