



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 05:19 pm BST

PDB ID : 5VA8
Title : Crystal Structure of Short-chain dehydrogenase/reductase SDR from Burkholderia phymatum in complex with NADP
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2017-03-24
Resolution : 1.60 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

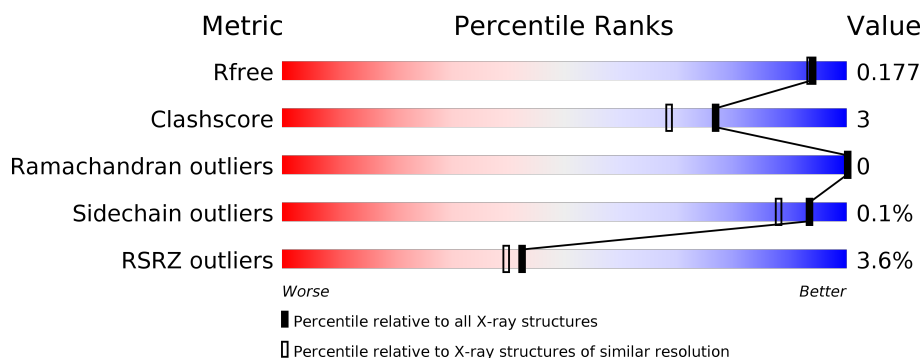
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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	267	<div> <div>4%</div> <div>91%</div> <div>6%</div> <div>.</div> </div>
1	B	267	<div> <div>3%</div> <div>90%</div> <div>8%</div> <div>.</div> </div>
1	C	267	<div> <div>3%</div> <div>93%</div> <div>.</div> <div>.</div> </div>
1	D	267	<div> <div>4%</div> <div>94%</div> <div>.</div> <div>.</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9211 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 Short-chain dehydrogenase/reductase SDR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	261	Total	C	N	O	S	0	8	0
			1906	1189	342	365	10			
1	B	261	Total	C	N	O	S	0	5	0
			1895	1182	338	365	10			
1	C	261	Total	C	N	O	S	0	9	0
			1918	1198	341	369	10			
1	D	261	Total	C	N	O	S	0	4	0
			1888	1177	337	364	10			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	expression tag	UNP B2JE32
A	-6	ALA	-	expression tag	UNP B2JE32
A	-5	HIS	-	expression tag	UNP B2JE32
A	-4	HIS	-	expression tag	UNP B2JE32
A	-3	HIS	-	expression tag	UNP B2JE32
A	-2	HIS	-	expression tag	UNP B2JE32
A	-1	HIS	-	expression tag	UNP B2JE32
A	0	HIS	-	expression tag	UNP B2JE32
B	-7	MET	-	expression tag	UNP B2JE32
B	-6	ALA	-	expression tag	UNP B2JE32
B	-5	HIS	-	expression tag	UNP B2JE32
B	-4	HIS	-	expression tag	UNP B2JE32
B	-3	HIS	-	expression tag	UNP B2JE32
B	-2	HIS	-	expression tag	UNP B2JE32
B	-1	HIS	-	expression tag	UNP B2JE32
B	0	HIS	-	expression tag	UNP B2JE32
C	-7	MET	-	expression tag	UNP B2JE32
C	-6	ALA	-	expression tag	UNP B2JE32
C	-5	HIS	-	expression tag	UNP B2JE32
C	-4	HIS	-	expression tag	UNP B2JE32
C	-3	HIS	-	expression tag	UNP B2JE32

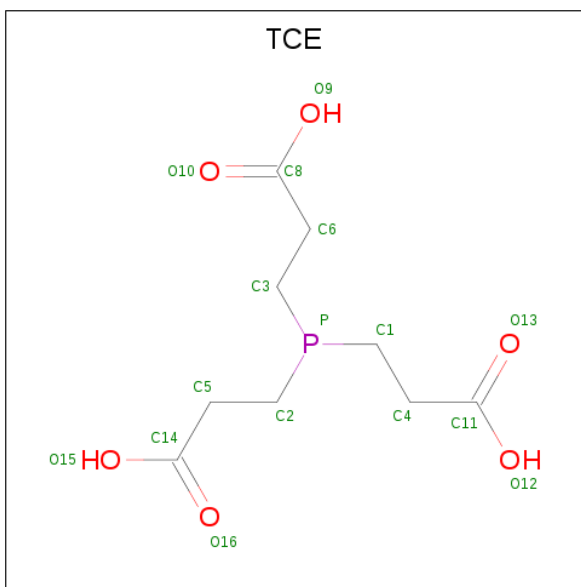
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Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	HIS	-	expression tag	UNP B2JE32
C	-1	HIS	-	expression tag	UNP B2JE32
C	0	HIS	-	expression tag	UNP B2JE32
D	-7	MET	-	expression tag	UNP B2JE32
D	-6	ALA	-	expression tag	UNP B2JE32
D	-5	HIS	-	expression tag	UNP B2JE32
D	-4	HIS	-	expression tag	UNP B2JE32
D	-3	HIS	-	expression tag	UNP B2JE32
D	-2	HIS	-	expression tag	UNP B2JE32
D	-1	HIS	-	expression tag	UNP B2JE32
D	0	HIS	-	expression tag	UNP B2JE32

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- The chemical structure of Naproxen (NAP) is shown, featuring a naphthalene ring system. The structure includes a chiral center at the 1-position, with a carboxylic acid group (-COOH) and a 6-methoxy group (-OCH₃). The stereochemistry is indicated by wedged and dashed bonds. The structure is labeled with 'NAP' and 'NAP' at the top, and 'NAP' and 'NAP' at the bottom, indicating the enantiomers.

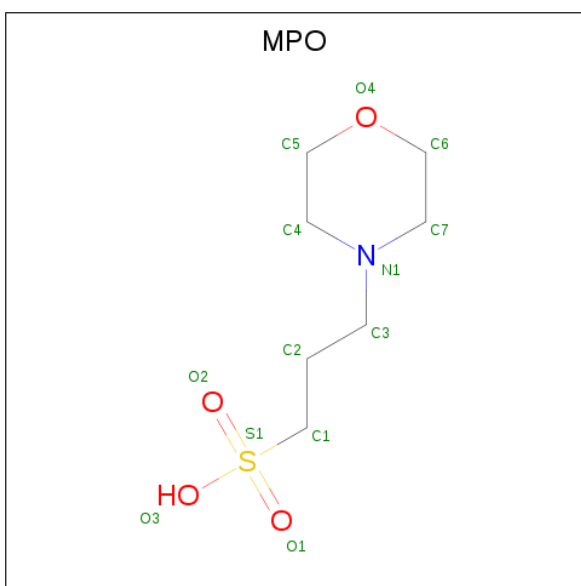
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	B	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	C	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	D	1	Total 48	C 21	N 7	O 17	P 3	0	0

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- WORLD WIDE
PDB
PROTEIN DATA BANK



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total 16	C 9	O 6	P 1	0	0
3	B	1	Total 16	C 9	O 6	P 1	0	0
3	C	1	Total 16	C 9	O 6	P 1	0	0
3	D	1	Total 16	C 9	O 6	P 1	0	0

- Molecule 4 is 3[N-MORPHOLINO]PROPANE SULFONIC ACID (three-letter code: MPO) (formula: $\text{C}_7\text{H}_{15}\text{NO}_4\text{S}$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			13	7	1	4	1		
4	C	1	Total	C	N	O	S	0	0
			13	7	1	4	1		
4	D	1	Total	C	N	O	S	0	0
			13	7	1	4	1		

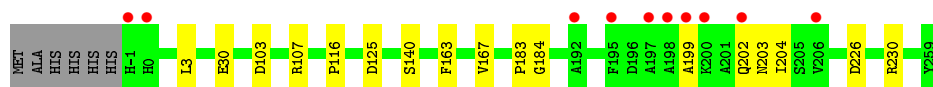
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	347	Total	O	0	3
			350	350		
5	B	293	Total	O	0	2
			295	295		
5	C	369	Total	O	0	8
			377	377		
5	D	284	Total	O	0	3
			287	287		

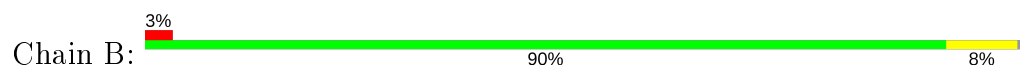
3 Residue-property plots [i](#)

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: Short-chain dehydrogenase/reductase SDR



- Molecule 1: Short-chain dehydrogenase/reductase SDR



- Molecule 1: Short-chain dehydrogenase/reductase SDR



- Molecule 1: Short-chain dehydrogenase/reductase SDR



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	133.56 Å 75.13 Å 109.88 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.35 – 1.60 44.35 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (44.35-1.60) 99.7 (44.35-1.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.46 (at 1.60 Å)	Xtriage
Refinement program	PHENIX (dev_2689: ???)	Depositor
R, R_{free}	0.149 , 0.177 0.149 , 0.177	Depositor DCC
R_{free} test set	1928 reflections (1.32%)	wwPDB-VP
Wilson B-factor (Å ²)	16.4	Xtriage
Anisotropy	0.538	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 52.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9211	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 52.62 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.7393e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MPO, NAP, TCE

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	0.34	0/1961	0.55	0/2672
1	B	0.32	0/1941	0.53	0/2645
1	C	0.34	0/1976	0.54	0/2692
1	D	0.32	0/1932	0.52	0/2633
All	All	0.33	0/7810	0.54	0/10642

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	1906	0	1895	13	0
1	B	1895	0	1887	15	0
1	C	1918	0	1919	10	0
1	D	1888	0	1869	6	0
2	A	48	0	25	2	0
2	B	48	0	25	3	0
2	C	48	0	25	2	0
2	D	48	0	25	2	0
3	A	16	0	12	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	16	0	12	1	0
3	C	16	0	12	1	0
3	D	16	0	12	1	0
4	A	13	0	15	0	0
4	C	13	0	15	0	0
4	D	13	0	15	0	0
5	A	350	0	0	5	4
5	B	295	0	0	4	0
5	C	377	0	0	6	3
5	D	287	0	0	1	0
All	All	9211	0	7763	47	4

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 3.

