



wwPDB X-ray Structure Validation Summary Report

Jan 3, 2024 – 02:33 am GMT

PDB ID : 4UOV
Title : The structure of a tetrameric alpha-carbonic anhydrase from *Thermovibrio ammonificans* reveals a core formed around intermolecular disulfides, which contribute to its thermostability.
Authors : James, P.; Isupov, M.; Sayer, C.; Berg, S.; Lioliou, M.; Kotlar, H.; Littlechild, J.
Deposited on : 2014-06-10
Resolution : 1.85 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

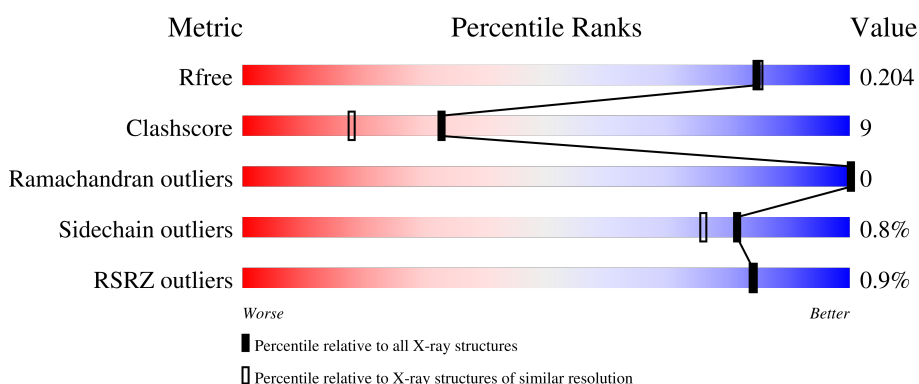
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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	247	
1	B	247	
1	C	247	
1	D	247	
1	E	247	

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Mol	Chain	Length	Quality of chain
1	F	247	 81% 10% 9%

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
11	PG5	B	312	-	-	X	-
5	B3P	C	301	-	X	-	X
5	B3P	D	301	-	-	-	X

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 14247 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.

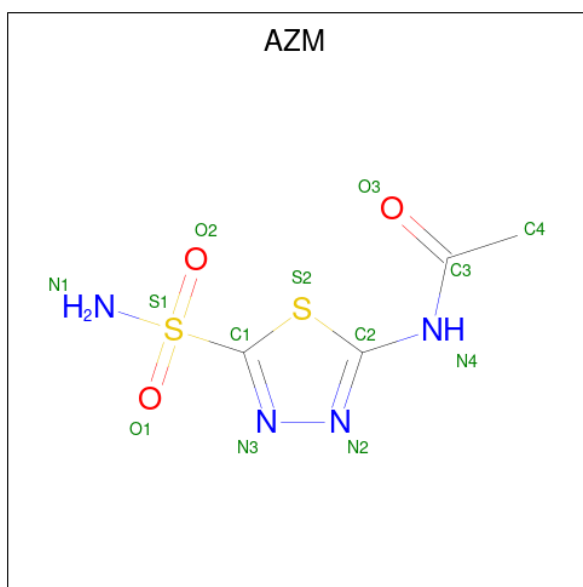
- Molecule 1 is a protein called CARBONATE DEHYDRATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	225	1994	1296	342	343	13	0	34	0
1	B	225	1997	1299	343	344	11	0	33	0
1	C	225	1979	1286	342	339	12	0	30	0
1	D	225	1991	1293	346	341	11	0	31	0
1	E	225	1961	1273	337	339	12	0	27	0
1	F	225	1962	1272	342	337	11	0	26	0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

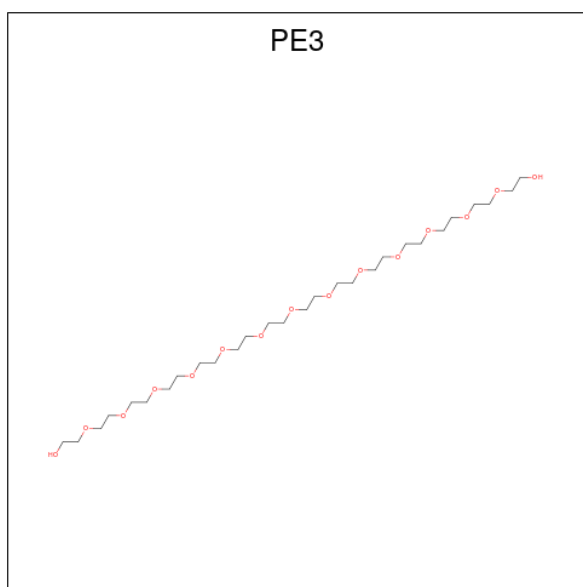
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total 1	Zn 1	0	0
2	B	1	Total 1	Zn 1	0	0
2	C	1	Total 1	Zn 1	0	0
2	D	1	Total 1	Zn 1	0	0
2	E	1	Total 1	Zn 1	0	0
2	F	1	Total 1	Zn 1	0	0

- Molecule 3 is 5-ACETAMIDO-1,3,4-THIADIAZOLE-2-SULFONAMIDE (three-letter code: AZM) (formula: C₄H₆N₄O₃S₂).



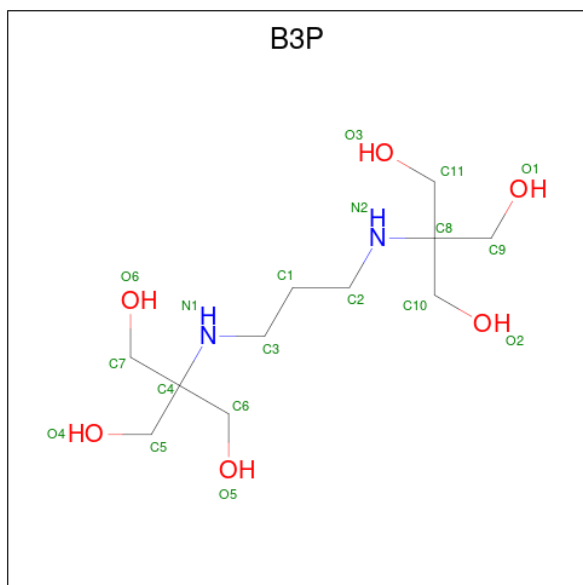
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
3	A	1	Total 13	C 4	N 4	O 3	S 2	0	0
3	B	1	Total 13	C 4	N 4	O 3	S 2	0	0
3	C	1	Total 13	C 4	N 4	O 3	S 2	0	0
3	D	1	Total 13	C 4	N 4	O 3	S 2	0	0
3	E	1	Total 13	C 4	N 4	O 3	S 2	0	0
3	F	1	Total 13	C 4	N 4	O 3	S 2	0	0

- Molecule 4 is 3,6,9,12,15,18,21,24,27,30,33,36,39-TRIDECAXAHENTETRACONTANE-1,41-DIOL (three-letter code: PE3) (formula: C₂₈H₅₈O₁₅).



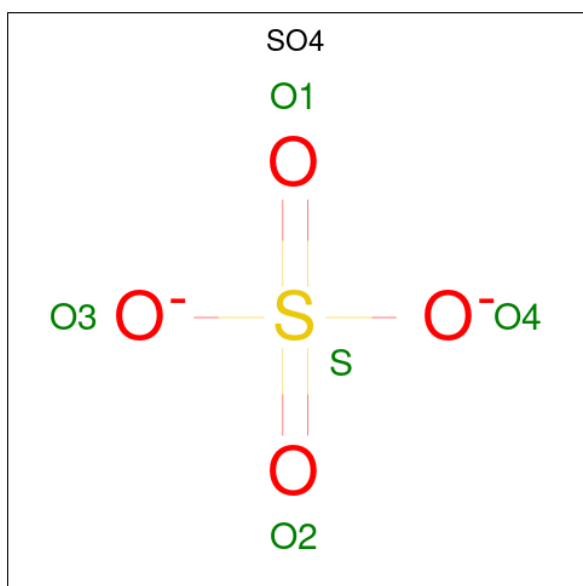
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	C O	0	0
			21	14 7		
4	B	1	Total	C O	0	0
			21	14 7		
4	F	1	Total	C O	0	0
			21	14 7		

- Molecule 5 is 2-[3-(2-HYDROXY-1,1-DIHYDROXYMETHYL-ETHYLAMINO)-PROPYL AMINO]-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: B3P) (formula: $C_{11}H_{26}N_2O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	1
			24	14	2	8		
5	C	1	Total	C	N	O	0	0
			19	11	2	6		
5	D	1	Total	C	N	O	0	0
			19	11	2	6		
5	E	1	Total	C	N	O	0	1
			22	11	2	9		
5	F	1	Total	C	N	O	0	1
			21	11	2	8		

- Molecule 6 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).

