



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 03:50 PM UTC

PDB ID : 4YLJ / pdb_00004ylj
Title : Crystal structure of DYRK1A in complex with 10-Iodo-substituted 11H-indol
o[3,2-c]quinoline-6-carboxylic acid inhibitor 5j
Authors : Chaikuad, A.; Falke, H.; Nowak, R.; von Delft, F.; Arrowsmith, C.H.; Ed-
wards, A.M.; Bountra, C.; Kunick, C.; Knapp, S.; Structural Genomics Con-
sortium (SGC)
Deposited on : 2015-03-05
Resolution : 2.58 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	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
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)	:	2.48.1

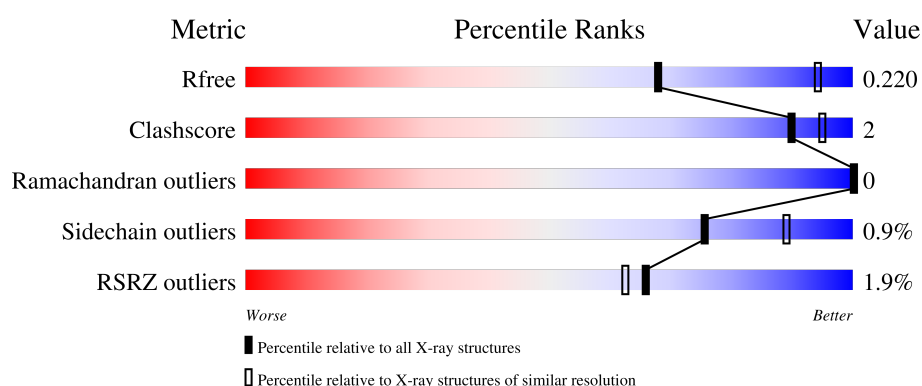
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.58 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}	164625	4456 (2.60-2.56)
Clashscore	180529	4905 (2.60-2.56)
Ramachandran outliers	177936	4847 (2.60-2.56)
Sidechain outliers	177891	4847 (2.60-2.56)
RSRZ outliers	164620	4456 (2.60-2.56)

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	361	0% 93% 5% •
1	B	361	2% 91% 5% •
1	C	361	2% 91% 5% •
1	D	361	2% 91% 5% •

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11988 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 Dual specificity tyrosine-phosphorylation-regulated kinase 1A.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	347	Total	C	N	O	P	S	0	3	0
			2860	1838	496	508	1	17			
1	B	348	Total	C	N	O	P	S	0	3	0
			2860	1838	489	514	2	17			
1	C	347	Total	C	N	O	P	S	0	3	0
			2857	1835	488	515	2	17			
1	D	348	Total	C	N	O	P	S	0	3	0
			2862	1838	489	516	2	17			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	125	SER	-	expression tag	UNP Q13627
A	126	MET	-	expression tag	UNP Q13627
B	125	SER	-	expression tag	UNP Q13627
B	126	MET	-	expression tag	UNP Q13627
C	125	SER	-	expression tag	UNP Q13627
C	126	MET	-	expression tag	UNP Q13627
D	125	SER	-	expression tag	UNP Q13627
D	126	MET	-	expression tag	UNP Q13627

- Molecule 2 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



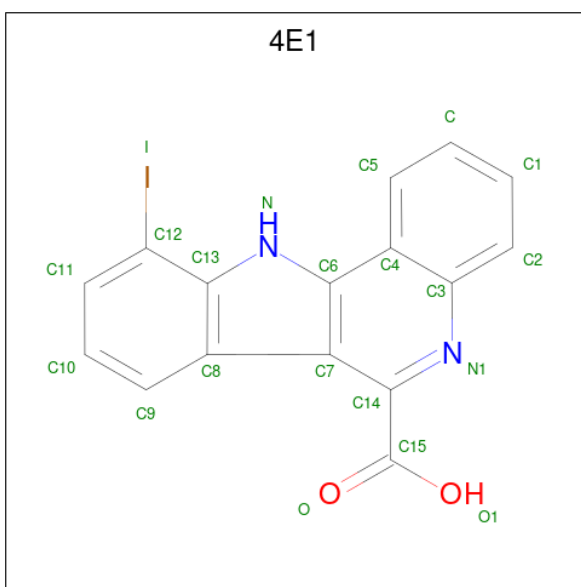
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	8	5		
2	A	1	Total	C	O	0	0
			10	6	4		
2	B	1	Total	C	O	0	0
			13	8	5		
2	B	1	Total	C	O	0	0
			13	8	5		
2	B	1	Total	C	O	0	0
			13	8	5		
2	C	1	Total	C	O	0	0
			13	8	5		
2	D	1	Total	C	O	0	0
			13	8	5		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is 10-iodo-11H-indolo[3,2-c]quinoline-6-carboxylic acid (CCD ID: 4E1) (formula: C₁₆H₉IN₂O₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 41	C 32	I 1	N 4	O 4	0	1
4	B	1	Total 41	C 32	I 1	N 4	O 4	0	1
4	C	1	Total 41	C 32	I 1	N 4	O 4	0	1
4	D	1	Total 41	C 32	I 1	N 4	O 4	0	1

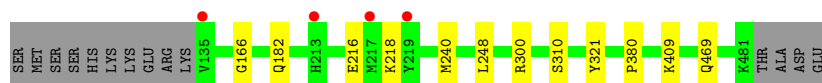
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	110	Total	O	0	0
			110	110		
5	B	59	Total	O	0	0
			59	59		
5	C	39	Total	O	0	0
			39	39		
5	D	39	Total	O	0	0
			39	39		

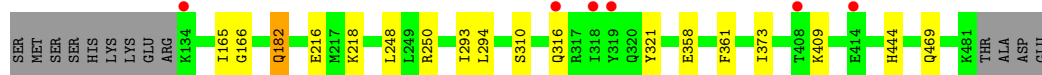
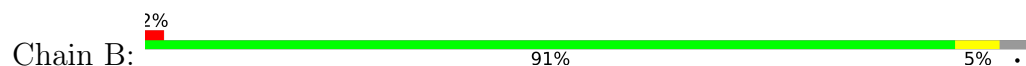
3 Residue-property plots [i](#)

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.

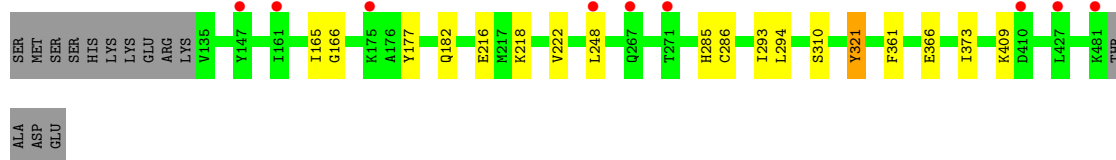
- Molecule 1: Dual specificity tyrosine-phosphorylation-regulated kinase 1A



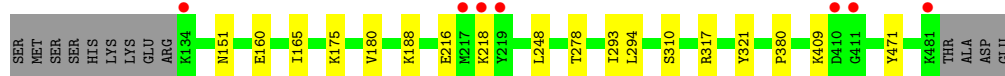
- Molecule 1: Dual specificity tyrosine-phosphorylation-regulated kinase 1A



- Molecule 1: Dual specificity tyrosine-phosphorylation-regulated kinase 1A



- Molecule 1: Dual specificity tyrosine-phosphorylation-regulated kinase 1A



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	265.47Å 65.53Å 139.37Å 90.00° 114.60° 90.00°	Depositor
Resolution (Å)	48.41 – 2.58 48.41 – 2.58	Depositor EDS
% Data completeness (in resolution range)	99.5 (48.41-2.58) 99.5 (48.41-2.58)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.192 , 0.222 0.195 , 0.220	Depositor DCC
R_{free} test set	3484 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	65.1	Xtriage
Anisotropy	0.337	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11988	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% 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: SO4, PTR, PG4, 4E1

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	0.84	0/2919	0.77	1/3932 (0.0%)
1	B	0.82	1/2902 (0.0%)	0.79	1/3913 (0.0%)
1	C	0.73	0/2899	0.76	1/3909 (0.0%)
1	D	0.75	0/2904	0.76	4/3916 (0.1%)
All	All	0.79	1/11624 (0.0%)	0.77	7/15670 (0.0%)

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	A	0	1
1	B	0	1
1	C	0	1
1	D	0	1
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	358	GLU	CA-C	-5.54	1.47	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	380	PRO	O-C-N	5.90	123.92	121.15
1	D	165	ILE	CB-CA-C	-5.73	104.72	111.94
1	B	165	ILE	CB-CA-C	-5.49	105.02	111.94
1	D	380	PRO	O-C-N	5.25	123.62	121.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	278	THR	CA-C-N	5.14	125.18	119.32
1	D	278	THR	C-N-CA	5.14	125.18	119.32
1	C	165	ILE	CB-CA-C	-5.09	105.52	111.94