All (47) 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:B:103[A]:ASP:OD2	1:B:107:ARG:NH1	2.05	0.90
1:C:103[B]:ASP:OD1	5:C:401:HOH:O	1.95	0.85
1:A:125:ASP:OD1	5:A:401:HOH:O	2.00	0.79
1:A:103[B]:ASP:OD2	5:A:402:HOH:O	2.02	0.78
1:A:103[A]:ASP:OD2	1:A:107[A]:ARG:NH1	2.17	0.78
1:C:125:ASP:OD1	5:C:402:HOH:O	2.02	0.77
1:B:259:TYR:O	5:B:401:HOH:O	2.04	0.76
2:C:301:NAP:C4N	3:C:302:TCE:P	2.76	0.74
2:B:301:NAP:C4N	3:B:302:TCE:P	2.78	0.71
2:A:301:NAP:C4N	3:A:302:TCE:P	2.78	0.71
1:C:203:ASN:OD1	5:C:403:HOH:O	2.08	0.71
1:B:125:ASP:OD1	5:B:402:HOH:O	2.09	0.70
1:A:202:GLN:O	5:A:403:HOH:O	2.09	0.70
1:C:202:GLN:O	5:C:404:HOH:O	2.11	0.68
1:B:202:GLN:NE2	5:B:404:HOH:O	2.22	0.67
2:D:301:NAP:C4N	3:D:302:TCE:P	2.85	0.65
1:A:3:LEU:HD12	1:A:30:GLU:HG3	1.80	0.61
1:D:3:LEU:HD12	1:D:30:GLU:HG3	1.83	0.60
1:C:3:LEU:HD12	1:C:30:GLU:HG3	1.84	0.58
1:A:203:ASN:OD1	5:A:404:HOH:O	2.18	0.57
1:C:116:PRO:HG3	5:C:633[B]:HOH:O	2.05	0.56
1:A:116:PRO:HG3	5:A:607[B]:HOH:O	2.05	0.56
1:B:3:LEU:HD12	1:B:30:GLU:HG3	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:140:SER:HA	1:C:183:PRO:HD2	1.90	0.54
1:A:184:GLY:O	2:A:301:NAP:H4N	2.09	0.53
1:A:226:ASP:O	1:A:230[B]:ARG:HG2	2.11	0.51
1:C:184:GLY:O	2:C:301:NAP:H4N	2.11	0.51
1:D:184:GLY:O	2:D:301:NAP:H4N	2.10	0.50
1:A:140:SER:HA	1:A:183:PRO:HD2	1.94	0.49
1:D:226:ASP:O	1:D:230:ARG:HG2	2.13	0.47
1:B:163:PHE:CZ	1:B:167[B]:VAL:HG21	2.49	0.47
1:B:226:ASP:O	1:B:230:ARG:HG2	2.14	0.46
1:B:183:PRO:HG2	2:B:301:NAP:C5N	2.46	0.46
1:A:199:ALA:HB1	1:A:204:ILE:O	2.16	0.45
1:D:187:ASP:HA	1:D:191:ILE:HD11	1.98	0.44
1:B:103[B]:ASP:OD1	5:B:403:HOH:O	2.20	0.43
1:B:184:GLY:O	2:B:301:NAP:H4N	2.17	0.43
1:C:21:ARG:NH2	5:C:406:HOH:O	2.37	0.43
1:D:38:ALA:O	1:D:63:ALA:HA	2.18	0.43
1:A:163:PHE:CZ	1:A:167[B]:VAL:HG21	2.54	0.43
1:B:97:PHE:HA	1:B:100:PHE:CD2	2.54	0.42
1:B:199:ALA:HB1	1:B:204:ILE:O	2.20	0.42
1:A:226:ASP:O	1:A:230[A]:ARG:HG2	2.18	0.42
1:B:38:ALA:O	1:B:63:ALA:HA	2.19	0.42
1:B:102:HIS:CE1	1:D:110:GLU:OE1	2.74	0.41
1:B:98:ARG:HG2	5:D:446:HOH:O	2.20	0.41
1:C:163:PHE:CZ	1:C:167[B]:VAL:HG21	2.56	0.41

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:608:HOH:O	5:C:451:HOH:O[2_455]	2.08	0.12
5:A:633:HOH:O	5:A:671:HOH:O[2_455]	2.10	0.10
5:A:682:HOH:O	5:C:412:HOH:O[2_455]	2.18	0.02
5:A:468:HOH:O	5:C:430:HOH:O[2_455]	2.18	0.02

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	267/267 (100%)	259 (97%)	8 (3%)	0	100	100
1	B	264/267 (99%)	256 (97%)	8 (3%)	0	100	100
1	C	268/267 (100%)	261 (97%)	7 (3%)	0	100	100
1	D	263/267 (98%)	255 (97%)	8 (3%)	0	100	100
All	All	1062/1068 (99%)	1031 (97%)	31 (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	190/195 (97%)	190 (100%)	0	100	100
1	B	190/195 (97%)	190 (100%)	0	100	100
1	C	194/195 (100%)	193 (100%)	1 (0%)	88	80
1	D	188/195 (96%)	188 (100%)	0	100	100
All	All	762/780 (98%)	761 (100%)	1 (0%)	93	88

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

Mol	Chain	Res	Type
1	C	183	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

11 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	MPO	C	303	-	13,13,13	1.49	3 (23%)	17,17,17	2.27	6 (35%)
2	NAP	C	301	-	45,52,52	0.56	0	56,80,80	0.82	2 (3%)
3	TCE	B	302	-	6,15,15	0.62	0	9,18,18	3.66	7 (77%)
3	TCE	A	302	-	6,15,15	0.53	0	9,18,18	4.31	6 (66%)
3	TCE	D	302	-	6,15,15	0.75	0	9,18,18	3.68	6 (66%)
3	TCE	C	302	-	6,15,15	0.67	0	9,18,18	3.68	6 (66%)
2	NAP	B	301	-	45,52,52	0.59	0	56,80,80	0.75	1 (1%)
4	MPO	D	303	-	13,13,13	1.45	3 (23%)	17,17,17	1.87	5 (29%)
2	NAP	D	301	-	45,52,52	0.62	0	56,80,80	0.77	1 (1%)
4	MPO	A	303	-	13,13,13	1.51	3 (23%)	17,17,17	1.98	4 (23%)
2	NAP	A	301	-	45,52,52	0.59	0	56,80,80	0.78	1 (1%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MPO	C	303	-	-	4/7/15/15	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAP	C	301	-	-	7/31/67/67	0/5/5/5
3	TCE	B	302	-	-	1/9/15/15	-
3	TCE	A	302	-	-	0/9/15/15	-
3	TCE	D	302	-	-	0/9/15/15	-
3	TCE	C	302	-	-	0/9/15/15	-
2	NAP	B	301	-	-	4/31/67/67	0/5/5/5
4	MPO	D	303	-	-	0/7/15/15	0/1/1/1
2	NAP	D	301	-	-	6/31/67/67	0/5/5/5
4	MPO	A	303	-	-	1/7/15/15	0/1/1/1
2	NAP	A	301	-	-	7/31/67/67	0/5/5/5

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	303	MPO	C1-S1	3.65	1.82	1.77
4	C	303	MPO	C1-S1	3.49	1.82	1.77
4	D	303	MPO	C1-S1	3.24	1.82	1.77
4	D	303	MPO	O2-S1	2.23	1.51	1.45
4	A	303	MPO	O1-S1	2.23	1.51	1.45
4	A	303	MPO	O2-S1	2.22	1.51	1.45
4	C	303	MPO	O1-S1	2.21	1.51	1.45
4	D	303	MPO	O1-S1	2.20	1.51	1.45
4	C	303	MPO	O2-S1	2.19	1.51	1.45

