



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 12:59 AM BST

PDB ID : 2R4T
Title : Crystal Structure of Wild-type E.coli GS in Complex with ADP and Glucose(wtGSc)
Authors : Sheng, F.; Geiger, J.
Deposited on : 2007-09-01
Resolution : 2.26 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 : 2.13.1
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) : 2.13.1

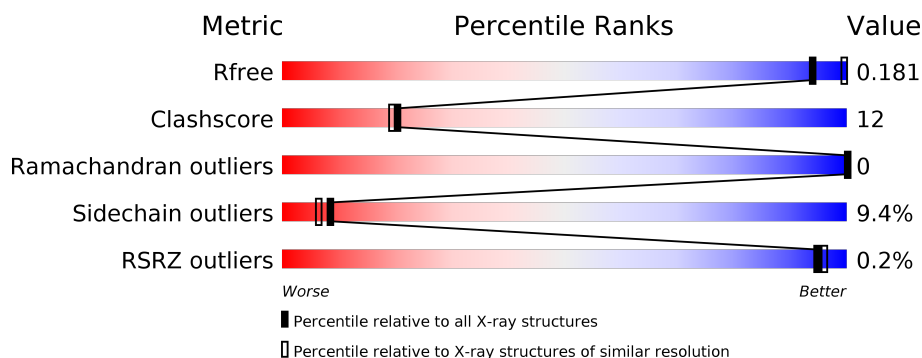
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.26 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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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	485	 78% 17% . .

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	PE3	A	509	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 4223 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.

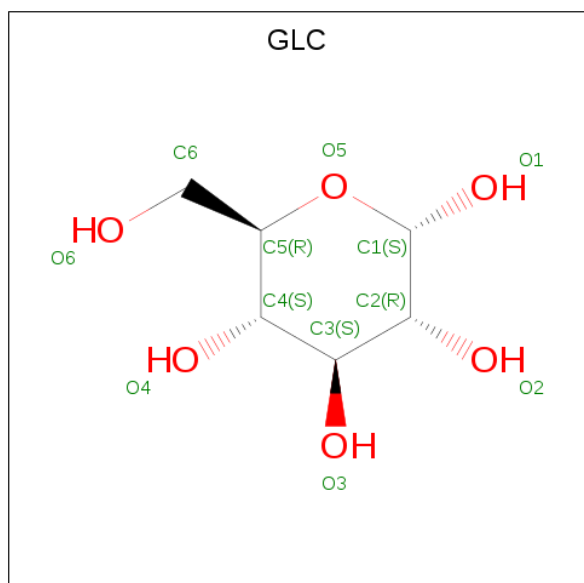
- Molecule 1 is a protein called Glycogen synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	477	3738	2403	652	669	14	0	1	0

There are 8 discrepancies between the modelled and reference sequences:

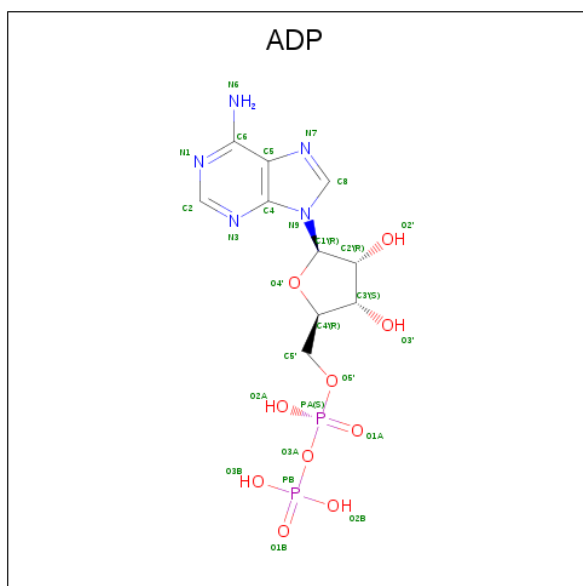
Chain	Residue	Modelled	Actual	Comment	Reference
A	478	LEU	-	expression tag	UNP P0A6U8
A	479	GLU	-	expression tag	UNP P0A6U8
A	480	HIS	-	expression tag	UNP P0A6U8
A	481	HIS	-	expression tag	UNP P0A6U8
A	482	HIS	-	expression tag	UNP P0A6U8
A	483	HIS	-	expression tag	UNP P0A6U8
A	484	HIS	-	expression tag	UNP P0A6U8
A	485	HIS	-	expression tag	UNP P0A6U8

- Molecule 2 is alpha-D-glucopyranose (three-letter code: GLC) (formula: C₆H₁₂O₆).



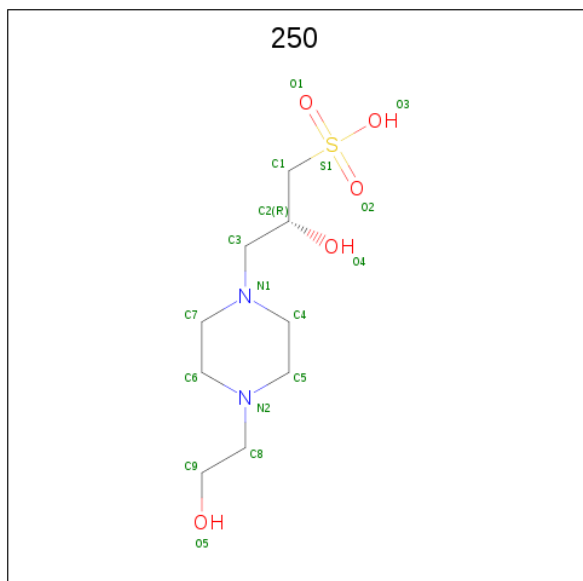
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			12	6	6		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



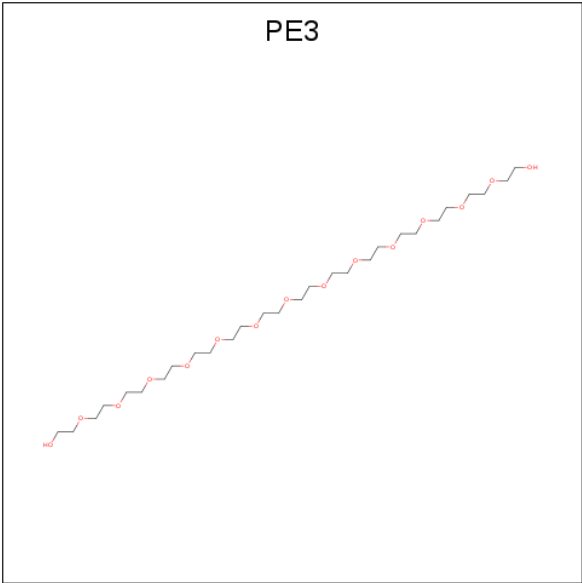
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 4 is (2R)-2-hydroxy-3-[4-(2-hydroxyethyl)piperazin-1-yl]propane-1-sulfonic acid (three-letter code: 250) (formula: $C_9H_{20}N_2O_5S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			17	9	2	5	1		

- Molecule 5 is 3,6,9,12,15,18,21,24,27,30,33,36,39-TRIDECAOXAHENTETRACONTANE-1,41-DIOL (three-letter code: PE3) (formula: C₂₈H₅₈O₁₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			13	9	4		
5	A	1	Total	C	O	0	0
			11	7	4		
5	A	1	Total	C	O	0	0
			8	5	3		
5	A	1	Total	C	O	0	0
			10	7	3		
5	A	1	Total	C	O	0	0
			8	5	3		
5	A	1	Total	C	O	0	0
			4	3	1		
5	A	1	Total	C	O	0	0
			6	4	2		
5	A	1	Total	C	O	0	0
			7	5	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			10	6	4		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			3	2	1		
5	A	1	Total	C	O	0	0
			6	4	2		
5	A	1	Total	C	O	0	0
			4	2	2		