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	409	LYS	Peptide
1	B	409	LYS	Peptide
1	C	409	LYS	Peptide
1	D	409	LYS	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	2860	0	2871	8	0
1	B	2860	0	2848	8	1
1	C	2857	0	2846	10	0
1	D	2862	0	2848	9	1
2	A	23	0	31	2	0
2	B	39	0	54	0	0
2	C	13	0	18	0	0
2	D	13	0	18	2	0
3	A	10	0	0	1	0
3	B	20	0	0	1	0
3	C	5	0	0	0	0
3	D	15	0	0	0	0
4	A	41	0	16	3	0
4	B	41	0	16	2	0
4	C	41	0	16	4	0
4	D	41	0	16	4	0
5	A	110	0	0	0	0
5	B	59	0	0	1	0
5	C	39	0	0	0	0
5	D	39	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	11988	0	11598	40	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 (40) 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)
4:D:505[B]:4E1:O	4:D:505[B]:4E1:H2	1.52	1.08
4:B:508[B]:4E1:H2	4:B:508[B]:4E1:O	1.65	0.97
1:A:300:ARG:HD3	1:C:177:TYR:CZ	2.12	0.84
4:A:505[B]:4E1:O1	4:A:505[B]:4E1:H2	1.86	0.76
1:A:240:MET:HE2	2:A:501:PG4:H22	1.71	0.72
1:D:218:LYS:NZ	5:D:613:HOH:O	2.23	0.70
4:C:503[B]:4E1:H2	4:C:503[B]:4E1:O	1.91	0.69
4:D:505[B]:4E1:O	4:D:505[B]:4E1:C9	2.38	0.69
1:B:166:GLY:HA2	1:D:180:VAL:O	1.92	0.68
1:D:151:ASN:HB3	5:D:614:HOH:O	1.93	0.67
1:B:250:ARG:NE	1:D:160:GLU:OE1	2.26	0.66
1:A:300:ARG:CD	1:C:177:TYR:CZ	2.81	0.63
4:B:508[B]:4E1:O	4:B:508[B]:4E1:C9	2.48	0.56
4:C:503[A]:4E1:H2	4:C:503[A]:4E1:O	2.07	0.54
1:B:316:GLN:CG	1:B:316:GLN:O	2.56	0.52
1:A:300:ARG:HD3	1:C:177:TYR:CE1	2.44	0.52
1:A:469:GLN:HB2	3:A:504:SO4:O4	2.10	0.51
1:D:317:ARG:NH2	5:D:630:HOH:O	2.43	0.51
1:C:166:GLY:HA3	4:C:503[B]:4E1:H7	1.93	0.51
1:B:469:GLN:HB2	3:B:507:SO4:O2	2.12	0.50
1:A:240:MET:HE2	2:A:501:PG4:C2	2.40	0.49
1:A:166:GLY:HA3	4:A:505[B]:4E1:H7	1.95	0.47
1:D:175:LYS:NZ	2:D:501:PG4:H72	2.28	0.47
1:B:444:HIS:CE1	5:B:650:HOH:O	2.68	0.47
1:D:188:LYS:NZ	4:D:505[B]:4E1:O1	2.45	0.47
4:D:505[A]:4E1:H2	4:D:505[A]:4E1:O	2.15	0.47
2:D:501:PG4:H31	5:D:626:HOH:O	2.15	0.46
4:A:505[A]:4E1:O1	4:A:505[A]:4E1:H2	2.16	0.45
1:A:216:GLU:HB2	1:A:218:LYS:HE3	1.99	0.44
1:B:293:ILE:C	1:B:294:LEU:HD12	2.43	0.44
1:C:216:GLU:HB2	1:C:218:LYS:HE3	1.99	0.43
1:D:216:GLU:HB2	1:D:218:LYS:HE3	2.00	0.43
1:C:321[B]:PTR:O2P	1:C:366:GLU:HG3	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:216:GLU:HB2	1:B:218:LYS:HE3	2.00	0.43
1:C:222:VAL:HG21	4:C:503[A]:4E1:H3	2.01	0.43
1:C:361:PHE:CE1	1:C:373:ILE:HA	2.54	0.42
1:B:361:PHE:CE1	1:B:373:ILE:HA	2.54	0.42
1:D:293:ILE:C	1:D:294:LEU:HD12	2.45	0.42
1:C:285:HIS:O	1:C:286:CYS:HB2	2.20	0.42
1:C:293:ILE:C	1:C:294:LEU:HD12	2.45	0.41

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:182:GLN:OE1	1:D:471:TYR:OH[4_445]	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	347/361 (96%)	336 (97%)	11 (3%)	0	100	100
1	B	347/361 (96%)	336 (97%)	11 (3%)	0	100	100
1	C	346/361 (96%)	335 (97%)	11 (3%)	0	100	100
1	D	347/361 (96%)	336 (97%)	11 (3%)	0	100	100
All	All	1387/1444 (96%)	1343 (97%)	44 (3%)	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	308/320 (96%)	305 (99%)	3 (1%)	73	87
1	B	305/320 (95%)	302 (99%)	3 (1%)	73	87
1	C	306/320 (96%)	303 (99%)	3 (1%)	73	87
1	D	306/320 (96%)	304 (99%)	2 (1%)	81	92
All	All	1225/1280 (96%)	1214 (99%)	11 (1%)	75	89

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

Mol	Chain	Res	Type
1	A	182	GLN
1	A	248	LEU
1	A	310	SER
1	B	182	GLN
1	B	248	LEU
1	B	310	SER
1	C	182	GLN
1	C	248	LEU
1	C	310	SER
1	D	248	LEU
1	D	310	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	201	GLN
1	B	199	GLN
1	B	223	HIS
1	B	232	ASN
1	B	383	HIS
1	C	199	GLN
1	C	213	HIS
1	C	223	HIS
1	C	232	ASN
1	D	199	GLN
1	D	213	HIS
1	D	383	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

7 non-standard protein/DNA/RNA residues are modelled in this entry.

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
1	PTR	B	321[B]	-	15,16,17	0.75	0	17,22,24	1.32	3 (17%)
1	PTR	C	321[A]	-	15,16,17	0.90	0	17,22,24	1.24	3 (17%)
1	PTR	B	321[A]	-	15,16,17	2.81	8 (53%)	17,22,24	2.44	6 (35%)
1	PTR	D	321[B]	-	15,16,17	1.20	0	17,22,24	1.07	1 (5%)
1	PTR	A	321	1	15,16,17	0.93	0	17,22,24	1.18	2 (11%)
1	PTR	C	321[B]	-	15,16,17	1.52	2 (13%)	17,22,24	1.33	2 (11%)
1	PTR	D	321[A]	-	15,16,17	1.03	1 (6%)	17,22,24	1.14	2 (11%)

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
1	PTR	B	321[B]	-	-	0/10/11/13	0/1/1/1
1	PTR	C	321[A]	-	-	0/10/11/13	0/1/1/1
1	PTR	B	321[A]	-	-	3/10/11/13	0/1/1/1
1	PTR	D	321[B]	-	-	3/10/11/13	0/1/1/1
1	PTR	A	321	1	-	0/10/11/13	0/1/1/1
1	PTR	C	321[B]	-	-	4/10/11/13	0/1/1/1
1	PTR	D	321[A]	-	-	0/10/11/13	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	321[A]	PTR	P-OH	5.87	1.70	1.59
1	B	321[A]	PTR	CE1-CD1	4.25	1.45	1.38
1	B	321[A]	PTR	CB-CA	3.71	1.61	1.53
1	B	321[A]	PTR	CB-CG	3.68	1.60	1.51
1	C	321[B]	PTR	P-OH	3.34	1.66	1.59
1	B	321[A]	PTR	CE2-CD2	2.97	1.43	1.38
1	B	321[A]	PTR	OH-CZ	2.83	1.47	1.40
1	B	321[A]	PTR	CE1-CZ	2.49	1.43	1.38
1	C	321[B]	PTR	CB-CG	2.37	1.56	1.51
1	B	321[A]	PTR	CA-N	2.26	1.54	1.48
1	D	321[A]	PTR	OH-CZ	-2.06	1.36	1.40

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	321[A]	PTR	P-OH-CZ	5.46	143.31	123.88
1	B	321[A]	PTR	CE2-CZ-CE1	-4.71	113.29	120.16
1	B	321[A]	PTR	OH-CZ-CE1	3.28	129.04	119.22
1	B	321[A]	PTR	CD2-CE2-CZ	3.23	123.42	119.73
1	B	321[A]	PTR	CD2-CG-CD1	-3.04	113.72	118.23
1	B	321[A]	PTR	CD1-CE1-CZ	2.63	122.73	119.73
1	C	321[B]	PTR	OH-P-O1P	-2.60	100.79	109.48
1	D	321[A]	PTR	O3P-P-O2P	2.56	117.42	107.80
1	A	321	PTR	O2P-P-O1P	2.51	120.62	110.83
1	A	321	PTR	OH-P-O1P	-2.38	101.55	109.48
1	B	321[B]	PTR	O3P-P-O2P	2.34	116.56	107.80
1	B	321[B]	PTR	O2P-P-OH	-2.30	98.50	105.32
1	C	321[A]	PTR	O2P-P-OH	-2.29	98.56	105.32
1	C	321[A]	PTR	O3P-P-O2P	2.24	116.19	107.80
1	C	321[B]	PTR	CE2-CZ-CE1	-2.23	116.91	120.16
1	D	321[A]	PTR	O3P-P-O1P	2.21	119.46	110.83
1	B	321[B]	PTR	O3P-P-O1P	2.19	119.37	110.83
1	C	321[A]	PTR	O3P-P-O1P	2.17	119.28	110.83
1	D	321[B]	PTR	P-OH-CZ	2.07	131.26	123.88

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	321[A]	PTR	O-C-CA-CB
1	B	321[A]	PTR	N-CA-CB-CG
1	B	321[A]	PTR	C-CA-CB-CG

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Mol	Chain	Res	Type	Atoms
1	C	321[B]	PTR	O-C-CA-CB
1	C	321[B]	PTR	N-CA-CB-CG
1	C	321[B]	PTR	C-CA-CB-CG
1	D	321[B]	PTR	O-C-CA-CB
1	D	321[B]	PTR	N-CA-CB-CG
1	D	321[B]	PTR	C-CA-CB-CG
1	C	321[B]	PTR	CZ-OH-P-O2P

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	321[B]	PTR	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

25 ligands are modelled in this entry.

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	4E1	C	503[B]	-	21,23,24	1.87	4 (19%)	27,34,36	1.35	4 (14%)
3	SO4	D	503	-	4,4,4	0.53	0	6,6,6	0.43	0
2	PG4	A	502	-	9,9,12	0.50	0	8,8,11	0.45	0
2	PG4	B	503	-	12,12,12	0.72	0	11,11,11	0.49	0
3	SO4	B	507	-	4,4,4	0.47	0	6,6,6	0.60	0
4	4E1	C	503[A]	-	21,24,24	1.82	4 (19%)	24,36,36	1.33	4 (16%)
3	SO4	A	503	-	4,4,4	0.63	0	6,6,6	0.29	0
3	SO4	A	504	-	4,4,4	0.58	0	6,6,6	0.54	0
3	SO4	B	506	-	4,4,4	0.54	0	6,6,6	0.23	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	4E1	A	505[B]	-	21,23,24	2.08	7 (33%)	27,34,36	1.35	5 (18%)
2	PG4	B	501	-	12,12,12	0.70	0	11,11,11	1.09	0
4	4E1	A	505[A]	-	21,24,24	2.09	8 (38%)	24,36,36	1.30	2 (8%)
3	SO4	C	502	-	4,4,4	0.57	0	6,6,6	0.24	0
2	PG4	C	501	-	12,12,12	0.67	0	11,11,11	0.73	0
2	PG4	D	501	-	12,12,12	0.56	0	11,11,11	0.64	0
2	PG4	B	502	-	12,12,12	0.71	0	11,11,11	0.31	0
4	4E1	B	508[B]	-	21,23,24	1.64	3 (14%)	27,34,36	1.42	4 (14%)
4	4E1	B	508[A]	-	21,24,24	1.76	5 (23%)	24,36,36	1.41	3 (12%)
2	PG4	A	501	-	12,12,12	0.85	0	11,11,11	0.62	0
3	SO4	B	504	-	4,4,4	0.50	0	6,6,6	0.60	0
3	SO4	D	504	-	4,4,4	0.65	0	6,6,6	0.33	0
4	4E1	D	505[B]	-	21,23,24	1.95	5 (23%)	27,34,36	1.48	5 (18%)
3	SO4	B	505	-	4,4,4	0.52	0	6,6,6	0.36	0
3	SO4	D	502	-	4,4,4	0.49	0	6,6,6	0.54	0
4	4E1	D	505[A]	-	21,24,24	1.91	7 (33%)	24,36,36	1.16	3 (12%)