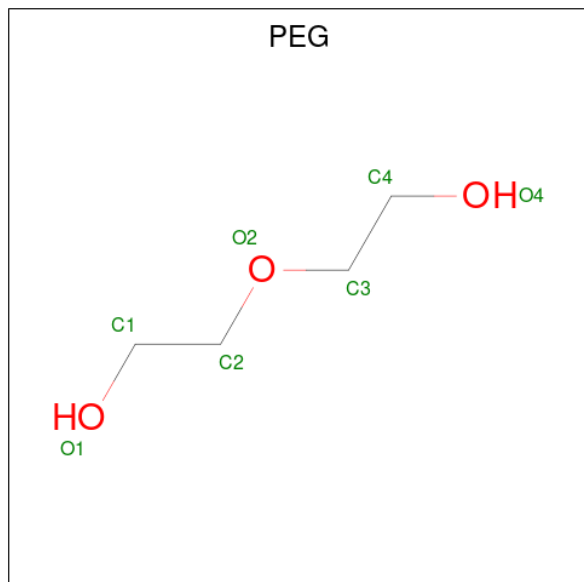


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	B	1	Total	O	S	0	0
			5	4	1		
6	C	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		
6	E	1	Total	O	S	0	0
			5	4	1		
6	F	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	3	Total Cl 3 3	0	0
7	B	3	Total Cl 3 3	0	0
7	C	3	Total Cl 3 3	0	0
7	D	2	Total Cl 2 2	0	0
7	E	2	Total Cl 2 2	0	0
7	F	2	Total Cl 2 2	0	0

- Molecule 8 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



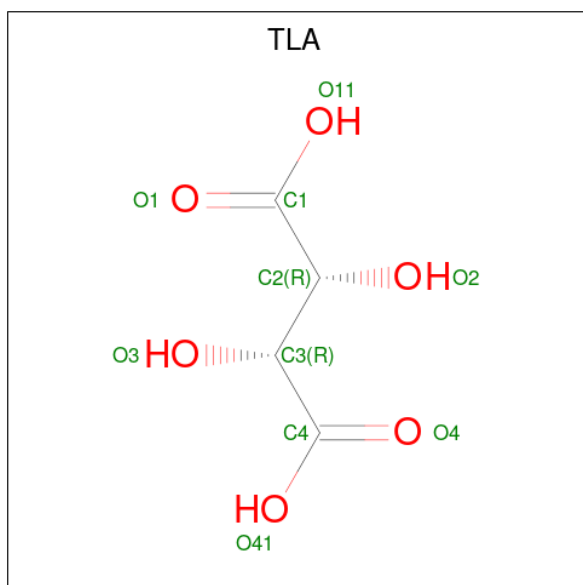
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 7 4 3	0	0
8	B	1	Total C O 7 4 3	0	0
8	B	1	Total C O 7 4 3	0	0
8	C	1	Total C O 7 4 3	0	0
8	E	1	Total C O 7 4 3	0	0
8	E	1	Total C O 7 4 3	0	0

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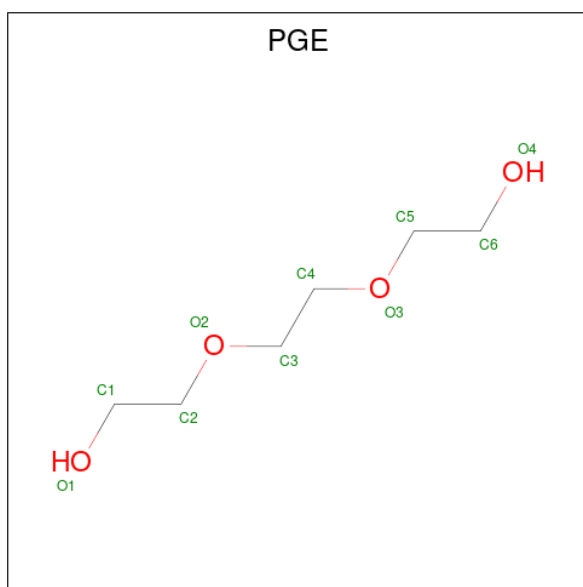
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	F	1	Total C O 7 4 3	0	0
8	F	1	Total C O 7 4 3	0	0

- Molecule 9 is L(+)-TARTARIC ACID (three-letter code: TLA) (formula: C₄H₆O₆).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	1	Total C O 10 4 6	0	0
9	B	1	Total C O 10 4 6	0	0
9	F	1	Total C O 10 4 6	0	0

- Molecule 10 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	B	1	Total C O 10 6 4	0	0
10	C	1	Total C O 10 6 4	0	0
10	D	1	Total C O 10 6 4	0	0
10	E	1	Total C O 10 6 4	0	0
10	F	1	Total C O 10 6 4	0	0

- Molecule 11 is 1-METHOXY-2-[2-(2-METHOXY-ETHOXY)]-ETHANE (three-letter code: PG5) (formula: C₈H₁₈O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	B	1	Total	C	O	0	0
			12	8	4		

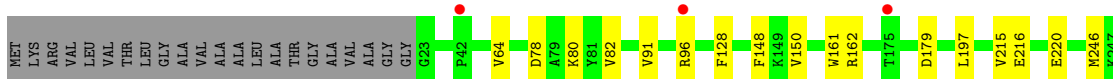
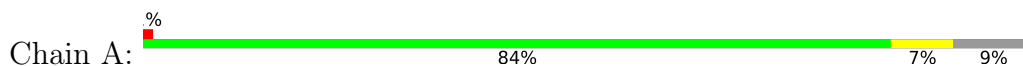
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	360	Total	O	0	0
			360	360		
12	B	363	Total	O	0	0
			363	363		
12	C	291	Total	O	0	0
			291	291		
12	D	242	Total	O	0	0
			242	242		
12	E	339	Total	O	0	0
			339	339		
12	F	323	Total	O	0	0
			323	323		

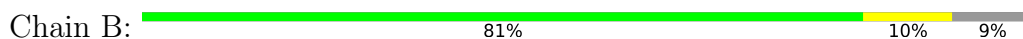
3 Residue-property plots [i](#)

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.

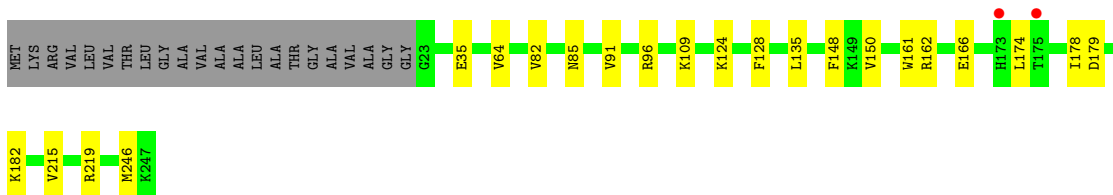
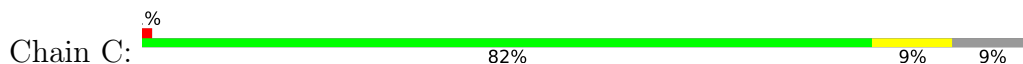
- Molecule 1: CARBONATE DEHYDRATASE



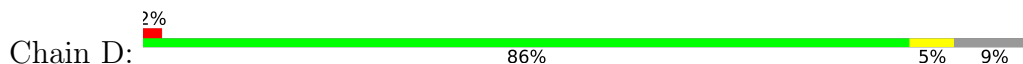
- Molecule 1: CARBONATE DEHYDRATASE



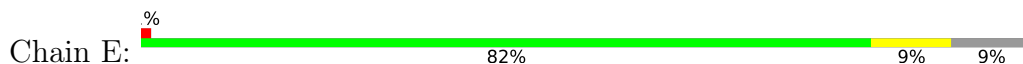
- Molecule 1: CARBONATE DEHYDRATASE



- Molecule 1: CARBONATE DEHYDRATASE



- Molecule 1: CARBONATE DEHYDRATASE





- Molecule 1: CARBONATE DEHYDRATASE

Chain F: 81% 10% 9%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	163.75Å 283.02Å 52.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.01 – 1.85 42.01 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.5 (42.01-1.85) 99.5 (42.01-1.85)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 1.86Å)	Xtrriage
Refinement program	REFMAC 5.8.0071	Depositor
R, R_{free}	0.172 , 0.204 0.172 , 0.204	Depositor DCC
R_{free} test set	10399 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	22.9	Xtrriage
Anisotropy	0.365	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 60.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	14247	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.78% 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: SO4, TLA, B3P, PE3, ZN, CL, PGE, PEG, AZM, PG5

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	0.42	0/2152	0.54	0/2904
1	B	0.44	0/2153	0.56	0/2904
1	C	0.41	0/2126	0.53	0/2867
1	D	0.41	0/2141	0.54	0/2886
1	E	0.44	0/2097	0.55	0/2829
1	F	0.46	0/2096	0.56	0/2825
All	All	0.43	0/12765	0.55	0/17215

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	1994	0	2071	26	0
1	B	1997	0	2070	76	1
1	C	1979	0	2048	30	0
1	D	1991	0	2064	28	0
1	E	1961	0	2024	38	0
1	F	1962	0	2022	39	1
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	13	0	6	0	0
3	B	13	0	5	0	0
3	C	13	0	6	1	0
3	D	13	0	6	0	0
3	E	13	0	6	0	0
3	F	13	0	6	0	0
4	A	21	0	24	0	0
4	B	21	0	24	0	0
4	F	21	0	24	0	0
5	A	24	0	20	0	0
5	C	19	0	26	3	0
5	D	19	0	26	0	0
5	E	22	0	18	0	0
5	F	21	0	12	0	0
6	A	5	0	0	1	0
6	B	5	0	0	0	0
6	C	5	0	0	1	0
6	D	5	0	0	0	0
6	E	5	0	0	0	0
6	F	5	0	0	0	0
7	A	3	0	0	1	0
7	B	3	0	0	1	0
7	C	3	0	0	1	0
7	D	2	0	0	1	0
7	E	2	0	0	1	0
7	F	2	0	0	1	0
8	A	7	0	10	1	0
8	B	14	0	20	2	0
8	C	7	0	10	1	0
8	E	14	0	20	3	0
8	F	14	0	20	4	0
9	A	10	0	4	0	0
9	B	10	0	4	0	0
9	F	10	0	4	0	0
10	B	10	0	14	5	0
10	C	10	0	14	0	0
10	D	10	0	14	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	E	10	0	14	1	0
10	F	10	0	14	0	0
11	B	12	0	18	31	0
12	A	360	0	0	13	0
12	B	363	0	0	16	0
12	C	291	0	0	8	0
12	D	242	0	0	15	0
12	E	339	0	0	18	0
12	F	323	0	0	11	0
All	All	14247	0	12688	218	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 9.