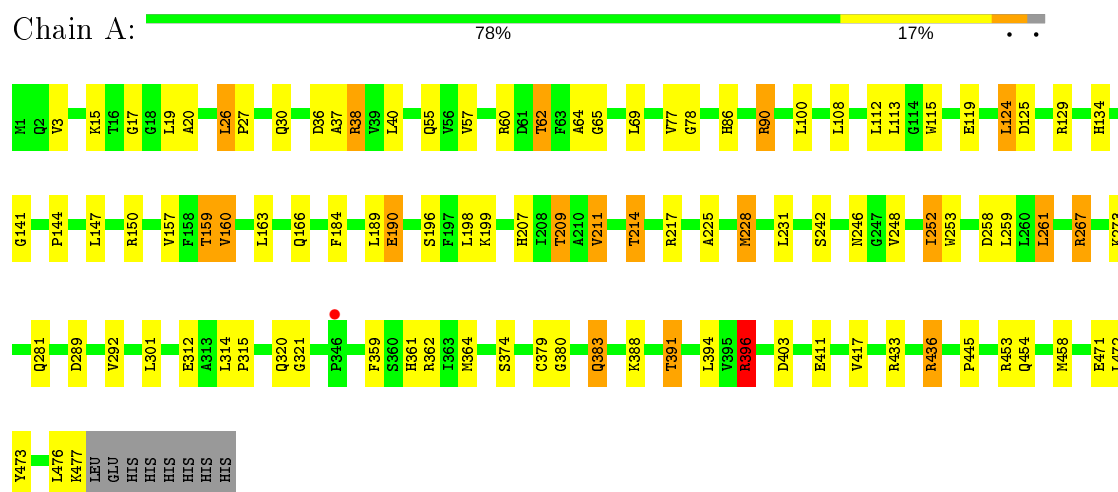
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	335	Total	O	0	0
			335	335		

3 Residue-property plots

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: Glycogen synthase



4 Data and refinement statistics

Property	Value	Source
Space group	I 41	Depositor
Cell constants a, b, c, α , β , γ	126.83Å 126.83Å 151.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	97.13 – 2.26 40.73 – 2.26	Depositor EDS
% Data completeness (in resolution range)	99.8 (97.13-2.26) 99.8 (40.73-2.26)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.09 (at 2.27Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.163 , 0.179 0.164 , 0.181	Depositor DCC
R_{free} test set	2845 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	36.1	Xtriage
Anisotropy	0.318	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 53.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.021 for -h,k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4223	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% 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: 250, GLC, ADP, PE3

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	0/3840	0.62	3/5221 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	396	ARG	NE-CZ-NH2	-6.22	117.19	120.30
1	A	38	ARG	NE-CZ-NH1	5.78	123.19	120.30
1	A	267	ARG	NE-CZ-NH2	-5.20	117.70	120.30

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	3738	0	3684	91	0
2	A	12	0	12	0	0
3	A	27	0	12	0	0
4	A	17	0	20	4	0
5	A	94	0	98	19	0
6	A	335	0	0	19	0
All	All	4223	0	3826	96	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (96) 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:A:36:ASP:OD1	5:A:506:PE3:H81	1.62	0.98
1:A:159:THR:HB	1:A:209:THR:HG22	1.49	0.93
1:A:78:GLY:H	5:A:506:PE3:C6	1.85	0.89
1:A:125:ASP:OD2	5:A:507:PE3:H81	1.73	0.88
1:A:150:ARG:HG3	6:A:674:HOH:O	1.73	0.88
1:A:273:LYS:NZ	1:A:391:THR:HG22	1.90	0.86
1:A:17:GLY:HA3	4:A:503:250:H92	1.59	0.84
1:A:62:THR:HG22	1:A:64:ALA:H	1.43	0.84
1:A:383:GLN:H	1:A:383:GLN:HE21	1.23	0.83
1:A:62:THR:CG2	1:A:64:ALA:H	1.92	0.81
1:A:134:HIS:NE2	1:A:159:THR:CG2	2.45	0.79
1:A:209:THR:CG2	6:A:530:HOH:O	2.31	0.79
1:A:211:VAL:HG13	1:A:379:CYS:HB2	1.67	0.76
1:A:364:MET:O	1:A:391:THR:HG21	1.85	0.75
1:A:38:ARG:HH11	5:A:506:PE3:H92	1.50	0.74
1:A:273:LYS:HZ1	1:A:391:THR:HG22	1.50	0.73
1:A:211:VAL:CG1	1:A:379:CYS:HB2	2.19	0.73
1:A:214:THR:CG2	1:A:403:ASP:OD1	2.37	0.72
1:A:252:ILE:HD11	6:A:526:HOH:O	1.89	0.72
1:A:69:LEU:HD22	1:A:124:LEU:HD13	1.72	0.72
1:A:38:ARG:HD2	5:A:506:PE3:H82	1.72	0.71
1:A:471:GLU:HG3	6:A:701:HOH:O	1.90	0.71
1:A:258:ASP:HB3	1:A:261:LEU:HD22	1.75	0.69
4:A:503:250:H32	6:A:688:HOH:O	1.91	0.69
1:A:163:LEU:HD23	1:A:225:ALA:HA	1.75	0.69
1:A:214:THR:HG23	1:A:403:ASP:CG	2.13	0.69
1:A:477:LYS:HE3	6:A:702:HOH:O	1.92	0.69
1:A:159:THR:HB	1:A:209:THR:CG2	2.24	0.67
1:A:209:THR:HG23	6:A:530:HOH:O	1.91	0.67
1:A:115:TRP:CH2	1:A:119:GLU:HG3	2.29	0.67
1:A:90:ARG:HD3	6:A:550:HOH:O	1.95	0.67
1:A:190:GLU:OE1	6:A:762:HOH:O	2.12	0.66
1:A:214:THR:HG23	1:A:403:ASP:OD1	1.95	0.65
1:A:383:GLN:N	1:A:383:GLN:HE21	1.95	0.65
1:A:380:GLY:C	1:A:383:GLN:HE22	2.01	0.64
1:A:252:ILE:HG13	1:A:253:TRP:CD1	2.33	0.63
1:A:62:THR:HG22	1:A:64:ALA:N	2.12	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:273:LYS:HZ2	1:A:391:THR:HG22	1.64	0.63
1:A:26:LEU:HD22	1:A:30:GLN:HG3	1.81	0.62
1:A:252:ILE:CD1	6:A:526:HOH:O	2.49	0.61
1:A:166:GLN:NE2	6:A:772:HOH:O	2.33	0.59
1:A:57:VAL:HG22	5:A:507:PE3:H112	1.84	0.59
1:A:199:LYS:HG2	1:A:228:MET:HE1	1.85	0.58
1:A:477:LYS:NZ	6:A:760:HOH:O	2.26	0.58
1:A:62:THR:HB	1:A:65:GLY:O	2.04	0.57
1:A:134:HIS:NE2	1:A:159:THR:HG21	2.20	0.56
1:A:312:GLU:OE2	6:A:828:HOH:O	2.18	0.56
5:A:507:PE3:H51	6:A:707:HOH:O	2.07	0.55
1:A:476:LEU:O	1:A:477:LYS:HB2	2.07	0.54
1:A:36:ASP:OD1	5:A:506:PE3:C8	2.46	0.54
1:A:57:VAL:HG13	5:A:507:PE3:C11	2.38	0.54
1:A:62:THR:HG23	1:A:64:ALA:H	1.72	0.52
1:A:214:THR:HB	1:A:217:ARG:NH1	2.24	0.52
1:A:383:GLN:H	1:A:383:GLN:NE2	2.00	0.52
1:A:57:VAL:HG13	5:A:507:PE3:H111	1.92	0.51
1:A:214:THR:HG23	1:A:403:ASP:OD2	2.11	0.51
1:A:477:LYS:HG2	6:A:836:HOH:O	2.11	0.51
1:A:267:ARG:HD2	1:A:453:ARG:O	2.12	0.50
1:A:60:ARG:NH1	6:A:720:HOH:O	2.44	0.50
1:A:3:VAL:HG21	1:A:473:TYR:CE2	2.47	0.49
1:A:196:SER:HB3	1:A:199:LYS:HB2	1.96	0.48
1:A:273:LYS:HZ1	1:A:391:THR:CG2	2.24	0.48
1:A:374:SER:O	1:A:396:ARG:HD3	2.13	0.48
1:A:477:LYS:CG	6:A:836:HOH:O	2.61	0.48
1:A:62:THR:HG21	1:A:112:LEU:HD13	1.95	0.48
1:A:36:ASP:HA	5:A:514:PE3:H122	1.95	0.48
1:A:38:ARG:HG3	5:A:506:PE3:C6	2.44	0.47
1:A:37:ALA:O	5:A:506:PE3:C6	2.63	0.47
1:A:228:MET:HE2	1:A:228:MET:HB3	1.64	0.47
1:A:252:ILE:HD12	1:A:361:HIS:CE1	2.51	0.46
1:A:26:LEU:HB3	1:A:27:PRO:HD3	1.98	0.46
1:A:211:VAL:HG22	1:A:246:ASN:HD21	1.81	0.45
1:A:86:HIS:HB2	5:A:512:PE3:H21	1.98	0.45
1:A:184:PHE:CE2	1:A:190:GLU:HG3	2.52	0.45
1:A:320:GLN:HE22	1:A:436:ARG:HG2	1.81	0.44
1:A:454:GLN:OE1	1:A:458:MET:CE	2.65	0.44
4:A:503:250:H42	4:A:503:250:H2	1.64	0.44
1:A:38:ARG:CD	5:A:506:PE3:H82	2.45	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:207:HIS:HE1	1:A:242:SER:OG	2.01	0.43
1:A:57:VAL:HG13	5:A:507:PE3:H112	2.00	0.42
1:A:15:LYS:HG3	1:A:20:ALA:HB3	2.01	0.42
1:A:454:GLN:OE1	1:A:458:MET:HE3	2.20	0.42
1:A:359:PHE:HD1	1:A:362:ARG:NH1	2.18	0.42
1:A:38:ARG:HD2	5:A:506:PE3:C8	2.47	0.41
1:A:160:VAL:HG21	1:A:163:LEU:HD13	2.02	0.41
1:A:445:PRO:HG3	5:A:504:PE3:H202	2.02	0.41
1:A:184:PHE:CZ	1:A:190:GLU:HG3	2.56	0.41
1:A:129:ARG:HD2	6:A:787:HOH:O	2.20	0.41
1:A:292:VAL:HG13	1:A:321:GLY:O	2.20	0.41
1:A:433:ARG:HA	1:A:433:ARG:HD2	1.80	0.41
1:A:141:GLY:O	1:A:144:PRO:HD2	2.21	0.41
1:A:394:LEU:C	1:A:394:LEU:HD23	2.42	0.40
4:A:503:250:H61	4:A:503:250:H91	1.62	0.40
5:A:504:PE3:H141	6:A:738:HOH:O	2.21	0.40
1:A:314:LEU:HB3	1:A:315:PRO:HD3	2.04	0.40
1:A:281:GLN:HE22	1:A:289:ASP:H	1.69	0.40