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	PG4	C	501	-	-	4/10/10/10	-
2	PG4	A	501	-	-	4/10/10/10	-
4	4E1	C	503[B]	-	-	2/4/4/4	0/4/4/4
2	PG4	D	501	-	-	4/10/10/10	-
4	4E1	A	505[B]	-	-	2/4/4/4	0/4/4/4
2	PG4	B	502	-	-	3/10/10/10	-
2	PG4	A	502	-	-	2/7/7/10	-
2	PG4	B	503	-	-	8/10/10/10	-
2	PG4	B	501	-	-	3/10/10/10	-
4	4E1	A	505[A]	-	-	3/4/4/4	0/4/4/4
4	4E1	B	508[B]	-	-	2/4/4/4	0/4/4/4
4	4E1	C	503[A]	-	-	2/4/4/4	0/4/4/4
4	4E1	D	505[B]	-	-	2/4/4/4	0/4/4/4
4	4E1	B	508[A]	-	-	2/4/4/4	0/4/4/4
4	4E1	D	505[A]	-	-	2/4/4/4	0/4/4/4

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	505[A]	4E1	C14-N1	4.17	1.36	1.32
4	A	505[A]	4E1	O-C15	4.13	1.34	1.22
4	C	503[A]	4E1	C14-N1	4.13	1.36	1.32
4	A	505[B]	4E1	O-C15	4.00	1.34	1.22
4	D	505[B]	4E1	C14-N1	4.00	1.36	1.32
4	A	505[B]	4E1	C14-N1	4.00	1.36	1.32
4	A	505[A]	4E1	C14-N1	3.96	1.36	1.32
4	D	505[B]	4E1	O-C15	3.93	1.34	1.22
4	D	505[A]	4E1	O-C15	3.92	1.34	1.22
4	C	503[A]	4E1	O-C15	3.81	1.33	1.22
4	C	503[B]	4E1	C14-N1	3.77	1.36	1.32
4	C	503[B]	4E1	O-C15	3.75	1.33	1.22
4	B	508[B]	4E1	C14-N1	3.64	1.36	1.32
4	A	505[B]	4E1	C4-C6	-3.63	1.38	1.45
4	B	508[B]	4E1	C4-C6	-3.62	1.38	1.45
4	B	508[A]	4E1	C14-N1	3.59	1.35	1.32
4	C	503[B]	4E1	C4-C6	-3.59	1.38	1.45
4	D	505[B]	4E1	C4-C6	-3.50	1.38	1.45
4	C	503[A]	4E1	C4-C6	-3.49	1.38	1.45
4	A	505[A]	4E1	C14-C15	-3.42	1.46	1.51
4	B	508[A]	4E1	C14-C15	-3.32	1.46	1.51
4	B	508[A]	4E1	C4-C6	-3.17	1.39	1.45
4	A	505[A]	4E1	C4-C6	-3.09	1.39	1.45
4	A	505[B]	4E1	C14-C15	-3.08	1.46	1.51
4	D	505[A]	4E1	C14-C15	-3.00	1.46	1.51
4	C	503[B]	4E1	O1-C15	-2.97	1.21	1.30
4	D	505[B]	4E1	O1-C15	-2.88	1.22	1.30
4	C	503[A]	4E1	O1-C15	-2.77	1.22	1.30
4	A	505[A]	4E1	O1-C15	-2.72	1.22	1.30
4	D	505[A]	4E1	C4-C6	-2.56	1.40	1.45
4	D	505[A]	4E1	C2-C3	-2.52	1.37	1.41
4	B	508[B]	4E1	C14-C7	-2.50	1.39	1.42
4	A	505[B]	4E1	O1-C15	-2.47	1.23	1.30
4	A	505[B]	4E1	C2-C3	-2.46	1.37	1.41
4	B	508[A]	4E1	C12-I	-2.44	2.06	2.10
4	D	505[B]	4E1	C14-C7	-2.41	1.39	1.42
4	A	505[A]	4E1	C10-C9	2.40	1.41	1.36
4	D	505[A]	4E1	O1-C15	-2.39	1.23	1.30
4	A	505[A]	4E1	C2-C3	-2.37	1.37	1.41
4	B	508[A]	4E1	C10-C9	2.26	1.41	1.36
4	A	505[A]	4E1	C12-I	-2.23	2.07	2.10
4	D	505[A]	4E1	C10-C9	2.14	1.41	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	505[B]	4E1	C14-C7	-2.05	1.39	1.42

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	508[B]	4E1	C9-C8-C13	3.59	122.92	118.17
4	D	505[B]	4E1	C9-C8-C13	3.57	122.90	118.17
4	B	508[A]	4E1	C7-C14-C15	3.32	129.52	123.30
4	C	503[B]	4E1	C9-C8-C13	3.24	122.46	118.17
4	B	508[A]	4E1	C15-C14-N1	-2.81	109.96	115.28
4	A	505[B]	4E1	C9-C8-C13	2.80	121.88	118.17
4	A	505[B]	4E1	O-C15-C14	-2.71	115.27	120.83
4	B	508[A]	4E1	C4-C3-N1	-2.70	120.47	122.98
4	D	505[B]	4E1	C5-C4-C3	2.68	121.34	117.58
4	A	505[A]	4E1	C4-C3-N1	-2.63	120.53	122.98
4	C	503[A]	4E1	C4-C3-N1	-2.62	120.54	122.98
4	A	505[A]	4E1	O-C15-C14	-2.58	115.55	120.83
4	B	508[B]	4E1	C5-C4-C3	2.51	121.10	117.58
4	A	505[B]	4E1	C5-C4-C3	2.50	121.08	117.58
4	D	505[A]	4E1	C4-C3-N1	-2.48	120.67	122.98
4	C	503[B]	4E1	C5-C4-C3	2.37	120.90	117.58
4	D	505[A]	4E1	C7-C14-C15	2.31	127.63	123.30
4	C	503[B]	4E1	C4-C3-N1	-2.31	120.83	122.98
4	C	503[A]	4E1	C14-N1-C3	2.29	121.77	117.28
4	C	503[A]	4E1	C5-C4-C3	2.25	120.74	117.58
4	B	508[B]	4E1	C4-C3-N1	-2.25	120.88	122.98
4	C	503[A]	4E1	C7-C14-C15	2.21	127.44	123.30
4	D	505[B]	4E1	C15-C14-N1	2.20	119.44	115.28
4	A	505[B]	4E1	C4-C3-N1	-2.18	120.95	122.98
4	C	503[B]	4E1	C14-N1-C3	2.17	121.53	117.28
4	D	505[A]	4E1	C14-N1-C3	2.16	121.52	117.28
4	A	505[B]	4E1	O1-C15-C14	2.15	121.16	114.35
4	B	508[B]	4E1	C9-C8-C7	-2.11	128.88	134.50
4	D	505[B]	4E1	C4-C3-N1	-2.06	121.06	122.98
4	D	505[B]	4E1	C9-C8-C7	-2.04	129.07	134.50

There are no chirality outliers.

All (45) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	505[A]	4E1	C7-C14-C15-O
4	A	505[A]	4E1	C7-C14-C15-O1

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Mol	Chain	Res	Type	Atoms
4	A	505[B]	4E1	C7-C14-C15-O
4	A	505[B]	4E1	C7-C14-C15-O1
4	B	508[A]	4E1	C7-C14-C15-O
4	B	508[A]	4E1	C7-C14-C15-O1
2	A	502	PG4	O2-C3-C4-O3
2	B	501	PG4	O3-C5-C6-O4
2	D	501	PG4	O3-C5-C6-O4
2	A	502	PG4	C3-C4-O3-C5
4	B	508[B]	4E1	C7-C14-C15-O
4	B	508[B]	4E1	C7-C14-C15-O1
4	C	503[A]	4E1	C7-C14-C15-O
4	C	503[A]	4E1	C7-C14-C15-O1
4	C	503[B]	4E1	C7-C14-C15-O
4	C	503[B]	4E1	C7-C14-C15-O1
4	D	505[A]	4E1	C7-C14-C15-O
4	D	505[A]	4E1	C7-C14-C15-O1
4	D	505[B]	4E1	C7-C14-C15-O
4	D	505[B]	4E1	C7-C14-C15-O1
2	B	503	PG4	O4-C7-C8-O5
2	A	501	PG4	O3-C5-C6-O4
2	B	502	PG4	O4-C7-C8-O5
2	B	501	PG4	O1-C1-C2-O2
2	B	501	PG4	O4-C7-C8-O5
2	D	501	PG4	O4-C7-C8-O5
2	A	501	PG4	O1-C1-C2-O2
2	C	501	PG4	C1-C2-O2-C3
2	C	501	PG4	O3-C5-C6-O4
2	C	501	PG4	O2-C3-C4-O3
2	B	502	PG4	C5-C6-O4-C7
2	B	503	PG4	O1-C1-C2-O2
2	B	503	PG4	C3-C4-O3-C5
2	A	501	PG4	C8-C7-O4-C6
2	D	501	PG4	C1-C2-O2-C3
2	B	503	PG4	C8-C7-O4-C6
2	B	503	PG4	C6-C5-O3-C4
2	A	501	PG4	C6-C5-O3-C4
2	B	502	PG4	C8-C7-O4-C6
2	D	501	PG4	C3-C4-O3-C5
4	A	505[A]	4E1	N1-C14-C15-O1
2	C	501	PG4	C4-C3-O2-C2
2	B	503	PG4	C5-C6-O4-C7
2	B	503	PG4	C4-C3-O2-C2

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Mol	Chain	Res	Type	Atoms
2	B	503	PG4	O2-C3-C4-O3

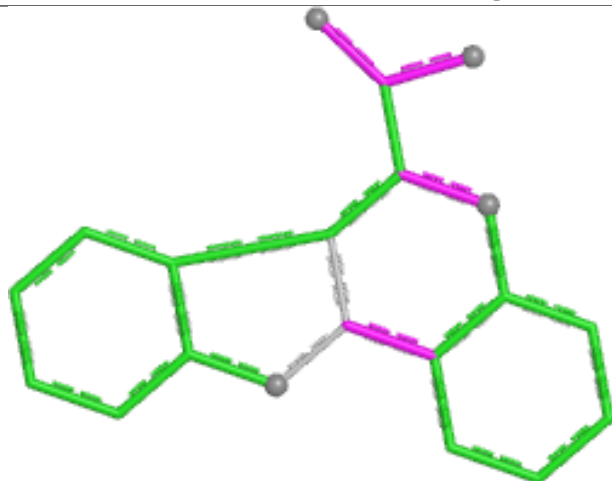
There are no ring outliers.

