



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:32 PM UTC

PDB ID : 7AT9 / pdb_00007at9
Title : Structure of protein kinase ck2 catalytic subunit (csnk2a2 gene product) in complex with the ATP-competitive inhibitor MB002 and the alphaD-pocket ligand 3,4-dichlorophenethylamine
Authors : Lindenblatt, D.; Applegate, V.; Nickelsen, A.; Klussmann, M.; Neundorf, I.; Goetz, C.; Jose, J.; Niefind, K.
Deposited on : 2020-10-29
Resolution : 1.05 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

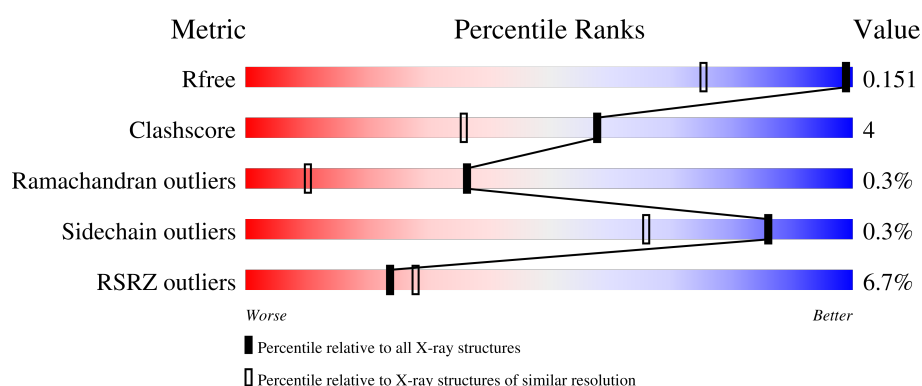
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.05 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	1182 (1.08-1.04)
Clashscore	190562	1204 (1.08-1.04)
Ramachandran outliers	187476	1175 (1.08-1.04)
Sidechain outliers	187428	1176 (1.08-1.04)
RSRZ outliers	180081	1181 (1.08-1.04)

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	364	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	4B0	A	511[A]	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6417 atoms, of which 2989 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.

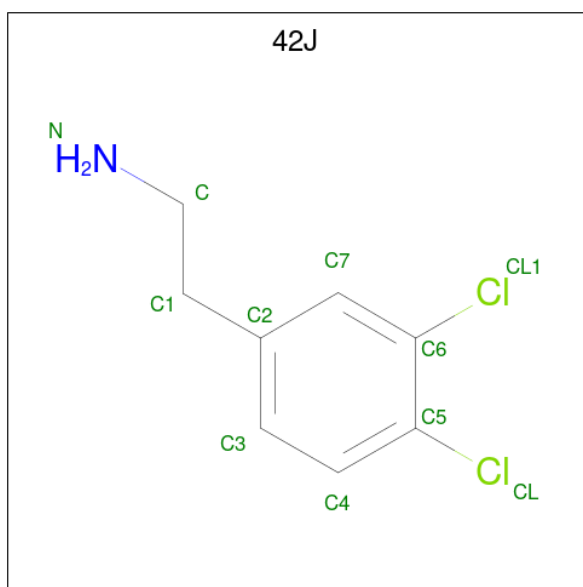
- Molecule 1 is a protein called Casein kinase II subunit alpha'.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	328	Total	C	H	N	O	S	0	21	0
			5851	1888	2915	509	526	13			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	initiating methionine	UNP P19784
A	-12	GLY	-	expression tag	UNP P19784
A	-11	SER	-	expression tag	UNP P19784
A	-10	SER	-	expression tag	UNP P19784
A	-9	HIS	-	expression tag	UNP P19784
A	-8	HIS	-	expression tag	UNP P19784
A	-7	HIS	-	expression tag	UNP P19784
A	-6	HIS	-	expression tag	UNP P19784
A	-5	HIS	-	expression tag	UNP P19784
A	-4	HIS	-	expression tag	UNP P19784
A	-3	SER	-	expression tag	UNP P19784
A	-2	GLN	-	expression tag	UNP P19784
A	-1	ASP	-	expression tag	UNP P19784
A	0	PRO	-	expression tag	UNP P19784
A	336	SER	CYS	conflict	UNP P19784

- Molecule 2 is 2-(3,4-dichlorophenyl)ethanamine (CCD ID: 42J) (formula: C₈H₉Cl₂N) (labeled as "Ligand of Interest" by depositor).