There are no symmetry-related clashes.

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	476/485 (98%)	467 (98%)	9 (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	383/390 (98%)	347 (91%)	36 (9%)	8 6

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	26	LEU
1	A	40	LEU
1	A	55	GLN
1	A	62	THR
1	A	77	VAL
1	A	90	ARG
1	A	100	LEU
1	A	108	LEU
1	A	113	LEU
1	A	124	LEU
1	A	147	LEU
1	A	157	VAL
1	A	159	THR
1	A	160	VAL
1	A	189	LEU
1	A	190	GLU
1	A	198	LEU
1	A	209	THR
1	A	211	VAL
1	A	214	THR
1	A	228	MET
1	A	231	LEU
1	A	248	VAL
1	A	252	ILE
1	A	259	LEU
1	A	261	LEU
1	A	301	LEU
1	A	383	GLN
1	A	388	LYS
1	A	391	THR
1	A	396	ARG
1	A	411	GLU
1	A	417	VAL

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Mol	Chain	Res	Type
1	A	436	ARG
1	A	472	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	GLN
1	A	96	HIS
1	A	192	ASN
1	A	207	HIS
1	A	223	GLN
1	A	233	GLN
1	A	246	ASN
1	A	279	GLN
1	A	281	GLN
1	A	320	GLN
1	A	383	GLN

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](#)

16 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	$\# Z > 2$
5	PE3	A	514	-	5,5,42	0.48	0	4,4,41	0.17	0
2	GLC	A	502	-	12,12,12	0.55	0	17,17,17	1.03	1 (5%)
5	PE3	A	507	-	9,9,42	0.54	0	8,8,41	0.47	0
5	PE3	A	511	-	3,3,42	0.43	0	2,2,41	0.30	0
5	PE3	A	515	-	3,3,42	0.45	0	2,2,41	0.27	0
5	PE3	A	506	-	7,7,42	0.44	0	6,6,41	0.30	0
5	PE3	A	512	-	9,9,42	0.47	0	8,8,41	0.29	0
5	PE3	A	508	-	7,7,42	0.74	0	6,6,41	0.52	0
5	PE3	A	505	-	10,10,42	0.48	0	9,9,41	0.17	0
5	PE3	A	519	-	5,5,42	0.51	0	4,4,41	0.25	0
5	PE3	A	509	-	3,3,42	0.55	0	2,2,41	0.03	0
5	PE3	A	510	-	6,6,42	0.53	0	5,5,41	0.14	0
5	PE3	A	504	-	12,12,42	0.53	0	11,11,41	0.16	0
5	PE3	A	513	-	2,2,42	0.48	0	1,1,41	0.33	0
4	250	A	503	-	17,17,17	0.89	1 (5%)	21,23,23	2.99	11 (52%)
3	ADP	A	501	-	24,29,29	0.82	1 (4%)	29,45,45	1.28	1 (3%)

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
5	PE3	A	514	-	-	3/3/3/40	-
2	GLC	A	502	-	-	0/2/22/22	0/1/1/1
5	PE3	A	507	-	-	5/7/7/40	-
5	PE3	A	511	-	-	0/1/1/40	-
5	PE3	A	515	-	-	1/1/1/40	-
5	PE3	A	506	-	-	3/5/5/40	-
5	PE3	A	512	-	-	4/7/7/40	-
5	PE3	A	508	-	-	1/5/5/40	-
5	PE3	A	505	-	-	4/8/8/40	-
5	PE3	A	519	-	-	2/3/3/40	-
5	PE3	A	509	-	-	0/0/1/40	-
5	PE3	A	510	-	-	3/4/4/40	-
5	PE3	A	504	-	-	7/10/10/40	-
4	250	A	503	-	-	3/12/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ADP	A	501	-	-	4/12/32/32	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	503	250	C1-S1	-2.78	1.67	1.77
3	A	501	ADP	C2-N3	2.18	1.35	1.32

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	503	250	C4-N1-C7	5.32	120.81	108.83
4	A	503	250	C2-C3-N1	-5.17	102.85	112.23
4	A	503	250	C5-N2-C6	5.04	120.18	108.83
3	A	501	ADP	N3-C2-N1	-4.92	120.98	128.68
4	A	503	250	C4-C5-N2	-4.58	101.25	110.64
4	A	503	250	O1-S1-C1	4.32	112.07	106.94
4	A	503	250	C8-N2-C6	4.03	121.53	111.23
4	A	503	250	C5-C4-N1	-3.39	103.70	110.64
4	A	503	250	C7-N1-C3	3.36	120.96	111.20
4	A	503	250	C8-N2-C5	2.75	118.27	111.23
2	A	502	GLC	O5-C1-C2	2.60	114.93	110.28
4	A	503	250	C4-N1-C3	2.42	118.22	111.20
4	A	503	250	C9-C8-N2	-2.07	105.95	113.40

There are no chirality outliers.

All (40) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	503	250	C9-C8-N2-C6
4	A	503	250	N2-C8-C9-O5
3	A	501	ADP	PB-O3A-PA-O5'
5	A	504	PE3	C18-C17-O16-C15
5	A	504	PE3	O13-C14-C15-O16
5	A	507	PE3	C9-C8-O7-C6
5	A	508	PE3	O4-C5-C6-O7
5	A	506	PE3	O10-C11-C12-O13
5	A	505	PE3	O1-C2-C3-O4
5	A	505	PE3	O7-C8-C9-O10
3	A	501	ADP	C3'-C4'-C5'-O5'
5	A	519	PE3	O1-C2-C3-O4

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Mol	Chain	Res	Type	Atoms
5	A	506	PE3	O7-C8-C9-O10
4	A	503	250	C2-C3-N1-C7
3	A	501	ADP	O4'-C4'-C5'-O5'
5	A	507	PE3	O4-C5-C6-O7
5	A	504	PE3	O10-C11-C12-O13
5	A	512	PE3	O7-C8-C9-O10
5	A	514	PE3	O10-C11-C12-O13
5	A	507	PE3	C6-C5-O4-C3
5	A	506	PE3	C9-C8-O7-C6
5	A	515	PE3	O1-C2-C3-O4
3	A	501	ADP	PA-O3A-PB-O1B
5	A	514	PE3	C12-C11-O10-C9
5	A	514	PE3	C11-C12-O13-C14
5	A	512	PE3	C2-C3-O4-C5
5	A	507	PE3	C12-C11-O10-C9
5	A	504	PE3	C21-C20-O19-C18
5	A	519	PE3	C6-C5-O4-C3
5	A	505	PE3	O4-C5-C6-O7
5	A	505	PE3	C8-C9-O10-C11
5	A	510	PE3	C2-C3-O4-C5
5	A	510	PE3	C6-C5-O4-C3
5	A	507	PE3	C8-C9-O10-C11
5	A	504	PE3	C15-C14-O13-C12
5	A	512	PE3	O4-C5-C6-O7
5	A	510	PE3	C5-C6-O7-C8
5	A	512	PE3	C5-C6-O7-C8
5	A	504	PE3	C17-C18-O19-C20
5	A	504	PE3	O16-C17-C18-O19

