



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:31 AM UTC

PDB ID : 9UC2 / pdb\_00009uc2  
Title : Crystal structure of voltage-gated sodium channel NavAb N49K/S178A mutant  
Authors : Irie, K.; Maeda, Y.; Kojima, K.  
Deposited on : 2025-04-03  
Resolution : 2.80 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

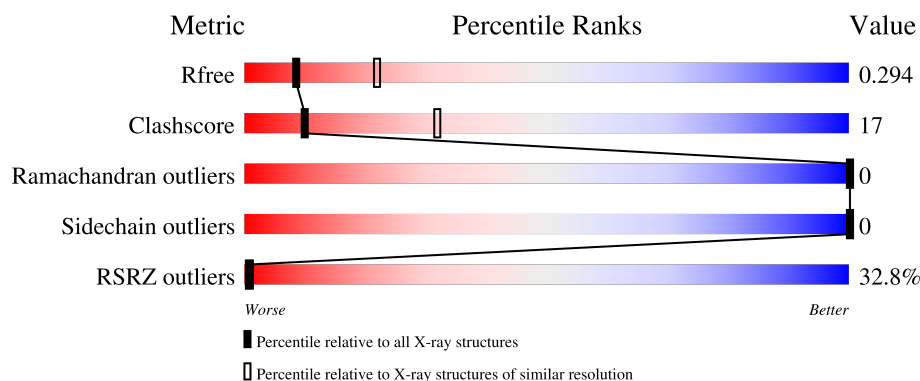
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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  $\geq 3$ , 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	271	<div> <div>28%</div> <div>58%</div> <div>24%</div> <div>•</div> <div>15%</div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 2346 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 Ion transport protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	229	Total	C	N	O	S	0	0	0
			1869	1273	281	303	12			

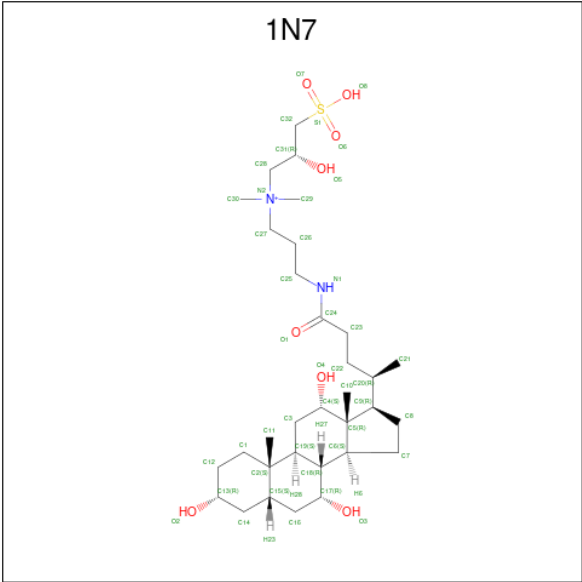
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	997	GLY	-	expression tag	UNP A8EVM5
A	998	SER	-	expression tag	UNP A8EVM5
A	999	GLY	-	expression tag	UNP A8EVM5
A	1000	SER	-	expression tag	UNP A8EVM5
A	1049	LYS	ASN	engineered mutation	UNP A8EVM5
A	1178	ALA	SER	engineered mutation	UNP A8EVM5

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

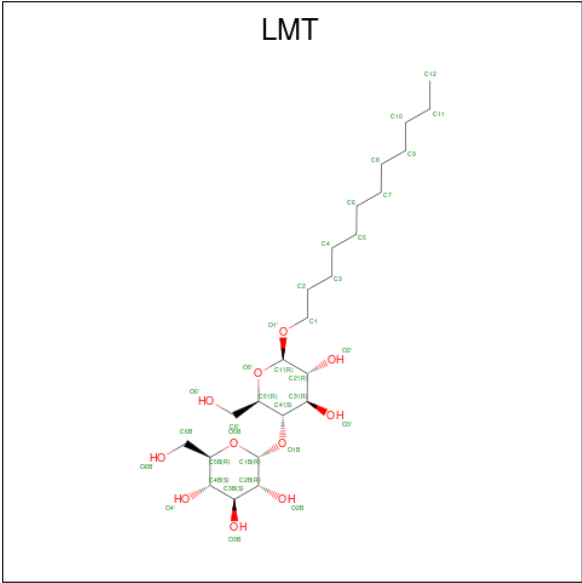
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Ca	0	0
			3	3		

