



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 11:39 am BST

PDB ID : 5TZ3
Title : CRYSTAL STRUCTURE OF HUMAN PHOSPHODIESTERASE 2A IN COMPLEX with [1,2,4]triazolo[1,5-a]pyrimidin-7-yl}-N-(naphthalene-2-yl)piperidine-3-carboxamide
Authors : Xu, R.; Aertgeerts, K.
Deposited on : 2016-11-21
Resolution : 1.72 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

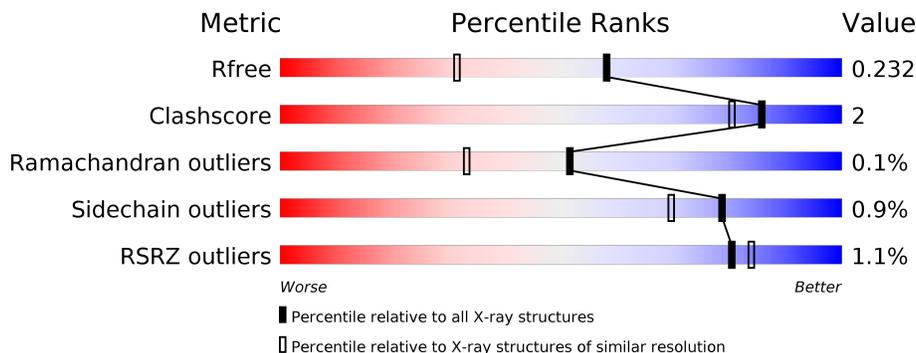
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

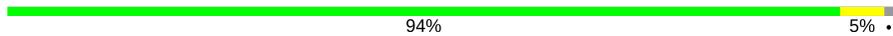
The reported resolution of this entry is 1.72 Å.

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	5722 (1.74-1.70)
Clashscore	141614	6152 (1.74-1.70)
Ramachandran outliers	138981	6051 (1.74-1.70)
Sidechain outliers	138945	6051 (1.74-1.70)
RSRZ outliers	127900	5629 (1.74-1.70)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	344	 94% 5% •
1	B	344	 93% 6% •
1	C	344	 89% 7% • •
1	D	344	 90% 5% 5%

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 11891 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.

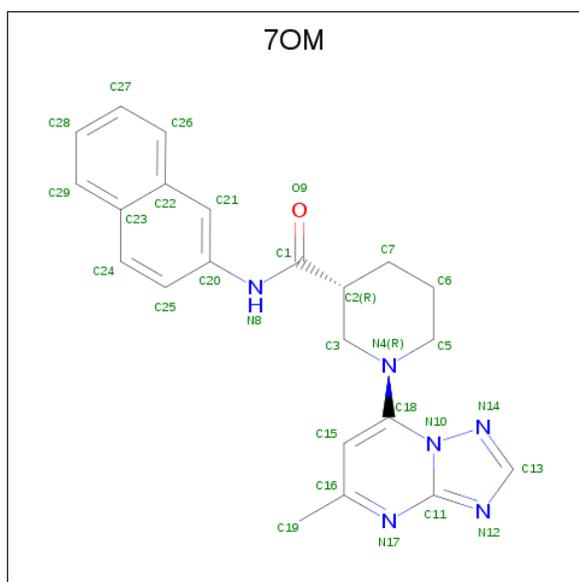
- Molecule 1 is a protein called cGMP-dependent 3',5'-cyclic phosphodiesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	341	2786	1774	477	509	26	0	0	0
1	B	340	2780	1771	476	507	26	0	0	0
1	C	334	2733	1743	468	497	25	0	0	0
1	D	327	2676	1707	460	484	25	0	0	0

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 (3 {R})-1-(5-methyl-[1,2,4]triazolo[1,5-a]pyrimidin-7-yl)- {N}-naphthalen-2-yl-piperidine-3-carboxamide (three-letter code: 7OM) (formula: C₂₂H₂₂N₆O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			29	22	6	1		
2	B	1	Total	C	N	O	0	0
			29	22	6	1		
2	C	1	Total	C	N	O	0	0
			29	22	6	1		
2	D	1	Total	C	N	O	0	0
			29	22	6	1		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	213	Total O 213 213	0	0
5	B	215	Total O 215 215	0	0
5	C	160	Total O 160 160	0	0
5	D	204	Total O 204 204	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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

Chain A: 



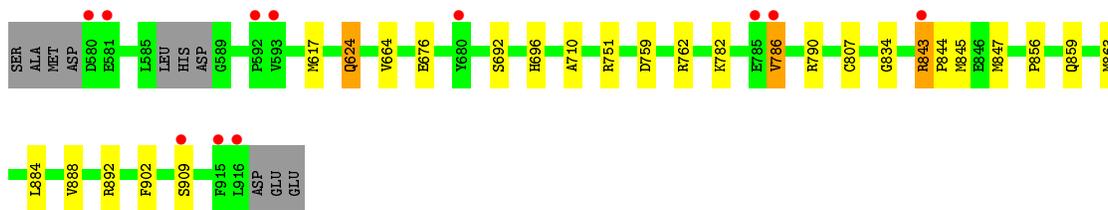
- Molecule 1: cGMP-dependent 3',5'-cyclic phosphodiesterase