11 monomers are involved in 19 short contacts:

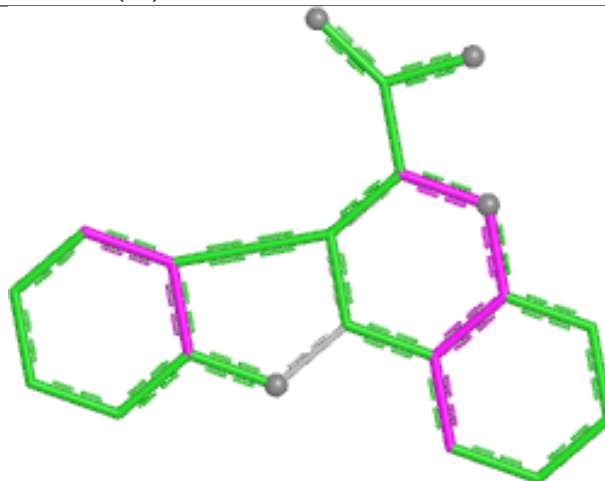
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	503[B]	4E1	2	0
3	B	507	SO4	1	0
4	C	503[A]	4E1	2	0
3	A	504	SO4	1	0
4	A	505[B]	4E1	2	0
4	A	505[A]	4E1	1	0
2	D	501	PG4	2	0
4	B	508[B]	4E1	2	0
2	A	501	PG4	2	0
4	D	505[B]	4E1	3	0
4	D	505[A]	4E1	1	0

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.

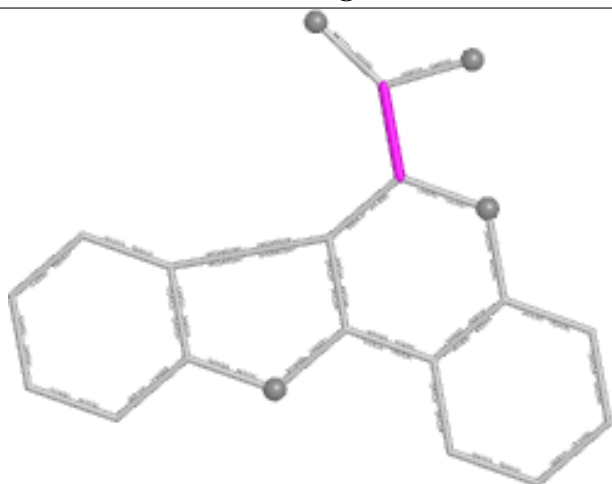
Ligand 4E1 C 503 (B)



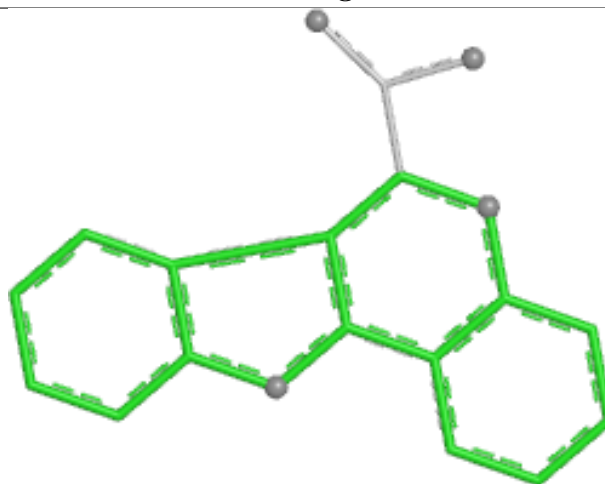
Bond lengths



Bond angles

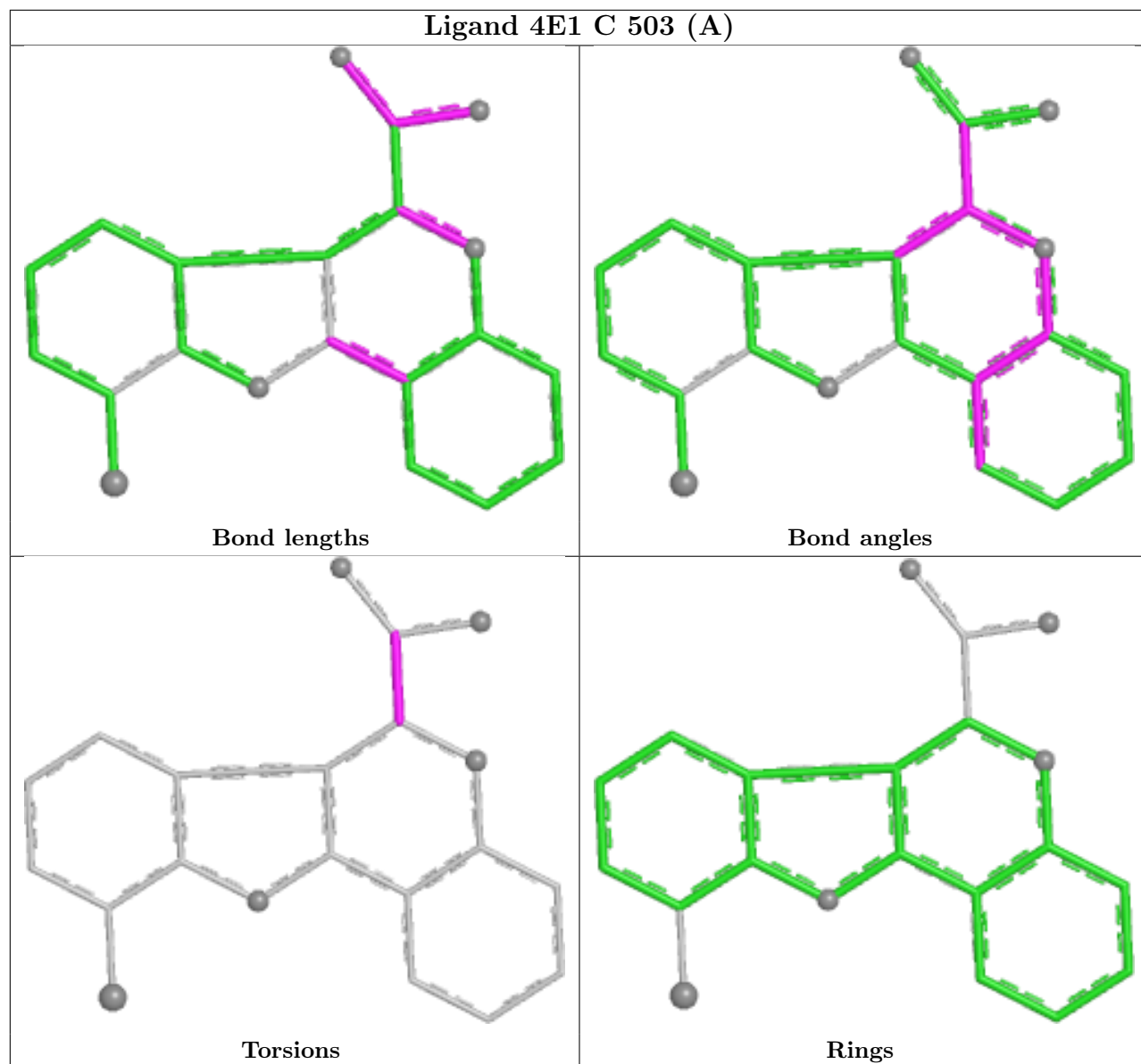


Torsions



Rings

Ligand 4E1 C 503 (A)



Ligand 4E1 A 505 (B)

