



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:04 PM UTC

PDB ID : 5TZA / pdb_00005tza
Title : CRYSTAL STRUCTURE OF HUMAN PHOSPHODIESTERASE 2A WITH
3-{5-methyl-[1,2,4]triazolo[1,5-a]pyrimidin-7-yl-1-[(naphthalene-2-yl)carbonyl]
piperidine
Authors : Xu, R.; Aertgeerts, K.
Deposited on : 2016-11-21
Resolution : 1.70 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

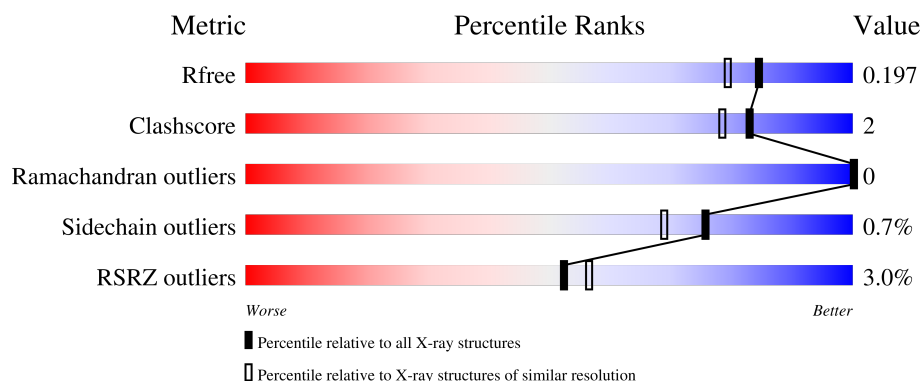
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	344	2% 94% 6%
1	B	344	4% 91% 6%
1	C	344	3% 88% 6%
1	D	344	3% 92% 5%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12265 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.

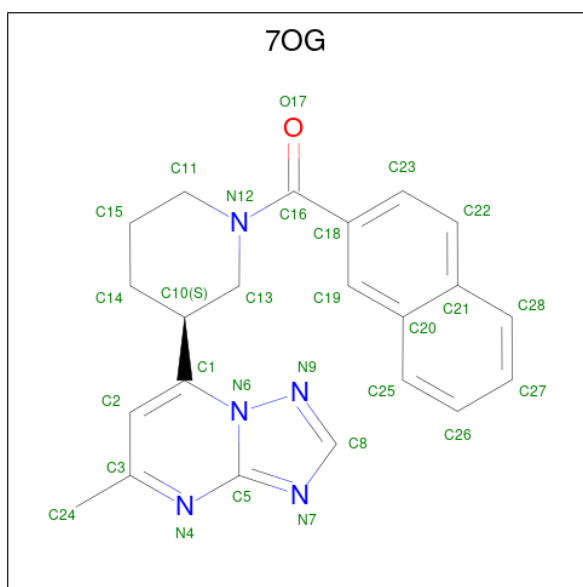
- Molecule 1 is a protein called cGMP-dependent 3',5'-cyclic phosphodiesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	0	0	0
			2786	1774	477	509	26			
1	B	338	Total	C	N	O	S	0	0	0
			2767	1762	474	505	26			
1	C	328	Total	C	N	O	S	0	0	0
			2680	1709	461	485	25			
1	D	333	Total	C	N	O	S	0	0	0
			2730	1744	469	492	25			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	576	SER	-	expression tag	UNP O00408
A	577	ALA	-	expression tag	UNP O00408
A	578	MET	-	expression tag	UNP O00408
B	576	SER	-	expression tag	UNP O00408
B	577	ALA	-	expression tag	UNP O00408
B	578	MET	-	expression tag	UNP O00408
C	576	SER	-	expression tag	UNP O00408
C	577	ALA	-	expression tag	UNP O00408
C	578	MET	-	expression tag	UNP O00408
D	576	SER	-	expression tag	UNP O00408
D	577	ALA	-	expression tag	UNP O00408
D	578	MET	-	expression tag	UNP O00408

- Molecule 2 is [(3S)-3-(5-methyl[1,2,4]triazolo[1,5-a]pyrimidin-7-yl)piperidin-1-yl](naphthalen-2-yl)methanone (CCD ID: 7OG) (formula: C₂₂H₂₁N₅O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			28	22	5	1		
2	B	1	Total	C	N	O	0	0
			28	22	5	1		
2	C	1	Total	C	N	O	0	0
			28	22	5	1		
2	D	1	Total	C	N	O	0	0
			28	22	5	1		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total 1	Mg 1	0	0
4	C	1	Total 1	Mg 1	0	0
4	D	1	Total 1	Mg 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	319	Total 319	O 319	0	0
5	B	298	Total 298	O 298	0	0
5	C	251	Total 251	O 251	0	0
5	D	314	Total 314	O 314	0	0