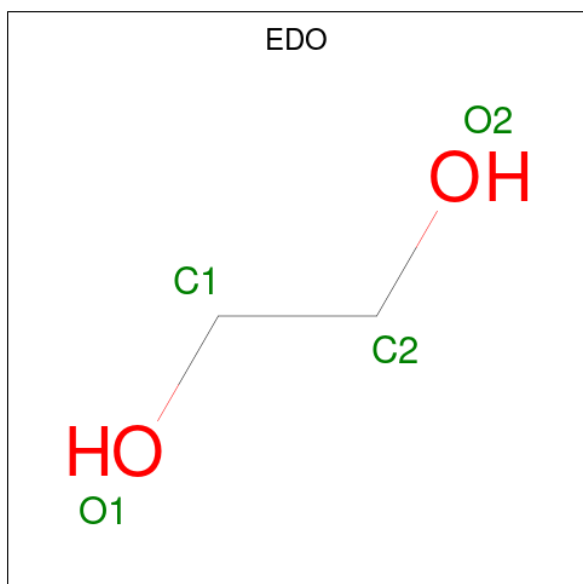


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Cl	H	N	0	0
			20	8	2	9	1		
2	A	1	Total	C	Cl	H	N	0	1
			20	8	2	9	1		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

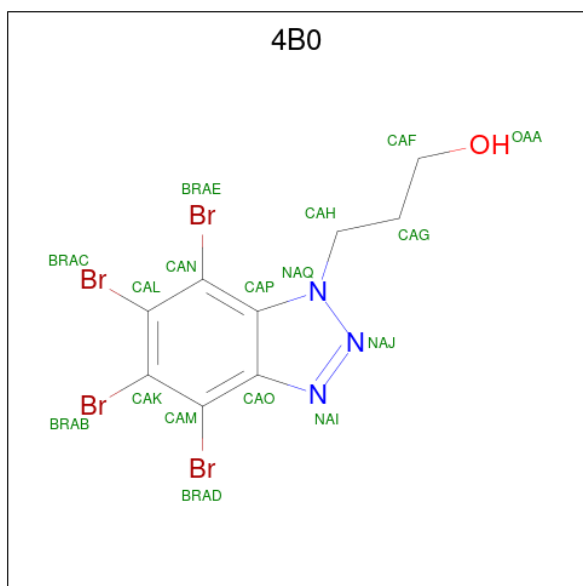
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Cl	0	0
			2	2		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			10	2	6	2		
4	A	1	Total	C	H	O	0	0
			10	2	6	2		
4	A	1	Total	C	H	O	0	0
			10	2	6	2		
4	A	1	Total	C	H	O	0	0
			10	2	6	2		
4	A	1	Total	C	H	O	0	0
			10	2	6	2		
4	A	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 5 is 3-(4,5,6,7-tetrabromo-1H-benzotriazol-1-yl)propan-1-ol (CCD ID: 4B0) (formula: C₉H₇Br₄N₃O).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total	Br	C	H	N	O	0	1
			48	8	18	14	6	2		

- Molecule 6 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	2	Total	Na	0	0
			2	2		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	390	Total 404	O 404	0	14

- Molecule 1: Casein kinase II subunit alpha'



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	46.48Å 47.66Å 50.73Å 66.61° 89.44° 88.66°	Depositor
Resolution (Å)	43.74 – 1.05 43.74 – 1.05	Depositor EDS
% Data completeness (in resolution range)	81.2 (43.74-1.05) 82.3 (43.74-1.05)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.26 (at 1.05Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.129 , 0.150 0.130 , 0.151	Depositor DCC
R_{free} test set	1555 reflections (0.82%)	wwPDB-VP
Wilson B-factor (Å ²)	11.5	Xtriage
Anisotropy	0.007	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.010 for h,-k,-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	6417	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.59% 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: NA, EDO, CL, 42J, 4B0

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	0.47	0/3012	0.65	0/4064

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	2936	2915	2897	22	0
2	A	22	18	18	0	0
3	A	2	0	0	0	0
4	A	28	42	42	1	0
5	A	34	14	14	14	0
6	A	2	0	0	0	0
7	A	404	0	0	6	0
All	All	3428	2989	2971	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 4.