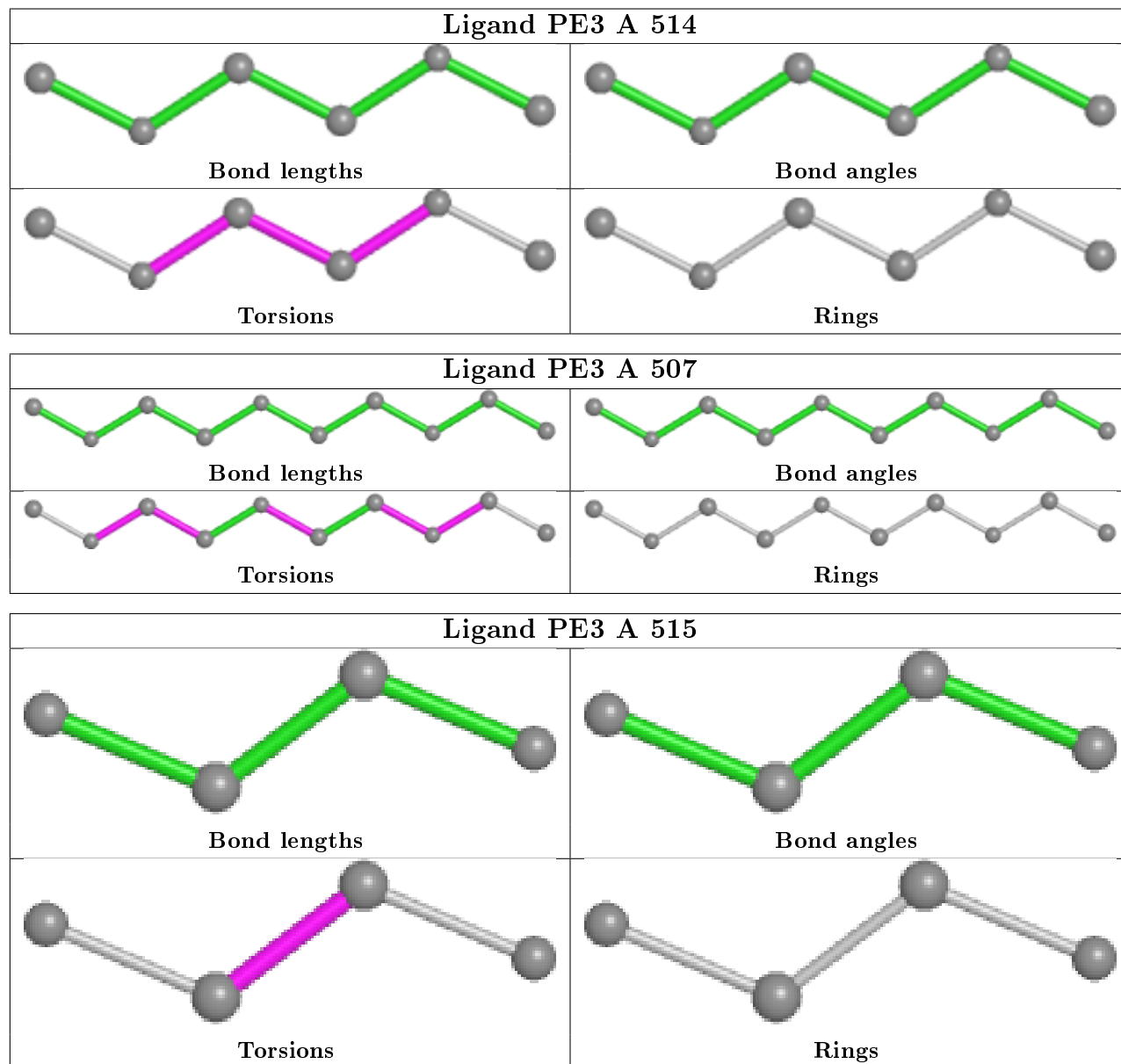
There are no ring outliers.

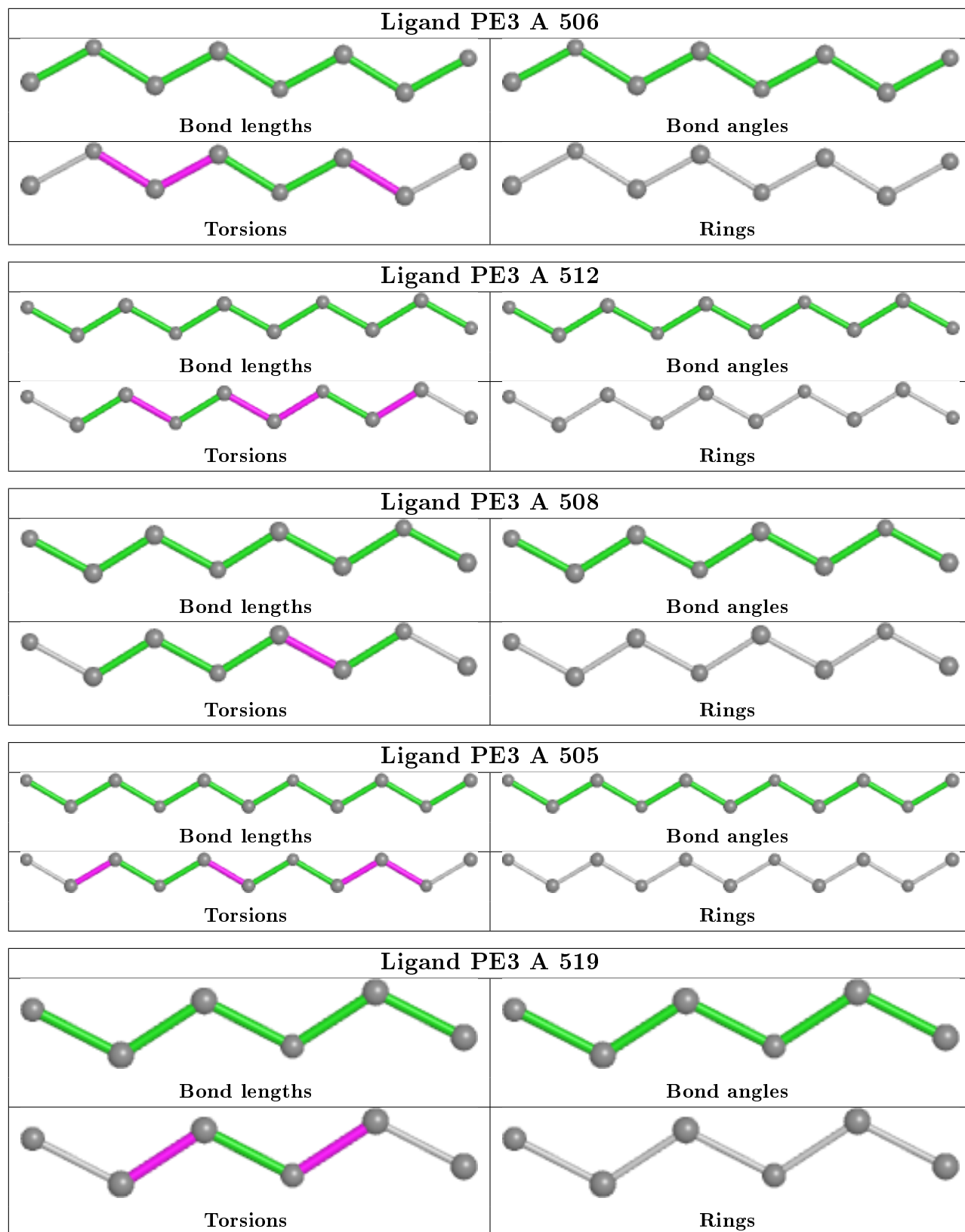
6 monomers are involved in 23 short contacts:

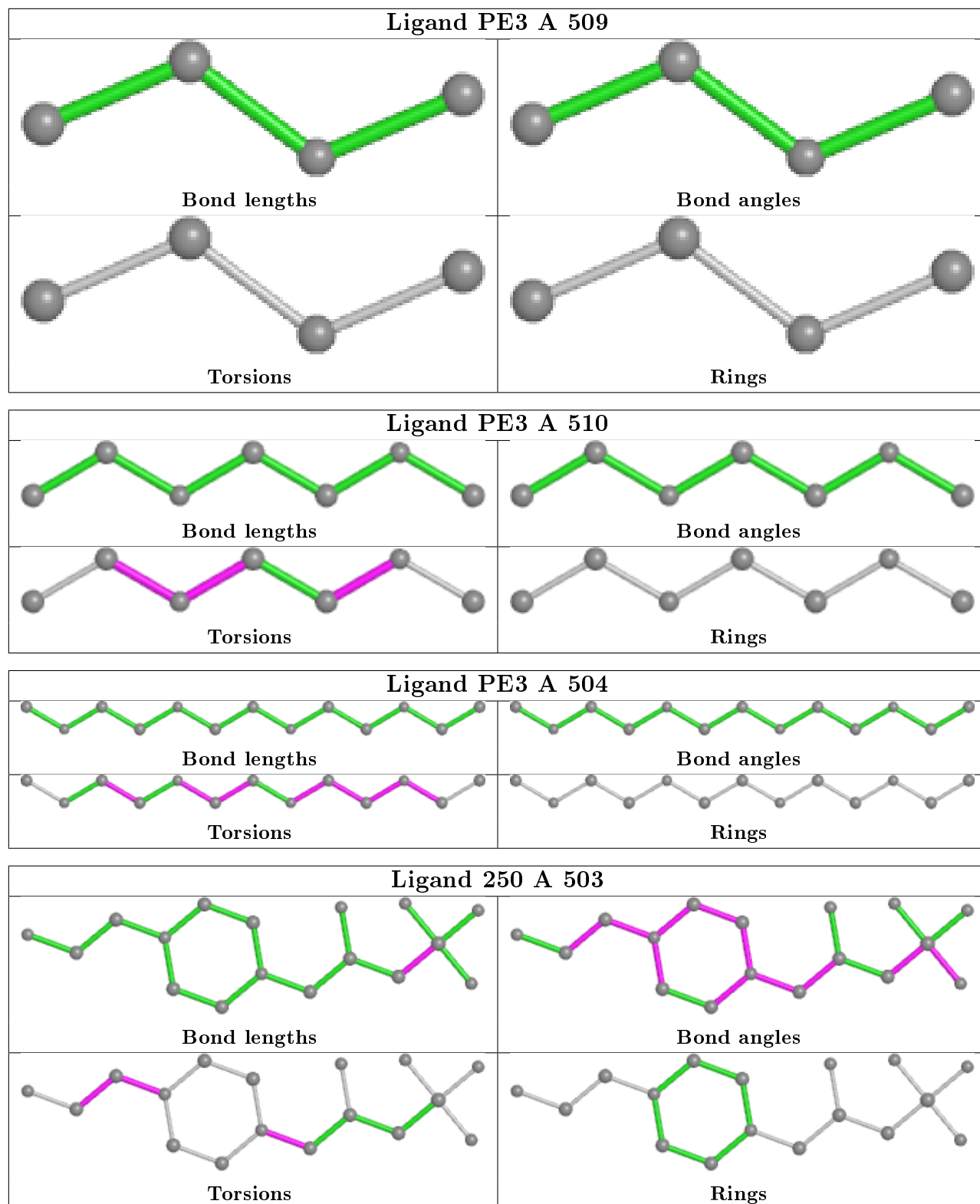
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	514	PE3	1	0
5	A	507	PE3	6	0
5	A	506	PE3	9	0
5	A	512	PE3	1	0
5	A	504	PE3	2	0
4	A	503	250	4	0

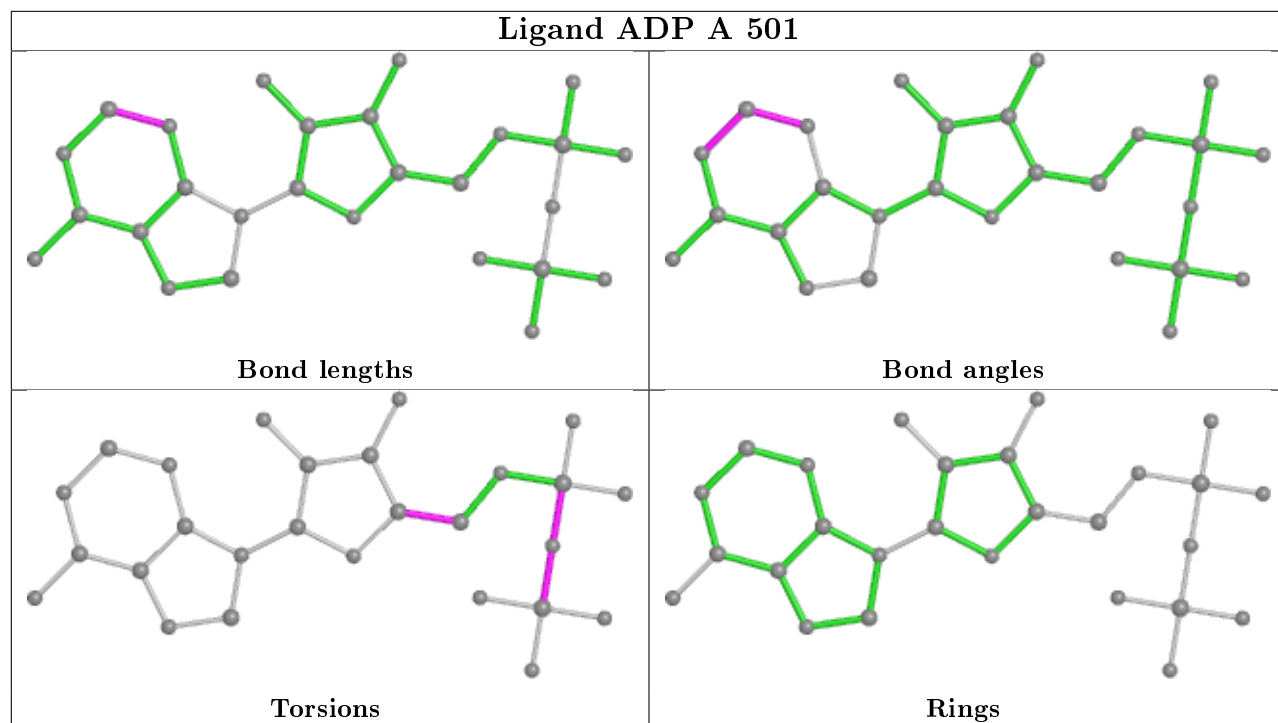
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

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	477/485 (98%)	-0.42	1 (0%) 95 96	26, 35, 50, 59	30 (6%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	346	PRO	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 monosaccharides 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.

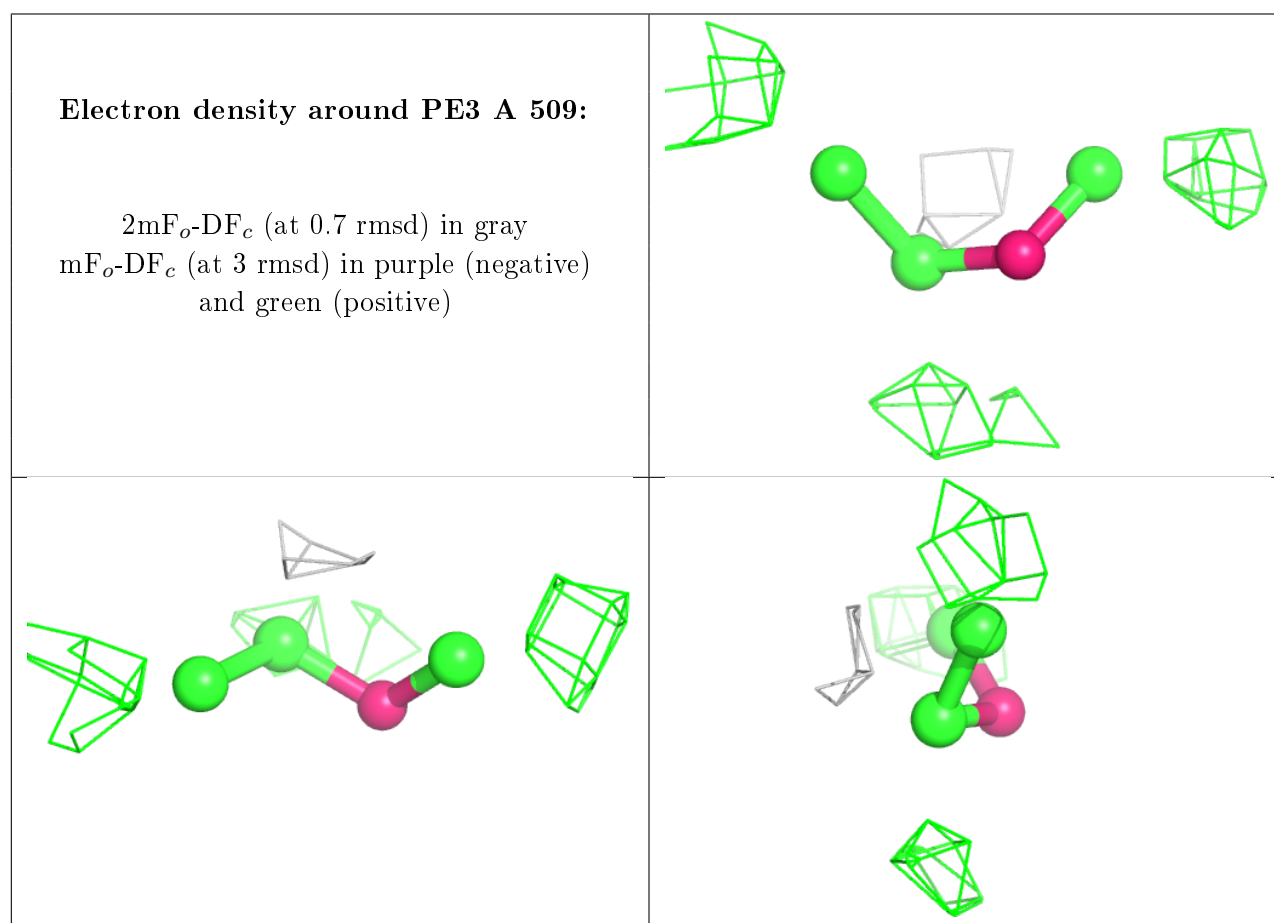
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	PE3	A	509	4/43	0.26	0.76	165,165,165,165	0
5	PE3	A	510	7/43	0.64	0.25	90,91,91,91	0
5	PE3	A	515	4/43	0.69	0.25	79,79,80,80	0
5	PE3	A	514	6/43	0.76	0.25	75,76,76,76	0
5	PE3	A	519	6/43	0.78	0.21	78,78,78,78	0
5	PE3	A	504	13/43	0.78	0.27	73,73,75,75	0