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



- | CLU | SFR |
|------|------|
| M578 | ALA |
| D579 | D580 |
| V693 | V693 |
| N599 | N599 |
| T604 | T604 |
| Y605 | Y605 |
| N630 | N630 |
| N631 | N631 |
| Y632 | Y632 |
| G636 | G636 |
| S692 | S692 |
| H696 | H696 |
| S721 | S721 |
| E722 | E722 |
| G723 | G723 |
| H748 | H748 |
| K782 | K782 |
| I826 | I826 |
| P856 | P856 |
| Q859 | Q859 |
| M863 | M863 |
| E864 | E864 |
| E883 | E883 |
| R887 | R887 |
| R892 | R892 |
| E893 | E893 |
| H894 | H894 |
| R897 | R897 |
| V898 | V898 |
| F902 | F902 |
| S909 | S909 |
| D914 | D914 |
| F915 | F915 |
| LEU | LEU |
| ASP | ASP |
| GLU | GLU |

-
- | Category | Count |
|----------|-------|
| SER | 1 |
| ALA | 1 |
| MET | 1 |
| ASP | 1 |
| ASP | 1 |
| GLU | 1 |
| TYR | 1 |
| THR | 1 |
| LYS | 1 |
| LEU | 1 |
| HIS | 1 |
| ASP | 1 |
| ASP | 1 |
| P603 | 1 |
| M617 | 1 |
| L624 | 1 |
| M627 | 1 |
| F628 | 1 |
| V632 | 1 |
| L670 | 1 |
| L671 | 1 |
| M679 | 1 |
| V680 | 1 |
| S692 | 1 |
| H696 | 1 |
| E722 | 1 |
| K752 | 1 |
| Q781 | 1 |
| K782 | 1 |
| E785 | 1 |
| V786 | 1 |
| Y787 | 1 |
| Y788 | 1 |
| D789 | 1 |
| R790 | 1 |
| L803 | 1 |
| K821 | 1 |
| E824 | 1 |

- [illegible]

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	55.80Å 73.44Å 91.61Å 109.36° 91.17° 91.14°	Depositor
Resolution (Å)	55.77 – 1.70 55.77 – 1.70	Depositor EDS
% Data completeness (in resolution range)	93.9 (55.77-1.70) 93.9 (55.77-1.70)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 1.38Å)	Xtriage
Refinement program	PHENIX phenix.refine: 1.9_1692	Depositor
R, R_{free}	0.165 , 0.197 0.166 , 0.197	Depositor DCC
R_{free} test set	8675 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	15.8	Xtriage
Anisotropy	0.257	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.014 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	12265	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.11% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, 7OG, ZN

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	0.55	3/2853 (0.1%)	0.72	0/3849
1	B	0.49	2/2834 (0.1%)	0.72	0/3823
1	C	0.50	1/2745 (0.0%)	0.69	0/3703
1	D	0.47	1/2796 (0.0%)	0.69	0/3771
All	All	0.50	7/11228 (0.1%)	0.71	0/15146

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	1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	845	MET	C-O	-5.96	1.16	1.23
1	A	636	CYS	C-O	-5.76	1.19	1.24
1	B	864	GLU	C-O	-5.73	1.17	1.24
1	B	782	LYS	C-O	-5.24	1.18	1.24
1	C	824	GLU	C-O	-5.23	1.18	1.24
1	D	584	LYS	C-O	-5.15	1.18	1.24
1	A	899	SER	C-O	-5.11	1.18	1.24

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	721	SER	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	A	2786	0	2728	10	0
1	B	2767	0	2707	16	0
1	C	2680	0	2633	18	0
1	D	2730	0	2687	10	0
2	A	28	0	0	0	0
2	B	28	0	0	1	0
2	C	28	0	0	0	0
2	D	28	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	319	0	0	0	1
5	B	298	0	0	4	1
5	C	251	0	0	2	0
5	D	314	0	0	2	0
All	All	12265	0	10755	52	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 2.