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	302	TCE	P-C3-C6	-9.25	102.81	113.67
3	C	302	TCE	P-C3-C6	-6.05	106.57	113.67
3	D	302	TCE	P-C2-C5	-5.51	107.21	113.67
3	A	302	TCE	P-C1-C4	-5.33	107.42	113.67
3	B	302	TCE	P-C2-C5	-5.19	107.58	113.67
3	D	302	TCE	P-C3-C6	-4.89	107.93	113.67
3	B	302	TCE	P-C1-C4	-4.80	108.04	113.67
3	B	302	TCE	P-C3-C6	-4.80	108.04	113.67
3	C	302	TCE	P-C1-C4	-4.75	108.10	113.67
3	D	302	TCE	P-C1-C4	-4.68	108.18	113.67
3	C	302	TCE	P-C2-C5	-4.66	108.20	113.67
4	C	303	MPO	O2-S1-C1	4.65	112.52	106.92
3	D	302	TCE	C1-P-C3	4.14	114.12	100.95
4	C	303	MPO	O3-S1-O1	-4.03	101.42	111.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	302	TCE	P-C2-C5	-3.93	109.06	113.67
4	A	303	MPO	O2-S1-C1	3.85	111.55	106.92
4	A	303	MPO	O3-S1-C1	3.83	111.96	105.77
3	B	302	TCE	C3-P-C2	3.81	113.09	100.95
3	B	302	TCE	C1-P-C3	3.77	112.95	100.95
4	A	303	MPO	O2-S1-O1	-3.60	101.48	113.95
3	C	302	TCE	C3-P-C2	3.50	112.11	100.95
3	A	302	TCE	C1-P-C3	3.47	112.02	100.95
4	D	303	MPO	O2-S1-O1	-3.47	101.94	113.95
4	D	303	MPO	O2-S1-C1	3.46	111.08	106.92
3	A	302	TCE	C3-P-C2	3.36	111.65	100.95
4	C	303	MPO	C6-C7-N1	3.33	115.15	110.10
3	D	302	TCE	C1-P-C2	3.32	111.53	100.95
4	A	303	MPO	O1-S1-C1	3.29	110.88	106.92
3	C	302	TCE	C1-P-C3	3.26	111.34	100.95
4	D	303	MPO	O3-S1-C1	3.26	111.04	105.77
4	C	303	MPO	O3-S1-C1	3.21	110.95	105.77
3	D	302	TCE	C3-P-C2	3.17	111.06	100.95
3	C	302	TCE	C1-P-C2	3.07	110.74	100.95
3	B	302	TCE	C1-P-C2	3.02	110.56	100.95
4	D	303	MPO	O1-S1-C1	2.98	110.51	106.92
4	C	303	MPO	O1-S1-C1	2.91	110.42	106.92
3	A	302	TCE	C1-P-C2	2.49	108.89	100.95
2	D	301	NAP	C5A-C6A-N6A	2.37	123.96	120.35
4	C	303	MPO	C7-N1-C4	2.27	113.94	108.83
2	A	301	NAP	C5A-C6A-N6A	2.25	123.78	120.35
2	C	301	NAP	C5A-C6A-N6A	2.24	123.76	120.35
4	D	303	MPO	C2-C1-S1	-2.24	109.81	113.25
2	C	301	NAP	C2N-C3N-C7N	2.19	125.81	119.46
3	B	302	TCE	C3-C6-C8	-2.17	106.22	112.65
2	B	301	NAP	C5A-C6A-N6A	2.15	123.62	120.35

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	303	MPO	C1-C2-C3-N1
2	C	301	NAP	C5D-O5D-PN-O1N
2	C	301	NAP	O4D-C1D-N1N-C2N
2	C	301	NAP	O4D-C1D-N1N-C6N
2	C	301	NAP	C2D-C1D-N1N-C6N
2	B	301	NAP	C5D-O5D-PN-O1N

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Mol	Chain	Res	Type	Atoms
2	B	301	NAP	O4D-C1D-N1N-C2N
2	D	301	NAP	O4D-C1D-N1N-C2N
2	D	301	NAP	O4D-C1D-N1N-C6N
2	A	301	NAP	C5D-O5D-PN-O1N
2	A	301	NAP	O4D-C1D-N1N-C2N
2	A	301	NAP	O4D-C1D-N1N-C6N
4	C	303	MPO	C2-C3-N1-C7
4	C	303	MPO	C2-C3-N1-C4
4	C	303	MPO	C2-C1-S1-O1
4	A	303	MPO	C1-C2-C3-N1
2	C	301	NAP	O4B-C4B-C5B-O5B
2	B	301	NAP	C5D-O5D-PN-O3
2	A	301	NAP	C5D-O5D-PN-O3
2	D	301	NAP	C5D-O5D-PN-O1N
3	B	302	TCE	C5-C2-P-C3
2	C	301	NAP	C5D-O5D-PN-O3
2	D	301	NAP	C5D-O5D-PN-O3
2	D	301	NAP	C2D-C1D-N1N-C6N
2	A	301	NAP	C2D-C1D-N1N-C6N
2	C	301	NAP	C3B-C4B-C5B-O5B
2	B	301	NAP	O4B-C4B-C5B-O5B
2	A	301	NAP	O4B-C4B-C5B-O5B
2	A	301	NAP	C5D-O5D-PN-O2N
2	D	301	NAP	O4B-C4B-C5B-O5B

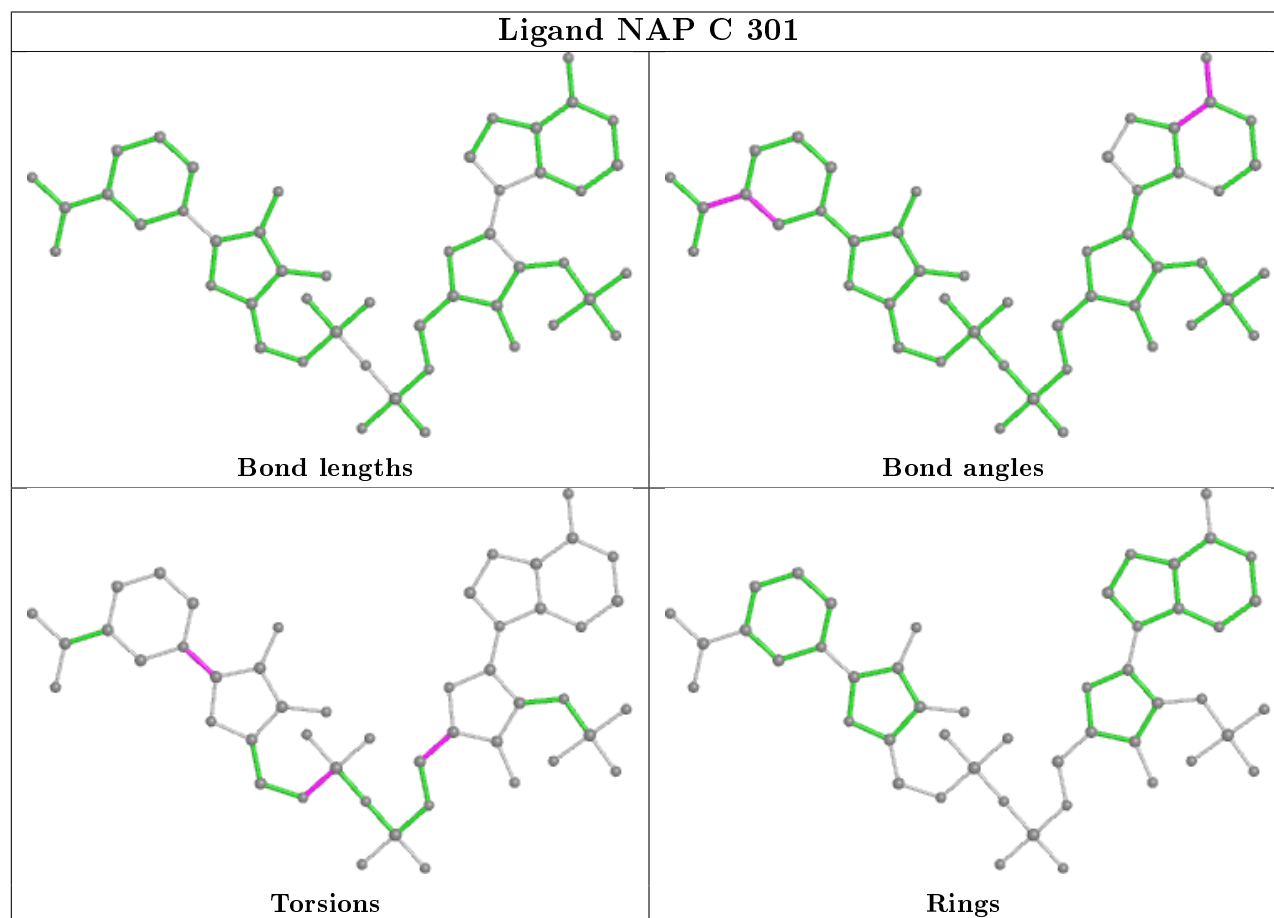
There are no ring outliers.