- Molecule 3 is CHAPSO (CCD ID: 1N7) (formula: C<sub>32</sub>H<sub>59</sub>N<sub>2</sub>O<sub>8</sub>S).



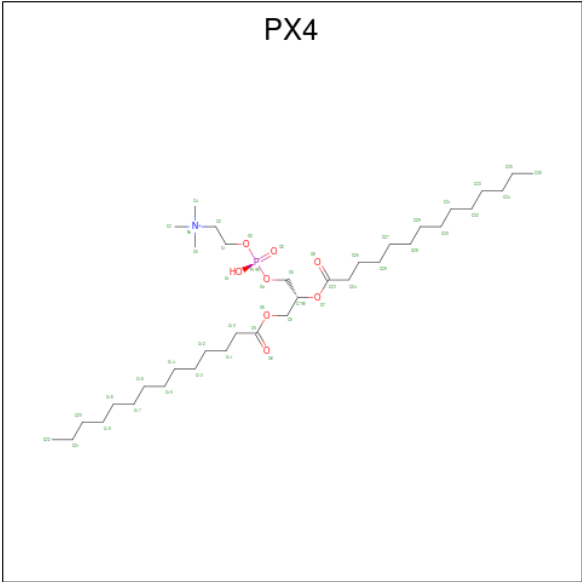
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			43	32	2	8	1		

- Molecule 4 is DODECYL-BETA-D-MALTOSIDE (CCD ID: LMT) (formula:  $C_{24}H_{46}O_{11}$ ).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	C O	0	0
			35	24 11		

- Molecule 5 is 1,2-DIMYRISTOYL-SN-GLYCERO-3-PHOSPHOCHOLINE (CCD ID: PX4) (formula:  $C_{36}H_{73}NO_8P$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			46	36	1	8	1		
5	A	1	Total	C	N	O	P	0	0
			46	36	1	8	1		
5	A	1	Total	C	N	O	P	0	0
			46	36	1	8	1		
5	A	1	Total	C	N	O	P	0	0
			46	36	1	8	1		
5	A	1	Total	C	N	O	P	0	0
			46	36	1	8	1		
5	A	1	Total	C	N	O	P	0	0
			46	36	1	8	1		
5	A	1	Total	C	N	O	P	0	0
			46	36	1	8	1		

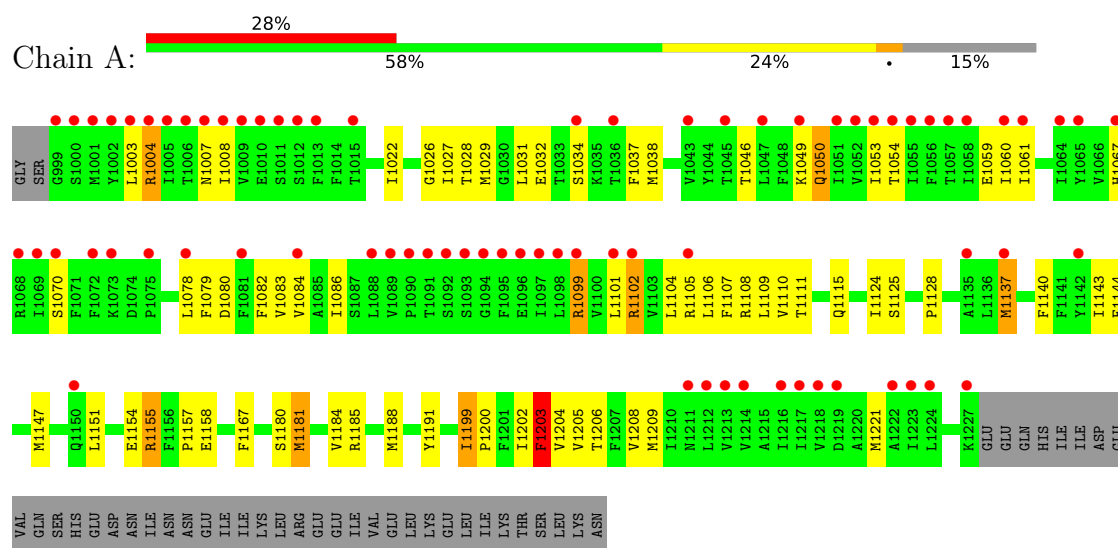
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	28	Total	O	0	0
			28	28		