All (52) 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:D:843:ARG:NH2	5:D:1101:HOH:O	2.05	0.90
1:C:679:ASN:O	1:C:680:TYR:CD1	2.38	0.76
1:C:679:ASN:O	1:C:680:TYR:HD1	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:679:ASN:C	1:C:680:TYR:CD1	2.73	0.67
1:C:790:ARG:CG	1:C:790:ARG:HH11	2.09	0.66
1:C:790:ARG:HH11	1:C:790:ARG:HG3	1.61	0.66
1:B:892:ARG:HH21	1:B:893:GLU:HG2	1.65	0.61
1:B:630:ASN:ND2	5:B:1101:HOH:O	2.34	0.60
1:C:722:GLU:OE1	1:D:722:GLU:OE1	2.20	0.59
1:C:781:GLN:O	1:C:785:GLU:HG3	2.03	0.59
1:B:856:PRO:HG3	1:B:902:PHE:CD1	2.38	0.58
1:A:864:GLU:OE2	1:A:892:ARG:NH2	2.36	0.58
1:B:883:GLU:OE1	1:B:887:ARG:NH1	2.38	0.57
1:C:782:LYS:HE3	5:C:1274:HOH:O	2.08	0.54
1:A:576:SER:HB3	1:A:579:ASP:OD2	2.10	0.52
1:B:580:ASP:HB3	5:B:1336:HOH:O	2.10	0.50
1:A:582:TYR:CE1	1:A:641:ARG:HG3	2.46	0.50
1:A:576:SER:CB	1:A:579:ASP:OD2	2.60	0.50
1:D:688:ALA:HA	1:D:757:MET:HE1	1.96	0.47
1:B:748:HIS:HB3	5:B:1132:HOH:O	2.13	0.47
1:C:603:PHE:HA	1:C:670:LEU:HD13	1.96	0.47
1:A:576:SER:OG	1:A:579:ASP:OD2	2.33	0.47
1:A:578:MET:HE3	1:A:578:MET:HB2	1.81	0.46
1:D:746:PHE:CZ	1:D:757:MET:HE2	2.51	0.46
1:B:721:SER:C	1:B:722:GLU:HG2	2.40	0.46
1:A:599:ASN:O	1:A:605:TYR:HB2	2.16	0.46
1:D:813:THR:O	1:D:887:ARG:HG2	2.16	0.45
1:A:859:GLN:O	1:A:863:MET:HG2	2.17	0.45
1:B:914:ASP:HB2	5:B:1303:HOH:O	2.17	0.45
1:C:897:LYS:HD3	1:C:897:LYS:HA	1.64	0.45
1:C:680:TYR:HB3	1:C:788:TYR:CE2	2.51	0.44
1:B:897:LYS:HB3	1:B:897:LYS:HE3	1.59	0.43
1:C:722:GLU:OE2	1:D:728:ARG:NH1	2.48	0.43
1:A:585:LEU:HG	1:A:620:LEU:HD11	2.01	0.42
1:C:671:LEU:HD13	1:C:803:LEU:HD22	2.02	0.42
1:B:826:ILE:HD13	2:B:1001:7OG:N6	2.34	0.42
1:C:617:MET:HG3	5:C:1293:HOH:O	2.19	0.42
1:B:892:ARG:HH21	1:B:893:GLU:CG	2.32	0.42
1:D:585:LEU:HD12	1:D:585:LEU:HA	1.85	0.42
1:D:664:VAL:HG13	1:D:807:CYS:HB3	2.02	0.41
1:B:599:ASN:O	1:B:605:TYR:HB2	2.20	0.41
1:B:859:GLN:O	1:B:863:MET:HG2	2.20	0.41
1:C:692:SER:O	1:C:696:HIS:HB3	2.20	0.41
1:C:836:LEU:O	1:C:840:MET:HG3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:837:GLU:OE1	1:D:844:PRO:HB3	2.21	0.41
1:A:603:PHE:HA	1:A:670:LEU:HD13	2.02	0.41
1:B:894:HIS:HD2	1:B:898:VAL:HG13	1.86	0.40
1:C:821:LYS:O	1:C:824:GLU:HB2	2.21	0.40
1:C:856:PRO:HG3	1:C:902:PHE:CE1	2.56	0.40
1:B:599:ASN:HD21	1:B:604:THR:HG1	1.68	0.40
1:D:797:ARG:HG2	5:D:1324:HOH:O	2.20	0.40
1:B:692:SER:O	1:B:696:HIS:HB3	2.21	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)
5:A:1251:HOH:O	5:B:1322:HOH:O[1_455]	2.15	0.05

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	A	339/344 (98%)	338 (100%)	1 (0%)	0	100	100
1	B	336/344 (98%)	330 (98%)	6 (2%)	0	100	100
1	C	326/344 (95%)	322 (99%)	4 (1%)	0	100	100
1	D	329/344 (96%)	328 (100%)	1 (0%)	0	100	100
All	All	1330/1376 (97%)	1318 (99%)	12 (1%)	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	A	307/310 (99%)	306 (100%)	1 (0%)	86	83
1	B	305/310 (98%)	303 (99%)	2 (1%)	76	69
1	C	295/310 (95%)	292 (99%)	3 (1%)	68	58
1	D	301/310 (97%)	299 (99%)	2 (1%)	76	69
All	All	1208/1240 (97%)	1200 (99%)	8 (1%)	76	69

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	596	ILE
1	B	893	GLU
1	B	897	LYS
1	C	752	LYS
1	C	790	ARG
1	C	843	ARG
1	D	584	LYS
1	D	897	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	708	GLN
1	A	812	GLN
1	A	875	GLN
1	B	624	GLN
1	B	708	GLN
1	B	741	HIS
1	B	781	GLN
1	B	812	GLN
1	B	865	HIS
1	B	894	HIS
1	C	624	GLN

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Mol	Chain	Res	Type
1	C	627	ASN
1	C	630	ASN
1	C	812	GLN
1	C	875	GLN
1	C	911	ASN
1	D	733	GLN
1	D	812	GLN
1	D	894	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 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	7OG	C	1001	-	31,32,32	1.30	3 (9%)	42,46,46	1.35	5 (11%)
2	7OG	D	1001	-	31,32,32	1.37	3 (9%)	42,46,46	1.44	6 (14%)
2	7OG	B	1001	-	31,32,32	1.28	4 (12%)	42,46,46	1.26	5 (11%)
2	7OG	A	1001	-	31,32,32	1.28	3 (9%)	42,46,46	1.25	4 (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	7OG	C	1001	-	-	0/12/22/22	0/5/5/5
2	7OG	D	1001	-	-	0/12/22/22	0/5/5/5
2	7OG	B	1001	-	-	0/12/22/22	0/5/5/5
2	7OG	A	1001	-	-	0/12/22/22	0/5/5/5