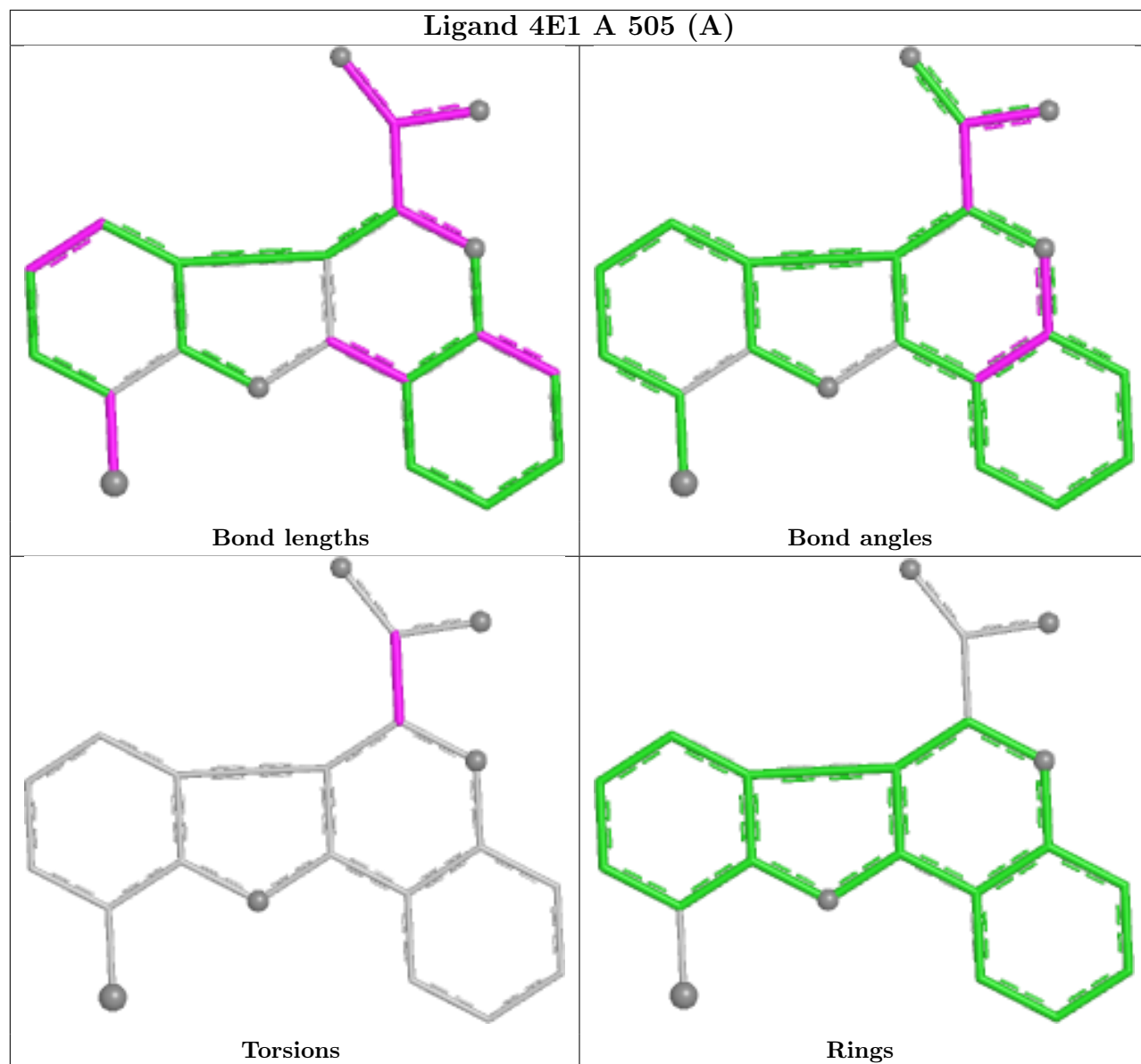
Bond lengths

Bond angles

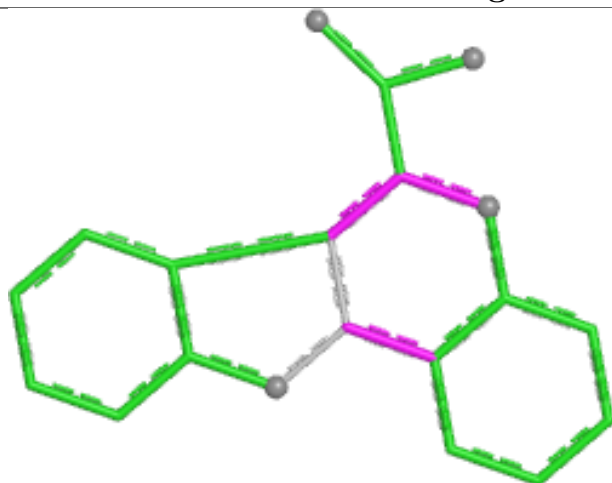
Torsions

Rings

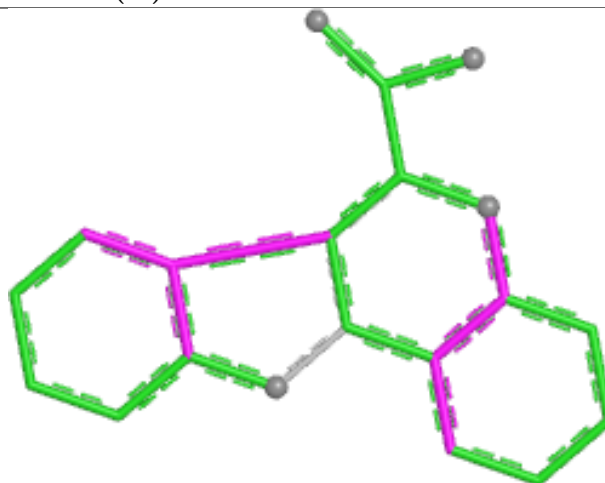
Ligand 4E1 A 505 (A)



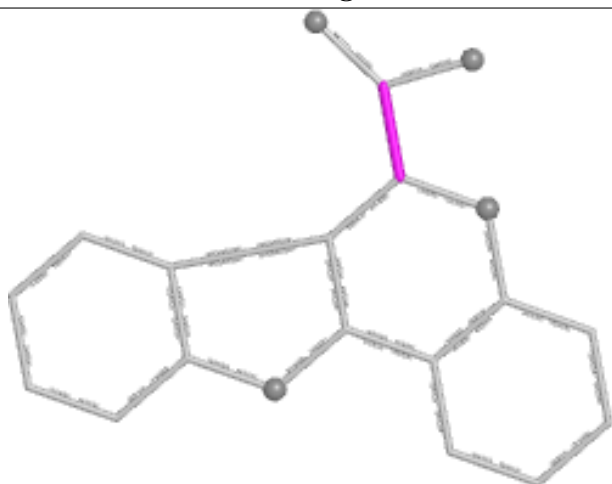
Ligand 4E1 B 508 (B)



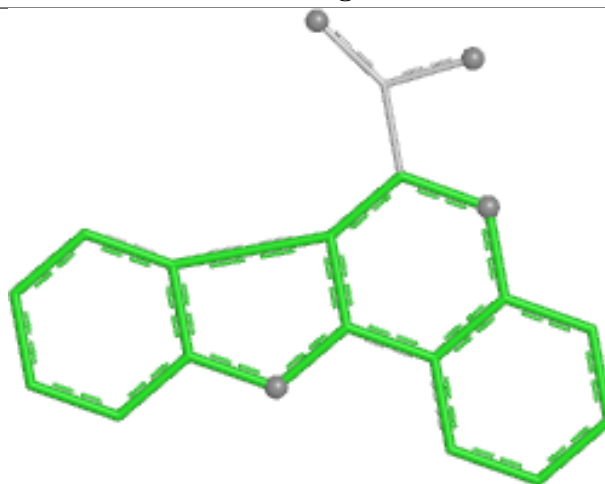
Bond lengths



Bond angles

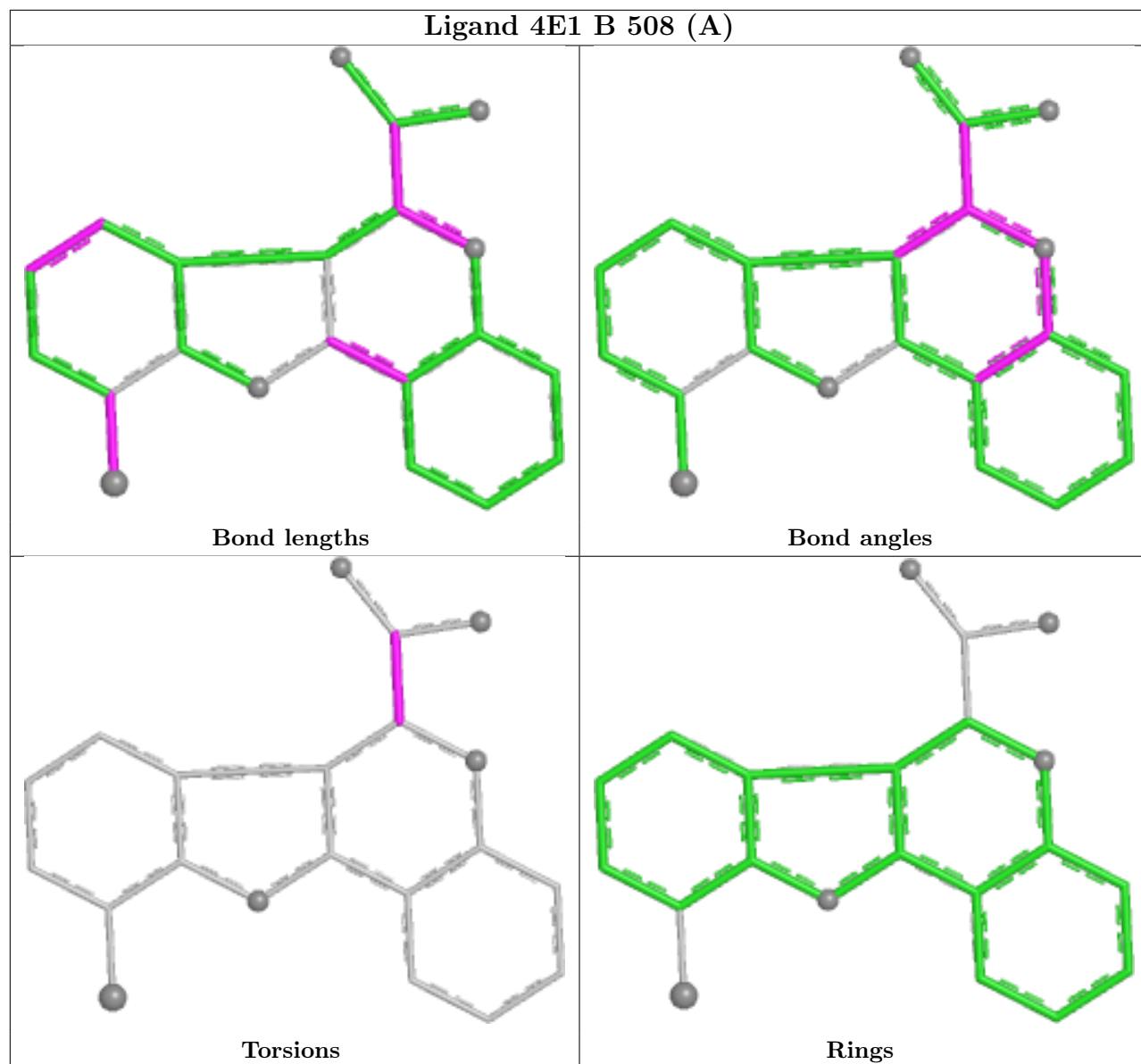


Torsions

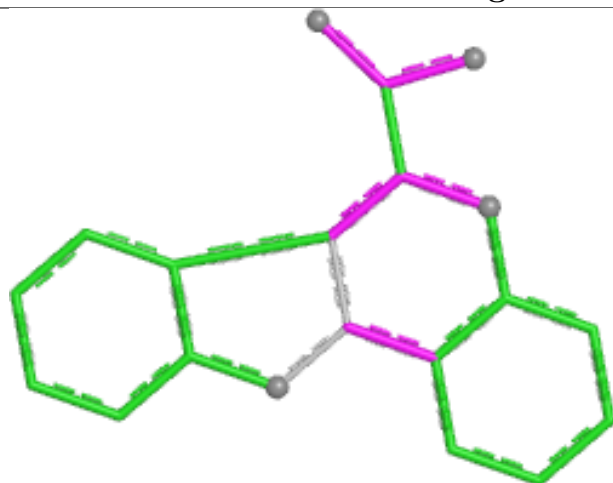


Rings

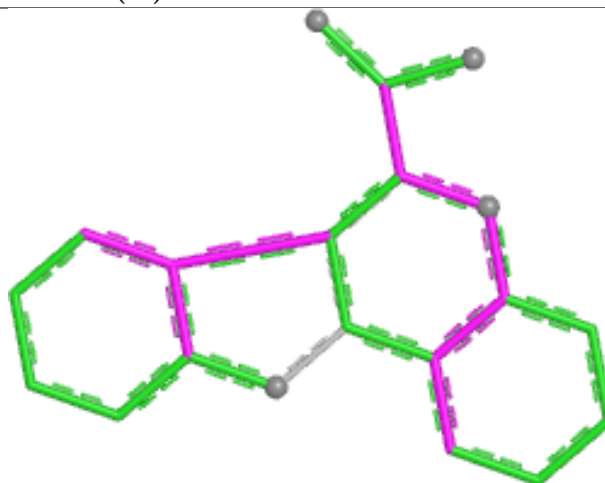
Ligand 4E1 B 508 (A)