All (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)
5:A:511[A]:4B0:BRAE	5:A:511[A]:4B0:HAGA	2.17	0.99
5:A:511[A]:4B0:BRAE	5:A:511[A]:4B0:CAG	2.69	0.96
5:A:511[A]:4B0:BRAE	5:A:511[A]:4B0:HAG	2.35	0.80
1:A:332:GLN:NE2	7:A:601:HOH:O	2.24	0.70
1:A:214:LEU:CD1	7:A:700:HOH:O	2.39	0.70
1:A:214:LEU:HD11	7:A:700:HOH:O	1.93	0.67
1:A:125:LEU:HD13	1:A:128:ILE:HA	1.83	0.60
1:A:117:ILE:O	5:A:511[A]:4B0:BRAB	2.77	0.58
1:A:164[A]:MET:SD	5:A:511[A]:4B0:BRAB	3.17	0.57
1:A:232:PRO:HA	4:A:509:EDO:H22	1.86	0.56
1:A:164[A]:MET:HE1	5:A:511[A]:4B0:BRAC	2.62	0.55
1:A:164[A]:MET:SD	5:A:511[A]:4B0:CAK	2.95	0.54
1:A:54[A]:VAL:HG21	5:A:511[A]:4B0:HAH	1.89	0.52
1:A:226[B]:MET:HE1	1:A:231:GLU:OE2	2.09	0.52
1:A:122:PHE:HD1	1:A:125:LEU:HD11	1.76	0.51
1:A:117:ILE:O	5:A:511[B]:4B0:BRAC	2.87	0.48
1:A:164[A]:MET:SD	5:A:511[A]:4B0:CAL	3.02	0.47
1:A:135:ARG:HG2	1:A:324:TYR:CZ	2.49	0.47
1:A:48:ARG:HD2	1:A:53:GLU:OE1	2.15	0.47
5:A:511[B]:4B0:HAHA	5:A:511[B]:4B0:BRAE	2.71	0.46
1:A:69:LYS:HD2	5:A:511[A]:4B0:HAF	1.98	0.46
1:A:117:ILE:HG13	1:A:164[A]:MET:CE	2.47	0.45
5:A:511[A]:4B0:HAHA	7:A:872:HOH:O	2.19	0.42
1:A:51:TYR:C	1:A:72:LYS:HD2	2.46	0.40
1:A:164[A]:MET:SD	5:A:511[A]:4B0:BRAC	3.34	0.40
1:A:36:ASN:ND2	7:A:611:HOH:O	2.54	0.40
1:A:78:LYS:CG	7:A:620:HOH:O	2.70	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	347/364 (95%)	336 (97%)	10 (3%)	1 (0%)	36 10

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	128	ILE

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	320/327 (98%)	319 (100%)	1 (0%)	86 66

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	128	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	208	GLN
1	A	237	GLN
1	A	263	HIS

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 15 ligands modelled in this entry, 4 are monoatomic - leaving 11 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	42J	A	501	-	11,11,11	0.24	0	14,14,14	0.49	0
4	EDO	A	507	-	3,3,3	0.43	0	2,2,2	0.39	0
4	EDO	A	508	-	3,3,3	0.61	0	2,2,2	0.35	0
4	EDO	A	510	-	3,3,3	0.41	0	2,2,2	0.39	0
4	EDO	A	509	-	3,3,3	0.44	0	2,2,2	0.34	0
5	4B0	A	511[A]	-	18,18,18	0.49	0	21,26,26	1.42	3 (14%)
4	EDO	A	505	-	3,3,3	0.37	0	2,2,2	0.37	0
5	4B0	A	511[B]	-	18,18,18	0.82	0	21,26,26	1.77	4 (19%)
4	EDO	A	506	-	3,3,3	0.71	0	2,2,2	0.39	0
2	42J	A	512[C]	-	11,11,11	0.22	0	14,14,14	0.49	0
4	EDO	A	504	-	3,3,3	0.49	0	2,2,2	0.09	0

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	42J	A	501	-	-	0/3/3/3	0/1/1/1
4	EDO	A	507	-	-	0/1/1/1	-
4	EDO	A	508	-	-	0/1/1/1	-
4	EDO	A	510	-	-	0/1/1/1	-
4	EDO	A	509	-	-	0/1/1/1	-
5	4B0	A	511[A]	-	-	2/4/4/4	0/2/2/2
4	EDO	A	505	-	-	0/1/1/1	-
5	4B0	A	511[B]	-	-	0/4/4/4	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	A	506	-	-	1/1/1/1	-
2	42J	A	512[C]	-	-	0/3/3/3	0/1/1/1
4	EDO	A	504	-	-	0/1/1/1	-

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	511[B]	4B0	BRAB-CAK-CAM	-5.08	112.50	120.56
5	A	511[B]	4B0	BRAB-CAK-CAL	4.32	127.41	120.56
5	A	511[A]	4B0	CAN-CAP-NAQ	3.24	136.89	132.14
5	A	511[B]	4B0	BRAD-CAM-CAK	-2.86	116.03	120.56
5	A	511[A]	4B0	BRAE-CAN-CAL	-2.22	117.04	120.56
5	A	511[B]	4B0	CAN-CAP-NAQ	2.16	135.30	132.14
5	A	511[A]	4B0	CAH-NAQ-NAJ	-2.02	116.85	120.14

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	511[A]	4B0	CAG-CAH-NAQ-NAJ
5	A	511[A]	4B0	CAG-CAH-NAQ-CAP
4	A	506	EDO	O1-C1-C2-O2