Chain B: 



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

Chain C: 



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

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	55.75Å 72.73Å 90.52Å 109.21° 90.70° 91.19°	Depositor
Resolution (Å)	68.65 – 1.72 68.65 – 1.64	Depositor EDS
% Data completeness (in resolution range)	95.0 (68.65-1.72) 89.7 (68.65-1.64)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 1.64Å)	Xtriage
Refinement program	PHENIX phenix.refine: 1.9_1692	Depositor
R, R_{free}	0.191 , 0.233 0.192 , 0.232	Depositor DCC
R_{free} test set	7091 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å ²)	17.7	Xtriage
Anisotropy	0.444	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 50.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.016 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11891	wwPDB-VP
Average B, all atoms (Å ²)	29.0	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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.39	0/2853	0.51	0/3849
1	B	0.35	0/2847	0.49	0/3841
1	C	0.36	0/2798	0.47	0/3773
1	D	0.37	0/2741	0.50	0/3698
All	All	0.37	0/11239	0.49	0/15161

There are no bond length outliers.

There are no bond angle outliers.

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	2786	0	2728	11	0
1	B	2780	0	2723	12	0
1	C	2733	0	2682	12	1
1	D	2676	0	2630	8	0
2	A	29	0	0	1	0
2	B	29	0	0	2	0
2	C	29	0	0	2	0
2	D	29	0	0	2	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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	213	0	0	3	0
5	B	215	0	0	3	0
5	C	160	0	0	1	0
5	D	204	0	0	0	0
All	All	11891	0	10763	49	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 (49) 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:C:759:ASP:OD1	1:C:762:ARG:NH2	2.22	0.72
2:D:1001:7OM:O9	2:D:1001:7OM:C25	2.43	0.65
1:D:654:PRO:HD2	1:D:829:GLU:HG3	1.82	0.62
2:A:1001:7OM:O9	2:A:1001:7OM:C25	2.47	0.61
1:D:593:VAL:HA	1:D:596:ILE:HD12	1.82	0.61
1:D:676:GLU:OE2	1:D:880:LYS:NZ	2.30	0.60
1:A:868:MET:SD	1:A:892:ARG:HD2	2.42	0.59
2:C:1001:7OM:C25	2:C:1001:7OM:O9	2.55	0.55
1:A:582:TYR:CE1	1:A:641:ARG:HG3	2.44	0.53
1:A:778:LYS:HG2	5:A:1119:HOH:O	2.09	0.53
1:C:884:LEU:O	1:C:888:VAL:HG23	2.09	0.52
1:A:712:LYS:NZ	5:A:1102:HOH:O	2.37	0.52
1:C:751:ARG:HG3	5:C:1225:HOH:O	2.09	0.52
2:D:1001:7OM:N14	2:D:1001:7OM:C3	2.73	0.51
1:B:856:PRO:HG3	1:B:902:PHE:CD1	2.45	0.51
1:B:887:ARG:NH1	5:B:1104:HOH:O	2.43	0.50
1:D:856:PRO:HG3	1:D:902:PHE:CD1	2.46	0.49
1:B:724:SER:N	5:B:1101:HOH:O	2.45	0.49
2:B:1001:7OM:C25	2:B:1001:7OM:O9	2.58	0.48
1:B:581:GLU:HA	1:B:584:LYS:HE3	1.95	0.48
1:B:591:GLN:HG2	1:B:595:ALA:HB3	1.95	0.47
1:B:821:LYS:HA	1:B:824:GLU:HG3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:843:ARG:HE	1:A:843:ARG:HB3	1.44	0.47
2:C:1001:7OM:C3	2:C:1001:7OM:N14	2.80	0.45
1:C:624:GLN:HE21	1:C:624:GLN:HA	1.81	0.45
1:A:636:CYS:HB2	1:A:637:PRO:HD3	1.99	0.44
1:C:664:VAL:HG13	1:C:807:CYS:HB3	2.00	0.44
1:A:671:LEU:HD13	1:A:803:LEU:HD22	2.00	0.44
1:B:581:GLU:OE2	1:B:647:LYS:NZ	2.51	0.44
1:A:797:ARG:HG2	5:A:1289:HOH:O	2.16	0.43
1:B:811:ASP:HB3	1:B:822:ILE:HG13	2.01	0.43
1:D:679:ASN:HD21	1:D:790:ARG:HH21	1.65	0.43
1:D:727:GLU:HA	1:D:730:HIS:CD2	2.54	0.43
1:C:692:SER:O	1:C:696:HIS:HB3	2.19	0.42
1:C:710:ALA:O	1:D:751:ARG:NH2	2.52	0.42
1:B:748:HIS:HB3	5:B:1124:HOH:O	2.19	0.42
1:B:664:VAL:HG13	1:B:807:CYS:HB3	2.01	0.42
1:C:834:GLY:HA2	1:C:844:PRO:HB3	2.01	0.42
1:C:856:PRO:HG3	1:C:902:PHE:CD1	2.54	0.42
1:C:845:MET:HG3	1:C:847:MET:HG2	2.02	0.42
1:C:859:GLN:O	1:C:863:MET:HG2	2.21	0.41
1:C:782:LYS:O	1:C:786:VAL:HG13	2.21	0.41
1:B:722:GLU:HB3	1:B:723:GLY:H	1.63	0.41
1:A:648:LYS:HD3	1:A:648:LYS:N	2.36	0.41
1:D:692:SER:O	1:D:696:HIS:HB3	2.21	0.41
2:B:1001:7OM:N14	2:B:1001:7OM:C3	2.85	0.40
1:A:727:GLU:HA	1:A:730:HIS:CD2	2.56	0.40
1:B:818:THR:O	1:B:822:ILE:HG12	2.22	0.40
1:A:859:GLN:O	1:A:863:MET:HG2	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)
1:C:676:GLU:OE2	1:C:843:ARG:NH2[1_655]	2.17	0.03