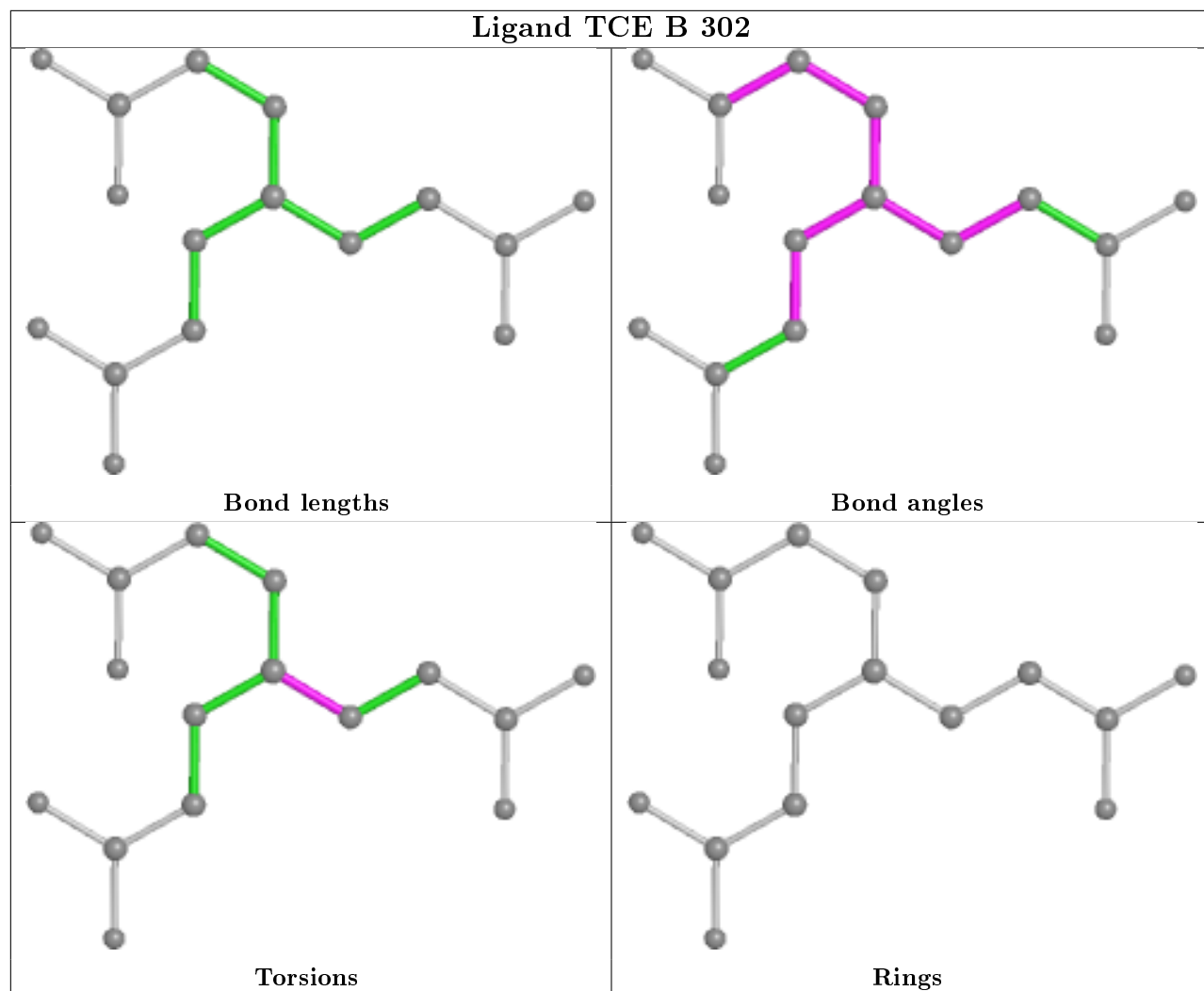
8 monomers are involved in 9 short contacts:

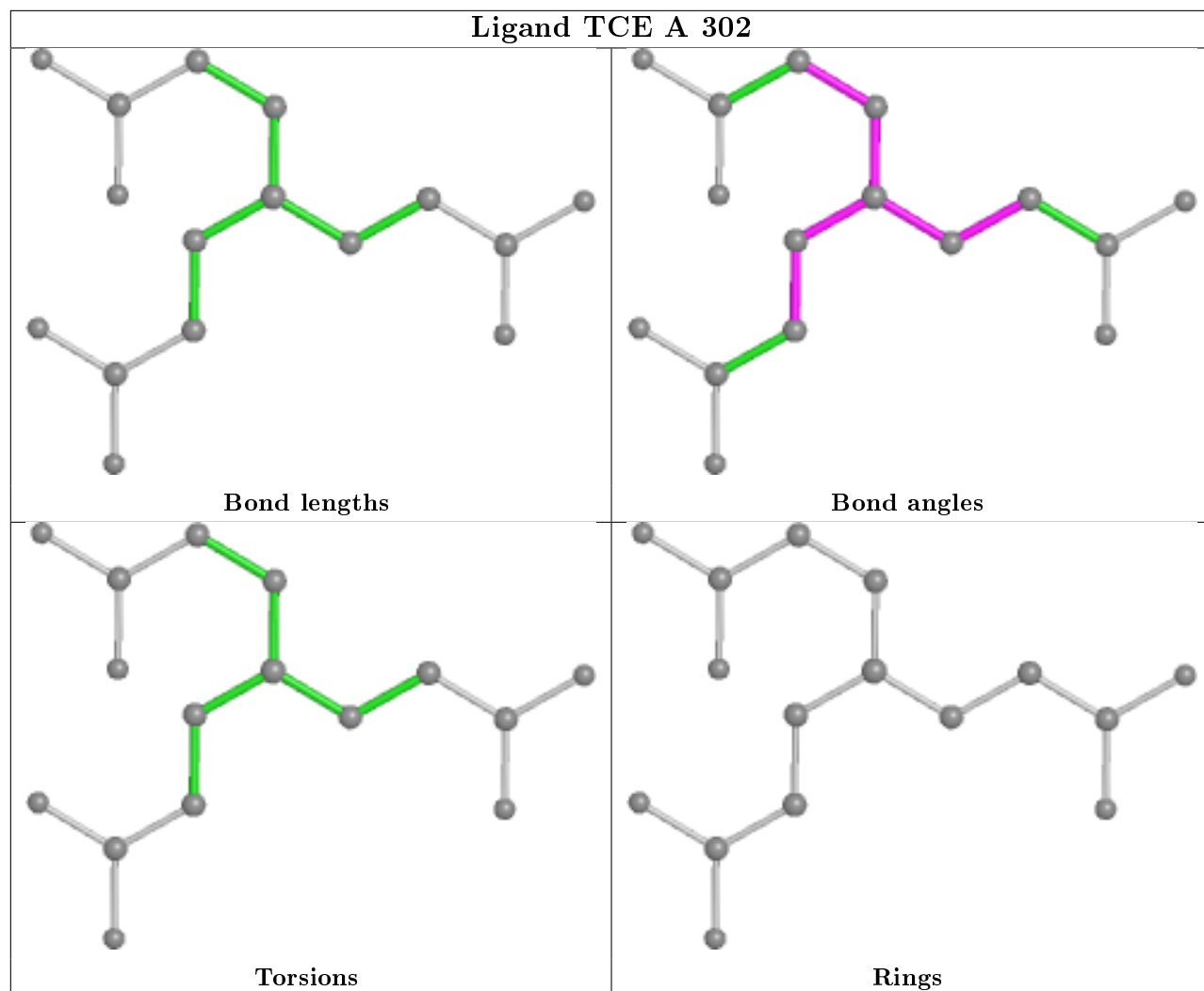
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	301	NAP	2	0
3	B	302	TCE	1	0
3	A	302	TCE	1	0
3	D	302	TCE	1	0
3	C	302	TCE	1	0
2	B	301	NAP	3	0
2	D	301	NAP	2	0
2	A	301	NAP	2	0