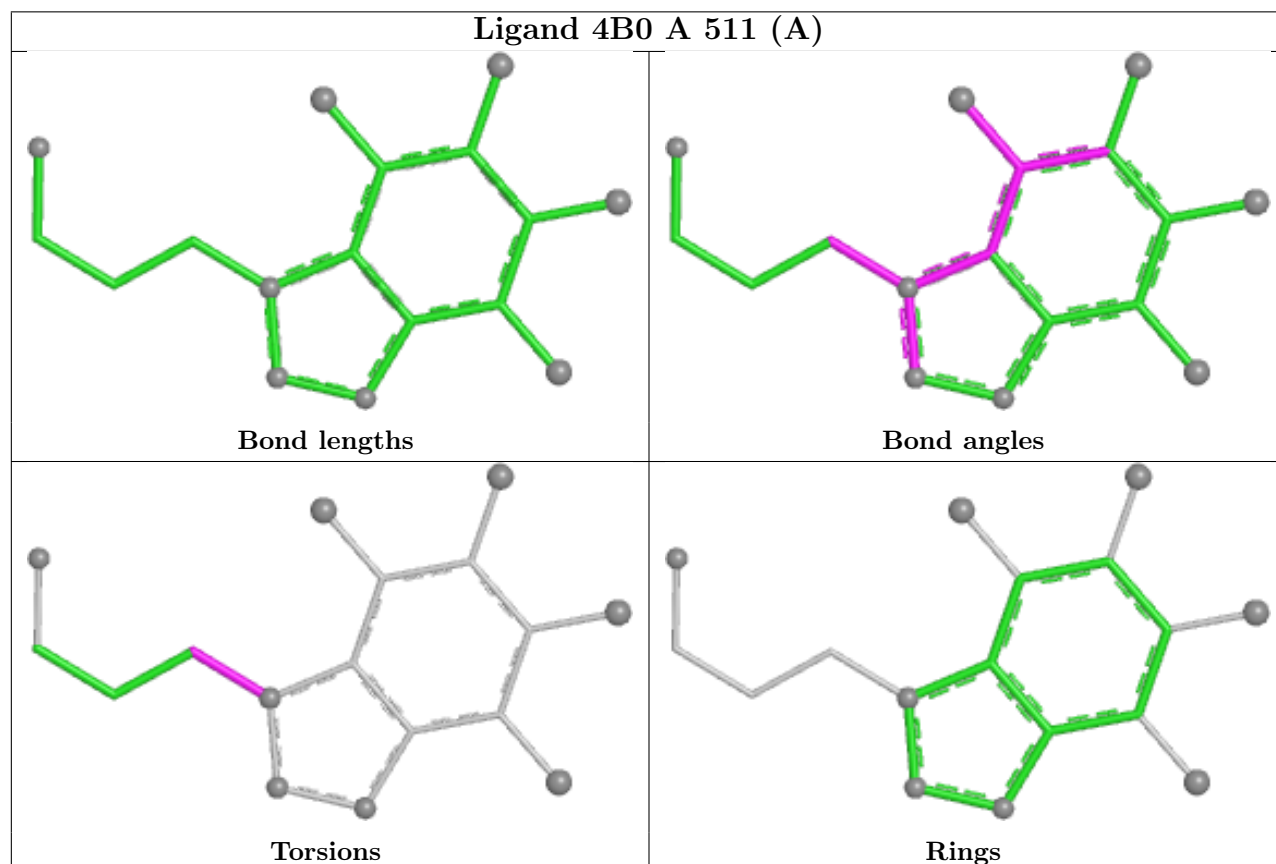
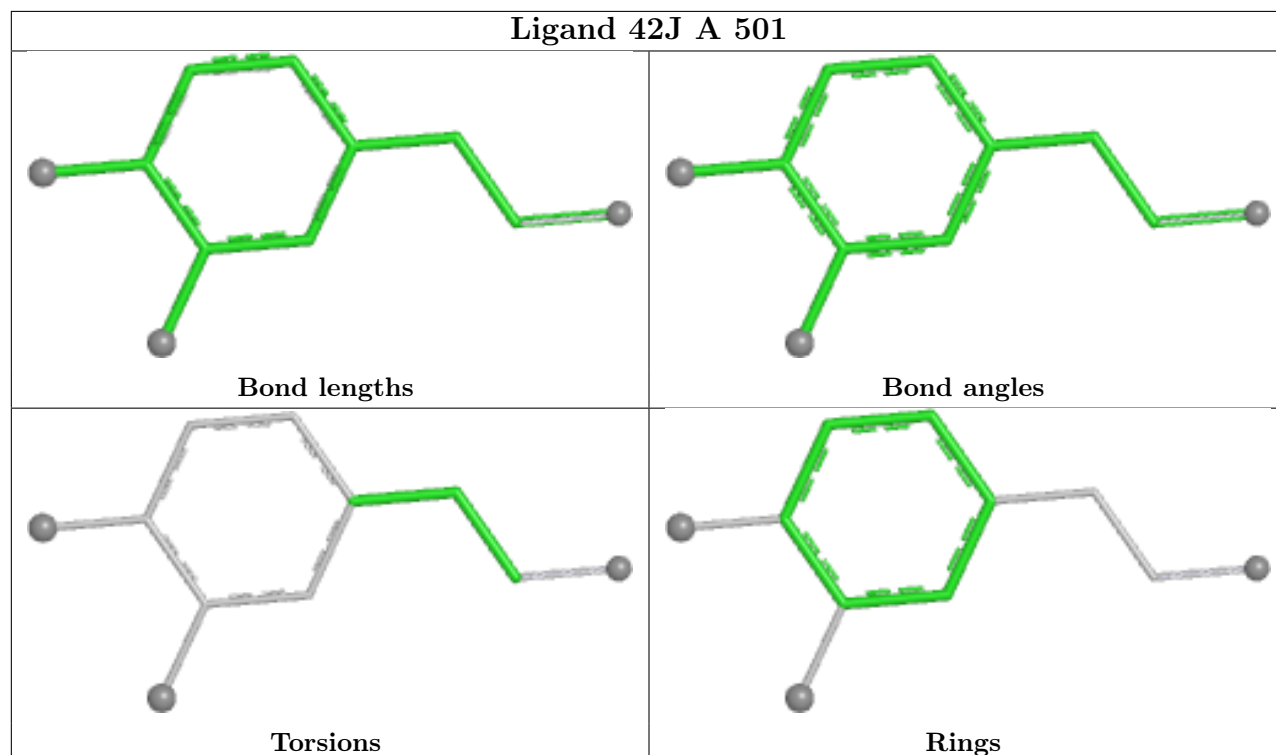
There are no ring outliers.