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	7OG	C2-C3	3.04	1.48	1.41
2	D	1001	7OG	C2-C3	3.02	1.47	1.41
2	C	1001	7OG	C2-C3	2.96	1.47	1.41
2	D	1001	7OG	C3-N4	-2.88	1.29	1.33
2	D	1001	7OG	C13-N12	-2.63	1.43	1.46
2	B	1001	7OG	C13-N12	-2.61	1.43	1.46
2	C	1001	7OG	C13-N12	-2.60	1.43	1.46
2	B	1001	7OG	C2-C3	2.45	1.46	1.41
2	A	1001	7OG	C3-N4	-2.39	1.30	1.33
2	C	1001	7OG	C3-N4	-2.33	1.30	1.33
2	A	1001	7OG	C13-N12	-2.24	1.44	1.46
2	B	1001	7OG	C3-N4	-2.18	1.30	1.33
2	B	1001	7OG	N6-N9	-2.07	1.34	1.37

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1001	7OG	C18-C16-N12	3.71	123.30	118.66
2	D	1001	7OG	O17-C16-C18	-3.46	113.43	120.29
2	C	1001	7OG	C24-C3-C2	-3.39	116.01	120.44
2	D	1001	7OG	C18-C16-N12	3.34	122.84	118.66
2	B	1001	7OG	C24-C3-C2	-3.21	116.24	120.44
2	D	1001	7OG	C2-C3-N4	2.80	125.47	122.71
2	D	1001	7OG	C24-C3-C2	-2.73	116.88	120.44
2	A	1001	7OG	O17-C16-C18	-2.65	115.04	120.29
2	C	1001	7OG	C24-C3-N4	2.60	119.74	116.84
2	D	1001	7OG	C11-N12-C13	2.48	117.86	113.07
2	B	1001	7OG	C24-C3-N4	2.43	119.55	116.84
2	D	1001	7OG	C8-N9-N6	2.38	102.80	100.86
2	B	1001	7OG	O17-C16-C18	-2.32	115.69	120.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1001	7OG	O17-C16-C18	-2.30	115.72	120.29
2	C	1001	7OG	C11-N12-C13	2.29	117.49	113.07
2	A	1001	7OG	C22-C23-C18	-2.29	118.07	120.78
2	A	1001	7OG	C23-C18-C19	2.19	121.86	119.21
2	B	1001	7OG	C11-N12-C13	2.14	117.20	113.07
2	A	1001	7OG	C24-C3-C2	-2.06	117.75	120.44
2	B	1001	7OG	C14-C10-C13	2.03	113.58	110.02

There are no chirality outliers.

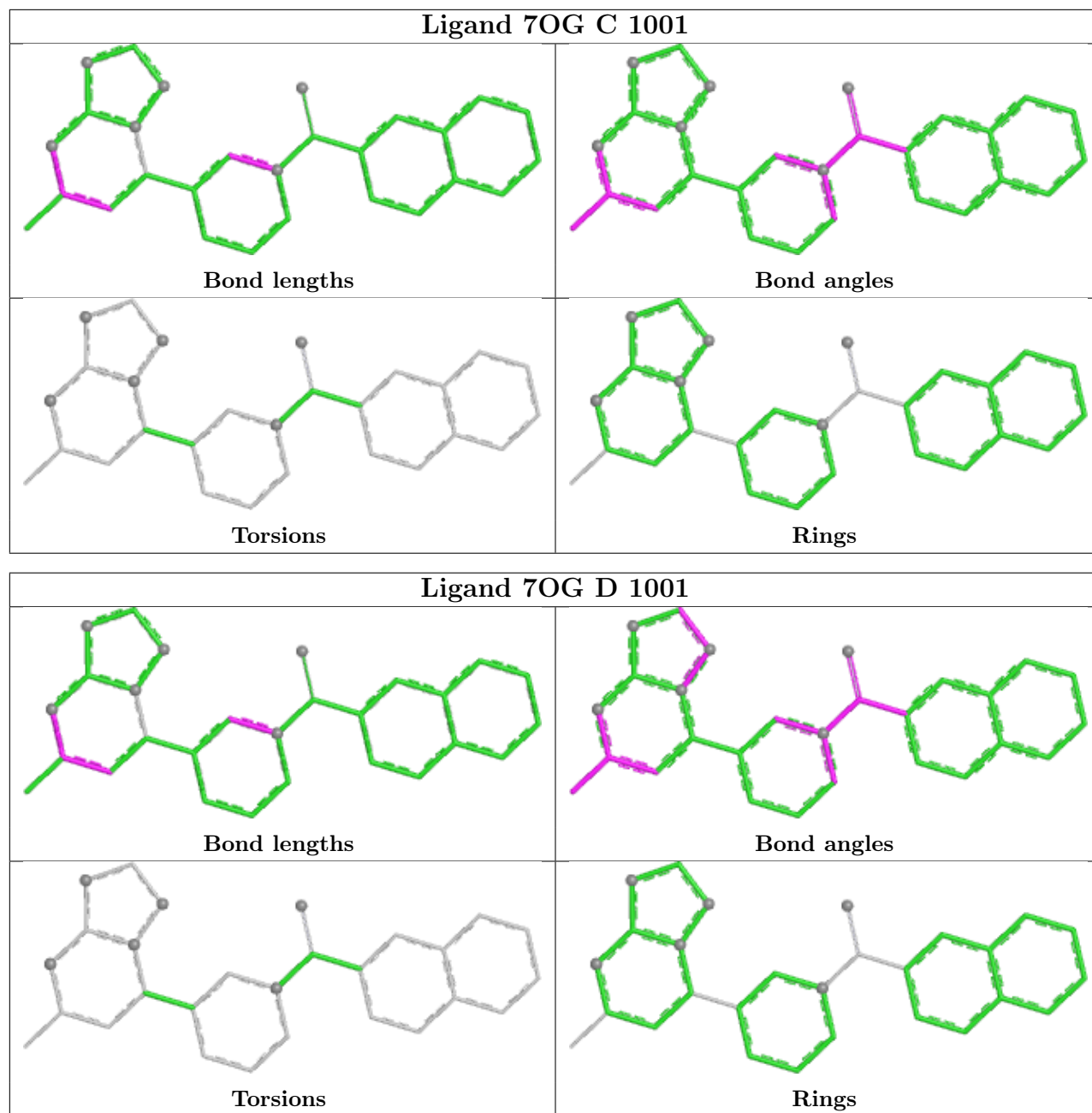
There are no torsion outliers.