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	339/344 (98%)	337 (99%)	2 (1%)	0	100	100
1	B	338/344 (98%)	334 (99%)	3 (1%)	1 (0%)	41	24
1	C	330/344 (96%)	328 (99%)	2 (1%)	0	100	100
1	D	325/344 (94%)	323 (99%)	2 (1%)	0	100	100
All	All	1332/1376 (97%)	1322 (99%)	9 (1%)	1 (0%)	51	33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	722	GLU

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	307/310 (99%)	305 (99%)	2 (1%)	84	76
1	B	306/310 (99%)	305 (100%)	1 (0%)	92	89
1	C	301/310 (97%)	294 (98%)	7 (2%)	50	31
1	D	295/310 (95%)	294 (100%)	1 (0%)	92	89
All	All	1209/1240 (98%)	1198 (99%)	11 (1%)	78	69

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	793	LYS
1	A	892	ARG
1	B	817	LYS
1	C	617	MET
1	C	624	GLN

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Mol	Chain	Res	Type
1	C	786	VAL
1	C	790	ARG
1	C	843	ARG
1	C	892	ARG
1	C	909	SER
1	D	756	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	842	ASN
1	C	624	GLN
1	C	812	GLN
1	C	894	HIS
1	D	679	ASN
1	D	708	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 carbohydrates 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	7OM	A	1001	-	28,33,33	1.44	3 (10%)	34,47,47	1.37	5 (14%)
2	7OM	B	1001	-	28,33,33	1.42	3 (10%)	34,47,47	1.56	5 (14%)
2	7OM	C	1001	-	28,33,33	1.43	4 (14%)	34,47,47	1.43	6 (17%)
2	7OM	D	1001	-	28,33,33	1.47	4 (14%)	34,47,47	1.62	7 (20%)

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	7OM	A	1001	-	-	4/12/22/22	0/5/5/5
2	7OM	B	1001	-	-	4/12/22/22	0/5/5/5
2	7OM	C	1001	-	-	3/12/22/22	0/5/5/5
2	7OM	D	1001	-	-	4/12/22/22	0/5/5/5

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1001	7OM	C15-C16	4.58	1.47	1.38
2	D	1001	7OM	C15-C16	4.54	1.47	1.38
2	A	1001	7OM	C15-C16	4.28	1.47	1.38
2	B	1001	7OM	C15-C16	3.96	1.46	1.38
2	A	1001	7OM	C15-C18	-2.96	1.33	1.38
2	B	1001	7OM	C15-C18	-2.86	1.33	1.38
2	D	1001	7OM	C11-N12	-2.84	1.31	1.35
2	D	1001	7OM	C15-C18	-2.54	1.34	1.38
2	A	1001	7OM	C11-N12	-2.54	1.32	1.35
2	B	1001	7OM	C11-N12	-2.51	1.32	1.35
2	C	1001	7OM	C15-C18	-2.42	1.34	1.38
2	C	1001	7OM	C11-N12	-2.39	1.32	1.35
2	C	1001	7OM	C16-N17	-2.07	1.29	1.33
2	D	1001	7OM	C16-N17	-2.04	1.29	1.33