3 monomers are involved in 15 short contacts:

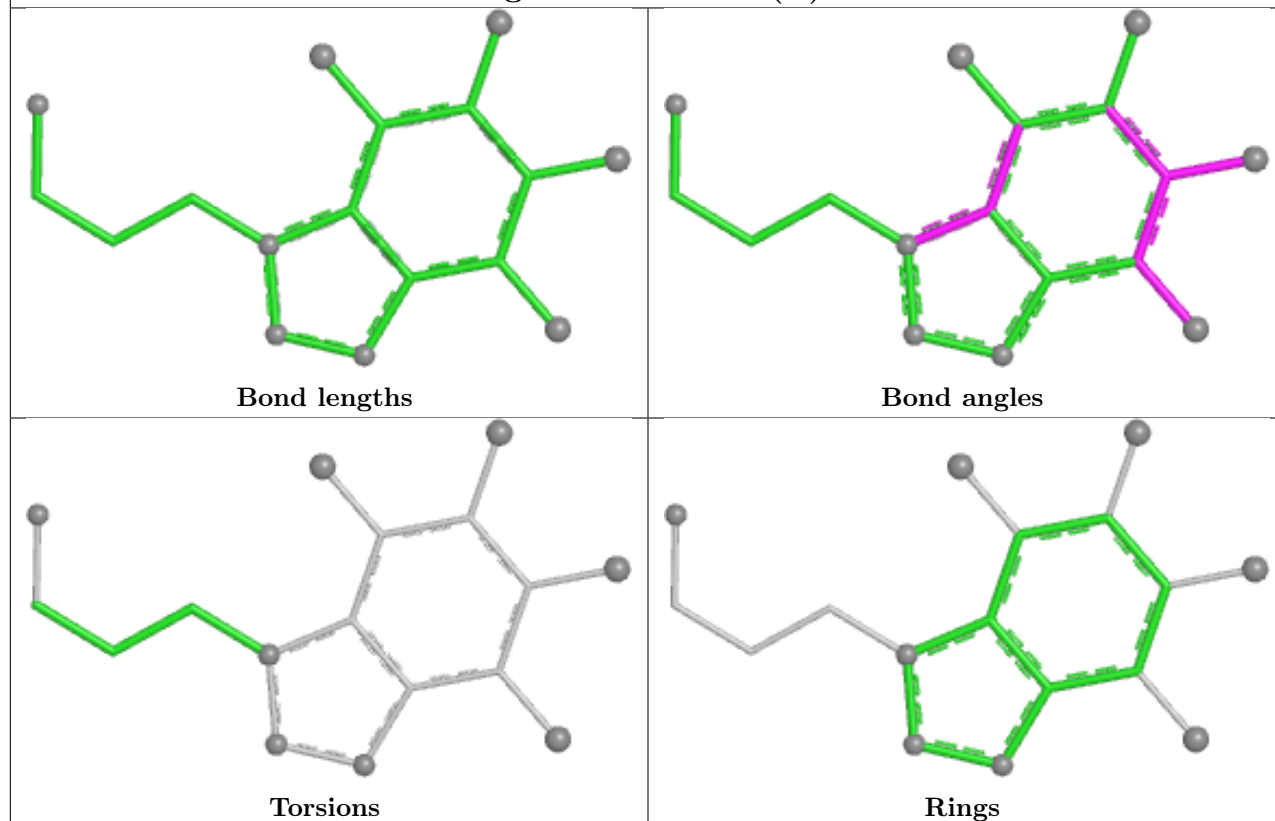
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	509	EDO	1	0
5	A	511[A]	4B0	12	0
5	A	511[B]	4B0	2	0