The worst 5 of 218 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:246[B]:MET:CE	1:D:64[B]:VAL:HG11	1.55	1.35
1:A:64[B]:VAL:HG11	1:C:246[B]:MET:CE	1.57	1.32
1:B:171[B]:LYS:NZ	11:B:312:PG5:H12	1.46	1.26
1:E:220[B]:GLU:CD	12:E:2279:HOH:O	1.73	1.24
1:B:173[B]:HIS:NE2	11:B:312:PG5:O1	1.71	1.23

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:175:THR:OG1	1:F:172[B]:ARG:NH2[3_444]	2.14	0.06

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	257/247 (104%)	253 (98%)	4 (2%)	0	100	100
1	B	256/247 (104%)	252 (98%)	4 (2%)	0	100	100
1	C	253/247 (102%)	249 (98%)	4 (2%)	0	100	100
1	D	254/247 (103%)	250 (98%)	4 (2%)	0	100	100
1	E	250/247 (101%)	245 (98%)	5 (2%)	0	100	100
1	F	249/247 (101%)	245 (98%)	4 (2%)	0	100	100
All	All	1519/1482 (102%)	1494 (98%)	25 (2%)	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	230/208 (111%)	229 (100%)	1 (0%)	91	89
1	B	229/208 (110%)	228 (100%)	1 (0%)	91	89
1	C	226/208 (109%)	225 (100%)	1 (0%)	91	89
1	D	227/208 (109%)	225 (99%)	2 (1%)	78	72
1	E	223/208 (107%)	220 (99%)	3 (1%)	69	58
1	F	222/208 (107%)	219 (99%)	3 (1%)	67	55
All	All	1357/1248 (109%)	1346 (99%)	11 (1%)	81	76

5 of 11 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	220[B]	GLU
1	F	152[A]	LYS
1	F	161	TRP
1	F	152[B]	LYS
1	D	236	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	140	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 61 ligands modelled in this entry, 21 are monoatomic - leaving 40 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	$\# Z > 2$
8	PEG	E	313	-	6,6,6	0.46	0	5,5,5	0.53	0
10	PGE	E	305	-	9,9,9	0.44	0	8,8,8	0.28	0
5	B3P	F	301[B]	-	18,18,18	0.42	0	21,23,23	1.82	3 (14%)
6	SO4	C	302	-	4,4,4	0.33	0	6,6,6	0.22	0
8	PEG	E	314	-	6,6,6	0.41	0	5,5,5	0.39	0
5	B3P	A	301[A]	-	18,18,18	0.52	0	21,23,23	1.97	2 (9%)
8	PEG	F	314	-	6,6,6	0.41	0	5,5,5	0.37	0
4	PE3	A	300	-	20,20,42	0.52	0	19,19,41	0.30	0
4	PE3	B	300	-	20,20,42	0.53	0	19,19,41	0.22	0
3	AZM	F	299	2	8,13,13	0.44	0	9,19,19	2.76	4 (44%)
6	SO4	A	302	-	4,4,4	0.28	0	6,6,6	0.07	0
8	PEG	A	305	-	6,6,6	0.40	0	5,5,5	0.37	0
9	TLA	B	306	-	9,9,9	1.10	0	12,12,12	1.11	2 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	SO4	B	302	-	4,4,4	0.26	0	6,6,6	0.08	0
5	B3P	F	301[A]	-	18,18,18	0.42	0	21,23,23	1.82	3 (14%)
6	SO4	E	302	-	4,4,4	0.25	0	6,6,6	0.06	0
10	PGE	D	305	-	9,9,9	0.44	0	8,8,8	0.22	0
10	PGE	F	309	-	9,9,9	0.44	0	8,8,8	0.30	0
4	PE3	F	300	-	20,20,42	0.48	0	19,19,41	0.23	0
5	B3P	E	301[B]	-	18,18,18	0.52	0	21,23,23	1.80	2 (9%)
5	B3P	C	301	-	18,18,18	0.65	0	21,23,23	5.54	13 (61%)
3	AZM	E	299	2	8,13,13	0.24	0	9,19,19	1.84	2 (22%)
9	TLA	A	306	-	9,9,9	1.10	0	12,12,12	0.89	0
8	PEG	B	310	-	6,6,6	0.49	0	5,5,5	0.21	0
10	PGE	B	305	-	9,9,9	0.47	0	8,8,8	0.34	0
3	AZM	A	299	2	8,13,13	0.35	0	9,19,19	2.16	2 (22%)
11	PG5	B	312	-	11,11,11	0.49	0	10,10,10	0.30	0
6	SO4	F	302	-	4,4,4	0.31	0	6,6,6	0.06	0
8	PEG	B	314	-	6,6,6	0.45	0	5,5,5	0.37	0
5	B3P	E	301[A]	-	18,18,18	0.51	0	21,23,23	1.80	2 (9%)
6	SO4	D	302	-	4,4,4	0.34	0	6,6,6	0.07	0
8	PEG	F	310	-	6,6,6	0.55	0	5,5,5	0.36	0
9	TLA	F	306	-	9,9,9	1.13	0	12,12,12	0.96	0
10	PGE	C	305	-	9,9,9	0.45	0	8,8,8	0.27	0
3	AZM	D	299	2	8,13,13	0.29	0	9,19,19	1.90	2 (22%)
3	AZM	B	299	2	8,13,13	0.32	0	9,19,19	2.44	3 (33%)
5	B3P	A	301[B]	-	18,18,18	0.56	0	21,23,23	1.84	2 (9%)
3	AZM	C	299	2	8,13,13	0.28	0	9,19,19	2.41	4 (44%)
8	PEG	C	314	-	6,6,6	0.46	0	5,5,5	0.32	0
5	B3P	D	301	-	18,18,18	0.45	0	21,23,23	2.14	2 (9%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	PEG	E	313	-	-	4/4/4/4	-
10	PGE	E	305	-	-	5/7/7/7	-
5	B3P	F	301[B]	-	-	2/28/28/28	-
8	PEG	E	314	-	-	1/4/4/4	-
5	B3P	A	301[A]	-	-	16/28/28/28	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	PEG	F	314	-	-	2/4/4/4	-
4	PE3	A	300	-	-	3/18/18/40	-
4	PE3	B	300	-	-	3/18/18/40	-
3	AZM	F	299	2	-	0/2/10/10	0/1/1/1
8	PEG	A	305	-	-	4/4/4/4	-
9	TLA	B	306	-	-	2/12/12/12	-
5	B3P	F	301[A]	-	-	1/28/28/28	-
10	PGE	F	309	-	-	4/7/7/7	-
10	PGE	D	305	-	-	3/7/7/7	-
4	PE3	F	300	-	-	1/18/18/40	-
5	B3P	E	301[B]	-	-	9/28/28/28	-
5	B3P	C	301	-	-	17/28/28/28	-
3	AZM	E	299	2	-	0/2/10/10	0/1/1/1
9	TLA	A	306	-	-	0/12/12/12	-
8	PEG	B	310	-	-	2/4/4/4	-
10	PGE	B	305	-	-	5/7/7/7	-
3	AZM	A	299	2	-	0/2/10/10	0/1/1/1
11	PG5	B	312	-	-	7/9/9/9	-
8	PEG	B	314	-	-	2/4/4/4	-
5	B3P	E	301[A]	-	-	5/28/28/28	-
8	PEG	F	310	-	-	2/4/4/4	-
9	TLA	F	306	-	-	0/12/12/12	-
10	PGE	C	305	-	-	4/7/7/7	-
3	AZM	D	299	2	-	0/2/10/10	0/1/1/1
3	AZM	B	299	2	-	2/2/10/10	0/1/1/1
5	B3P	A	301[B]	-	-	18/28/28/28	-
3	AZM	C	299	2	-	2/2/10/10	0/1/1/1
8	PEG	C	314	-	-	2/4/4/4	-
5	B3P	D	301	-	-	7/28/28/28	-

There are no bond length outliers.