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1001	7OM	C20-N8-C1	-4.87	115.73	127.40
2	A	1001	7OM	C20-N8-C1	-4.34	117.00	127.40
2	B	1001	7OM	C2-C1-N8	-4.32	108.67	115.02
2	B	1001	7OM	C19-C16-C15	-4.11	115.54	121.81
2	C	1001	7OM	C20-N8-C1	-3.94	117.97	127.40
2	C	1001	7OM	C7-C2-C3	3.81	116.75	109.92
2	D	1001	7OM	C18-C15-C16	-3.29	114.92	116.91
2	A	1001	7OM	C19-C16-C15	-3.28	116.81	121.81
2	D	1001	7OM	C19-C16-C15	-3.11	117.06	121.81
2	C	1001	7OM	C19-C16-C15	-3.03	117.18	121.81
2	B	1001	7OM	C20-N8-C1	-2.96	120.31	127.40
2	A	1001	7OM	C7-C2-C3	2.87	115.07	109.92
2	C	1001	7OM	C2-C3-N4	2.78	117.00	109.73
2	D	1001	7OM	C15-C16-N17	2.74	125.81	122.57
2	D	1001	7OM	C7-C2-C3	2.71	114.77	109.92
2	B	1001	7OM	C7-C2-C3	2.68	114.73	109.92
2	D	1001	7OM	C2-C3-N4	2.43	116.09	109.73
2	D	1001	7OM	C5-N4-C3	2.43	119.61	112.55
2	B	1001	7OM	O9-C1-C2	2.20	124.98	122.12
2	A	1001	7OM	C2-C3-N4	2.15	115.37	109.73
2	C	1001	7OM	C7-C2-C1	-2.07	106.14	110.69
2	A	1001	7OM	C2-C1-N8	-2.06	112.00	115.02
2	C	1001	7OM	C5-N4-C3	2.04	118.48	112.55

There are no chirality outliers.

All (15) torsion outliers are listed below:

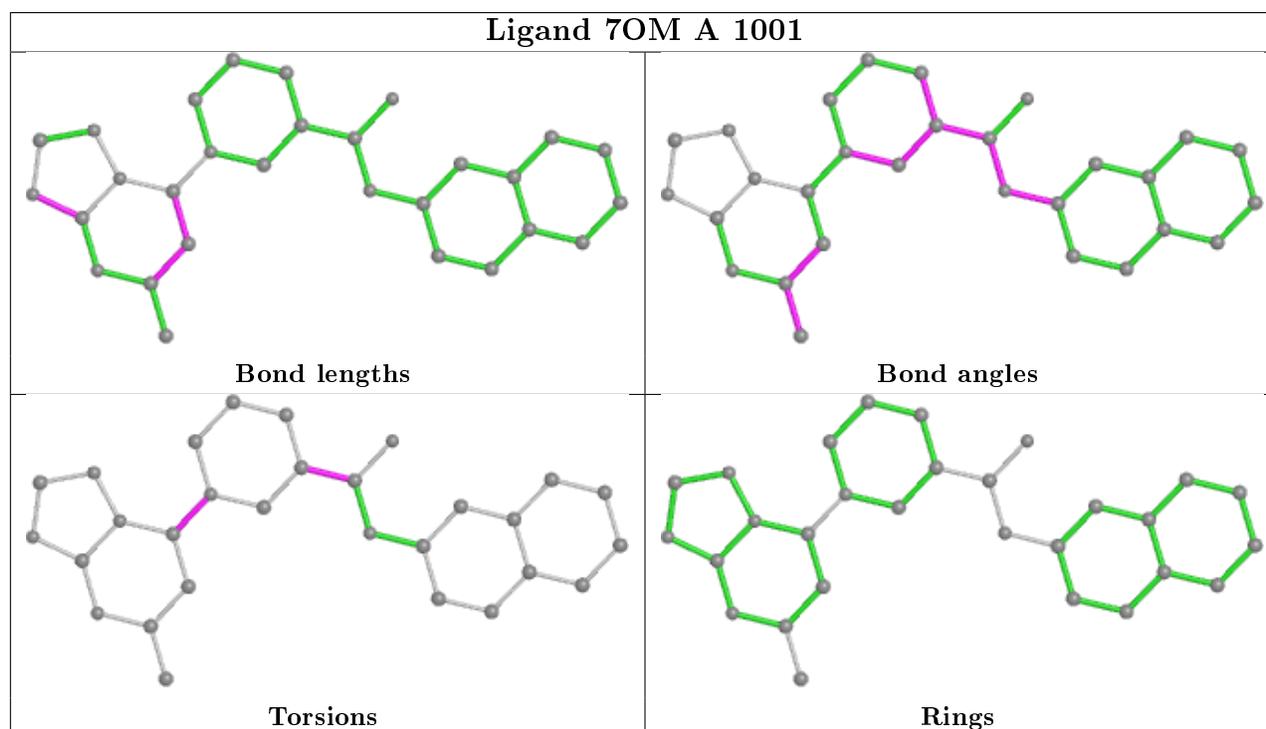
Mol	Chain	Res	Type	Atoms
2	A	1001	7OM	N8-C1-C2-C3
2	C	1001	7OM	N8-C1-C2-C3
2	B	1001	7OM	N8-C1-C2-C3
2	D	1001	7OM	N8-C1-C2-C3
2	A	1001	7OM	O9-C1-C2-C3
2	B	1001	7OM	O9-C1-C2-C3
2	C	1001	7OM	O9-C1-C2-C3
2	D	1001	7OM	O9-C1-C2-C3
2	A	1001	7OM	C15-C18-N4-C3
2	D	1001	7OM	C15-C18-N4-C3
2	B	1001	7OM	N10-C18-N4-C3
2	B	1001	7OM	C15-C18-N4-C3
2	C	1001	7OM	C15-C18-N4-C3
2	A	1001	7OM	N10-C18-N4-C3
2	D	1001	7OM	N10-C18-N4-C3