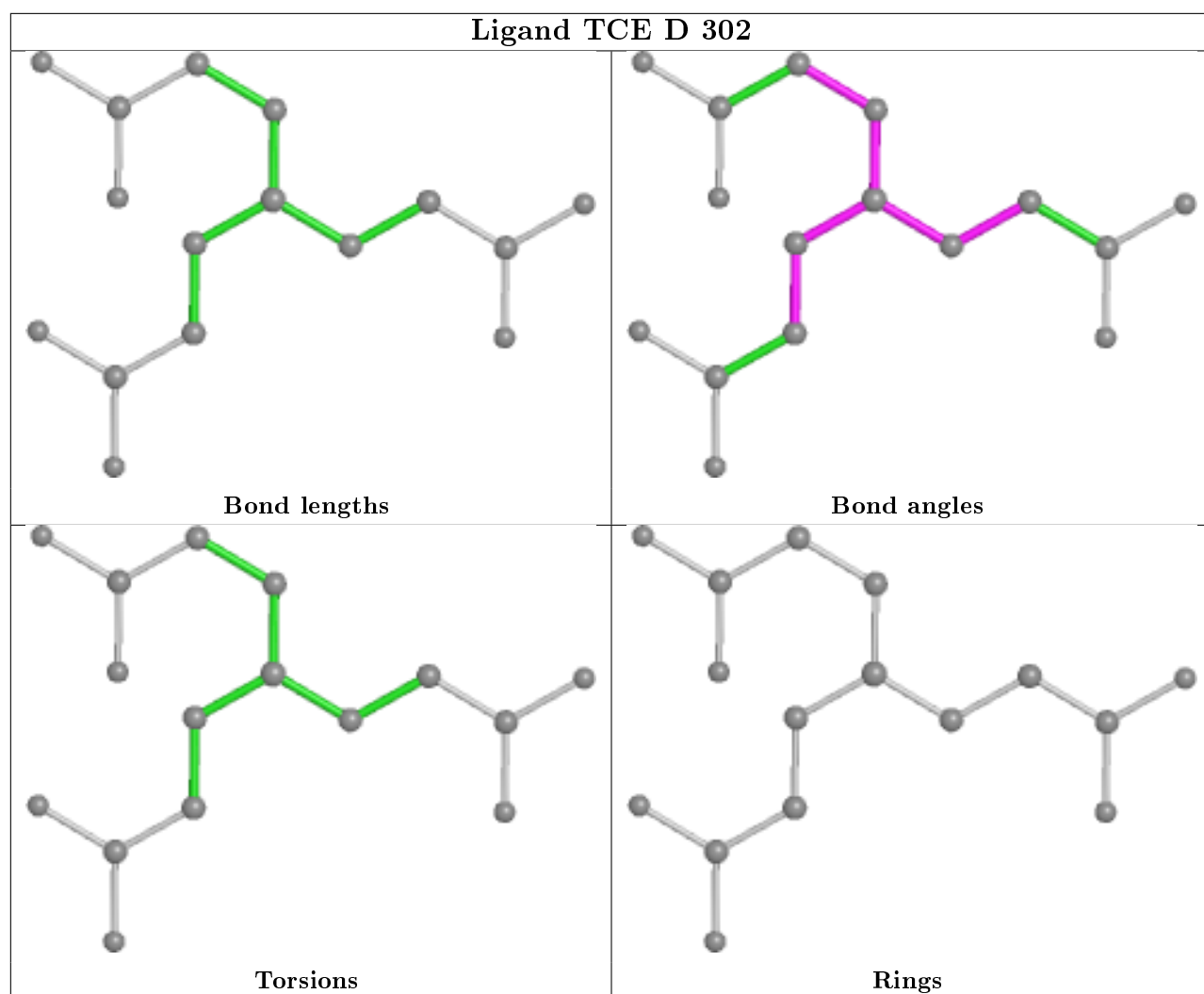
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is

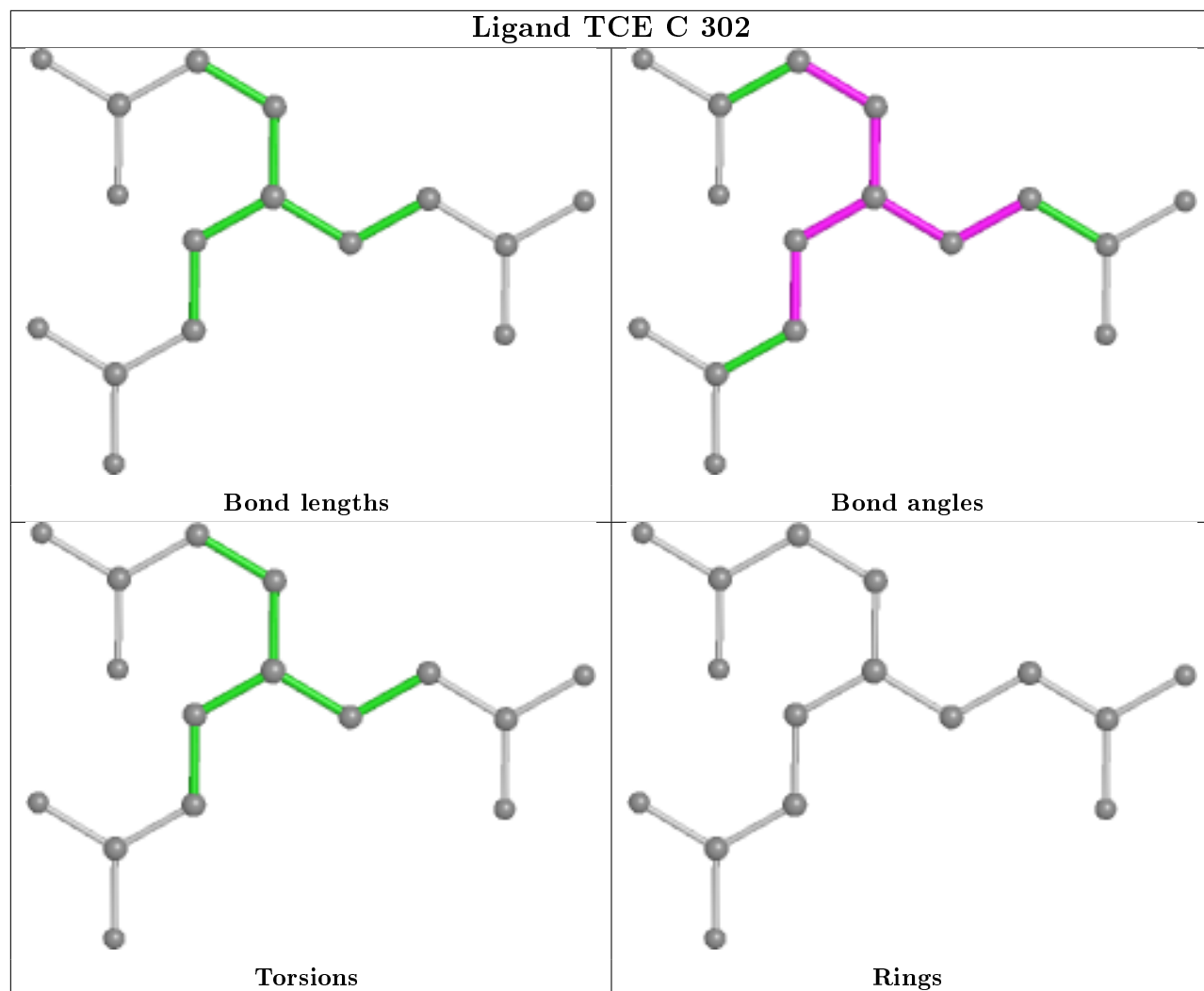
within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question 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.