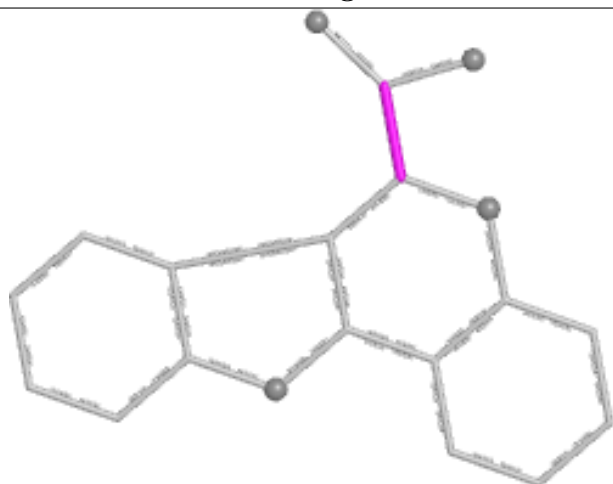
Ligand 4E1 D 505 (B)



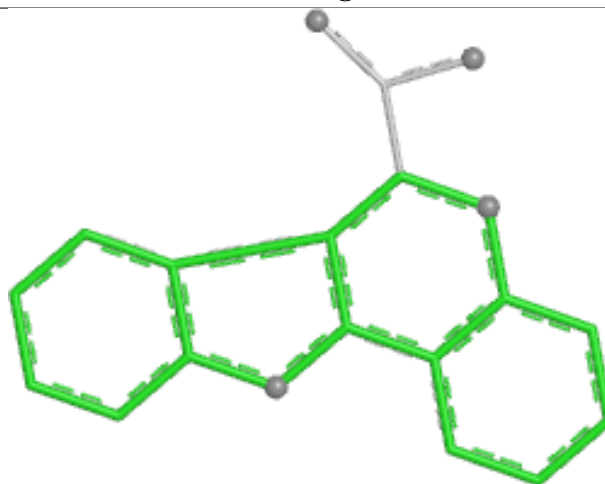
Bond lengths



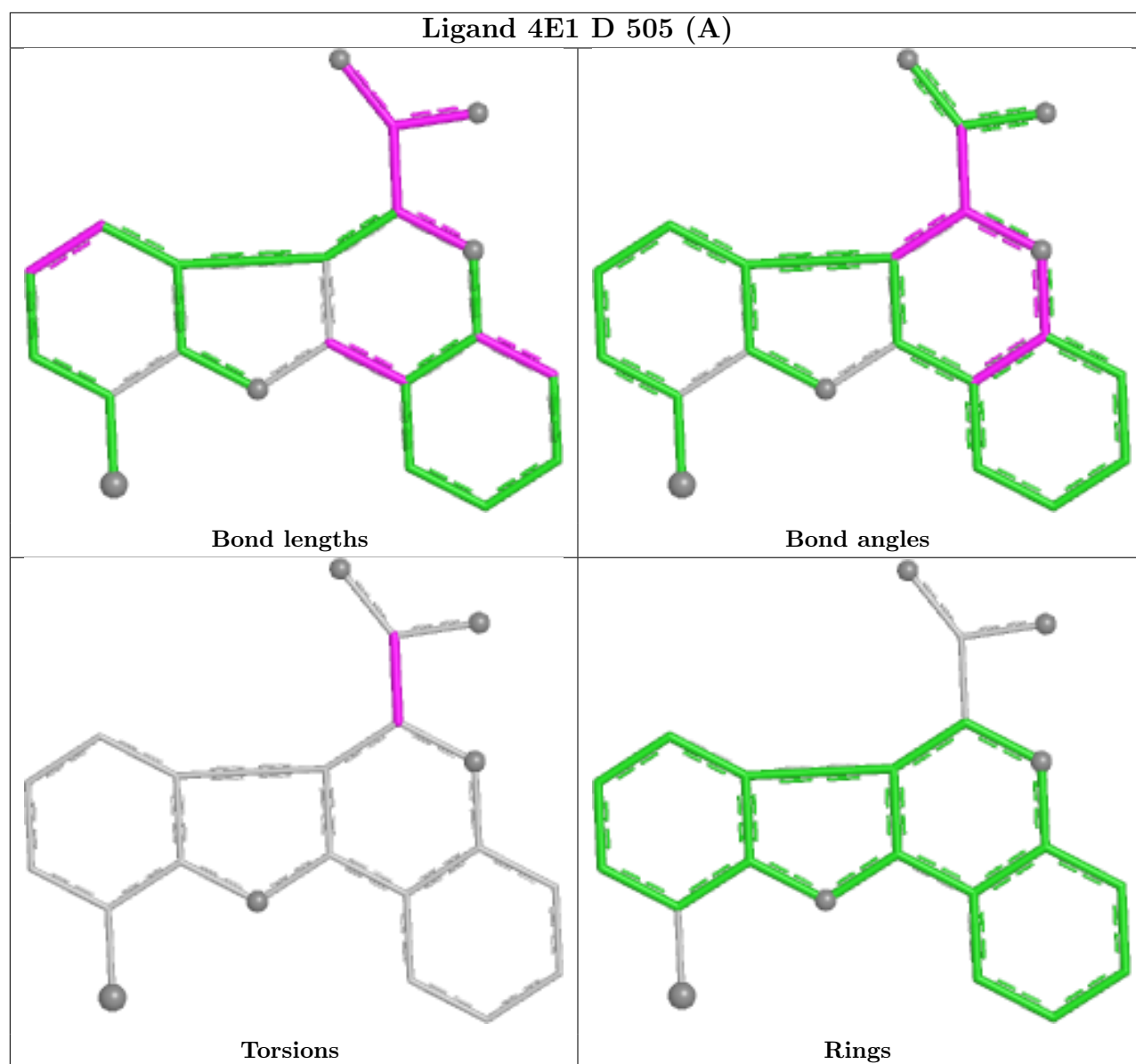
Bond angles



Torsions



Rings



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	346/361 (95%)	-0.45	4 (1%) 76 73	28, 57, 100, 161	3 (0%)
1	B	347/361 (96%)	-0.39	6 (1%) 69 65	27, 63, 111, 224	2 (0%)
1	C	346/361 (95%)	0.33	9 (2%) 57 53	53, 98, 143, 175	2 (0%)
1	D	347/361 (96%)	-0.20	7 (2%) 64 61	45, 76, 117, 151	2 (0%)
All	All	1386/1444 (95%)	-0.18	26 (1%) 66 62	27, 72, 126, 224	9 (0%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	217	MET	3.7
1	D	217	MET	3.3
1	A	219	TYR	3.3
1	D	481	LYS	3.3
1	C	481	LYS	3.2
1	C	248	LEU	3.2
1	D	134	LYS	3.0
1	C	161	ILE	2.9
1	C	410	ASP	2.9
1	B	318	ILE	2.8
1	C	147	TYR	2.6
1	D	219	TYR	2.6
1	D	410	ASP	2.4
1	A	213[A]	HIS	2.4
1	D	218	LYS	2.3
1	C	267	GLN	2.3
1	C	427	LEU	2.3
1	B	319	TYR	2.3
1	B	134	LYS	2.2
1	C	271	THR	2.2
1	B	316	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	408	THR	2.1
1	D	411	GLY	2.1
1	A	135	VAL	2.1
1	B	414	GLU	2.1
1	C	175	LYS	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	PTR	B	321[A]	16/17	0.91	0.13	62,80,87,87	13
1	PTR	B	321[B]	16/17	0.91	0.13	52,69,81,82	13
1	PTR	C	321[A]	16/17	0.91	0.11	65,80,85,86	13
1	PTR	C	321[B]	16/17	0.91	0.11	70,81,84,86	13
1	PTR	D	321[A]	16/17	0.91	0.11	62,81,87,91	13
1	PTR	D	321[B]	16/17	0.91	0.11	61,76,85,91	13
1	PTR	A	321	16/17	0.94	0.09	62,73,80,81	0

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
3	SO4	D	504	5/5	0.51	0.11	102,124,131,137	0
3	SO4	A	503	5/5	0.54	0.13	120,123,131,134	0
3	SO4	B	505	5/5	0.68	0.10	110,120,127,129	0
3	SO4	D	503	5/5	0.69	0.10	117,119,133,143	0
3	SO4	B	506	5/5	0.70	0.09	115,122,124,133	0
3	SO4	C	502	5/5	0.75	0.08	98,130,133,144	0