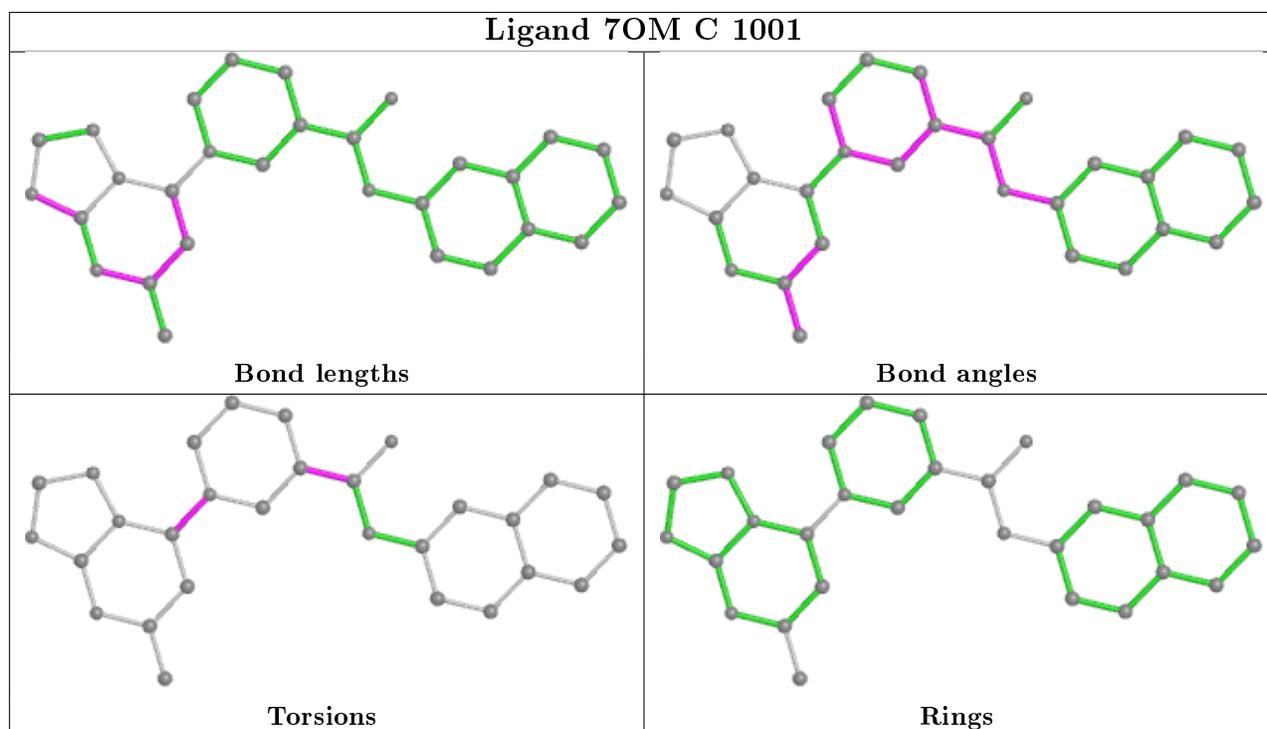
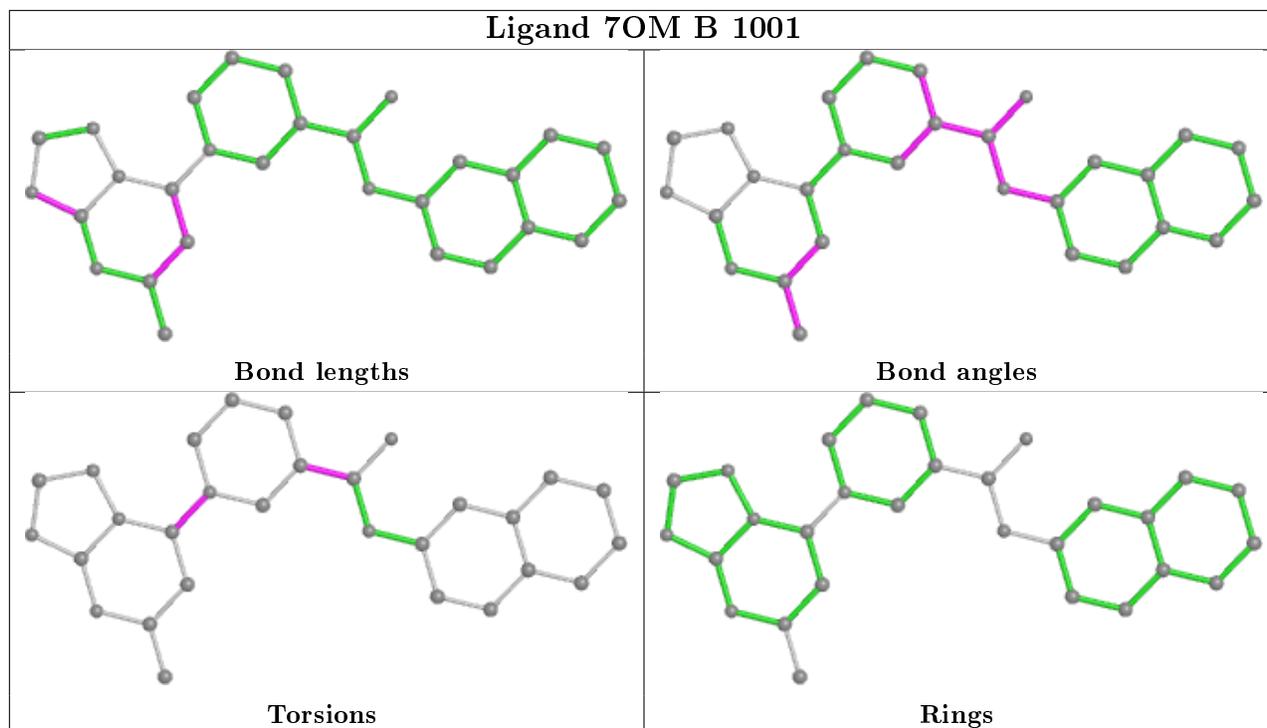
There are no ring outliers.