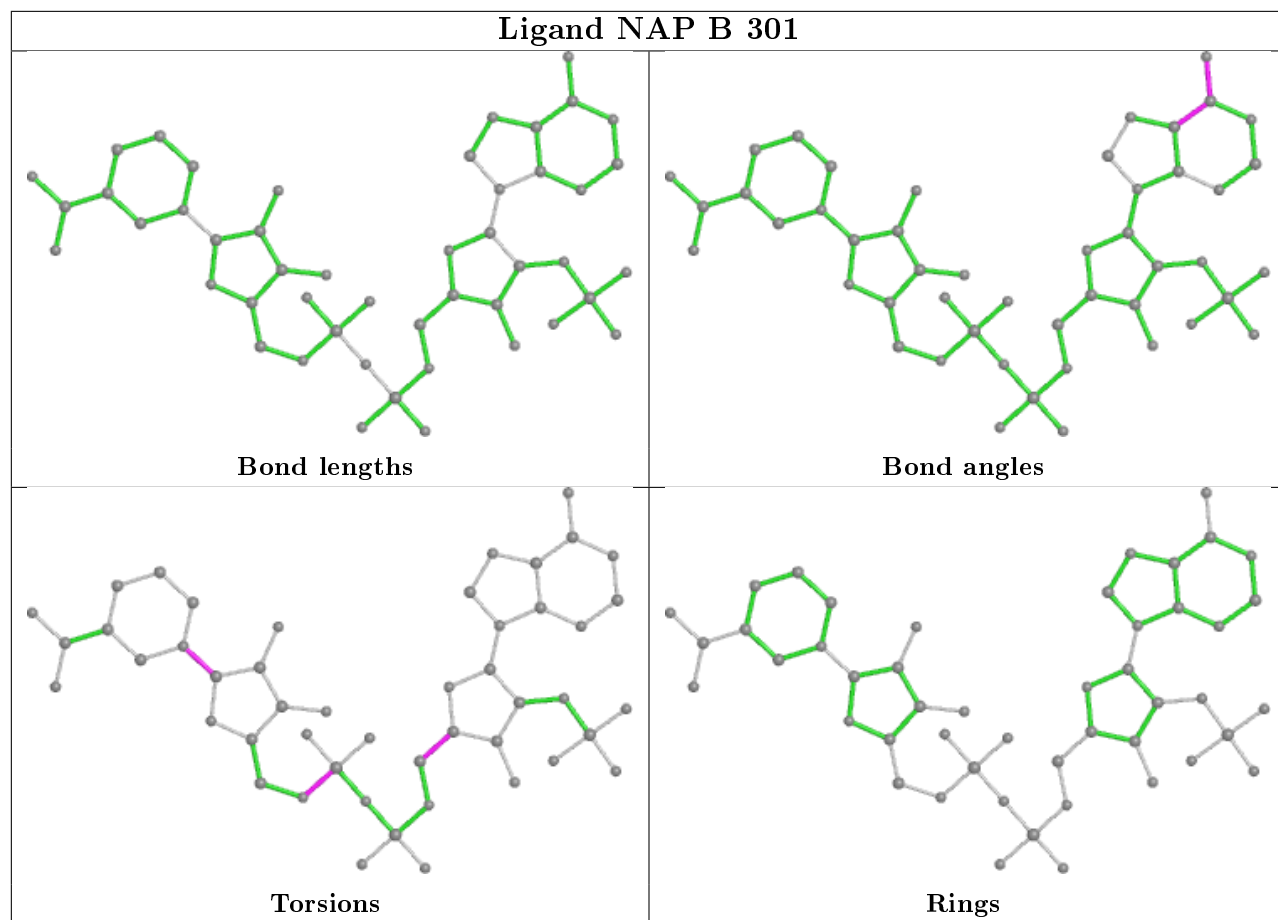


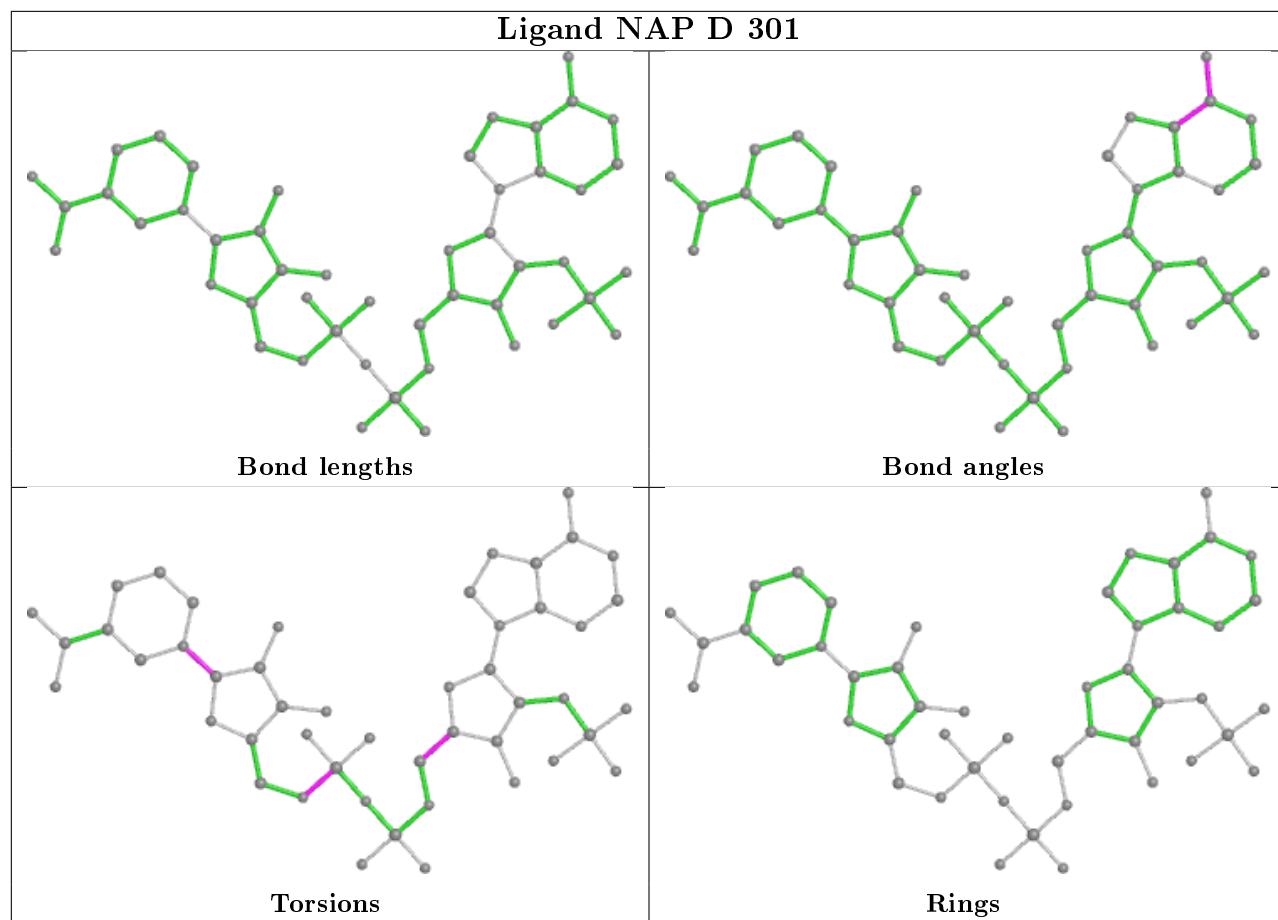


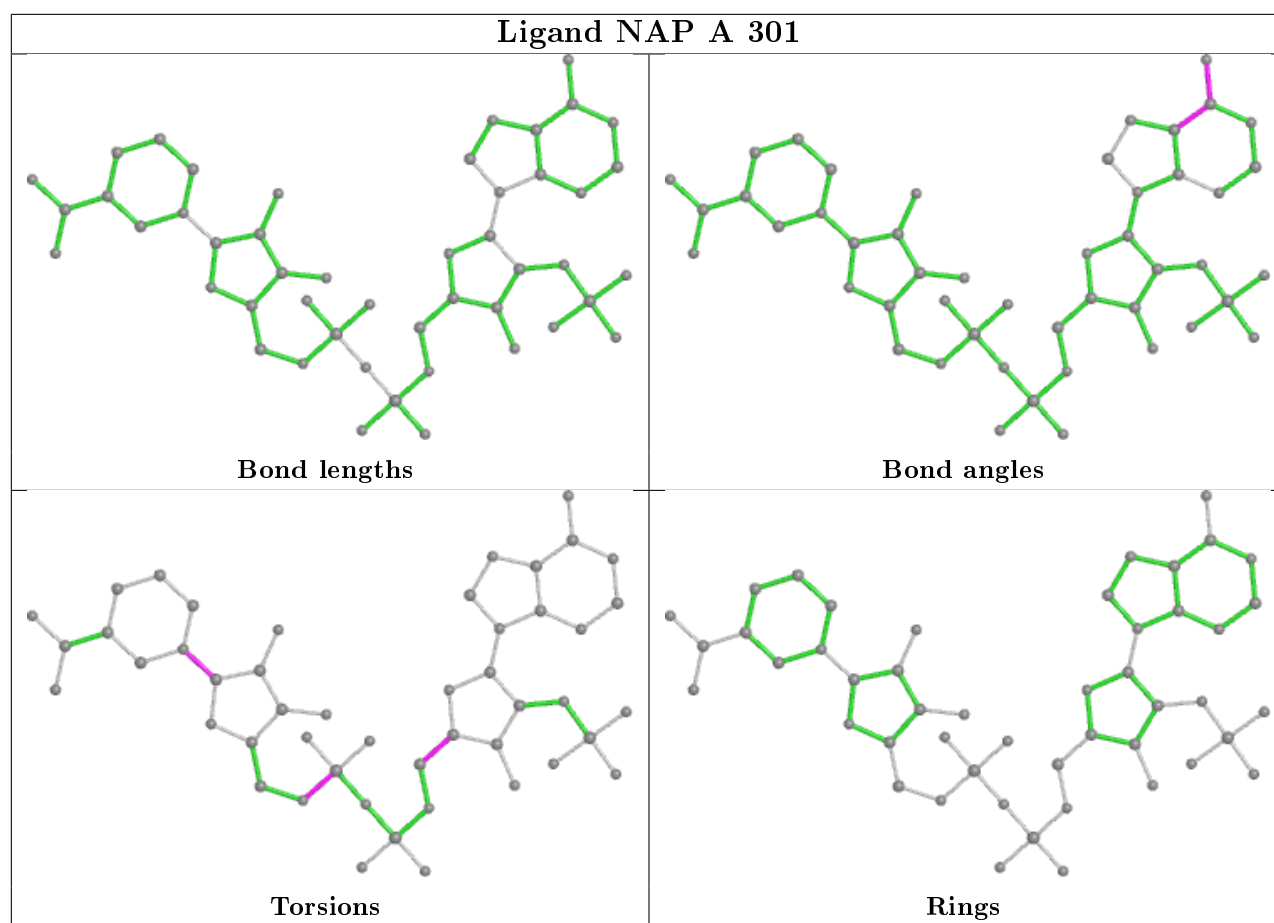












5.7 Other polymers [i](#)

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	261/267 (97%)	-0.07	10 (3%) 40 37	11, 17, 41, 79	0
1	B	261/267 (97%)	0.02	9 (3%) 45 42	13, 19, 40, 64	0
1	C	261/267 (97%)	-0.10	7 (2%) 54 52	11, 16, 37, 65	0
1	D	261/267 (97%)	0.10	12 (4%) 32 29	13, 21, 47, 72	0
All	All	1044/1068 (97%)	-0.01	38 (3%) 42 40	11, 18, 42, 79	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	193	VAL	5.9
1	A	-1	HIS	4.9
1	A	198	ALA	4.6
1	D	193	VAL	4.5
1	A	0	HIS	4.4
1	D	197	ALA	4.0
1	B	201	ALA	3.8
1	D	53	ALA	3.7
1	D	201	ALA	3.7
1	C	193	VAL	3.6
1	A	199	ALA	3.6
1	A	202	GLN	3.5
1	C	0	HIS	3.4
1	D	192	ALA	3.3
1	B	195	PHE	3.3
1	B	197	ALA	3.2
1	A	195	PHE	3.1
1	C	200	LYS	3.0
1	D	206	VAL	3.0
1	B	196	ASP	2.8
1	D	196	ASP	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	197	ALA	2.6
1	A	206	VAL	2.6
1	D	198	ALA	2.6
1	C	196[A]	ASP	2.5
1	B	192	ALA	2.5
1	D	207	ASP	2.5
1	D	54	ASN	2.4
1	B	56	GLY	2.4
1	B	54	ASN	2.4
1	A	197	ALA	2.2
1	A	200	LYS	2.2
1	C	201	ALA	2.1
1	B	0	HIS	2.1
1	C	-1	HIS	2.1
1	D	42	GLU	2.1
1	D	-1	HIS	2.1
1	A	192	ALA	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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.