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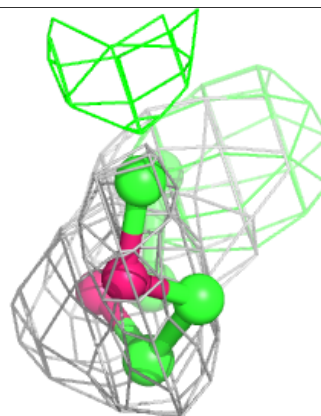
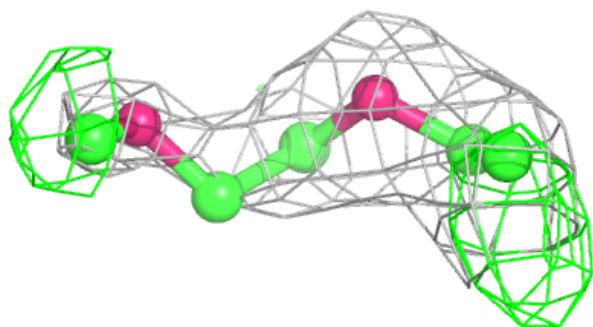
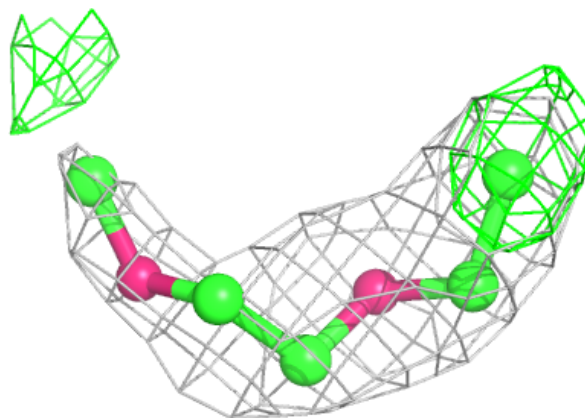
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	PE3	A	506	8/43	0.79	0.24	45,50,53,53	0
5	PE3	A	512	10/43	0.79	0.26	82,83,83,83	0
5	PE3	A	505	11/43	0.81	0.16	72,73,73,74	0
5	PE3	A	511	4/43	0.82	0.19	70,70,70,70	0
5	PE3	A	508	8/43	0.83	0.24	70,70,71,71	3
5	PE3	A	507	10/43	0.84	0.29	42,44,44,44	2
5	PE3	A	513	3/43	0.89	0.14	71,71,72,72	0
4	250	A	503	17/17	0.91	0.17	66,68,69,69	0
2	GLC	A	502	12/12	0.93	0.21	26,28,30,31	12
3	ADP	A	501	27/27	0.99	0.09	31,32,34,34	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.



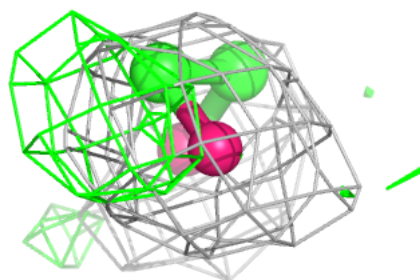
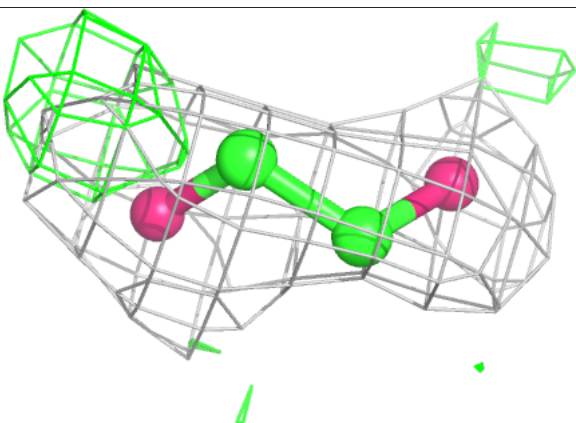
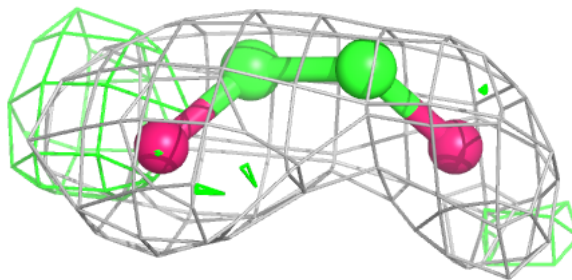
Electron density around PE3 A 510:

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

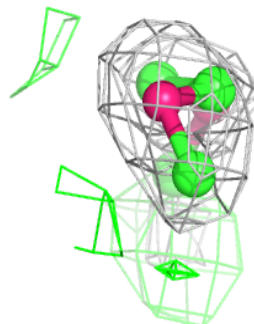
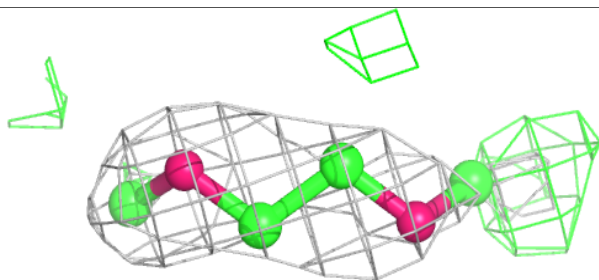
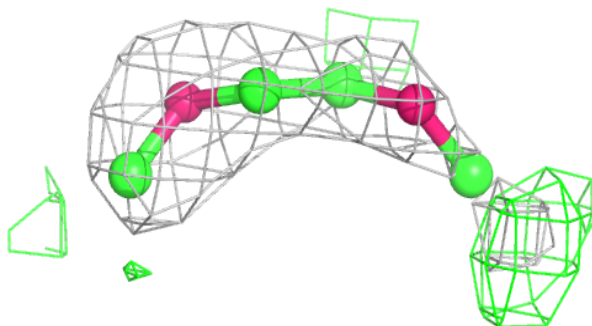


Electron density around PE3 A 515:

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

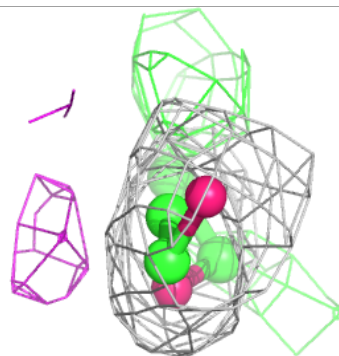
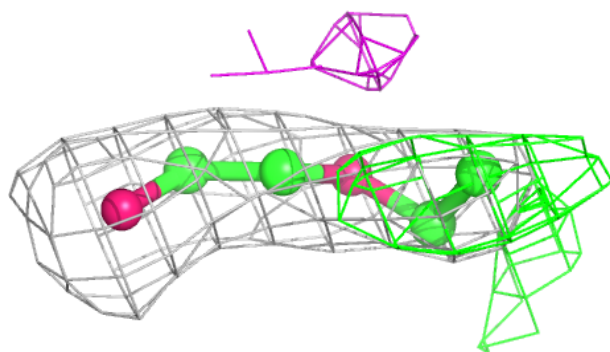
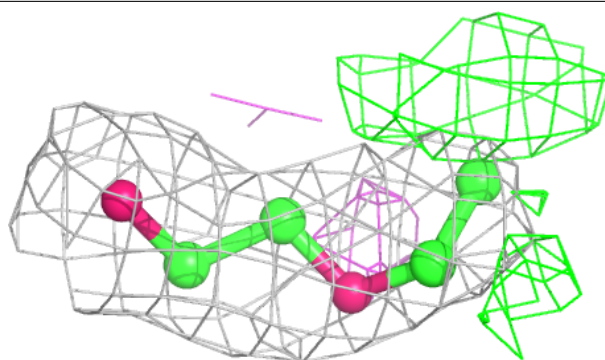
**Electron density around PE3 A 514:**

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

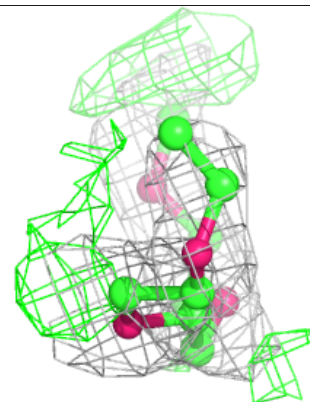
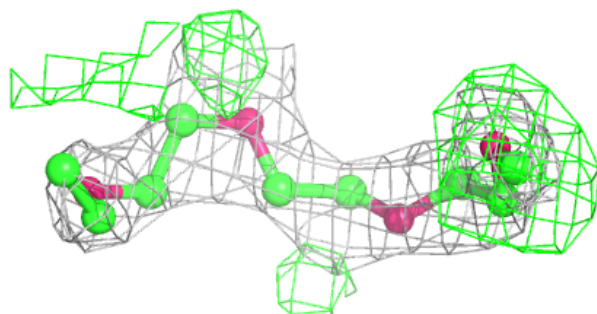
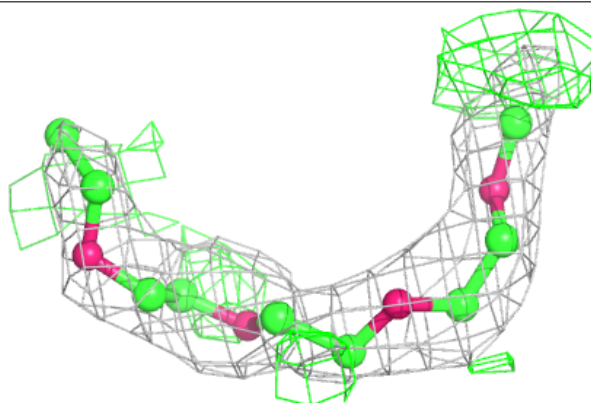


Electron density around PE3 A 519:

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

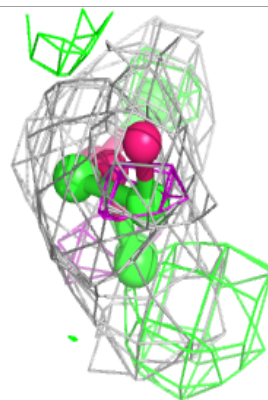
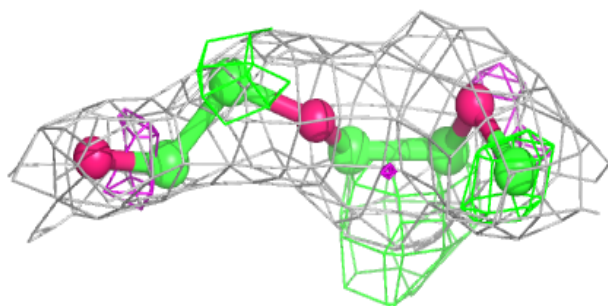
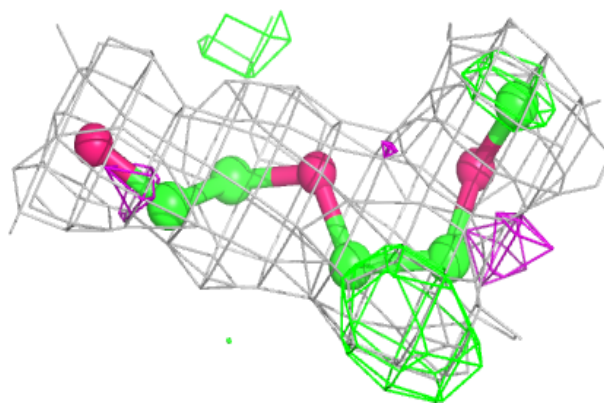
**Electron density around PE3 A 504:**

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

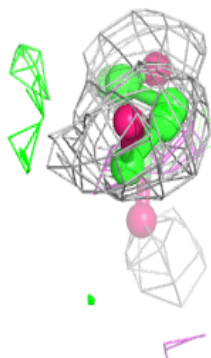
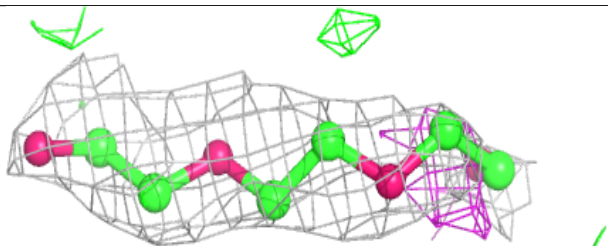
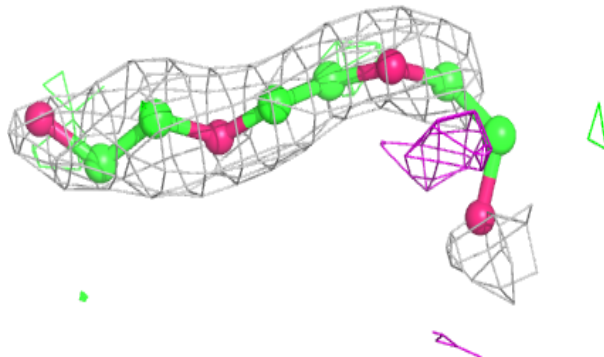


Electron density around PE3 A 506:

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

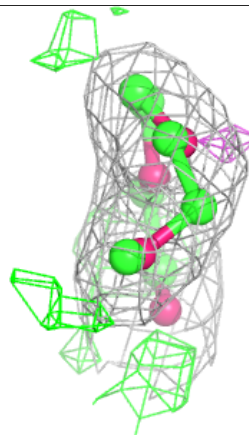
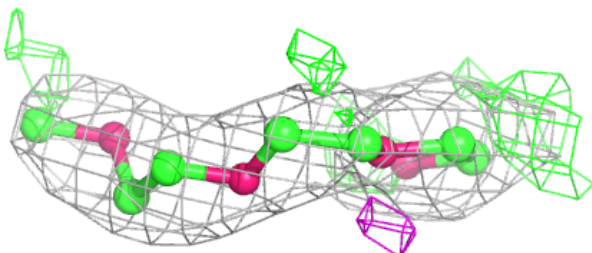
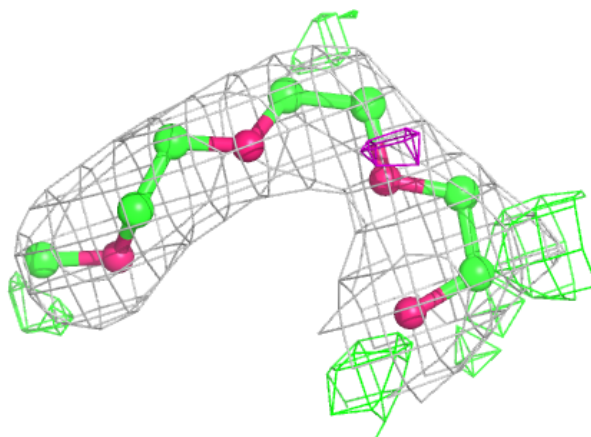
**Electron density around PE3 A 512:**