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

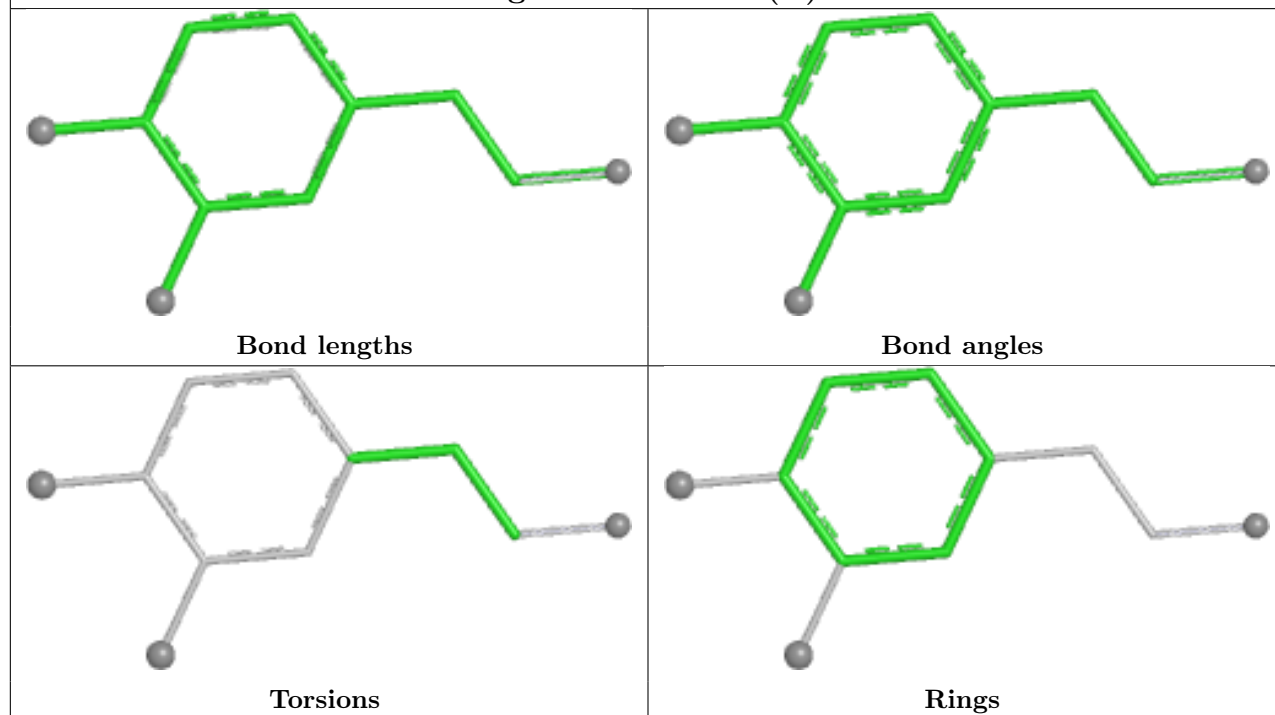
any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



Ligand 4B0 A 511 (B)



Ligand 42J A 512 (C)



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	328/364 (90%)	0.06	22 (6%) 24 28	6, 15, 41, 98	21 (6%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	74	VAL	10.5
1	A	128	ILE	10.3
1	A	125	LEU	8.7
1	A	126	TYR	7.9
1	A	6	ALA	7.4
1	A	73	PRO	5.4
1	A	129	LEU	4.2
1	A	51	TYR	4.0
1	A	332	GLN	3.9
1	A	124	GLN	3.9
1	A	7	GLY	3.7
1	A	127	GLN	3.0
1	A	333	SER	2.9
1	A	48	ARG	2.8
1	A	104	ASP	2.8
1	A	123	LYS	2.8
1	A	75	LYS	2.7
1	A	72	LYS	2.5
1	A	90	ARG	2.4
1	A	118	ASN	2.3
1	A	76	LYS	2.1
1	A	93	THR	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