The worst 5 of 48 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	301	B3P	C10-C8-C9	-11.56	85.60	110.04
5	C	301	B3P	C6-C4-C5	-10.84	87.11	110.04
5	C	301	B3P	C7-C4-C5	-9.77	89.39	110.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	301	B3P	C9-C8-N2	-7.48	86.52	109.03
5	C	301	B3P	C5-C4-N1	-7.35	86.92	109.03

There are no chirality outliers.

5 of 135 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	299	AZM	C4-C3-N4-C2
3	B	299	AZM	O3-C3-N4-C2
5	A	301[A]	B3P	C5-C4-N1-C3
5	A	301[A]	B3P	C6-C4-N1-C3
5	A	301[A]	B3P	C7-C4-N1-C3

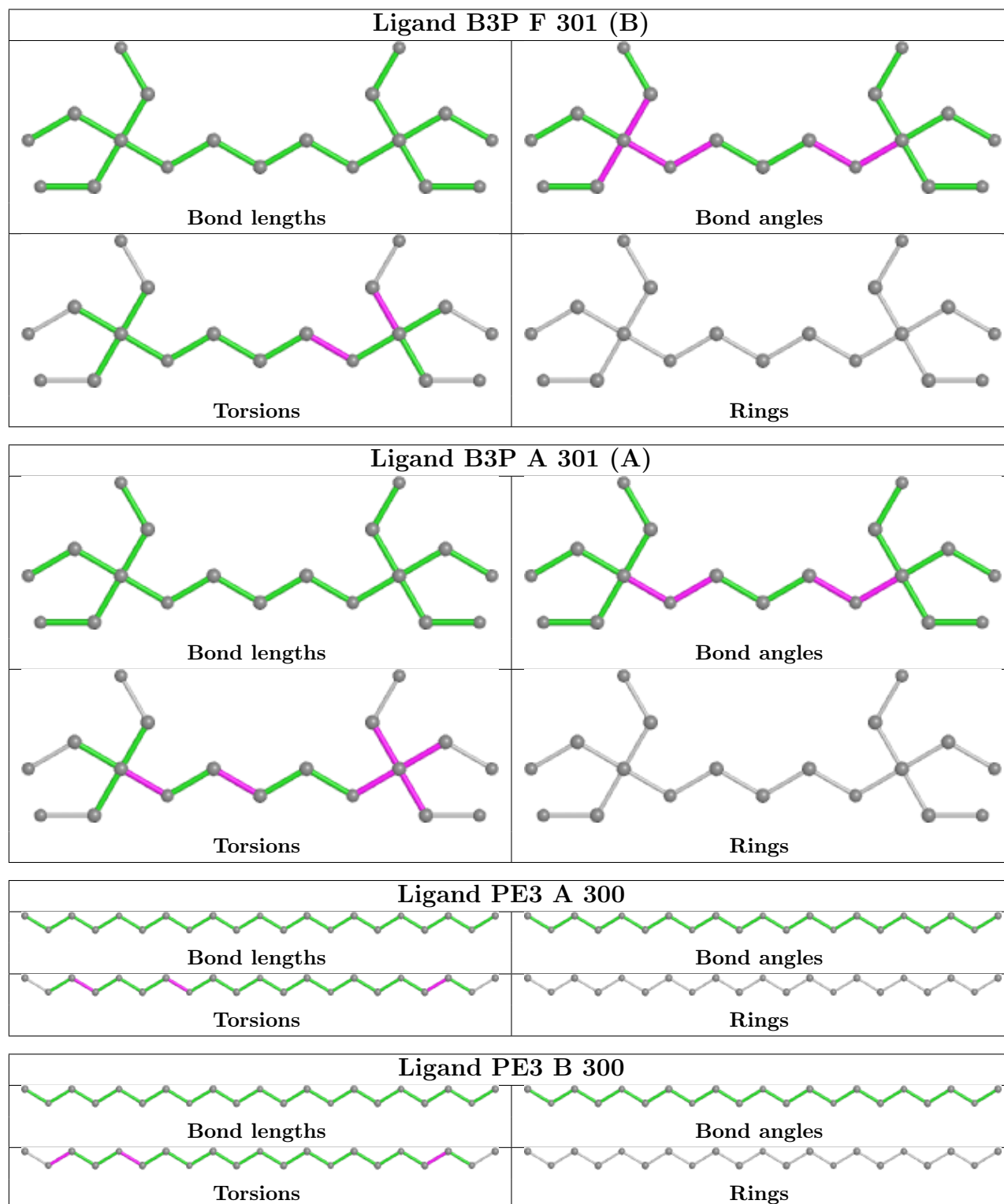
There are no ring outliers.

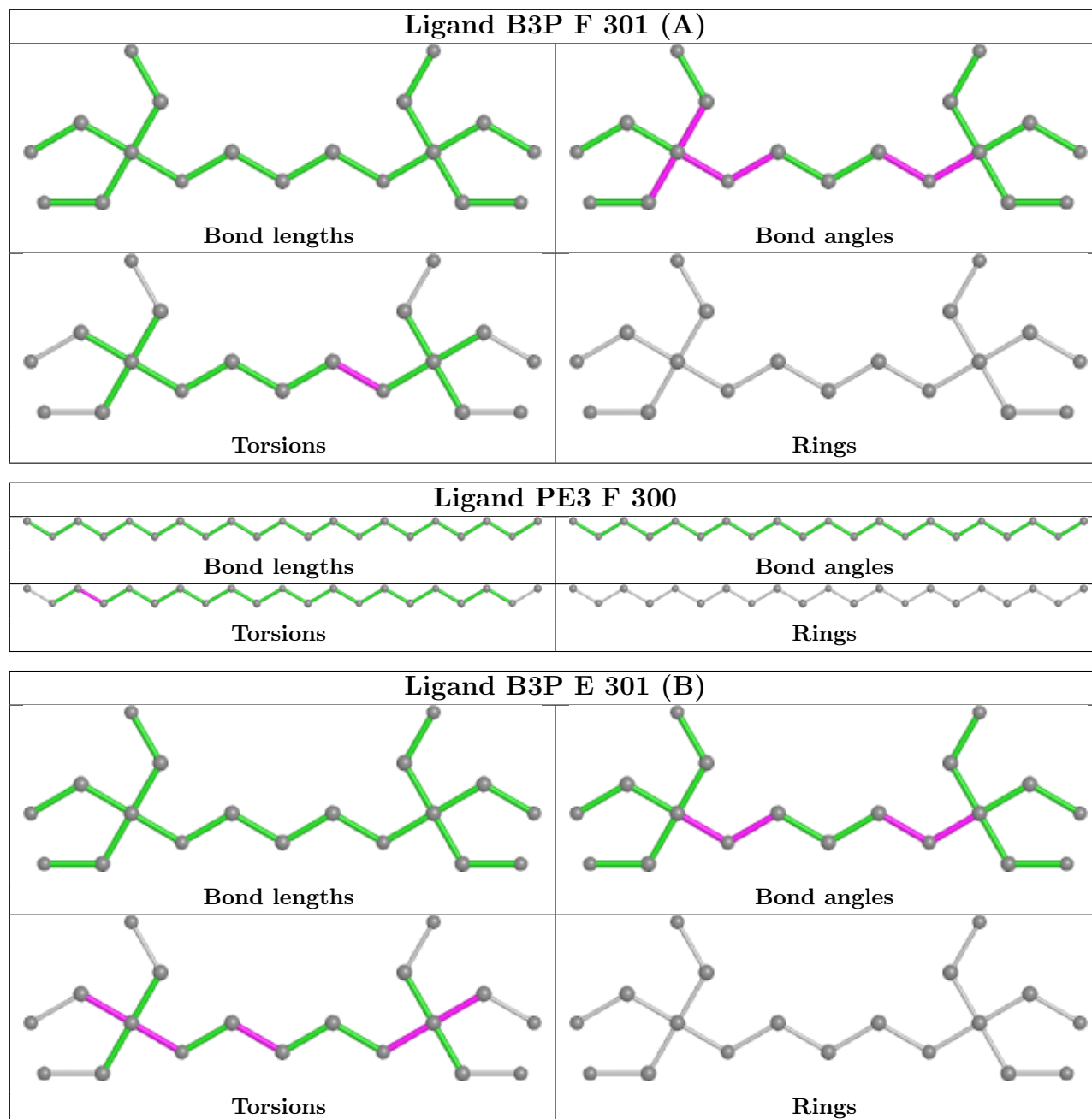
16 monomers are involved in 55 short contacts:

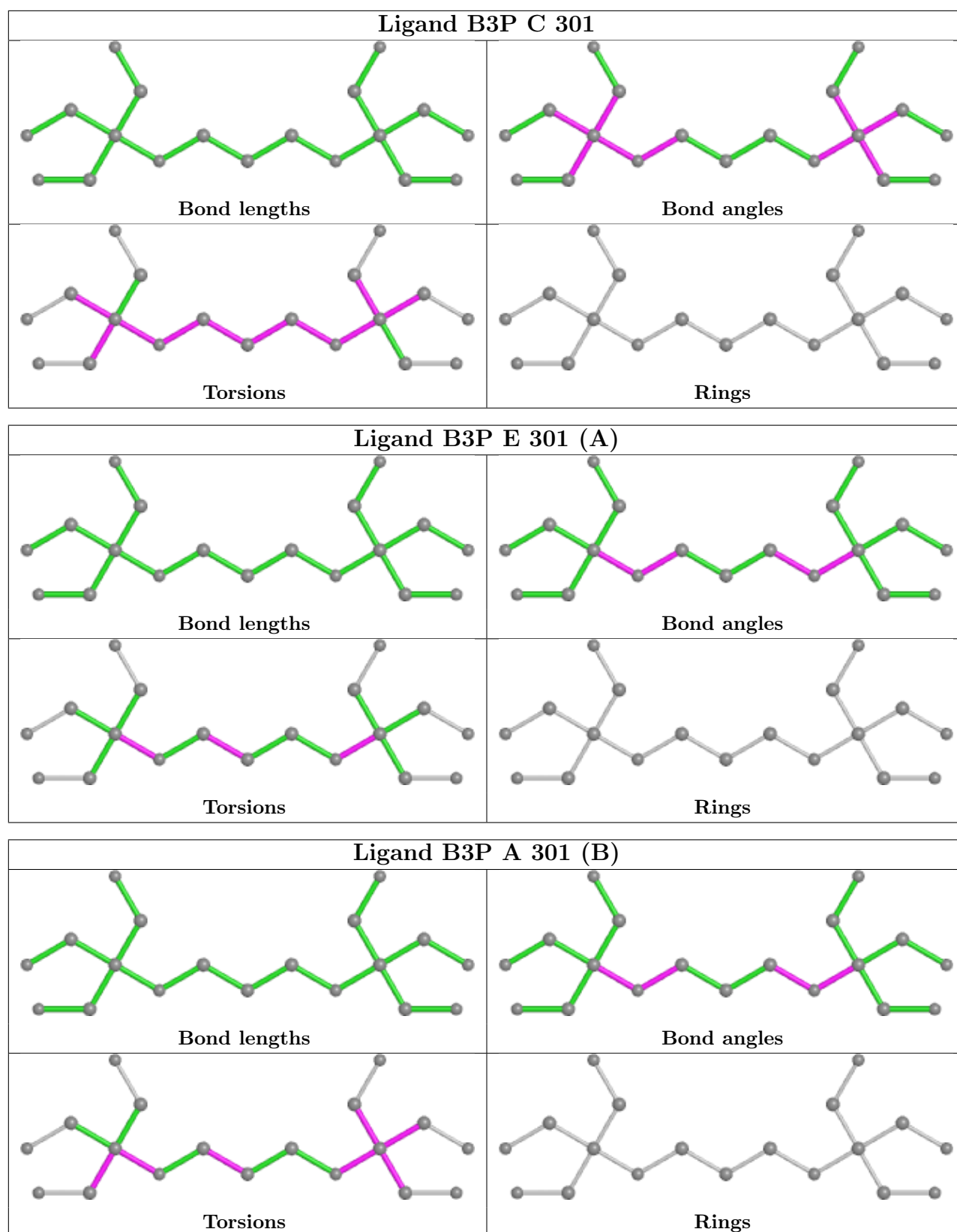
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	E	313	PEG	2	0
10	E	305	PGE	1	0
6	C	302	SO4	1	0
8	E	314	PEG	1	0
8	F	314	PEG	2	0
6	A	302	SO4	1	0
8	A	305	PEG	1	0
10	D	305	PGE	1	0
5	C	301	B3P	3	0
8	B	310	PEG	1	0
10	B	305	PGE	5	0
11	B	312	PG5	31	0
8	B	314	PEG	1	0
8	F	310	PEG	2	0
3	C	299	AZM	1	0
8	C	314	PEG	1	0

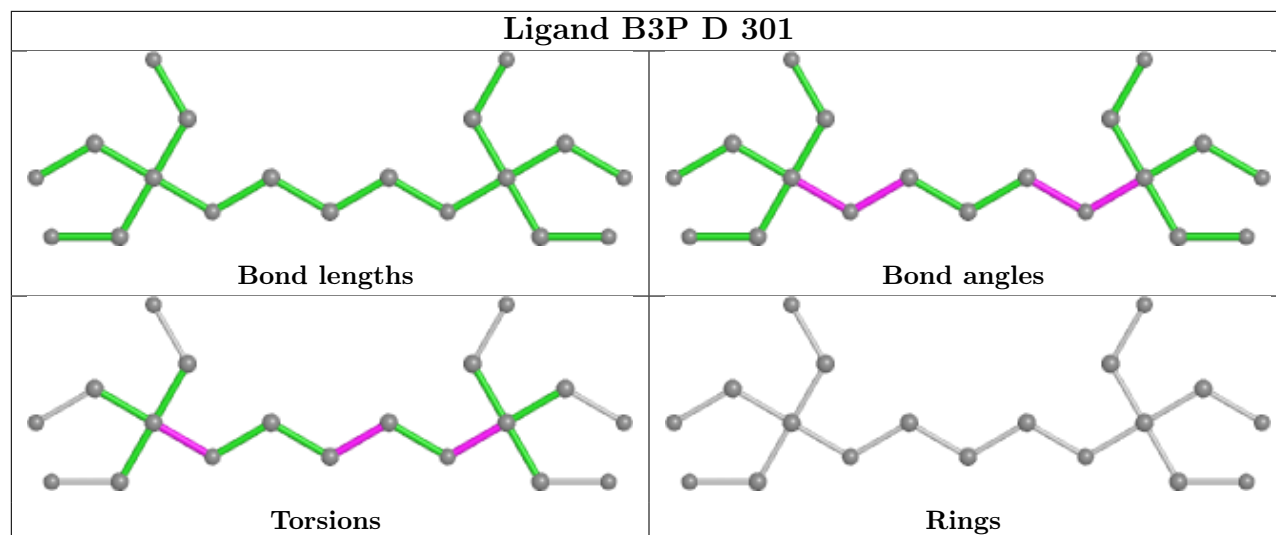
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	225/247 (91%)	-0.17	3 (1%) 77 78	14, 24, 42, 65	0
1	B	225/247 (91%)	-0.27	0 100 100	13, 22, 39, 61	0
1	C	225/247 (91%)	-0.20	2 (0%) 84 84	14, 26, 45, 71	0
1	D	225/247 (91%)	-0.14	4 (1%) 68 68	14, 26, 43, 67	0
1	E	225/247 (91%)	-0.23	2 (0%) 84 84	13, 24, 43, 58	0
1	F	225/247 (91%)	-0.36	1 (0%) 92 92	12, 21, 36, 61	0
All	All	1350/1482 (91%)	-0.23	12 (0%) 84 84	12, 24, 43, 71	0

The worst 5 of 12 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	175	THR	3.9
1	C	175	THR	3.2
1	C	173[A]	HIS	3.1
1	D	23	GLY	2.9
1	A	175	THR	2.5

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 monosaccharides in this entry.

6.4 Ligands

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
5	B3P	C	301	19/19	0.66	0.43	38,46,53,54	19
5	B3P	D	301	19/19	0.70	0.42	32,40,52,54	19
8	PEG	F	310	7/7	0.75	0.29	30,32,44,45	7
10	PGE	D	305	10/10	0.75	0.24	36,49,55,64	10
7	CL	D	307	1/1	0.76	0.13	48,48,48,48	1
10	PGE	F	309	10/10	0.77	0.18	34,38,42,42	10
10	PGE	B	305	10/10	0.78	0.17	34,43,54,59	10
8	PEG	B	310	7/7	0.79	0.26	24,29,32,38	7
10	PGE	E	305	10/10	0.81	0.17	33,42,47,51	10
8	PEG	C	314	7/7	0.85	0.18	46,61,79,85	0
9	TLA	B	306	10/10	0.85	0.22	20,25,27,30	10
11	PG5	B	312	12/12	0.85	0.21	33,40,44,47	12
5	B3P	E	301[A]	19/19	0.86	0.22	26,30,46,49	19
8	PEG	E	314	7/7	0.86	0.20	26,37,47,50	7
5	B3P	E	301[B]	19/19	0.86	0.22	23,34,46,49	19
8	PEG	A	305	7/7	0.87	0.17	33,35,43,52	7
10	PGE	C	305	10/10	0.87	0.21	43,47,56,56	10
8	PEG	B	314	7/7	0.87	0.19	36,53,64,68	0
5	B3P	A	301[A]	19/19	0.88	0.23	22,32,48,48	19
5	B3P	A	301[B]	19/19	0.88	0.23	25,33,48,48	19
8	PEG	E	313	7/7	0.89	0.17	39,41,45,50	7
8	PEG	F	314	7/7	0.89	0.18	25,30,41,43	7
4	PE3	A	300	21/43	0.89	0.17	35,41,63,66	0
7	CL	A	307	1/1	0.90	0.12	37,37,37,37	1
9	TLA	F	306	10/10	0.90	0.16	20,24,25,25	10
4	PE3	B	300	21/43	0.90	0.20	38,43,64,68	0
9	TLA	A	306	10/10	0.90	0.15	20,25,27,28	10
7	CL	C	307	1/1	0.91	0.10	33,33,33,33	1
5	B3P	F	301[B]	19/19	0.91	0.12	27,33,47,52	2
7	CL	E	307	1/1	0.91	0.11	37,37,37,37	1
5	B3P	F	301[A]	19/19	0.91	0.12	26,33,47,52	2
7	CL	C	303	1/1	0.91	0.17	37,37,37,37	1
4	PE3	F	300	21/43	0.92	0.15	33,40,62,67	0
7	CL	A	303	1/1	0.92	0.10	34,34,34,34	1
7	CL	F	308	1/1	0.94	0.11	25,25,25,25	1
6	SO4	A	302	5/5	0.95	0.17	37,41,46,51	5
6	SO4	D	302	5/5	0.95	0.13	71,73,74,79	0