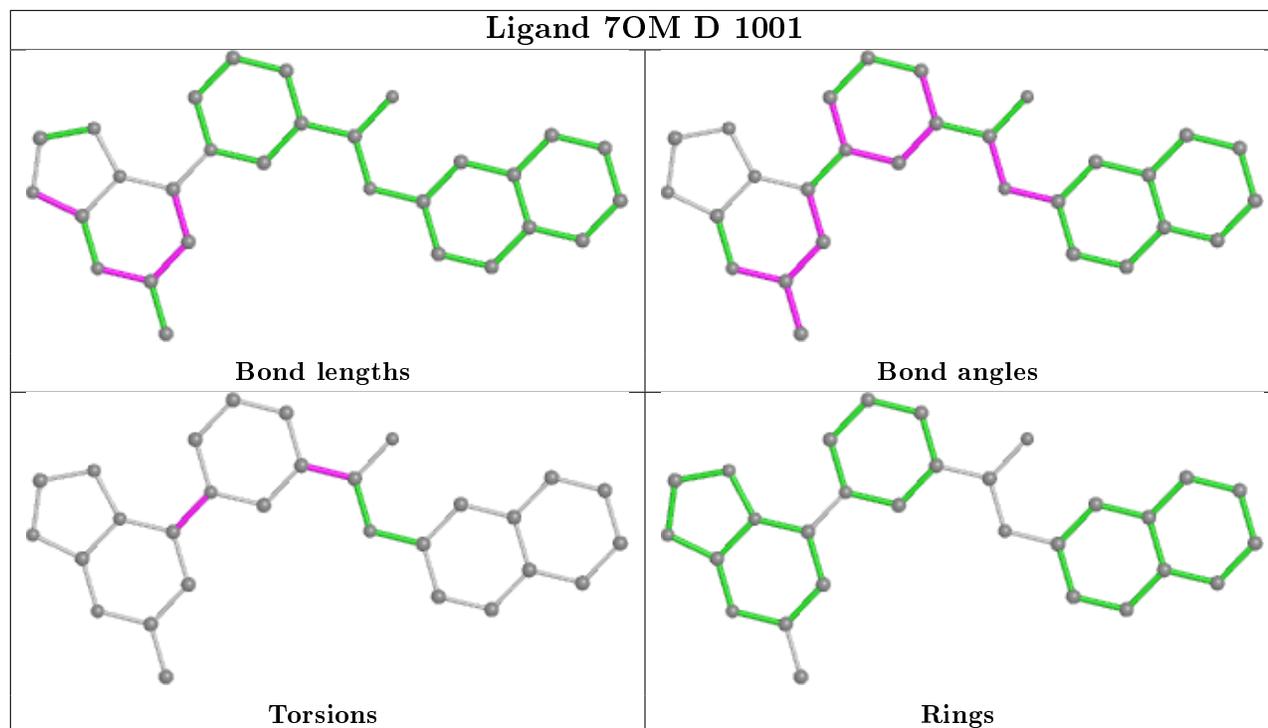
4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1001	7OM	1	0
2	B	1001	7OM	2	0
2	C	1001	7OM	2	0
2	D	1001	7OM	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	A	341/344 (99%)	-0.34	0 100 100	12, 25, 47, 65	0
1	B	340/344 (98%)	-0.39	3 (0%) 84 87	13, 26, 44, 62	0
1	C	334/344 (97%)	-0.07	11 (3%) 46 51	14, 32, 56, 73	0
1	D	327/344 (95%)	-0.37	1 (0%) 94 95	13, 26, 46, 64	0
All	All	1342/1376 (97%)	-0.29	15 (1%) 80 84	12, 27, 50, 73	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	909	SER	4.1
1	C	786	VAL	3.7
1	C	580	ASP	3.3
1	C	843	ARG	3.1
1	C	680	TYR	2.7
1	B	722	GLU	2.3
1	B	578	MET	2.3
1	C	581	GLU	2.2
1	C	915	PHE	2.2
1	B	723	GLY	2.2
1	C	785	GLU	2.1
1	C	916	LEU	2.1
1	D	592	PRO	2.1
1	C	593	VAL	2.1
1	C	592	PRO	2.1