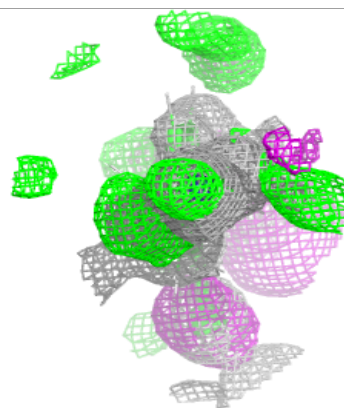
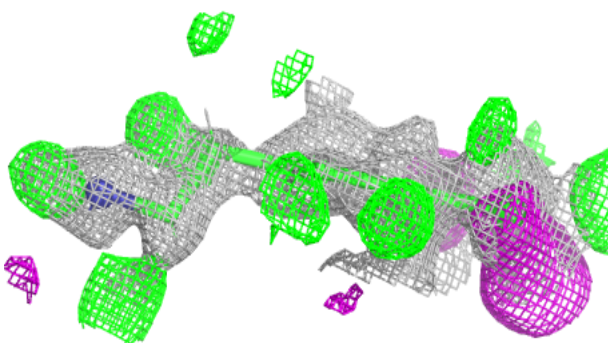
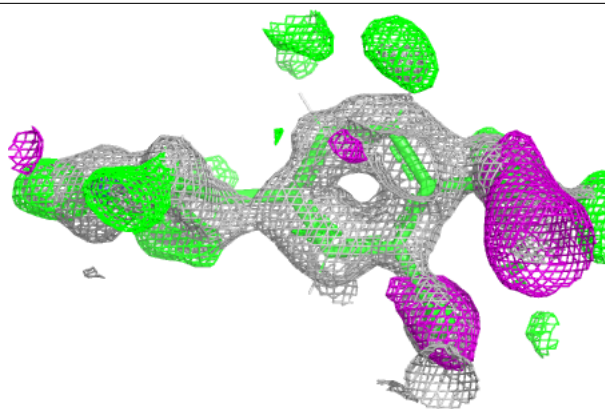
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	NA	A	513	1/1	0.48	0.27	44,44,44,44	0
4	EDO	A	507	4/4	0.76	0.18	51,61,61,62	0
2	42J	A	512[C]	11/11	0.79	0.22	16,20,24,25	20
4	EDO	A	510	4/4	0.87	0.12	35,42,43,43	0
5	4B0	A	511[B]	17/17	0.89	0.20	12,15,17,27	24
5	4B0	A	511[A]	17/17	0.89	0.20	7,14,20,33	24
4	EDO	A	509	4/4	0.90	0.14	12,22,33,40	0
4	EDO	A	504	4/4	0.91	0.13	22,36,40,43	0
4	EDO	A	506	4/4	0.91	0.11	17,21,25,30	0
4	EDO	A	508	4/4	0.96	0.07	27,33,37,37	0
4	EDO	A	505	4/4	0.96	0.07	22,27,29,33	0
6	NA	A	514	1/1	0.98	0.14	29,29,29,29	0
2	42J	A	501	11/11	0.99	0.06	13,15,18,18	0
3	CL	A	503	1/1	1.00	0.10	19,19,19,19	0
3	CL	A	502	1/1	1.00	0.10	18,18,18,18	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.

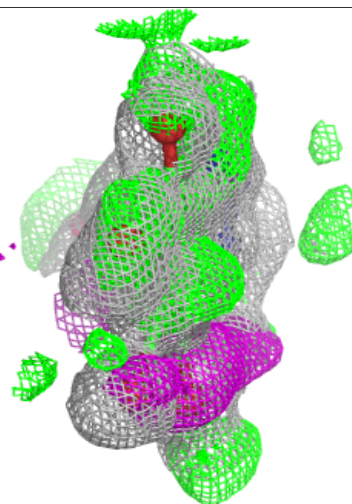
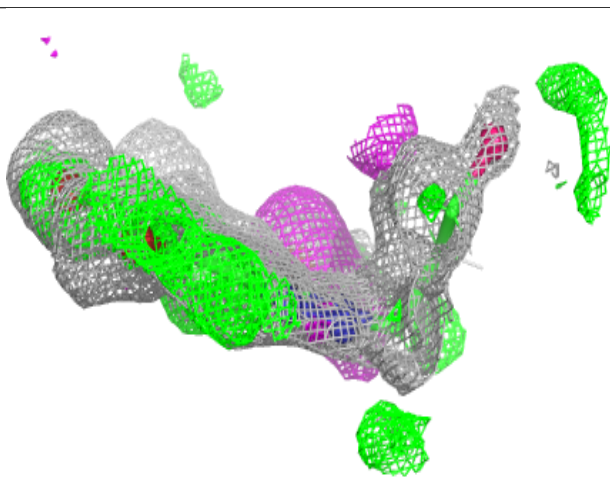
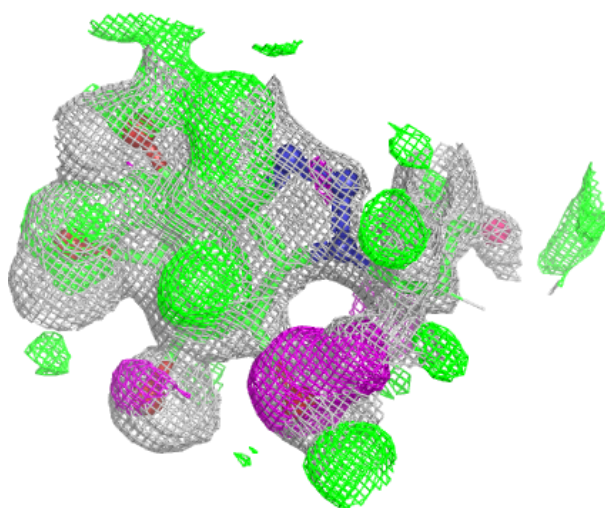
Electron density around 42J A 512 (C):