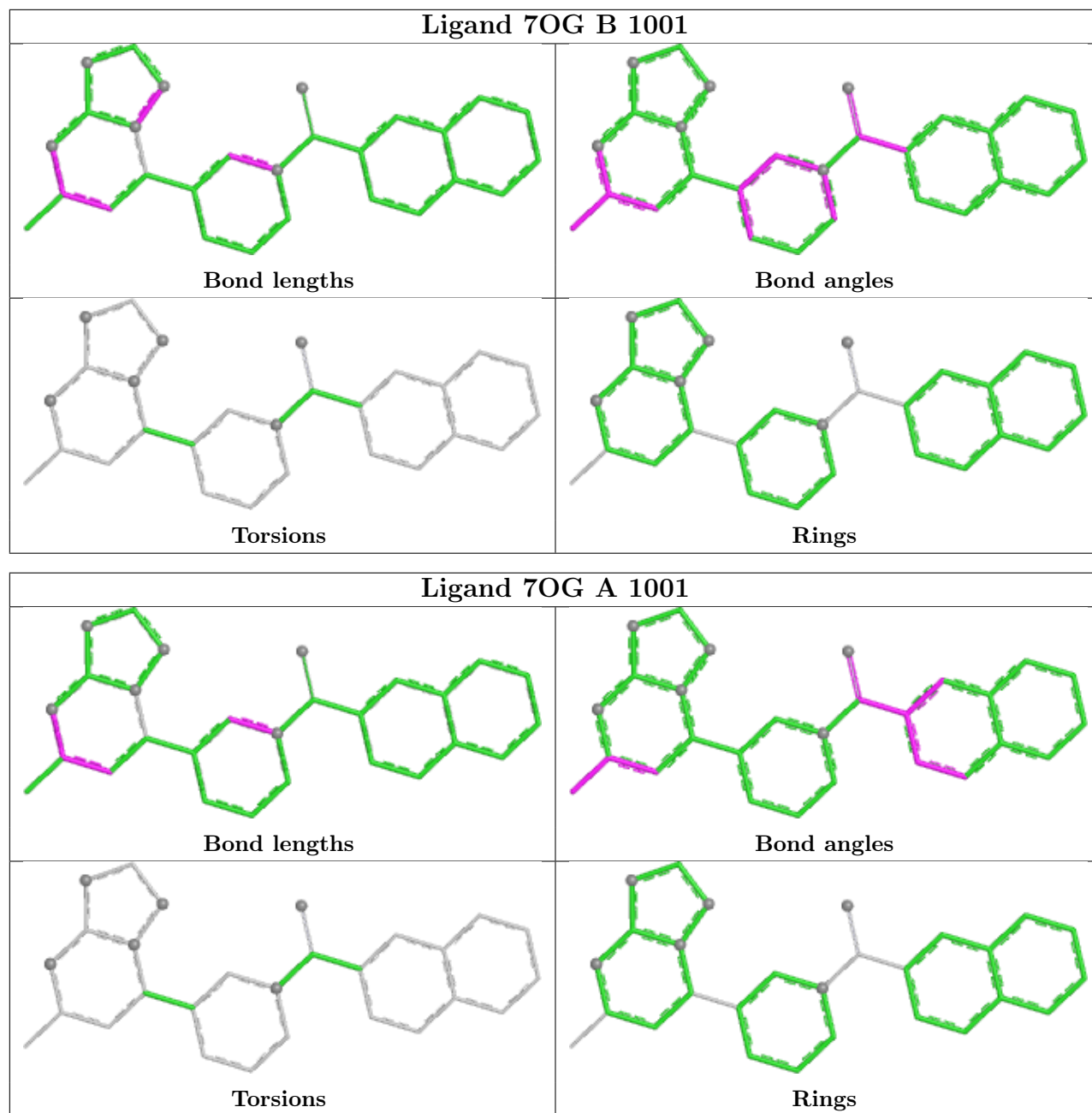
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1001	7OG	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	341/344 (99%)	-0.15	6 (1%) 67 72	8, 18, 36, 54	0
1	B	338/344 (98%)	0.06	13 (3%) 44 48	10, 22, 38, 56	0
1	C	328/344 (95%)	0.18	12 (3%) 45 49	10, 25, 45, 56	0
1	D	333/344 (96%)	-0.02	9 (2%) 56 60	9, 21, 39, 53	0
All	All	1340/1376 (97%)	0.02	40 (2%) 52 56	8, 21, 41, 56	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	680	TYR	4.7
1	B	723	GLY	3.7
1	B	909	SER	3.4
1	D	894	HIS	3.3
1	C	593	VAL	3.3
1	C	916	LEU	3.2
1	D	909	SER	3.2
1	C	624	GLN	2.9
1	B	593	VAL	2.8
1	B	914	ASP	2.8
1	B	579	ASP	2.8
1	C	786	VAL	2.8
1	B	894	HIS	2.7
1	B	636	CYS	2.7
1	C	909	SER	2.7
1	C	632	TYR	2.7
1	D	613	ASP	2.7
1	B	915	PHE	2.6
1	A	577	ALA	2.6
1	C	595	ALA	2.6
1	D	915	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	748	HIS	2.5
1	C	627	ASN	2.4
1	D	914	ASP	2.4
1	A	785	GLU	2.4
1	B	722	GLU	2.4
1	A	636	CYS	2.3
1	D	586	LEU	2.3
1	B	580	ASP	2.2
1	D	587	HIS	2.2
1	C	790	ARG	2.2
1	A	576	SER	2.2
1	D	585	LEU	2.2
1	C	785	GLU	2.2
1	B	632	TYR	2.2
1	A	578	MET	2.2
1	A	579	ASP	2.1
1	C	628	PHE	2.1
1	B	893	GLU	2.0
1	D	632	TYR	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.