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

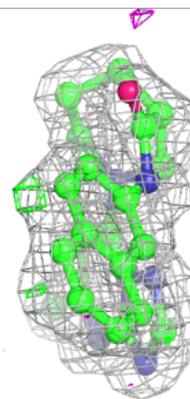
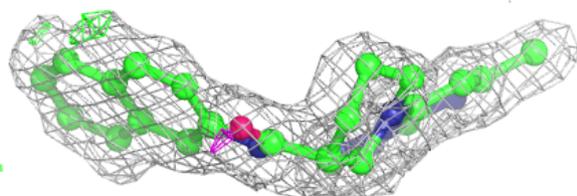
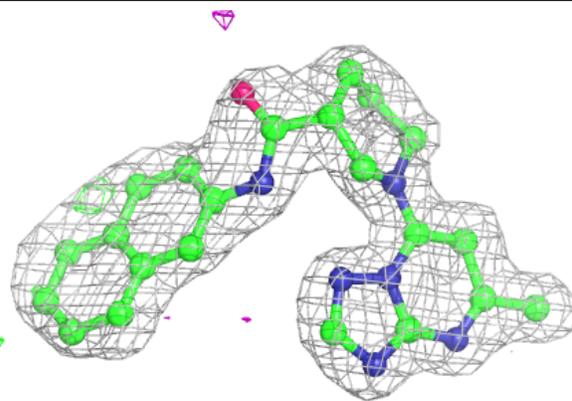
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	7OM	C	1001	29/29	0.90	0.11	18,30,44,48	0
2	7OM	B	1001	29/29	0.93	0.12	15,40,48,51	0
2	7OM	A	1001	29/29	0.93	0.11	12,32,39,46	0
2	7OM	D	1001	29/29	0.93	0.09	13,27,36,40	0
4	MG	D	1003	1/1	0.98	0.05	14,14,14,14	0
4	MG	B	1003	1/1	0.99	0.06	12,12,12,12	0
4	MG	C	1003	1/1	0.99	0.04	14,14,14,14	0
4	MG	A	1003	1/1	0.99	0.05	12,12,12,12	0
3	ZN	B	1002	1/1	1.00	0.06	16,16,16,16	0
3	ZN	C	1002	1/1	1.00	0.06	17,17,17,17	0
3	ZN	D	1002	1/1	1.00	0.06	18,18,18,18	0
3	ZN	A	1002	1/1	1.00	0.04	16,16,16,16	0