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

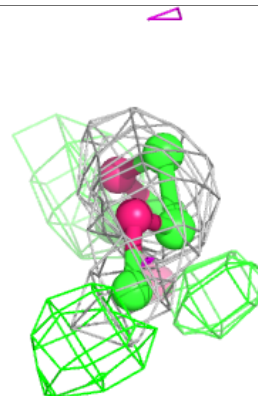
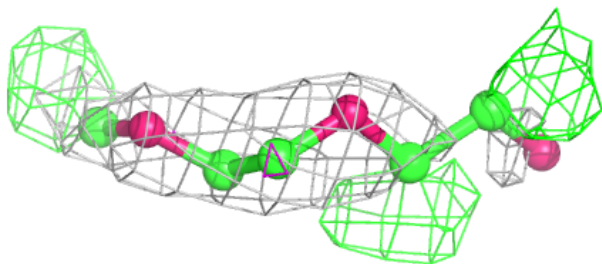
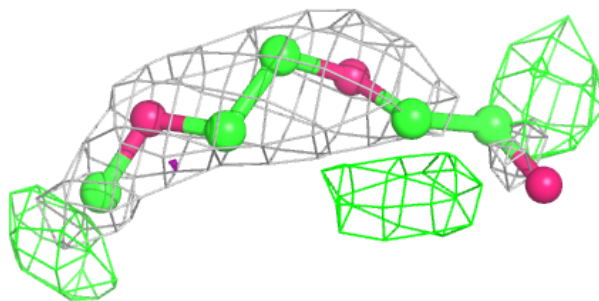


Electron density around PE3 A 505:

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and green (positive)

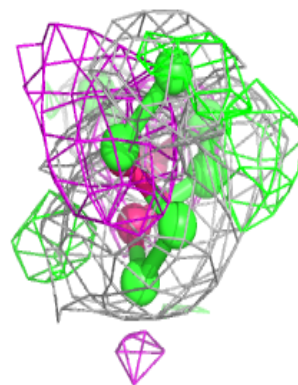
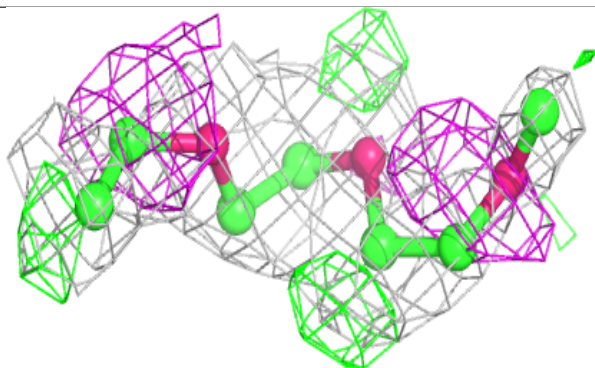
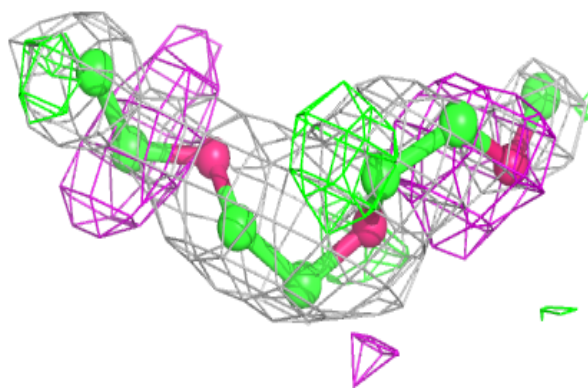
**Electron density around PE3 A 508:**

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

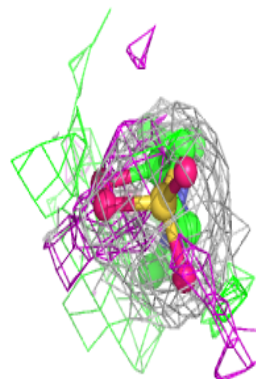
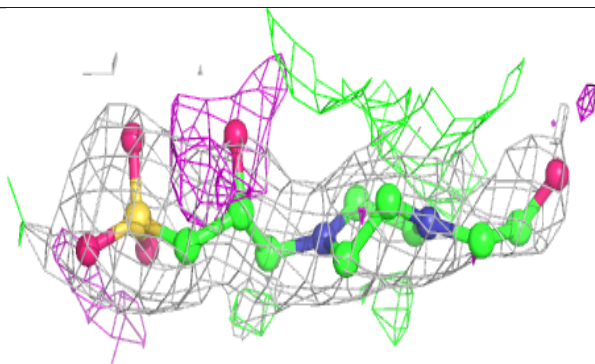
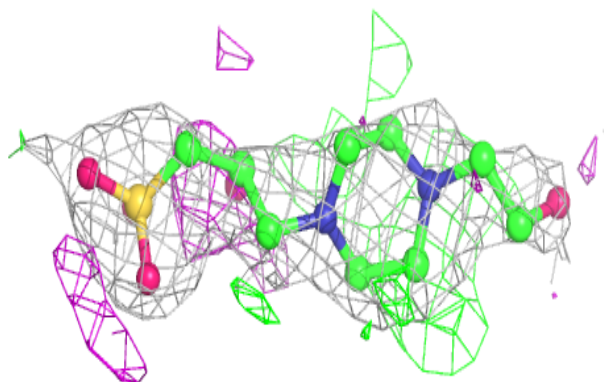


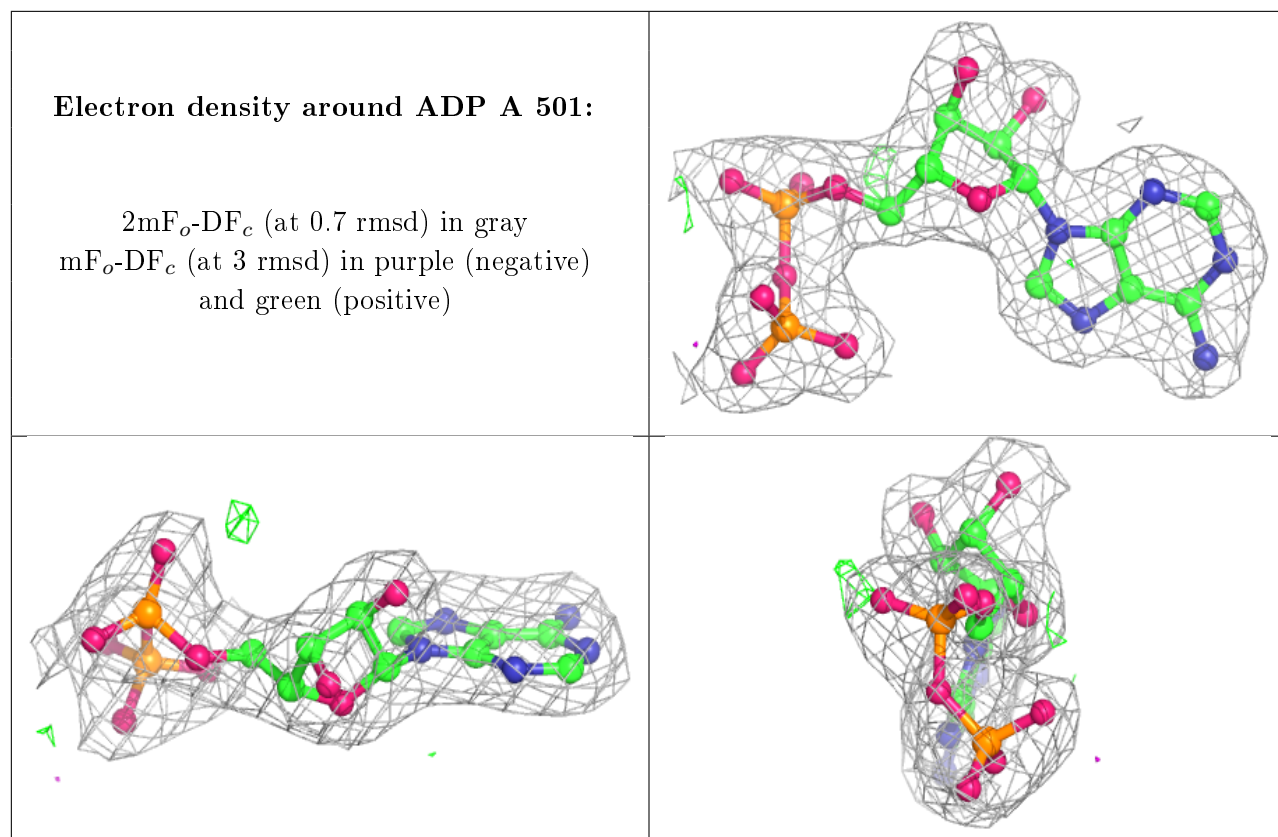
Electron density around PE3 A 507:

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

**Electron density around 250 A 503:**

$2mF_o-DF_c$ (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.