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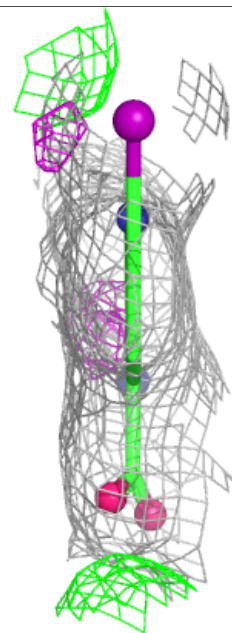
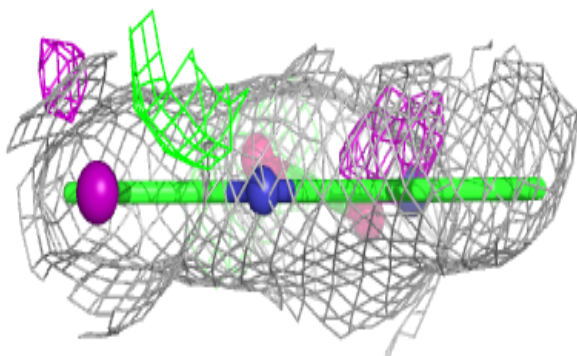
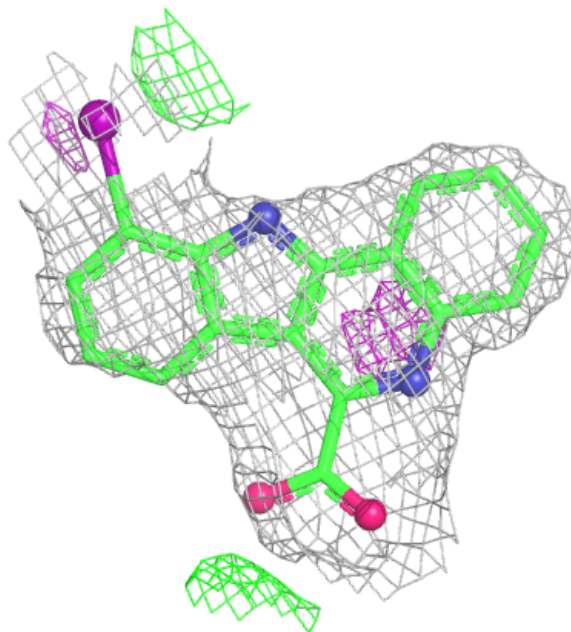
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	D	502	5/5	0.75	0.11	105,108,143,144	0
2	PG4	B	503	13/13	0.78	0.21	90,107,119,121	0
2	PG4	B	502	13/13	0.81	0.16	94,104,108,109	0
3	SO4	B	507	5/5	0.86	0.11	93,102,105,107	0
2	PG4	A	501	13/13	0.87	0.13	80,93,100,103	0
3	SO4	A	504	5/5	0.88	0.11	75,90,105,112	0
2	PG4	A	502	10/13	0.90	0.15	87,95,98,98	0
3	SO4	B	504	5/5	0.92	0.13	80,81,87,93	0
2	PG4	C	501	13/13	0.94	0.09	45,53,57,63	0
2	PG4	B	501	13/13	0.95	0.10	48,54,61,67	0
2	PG4	D	501	13/13	0.95	0.10	59,68,82,85	0
4	4E1	C	503[A]	21/21	0.97	0.11	68,77,90,92	21
4	4E1	C	503[B]	20/21	0.97	0.11	92,105,113,115	20
4	4E1	A	505[A]	21/21	0.98	0.09	49,60,74,80	21
4	4E1	A	505[B]	20/21	0.98	0.09	37,43,56,61	20
4	4E1	D	505[A]	21/21	0.98	0.09	60,70,76,82	21
4	4E1	D	505[B]	20/21	0.98	0.09	79,84,95,98	20
4	4E1	B	508[A]	21/21	0.99	0.07	54,58,70,74	21
4	4E1	B	508[B]	20/21	0.99	0.07	62,67,80,82	20

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.

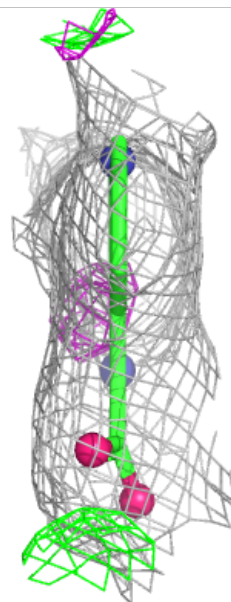
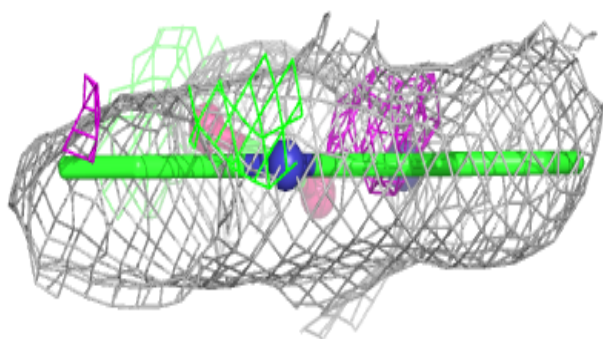
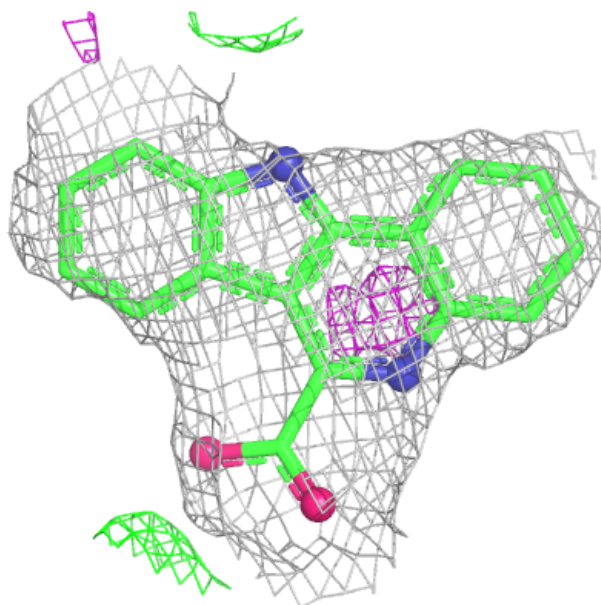
Electron density around 4E1 C 503 (A):

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



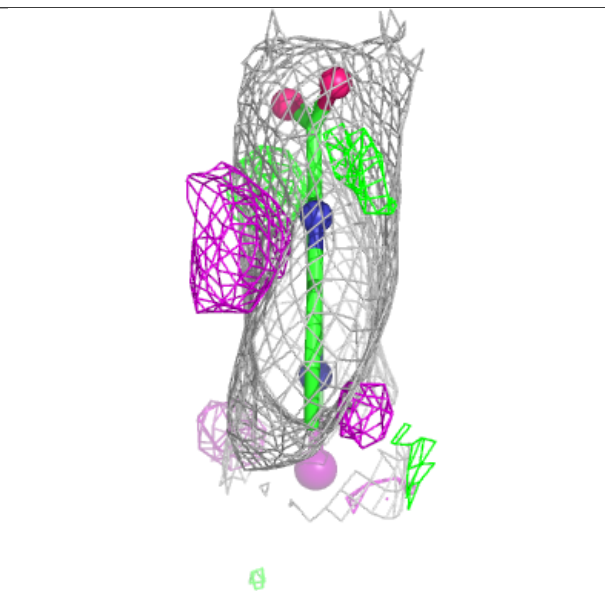
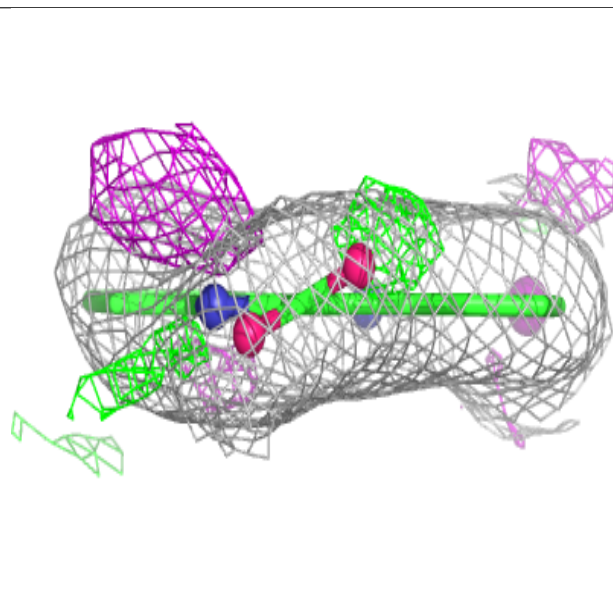
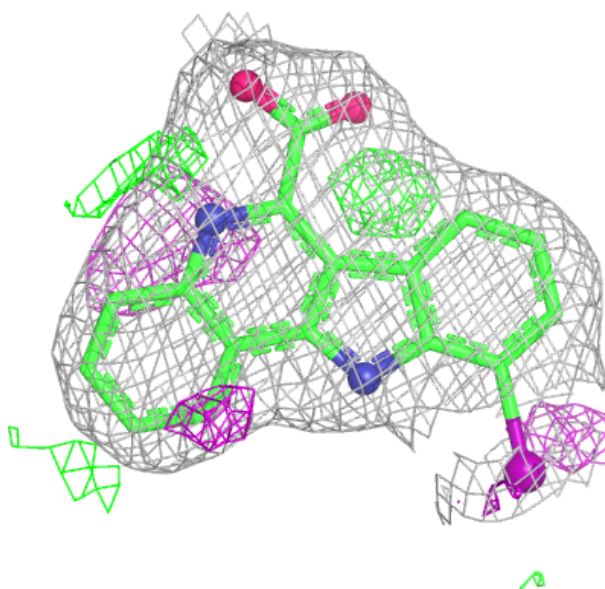
Electron density around 4E1 C 503 (B):

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



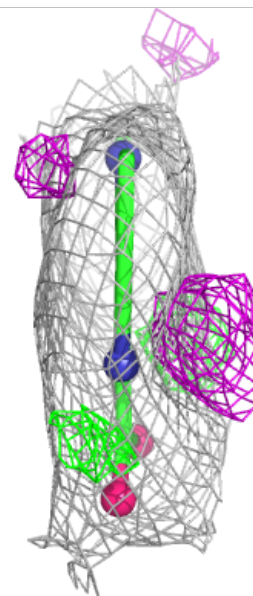
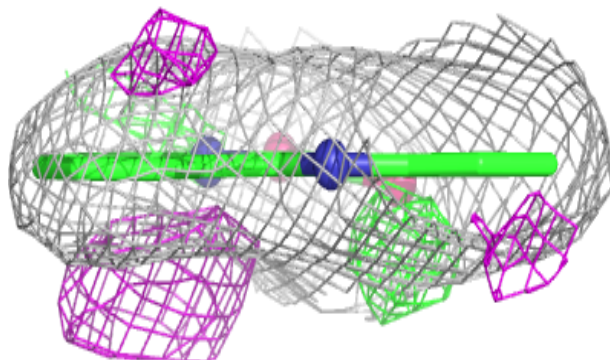
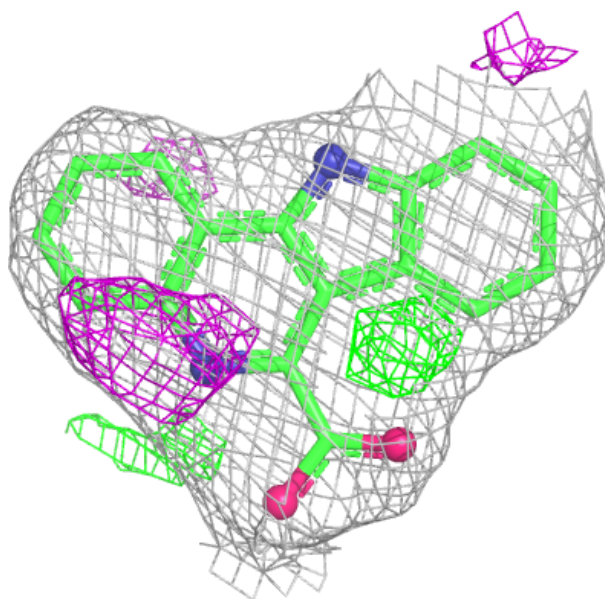
Electron density around 4E1 A 505 (A):