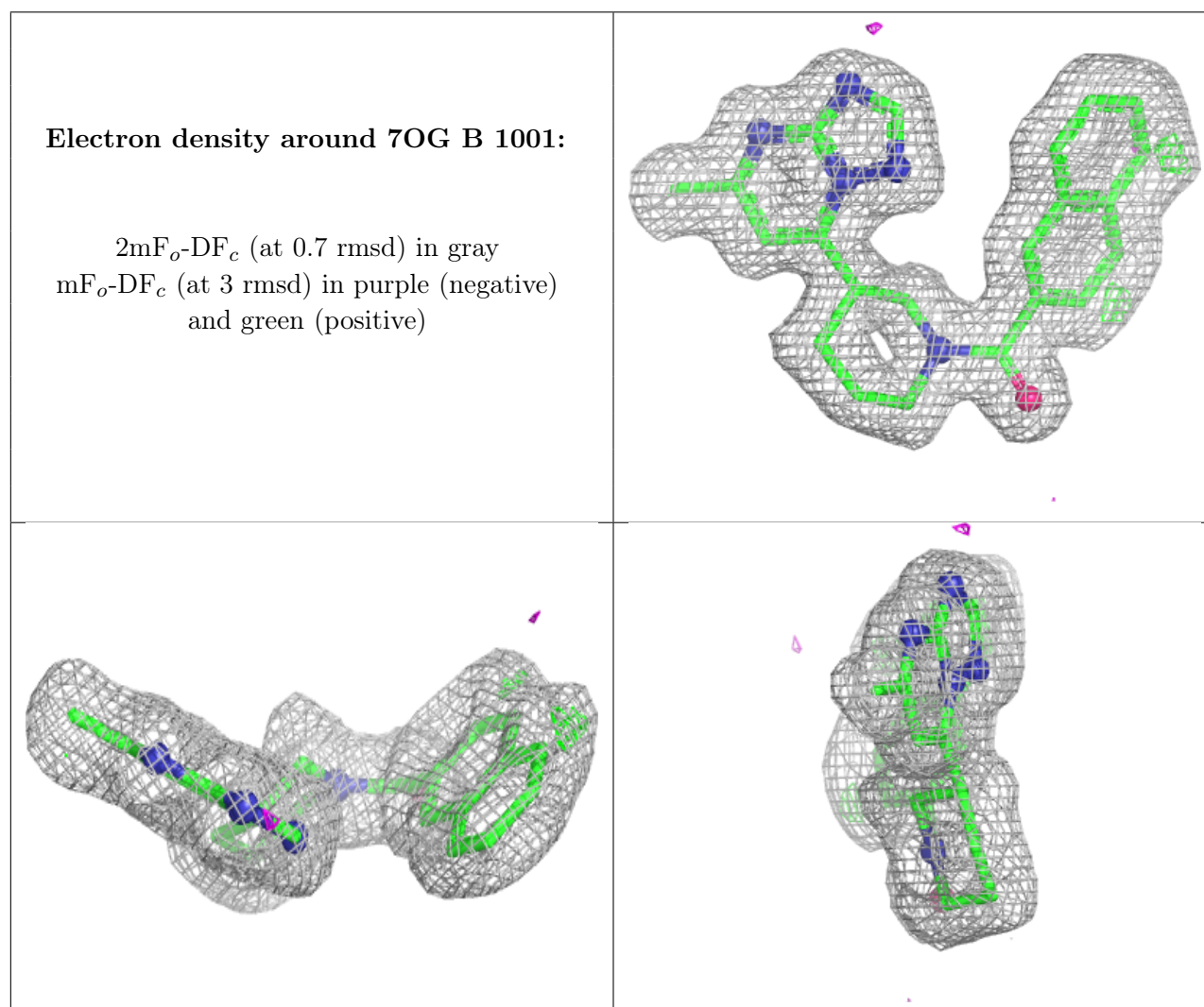
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	7OG	B	1001	28/28	0.94	0.07	12,16,19,20	0
2	7OG	C	1001	28/28	0.96	0.06	9,15,21,22	0
2	7OG	A	1001	28/28	0.97	0.05	8,14,19,20	0
2	7OG	D	1001	28/28	0.97	0.05	11,13,17,21	0