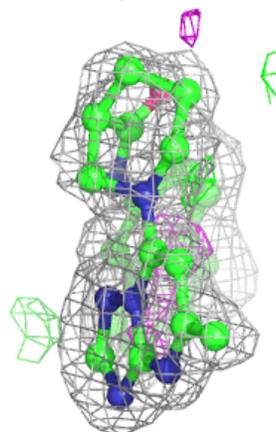
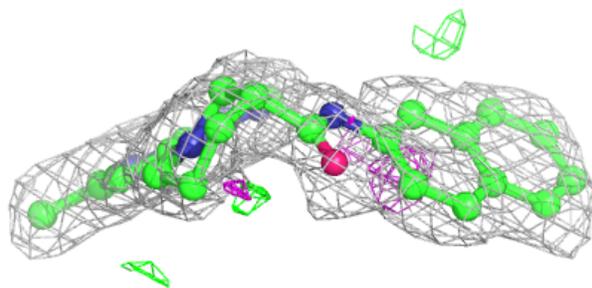
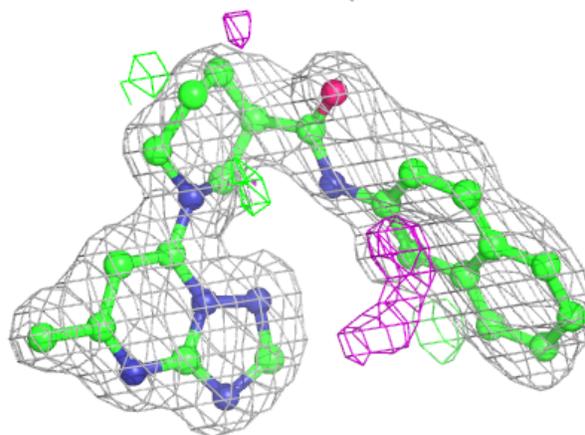
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 7OM 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)

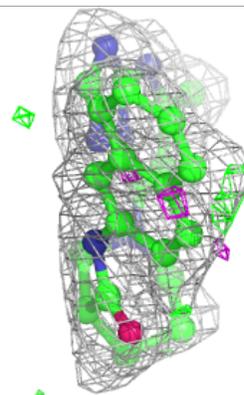
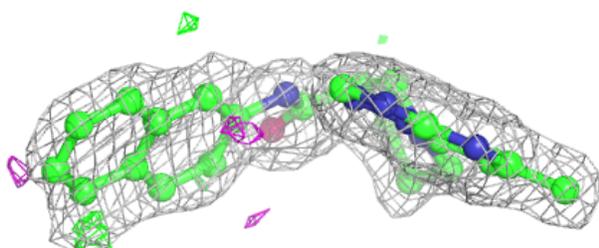
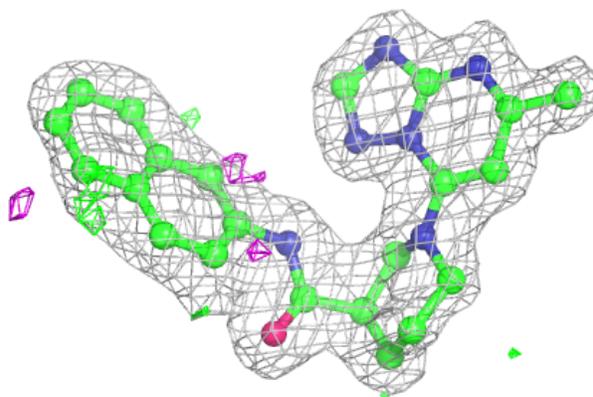
**Electron density around 7OM B 1001:**

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

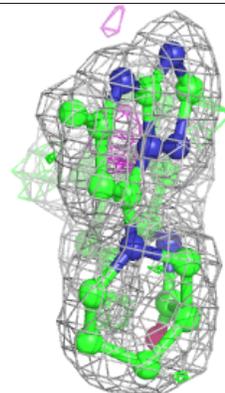
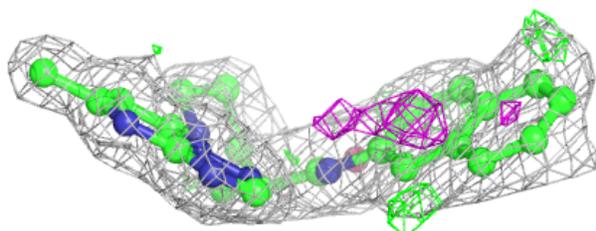
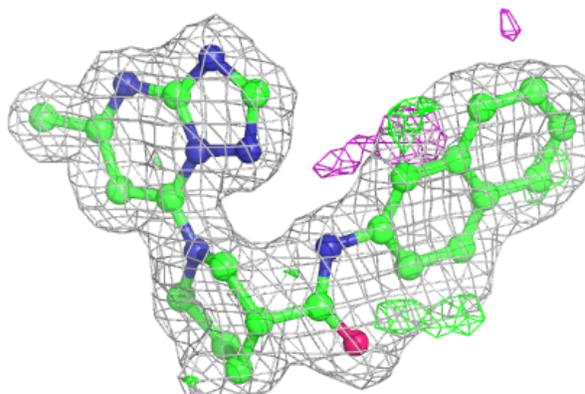


Electron density around 7OM A 1001:

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

**Electron density around 7OM D 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

There are no such residues in this entry.