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and green (positive)



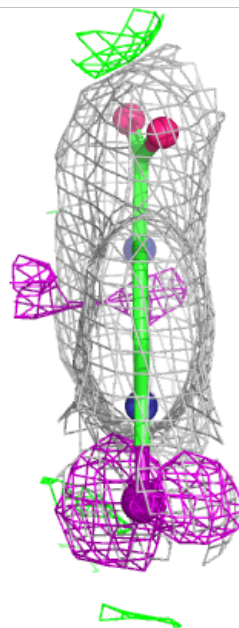
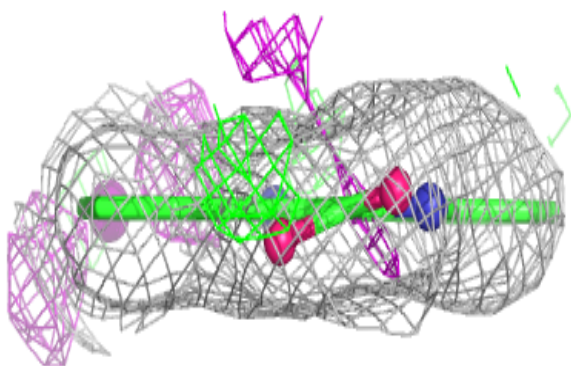
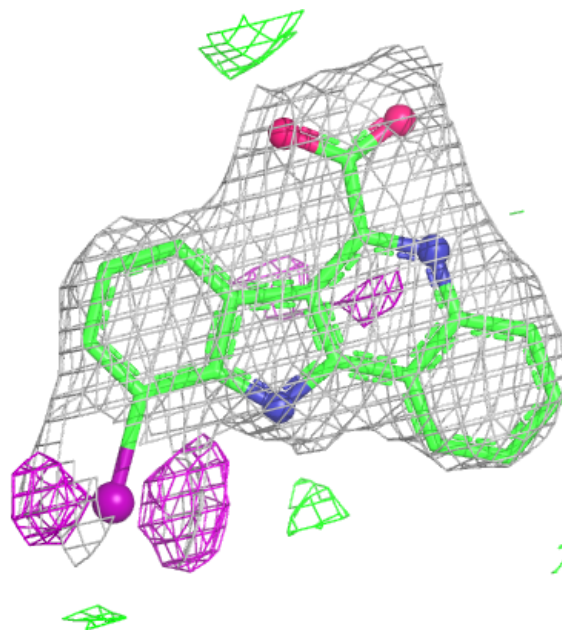
Electron density around 4E1 A 505 (B):

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



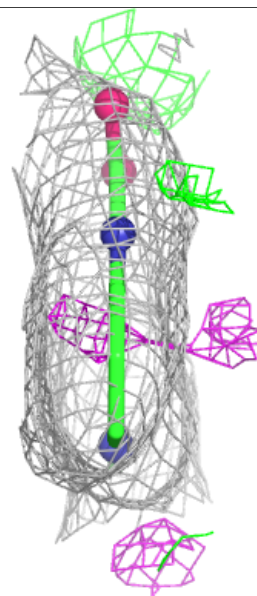
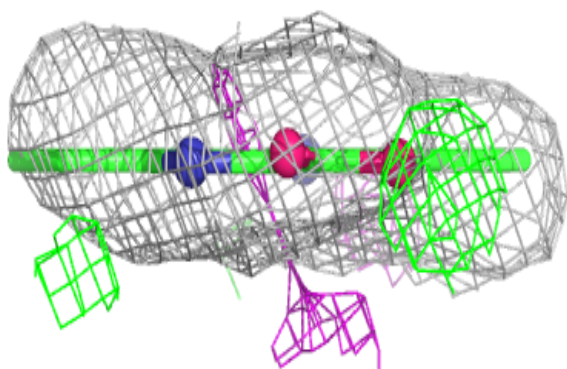
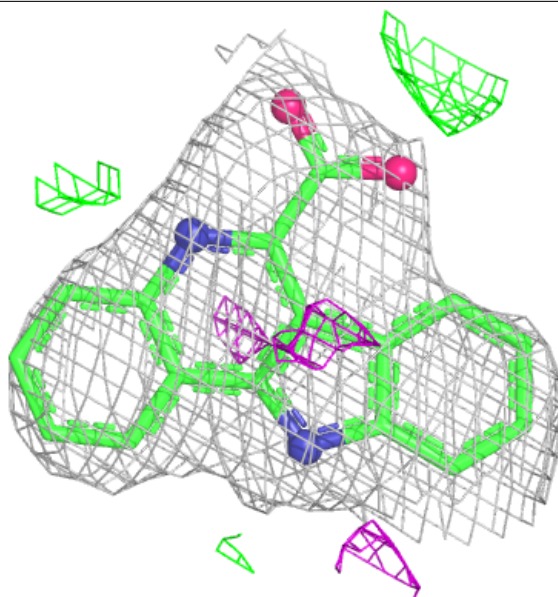
Electron density around 4E1 D 505 (A):

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and green (positive)



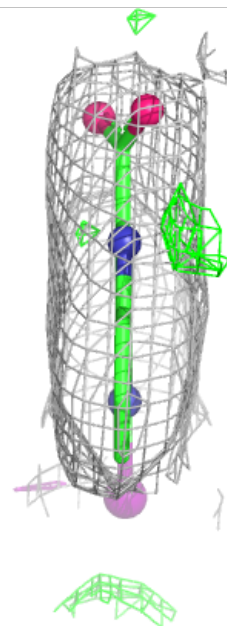
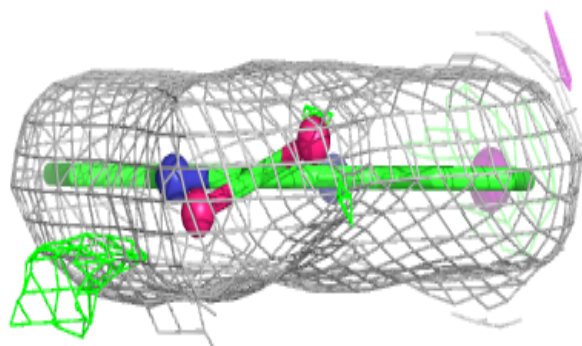
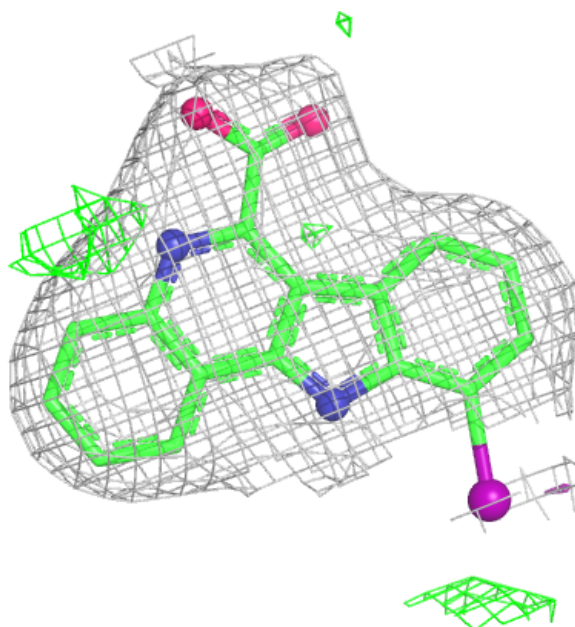
Electron density around 4E1 D 505 (B):

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and green (positive)



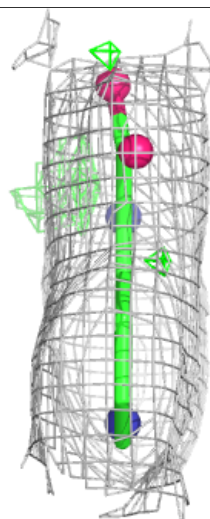
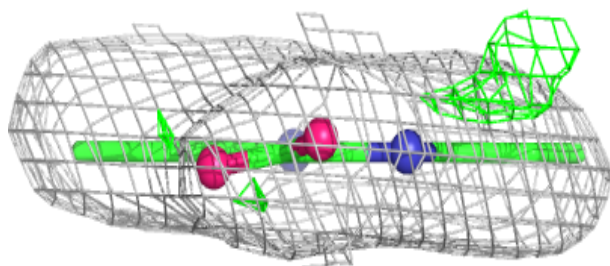
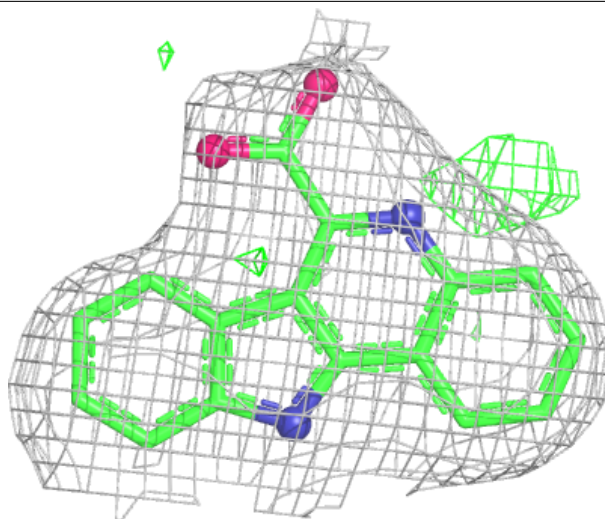
Electron density around 4E1 B 508 (A):

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



Electron density around 4E1 B 508 (B):

$2mF_o-DF_c$ (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.