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



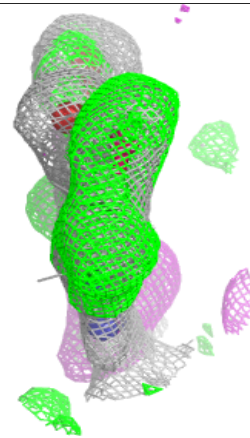
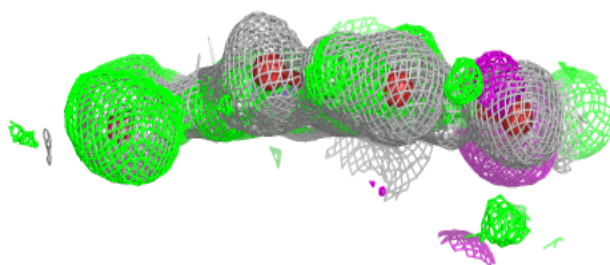
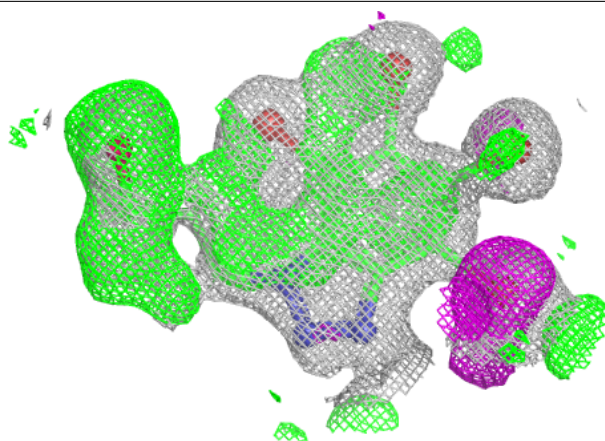
Electron density around 4B0 A 511 (B):

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



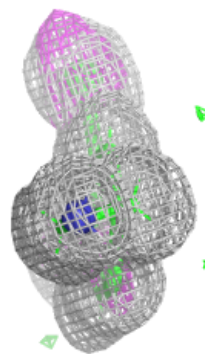
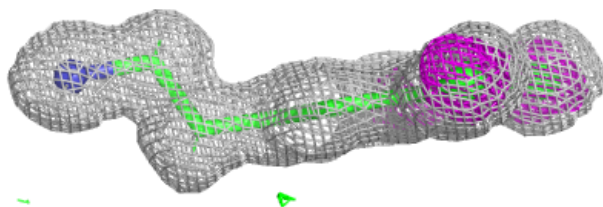
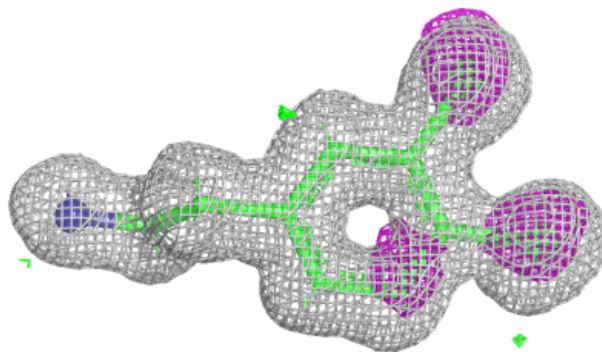
Electron density around 4B0 A 511 (A):

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



Electron density around 42J A 501:

$2mF_o-DF_c$ (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.