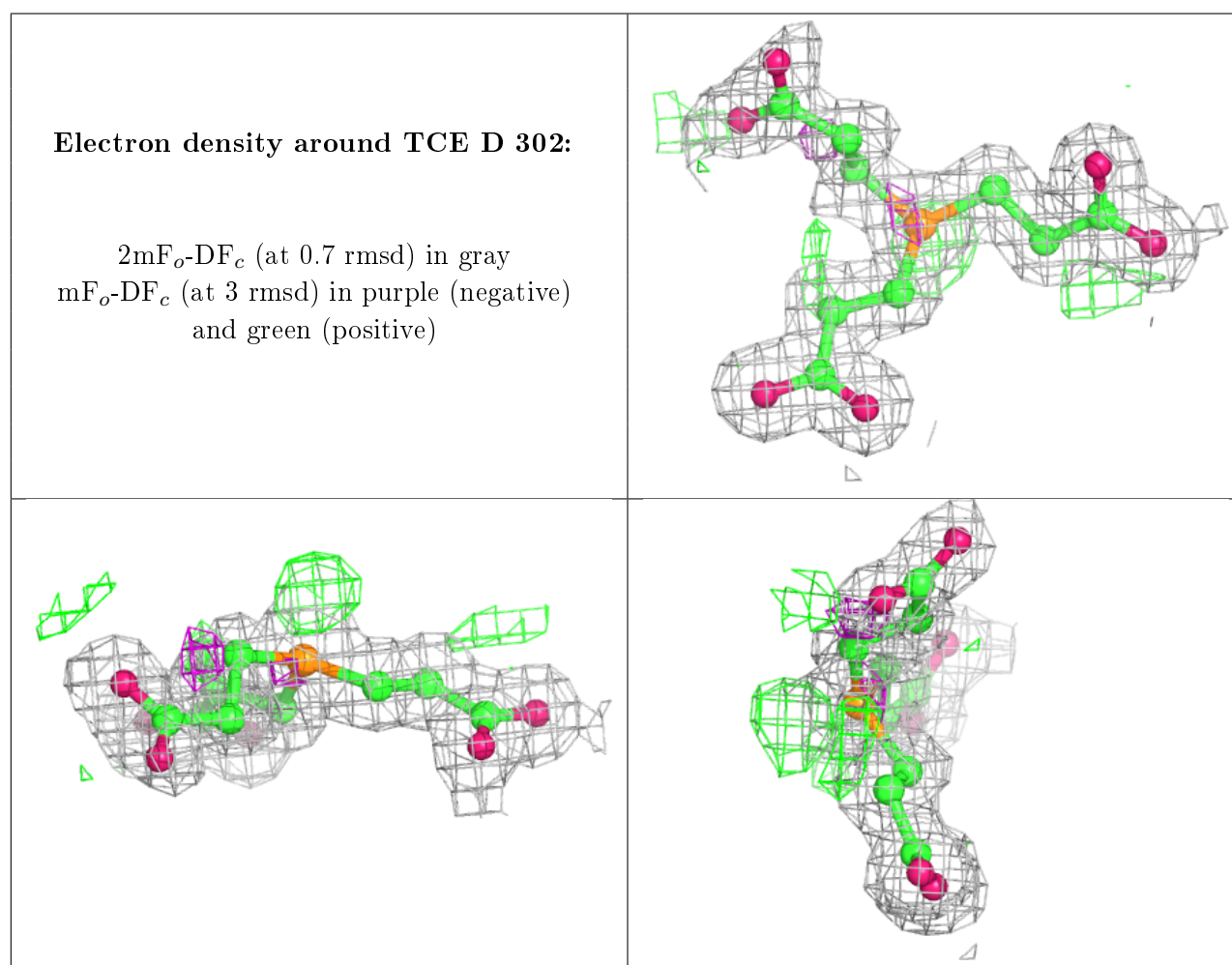
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	MPO	C	303	13/13	0.63	0.24	27,38,59,63	13
4	MPO	A	303	13/13	0.71	0.21	25,36,54,57	13
4	MPO	D	303	13/13	0.84	0.15	34,39,54,57	13
3	TCE	D	302	16/16	0.91	0.10	15,20,37,40	0
3	TCE	B	302	16/16	0.94	0.10	13,19,31,32	0
3	TCE	C	302	16/16	0.94	0.11	11,15,32,33	0

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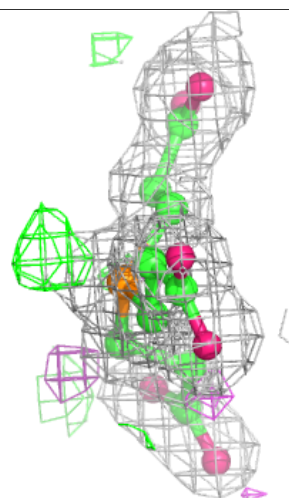
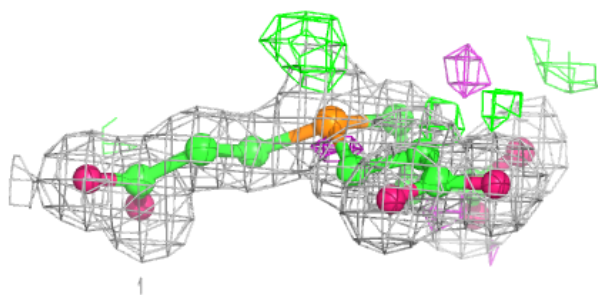
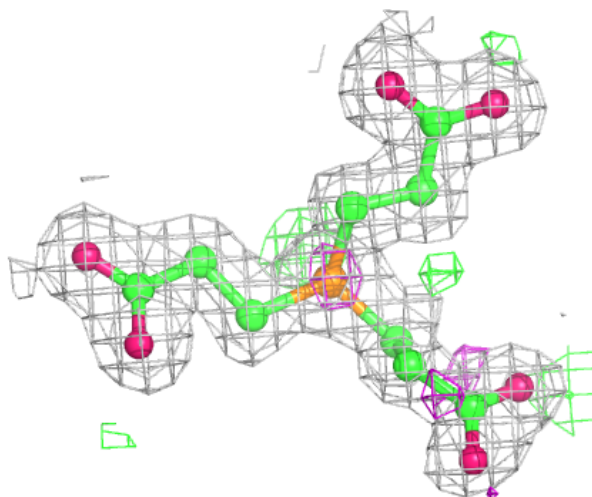
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	TCE	A	302	16/16	0.95	0.10	13,17,45,46	0
2	NAP	D	301	48/48	0.96	0.08	13,18,22,23	0
2	NAP	B	301	48/48	0.96	0.07	13,16,18,21	0
2	NAP	C	301	48/48	0.97	0.07	10,13,16,25	0
2	NAP	A	301	48/48	0.97	0.07	12,14,17,26	0

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.



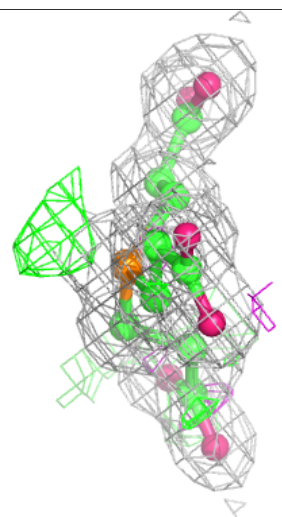
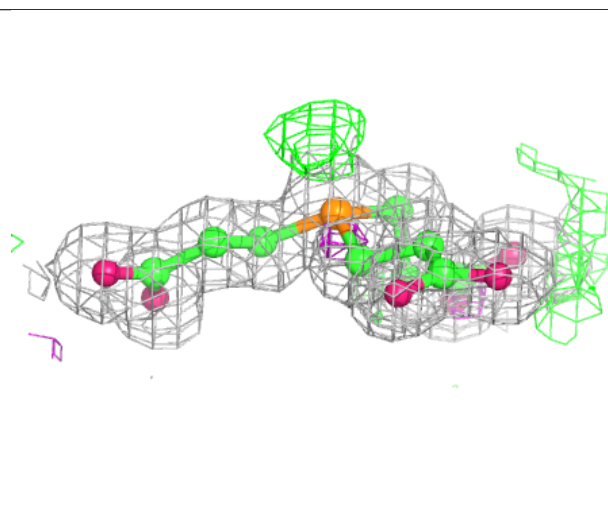
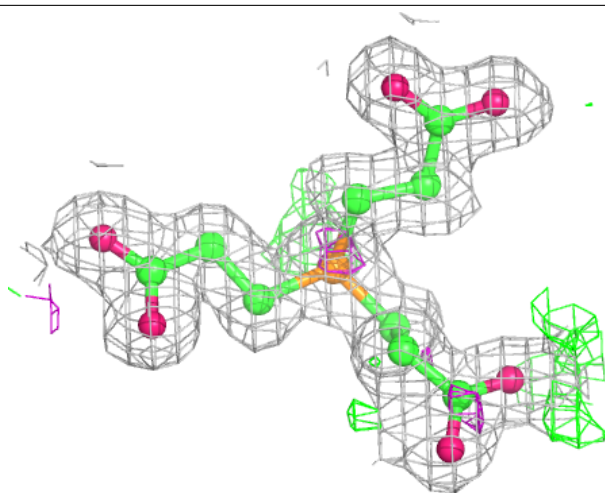
Electron density around TCE B 302:

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



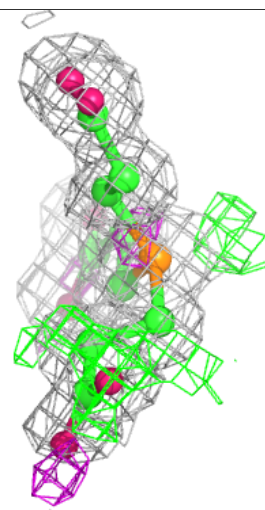
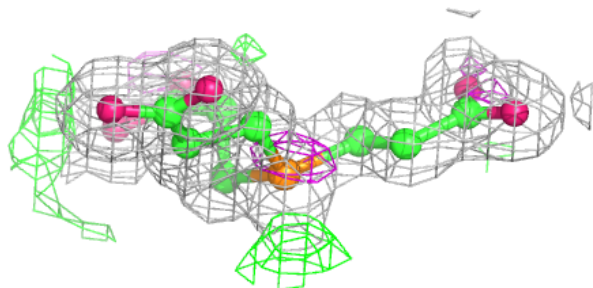
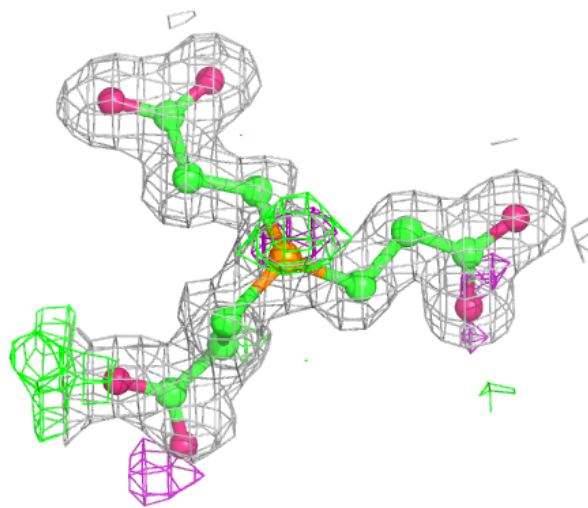
Electron density around TCE C 302:

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



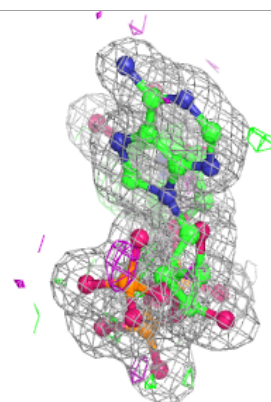
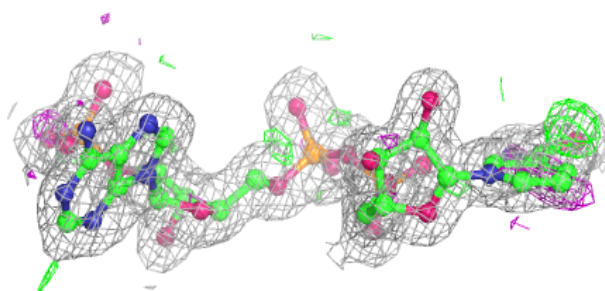
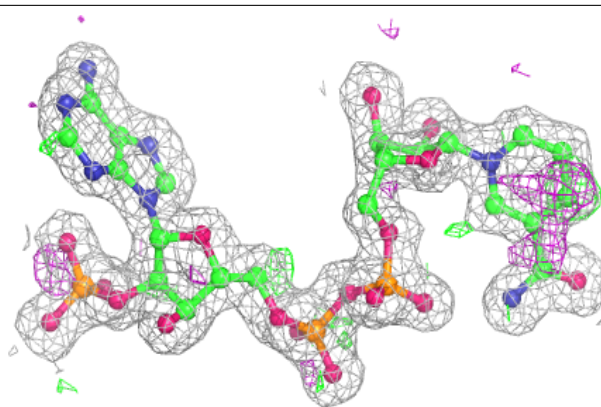
Electron density around TCE A 302:

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

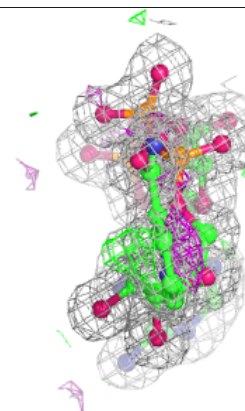
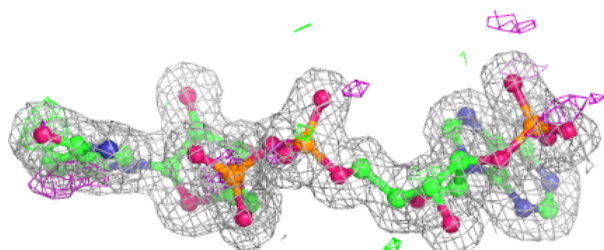
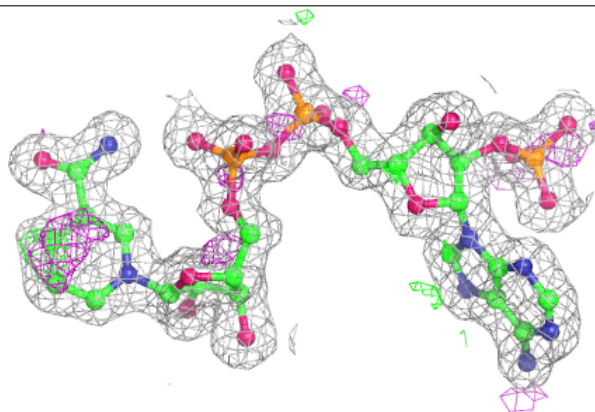


Electron density around NAP D 301:

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

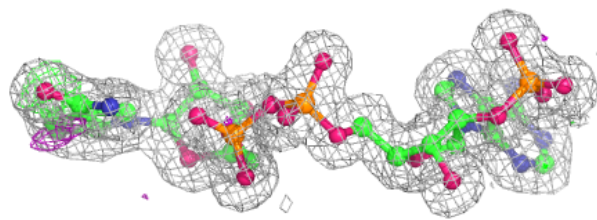
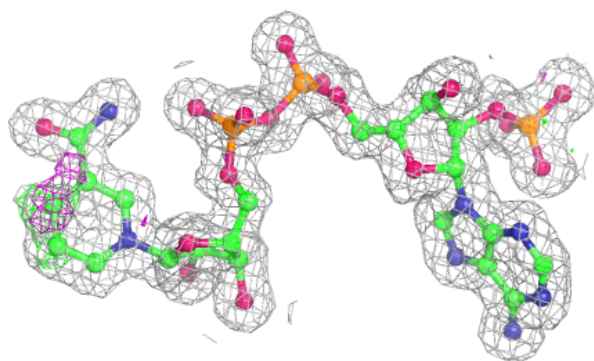
**Electron density around NAP B 301:**

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

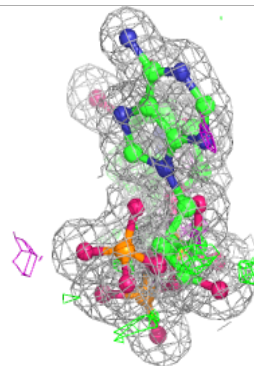
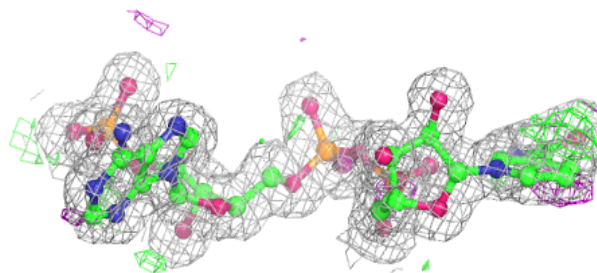
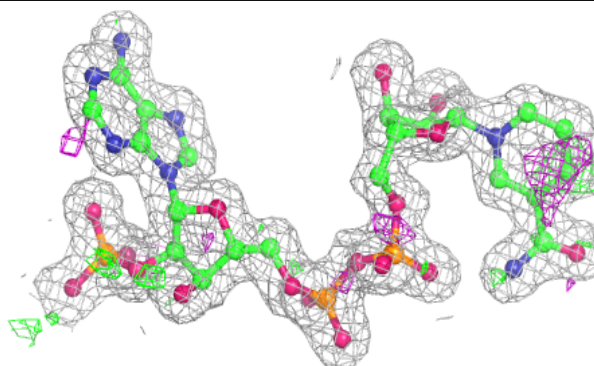


Electron density around NAP C 301:

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

**Electron density around NAP A 301:**

$2mF_o-DF_c$ (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.