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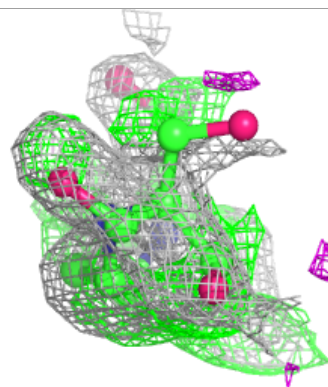
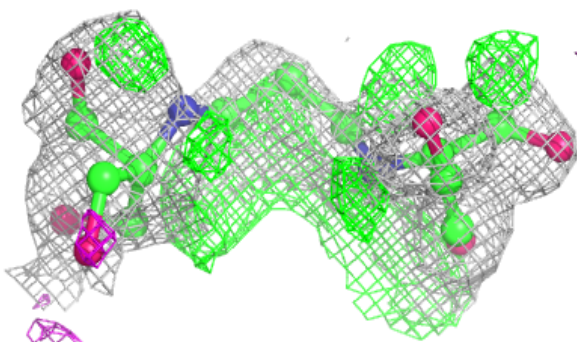
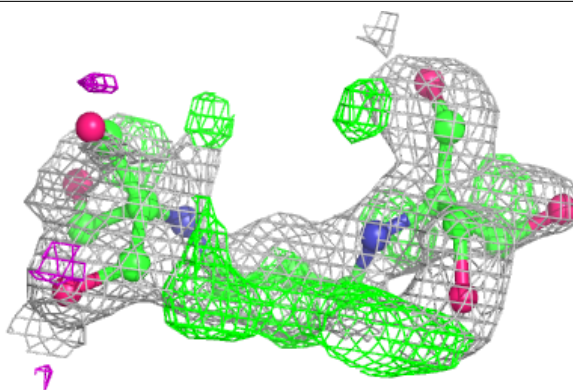
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	SO4	F	302	5/5	0.96	0.14	50,51,52,57	5
7	CL	B	303	1/1	0.96	0.09	29,29,29,29	0
6	SO4	C	302	5/5	0.96	0.11	44,50,52,53	5
6	SO4	E	302	5/5	0.97	0.11	37,41,42,47	5
7	CL	B	304	1/1	0.97	0.11	31,31,31,31	1
6	SO4	B	302	5/5	0.97	0.12	43,46,49,52	5
3	AZM	F	299	13/13	0.98	0.10	17,25,44,45	0
3	AZM	B	299	13/13	0.98	0.10	17,29,54,56	0
7	CL	F	304	1/1	0.98	0.07	32,32,32,32	0
7	CL	B	308	1/1	0.98	0.08	30,30,30,30	1
3	AZM	D	299	13/13	0.98	0.07	18,29,36,36	0
3	AZM	E	299	13/13	0.98	0.07	18,27,37,41	0
7	CL	D	304	1/1	0.98	0.05	34,34,34,34	0
2	ZN	B	298	1/1	0.99	0.07	17,17,17,17	0
3	AZM	C	299	13/13	0.99	0.08	17,27,36,41	0
7	CL	A	304	1/1	0.99	0.16	24,24,24,24	1
7	CL	E	304	1/1	0.99	0.05	34,34,34,34	0
3	AZM	A	299	13/13	0.99	0.10	21,28,42,43	0
7	CL	C	304	1/1	0.99	0.07	34,34,34,34	0
2	ZN	A	298	1/1	1.00	0.07	19,19,19,19	0
2	ZN	C	298	1/1	1.00	0.07	20,20,20,20	0
2	ZN	D	298	1/1	1.00	0.07	20,20,20,20	0
2	ZN	E	298	1/1	1.00	0.08	18,18,18,18	0
2	ZN	F	298	1/1	1.00	0.07	17,17,17,17	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.

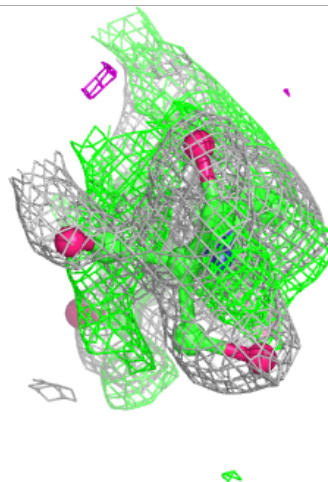
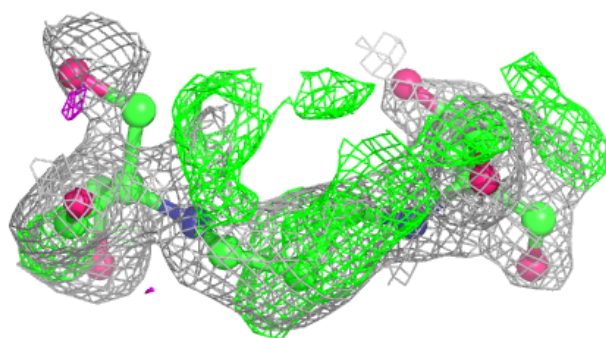
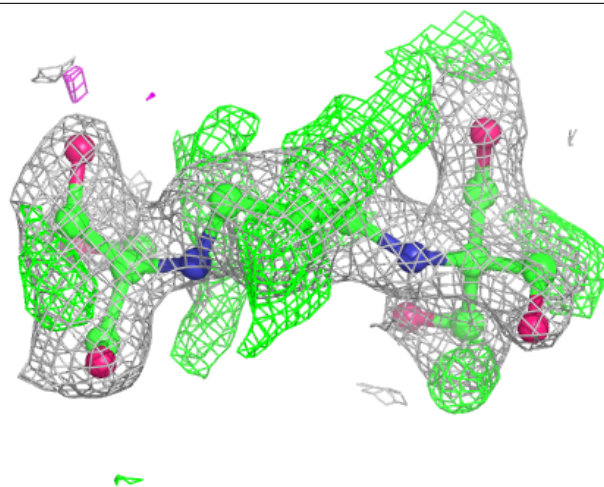
Electron density around B3P C 301:

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



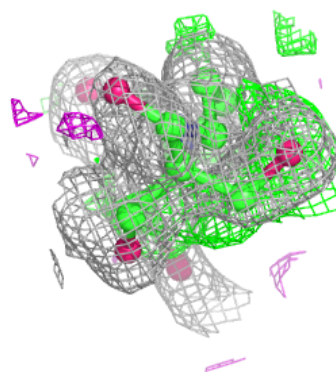
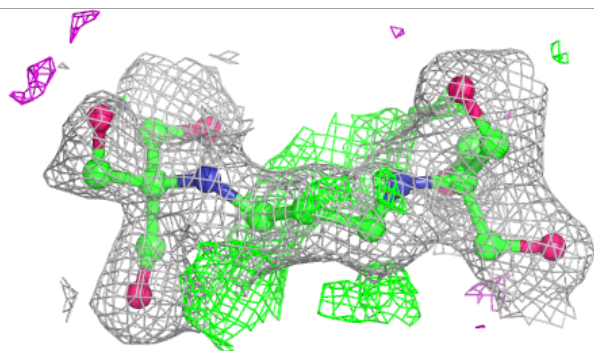
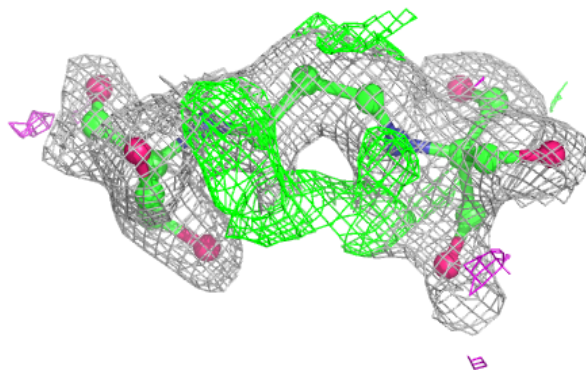
Electron density around B3P D 301:

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

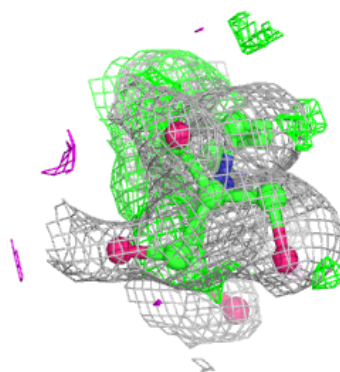
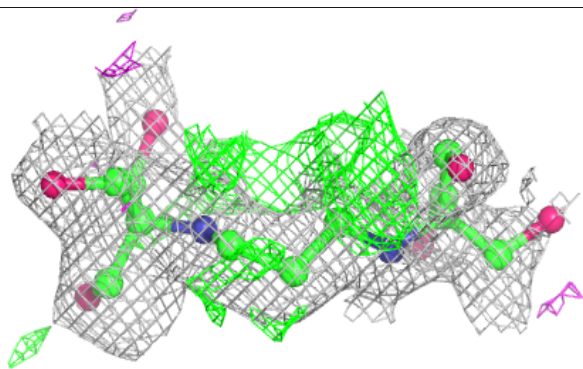
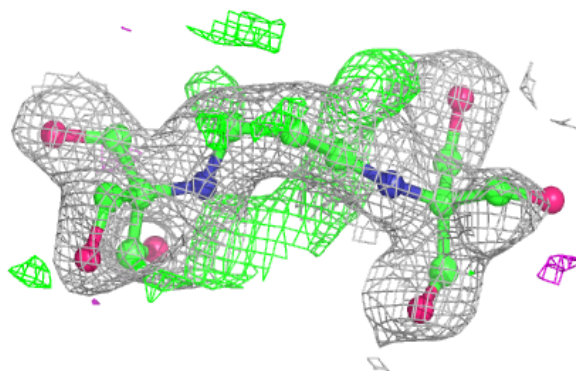


Electron density around B3P E 301 (A):

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

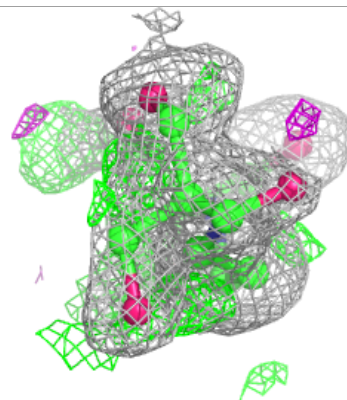
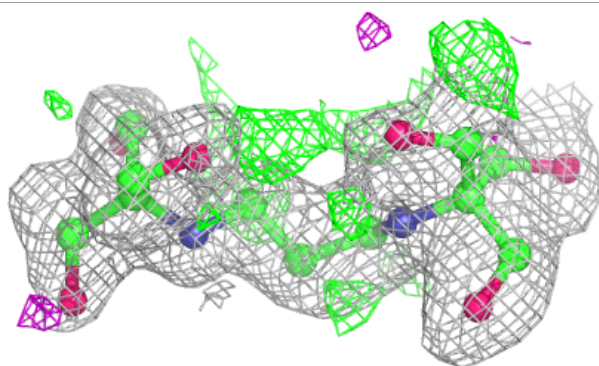
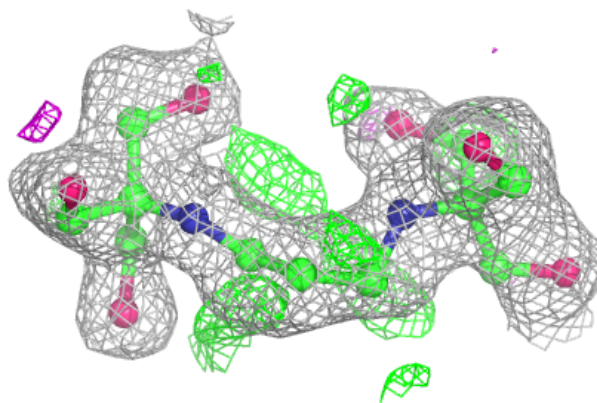
**Electron density around B3P E 301 (B):**

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

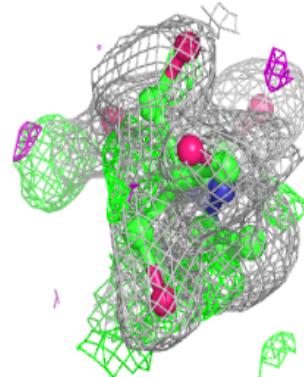
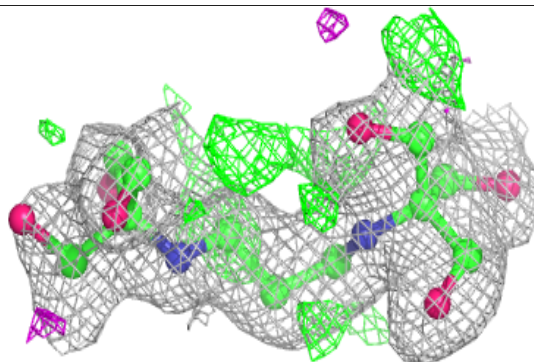
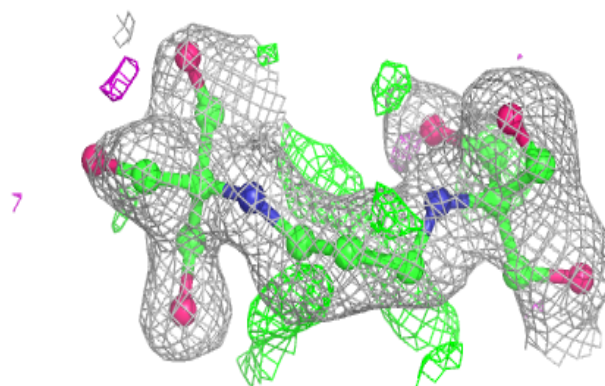


Electron density around B3P A 301 (A):

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

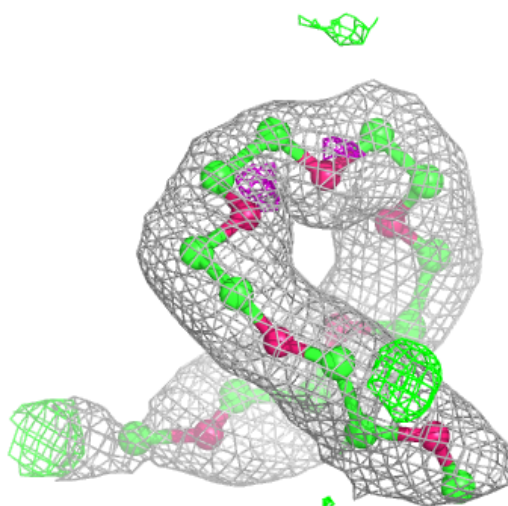
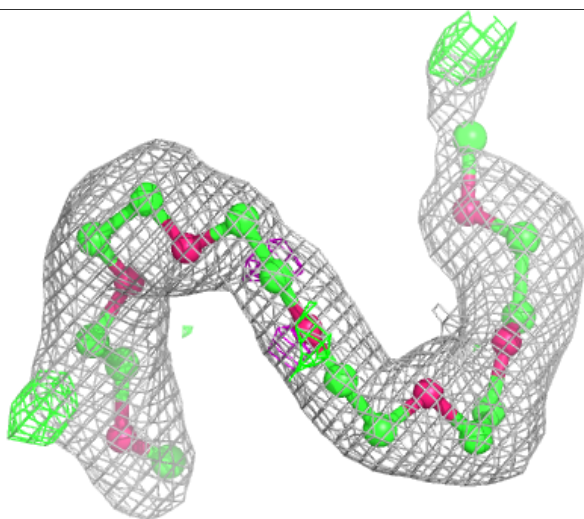
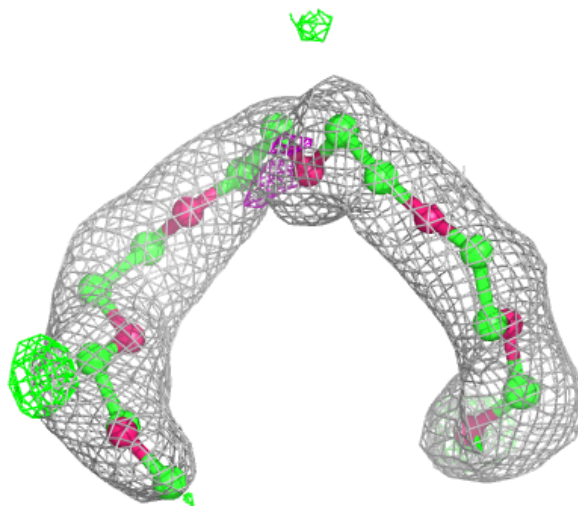
**Electron density around B3P A 301 (B):**