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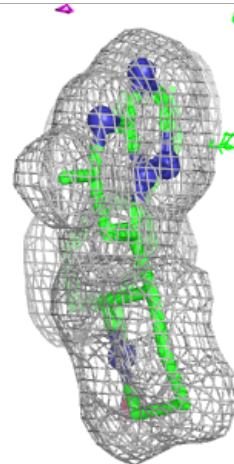
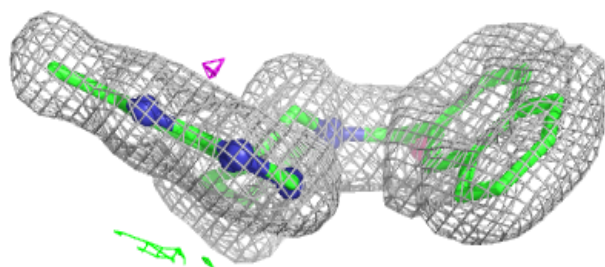
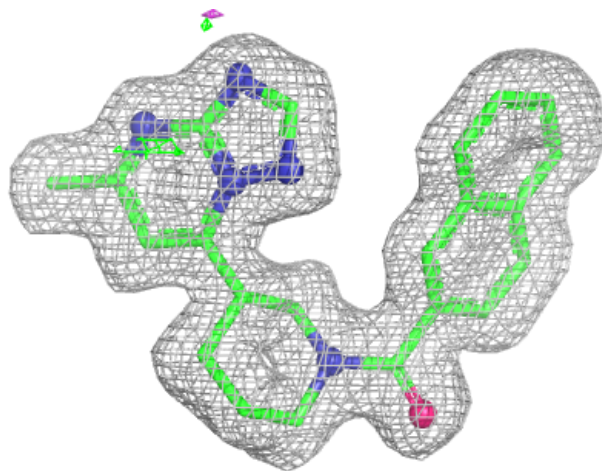
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MG	A	1003	1/1	0.99	0.02	8,8,8,8	0
4	MG	B	1003	1/1	0.99	0.03	10,10,10,10	0
4	MG	C	1003	1/1	0.99	0.03	11,11,11,11	0
4	MG	D	1003	1/1	0.99	0.03	9,9,9,9	0
3	ZN	A	1002	1/1	1.00	0.02	11,11,11,11	0
3	ZN	B	1002	1/1	1.00	0.01	12,12,12,12	0
3	ZN	C	1002	1/1	1.00	0.01	13,13,13,13	0
3	ZN	D	1002	1/1	1.00	0.01	12,12,12,12	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.



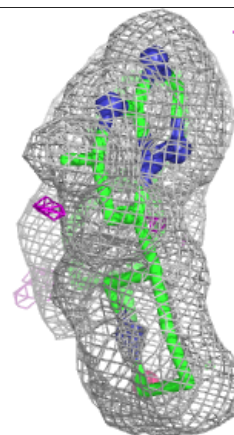
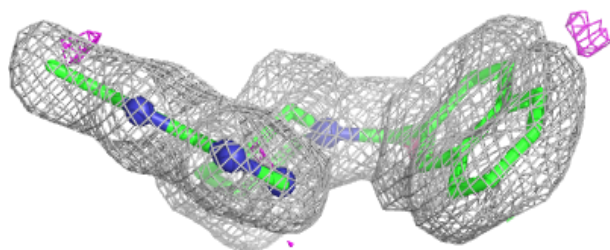
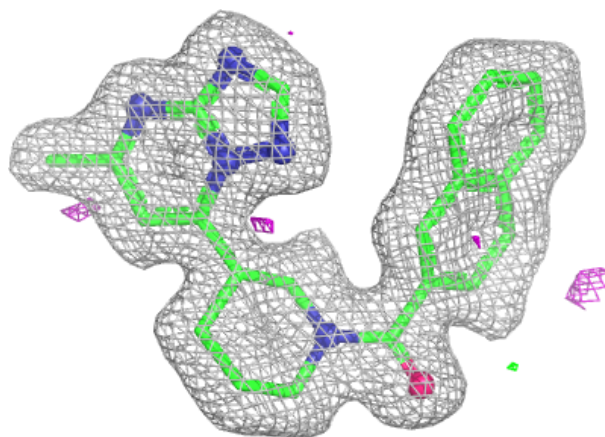
Electron density around 7OG C 1001:

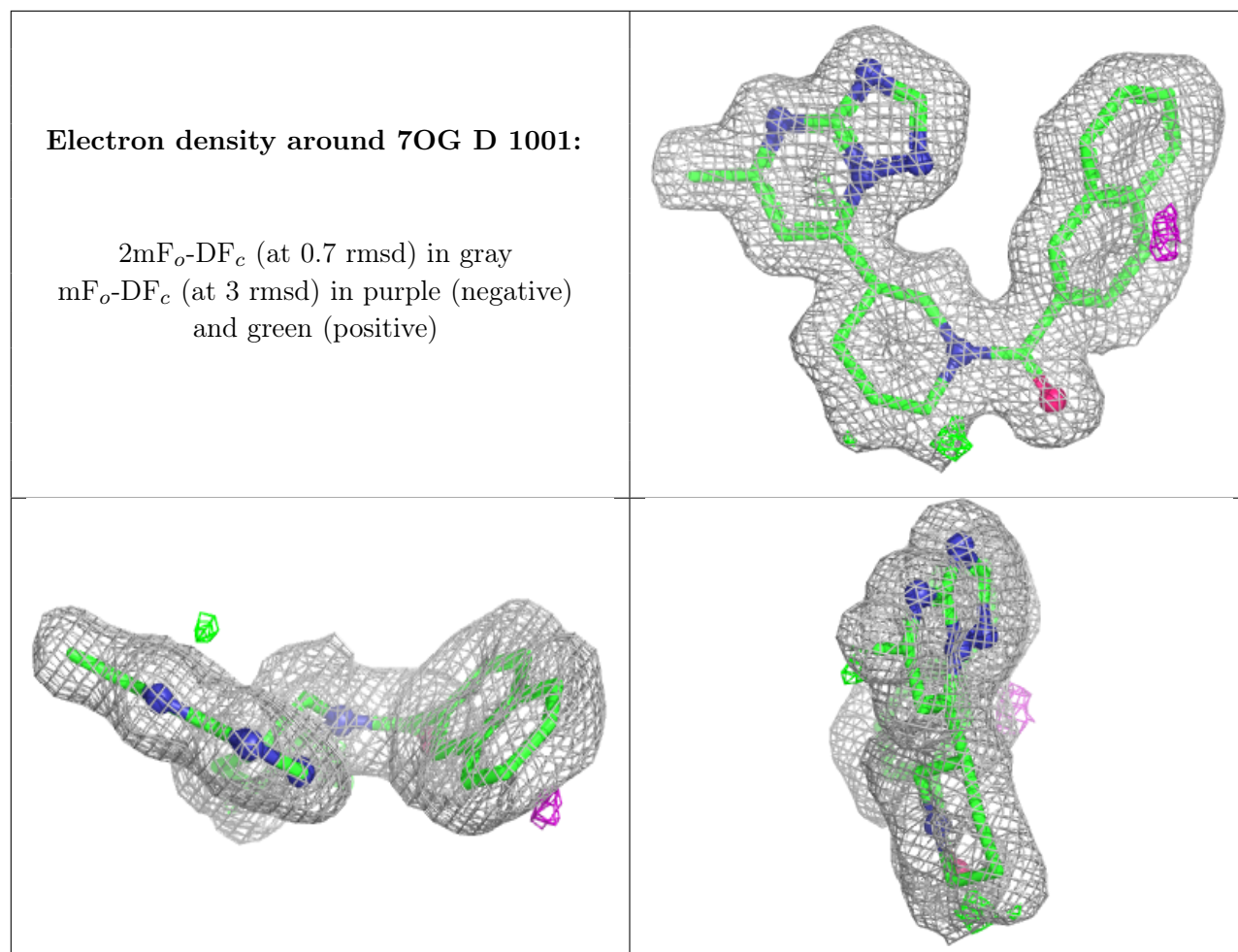
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 7OG A 1001:

$2mF_o-DF_c$ (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.