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



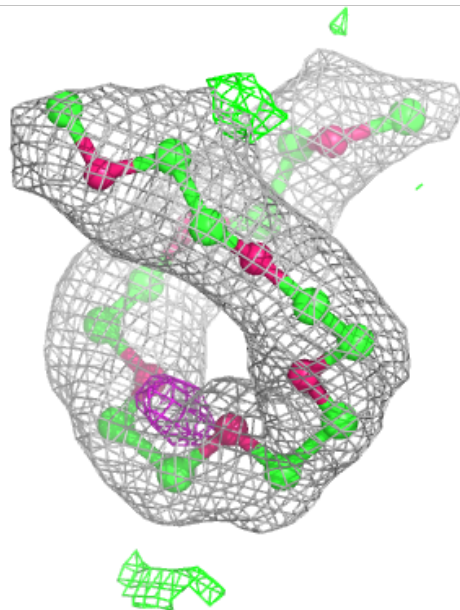
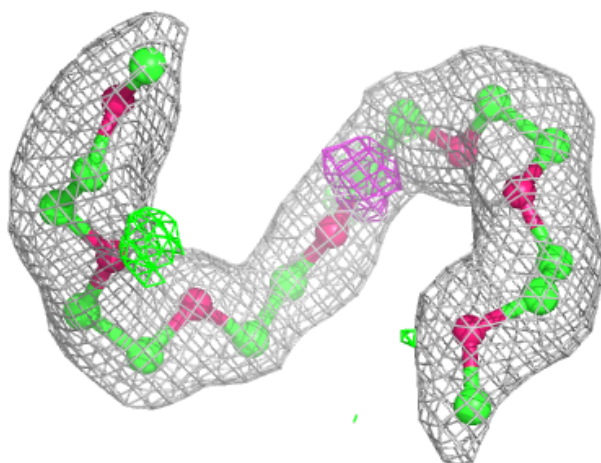
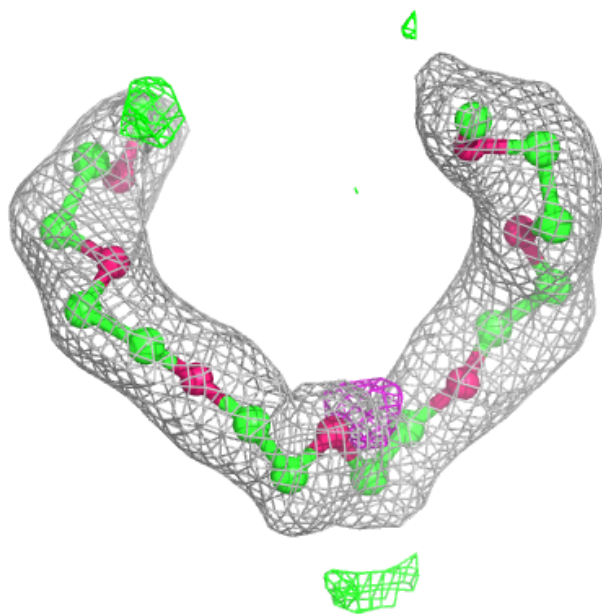
Electron density around PE3 A 300:

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



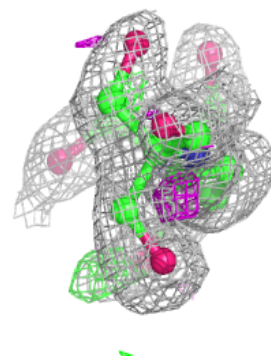
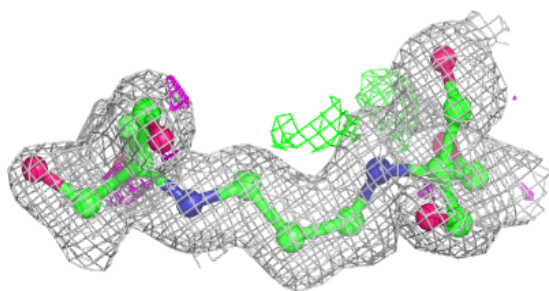
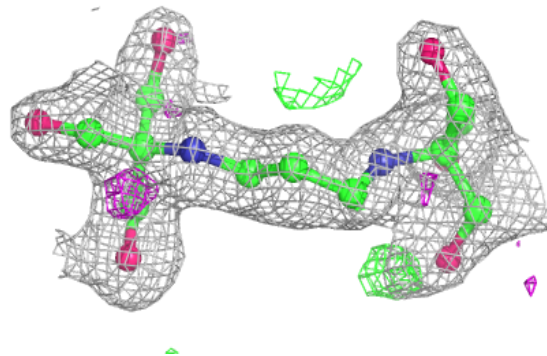
Electron density around PE3 B 300:

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

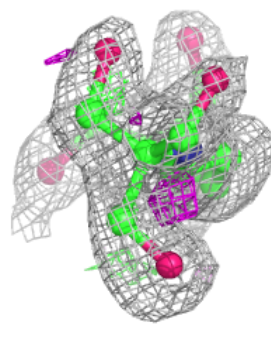
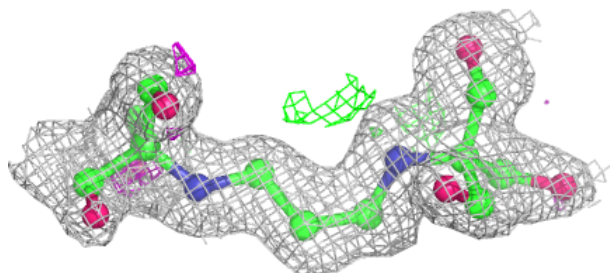
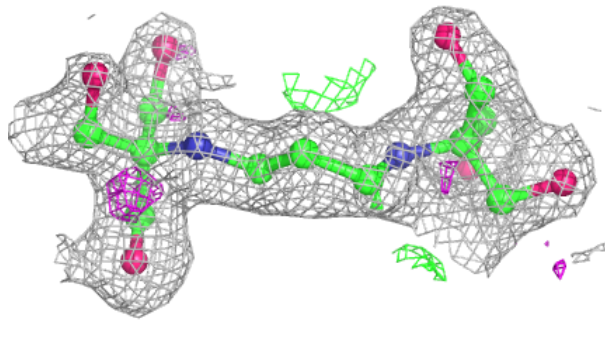


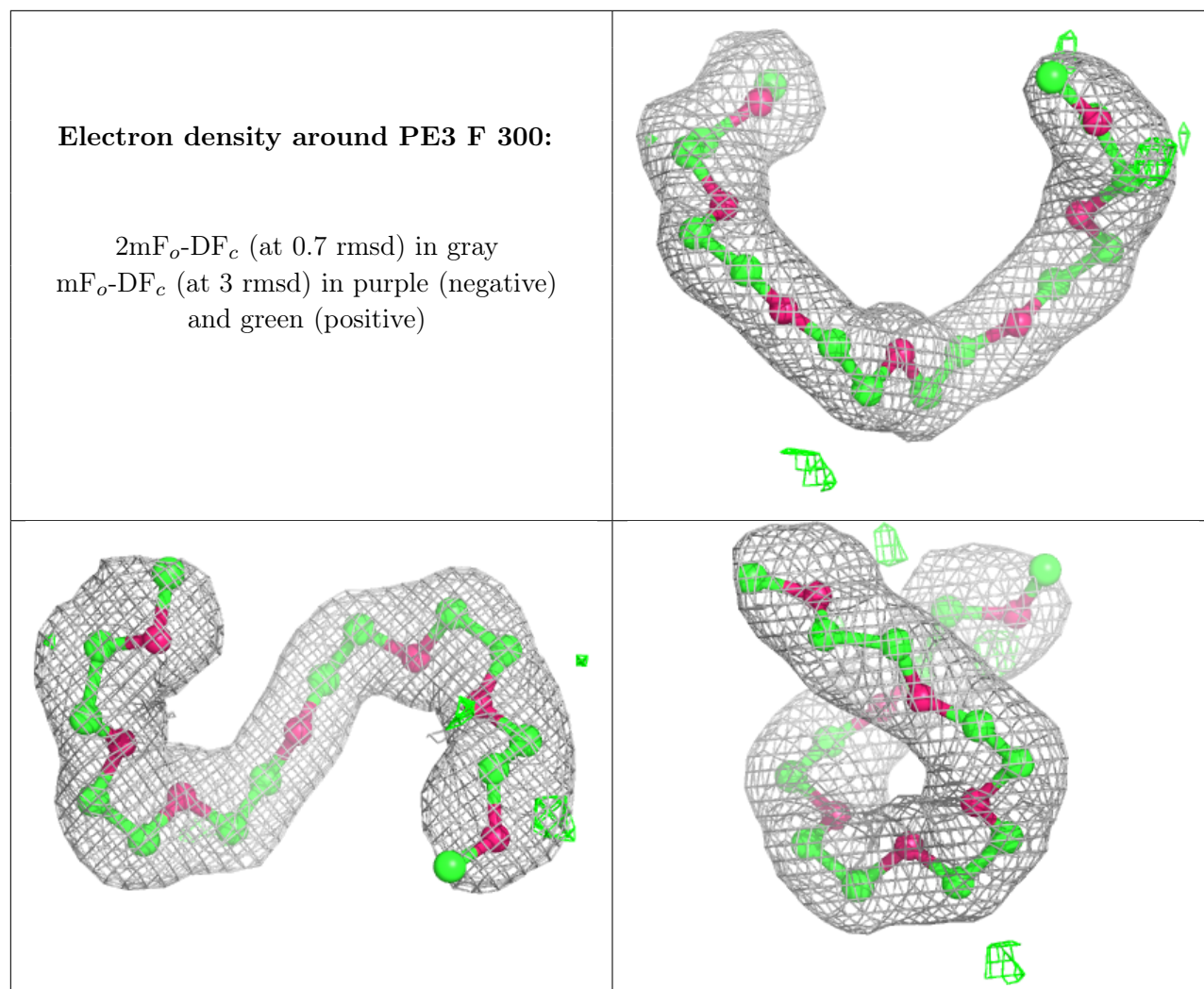
Electron density around B3P F 301 (B):

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

**Electron density around B3P F 301 (A):**

$2mF_o-DF_c$ (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.