### 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: Ion transport protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	126.59Å 126.59Å 201.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.76 – 2.80 44.76 – 2.80	Depositor EDS
% Data completeness (in resolution range)	84.6 (44.76-2.80) 84.5 (44.76-2.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.91 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, $R_{free}$	0.256 , 0.294 0.260 , 0.294	Depositor DCC
$R_{free}$ test set	855 reflections (3.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	62.8	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 70.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.83	EDS
Total number of atoms	2346	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.81% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: LMT, 1N7, PX4, CA

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	0.61	3/1920 (0.2%)	1.05	17/2609 (0.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1155	ARG	CD-NE	-8.15	1.34	1.46
1	A	1199	ILE	CA-CB	-5.96	1.51	1.54
1	A	1155	ARG	CB-CG	-5.79	1.35	1.52

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1155	ARG	NE-CZ-NH1	11.68	133.18	121.50
1	A	1155	ARG	NE-CZ-NH2	-8.65	111.41	119.20
1	A	1181	MET	CB-CG-SD	8.51	138.21	112.70
1	A	1004	ARG	CG-CD-NE	-7.53	95.44	112.00
1	A	1102	ARG	NE-CZ-NH2	6.47	125.02	119.20
1	A	1154	GLU	CA-C-N	-6.37	110.90	121.92
1	A	1154	GLU	C-N-CA	-6.37	110.90	121.92
1	A	1155	ARG	CD-NE-CZ	6.18	133.05	124.40
1	A	1181	MET	CA-CB-CG	-6.13	101.85	114.10
1	A	1167	PHE	N-CA-CB	6.07	119.05	110.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1099	ARG	NE-CZ-NH2	6.01	124.61	119.20
1	A	1203	PHE	N-CA-CB	5.88	118.77	110.12
1	A	1099	ARG	NE-CZ-NH1	-5.86	115.64	121.50
1	A	1102	ARG	NE-CZ-NH1	-5.81	115.69	121.50
1	A	1221	MET	CG-SD-CE	-5.55	88.69	100.90
1	A	1137	MET	CA-CB-CG	5.54	125.18	114.10
1	A	1050	GLN	CA-CB-CG	5.07	124.24	114.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1203	PHE	Sidechain

## 5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1869	0	1958	62	0
2	A	3	0	0	0	0
3	A	43	0	59	2	0
4	A	35	0	45	7	0
5	A	368	0	576	27	0
6	A	28	0	0	2	1
All	All	2346	0	2638	86	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 17.

All (86) 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:1059:GLU:OE2	1:A:1108:ARG:NH1	1.99	0.95
1:A:1137:MET:HE1	1:A:1205:VAL:HG12	1.69	0.75
1:A:1206:THR:HG23	4:A:1305:LMT:H6'2	1.71	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1004:ARG:HH22	1:A:1008:ILE:HG23	1.54	0.72
1:A:1083:VAL:HG21	1:A:1105:ARG:HA	1.71	0.71
1:A:1099:ARG:O	1:A:1102:ARG:NH2	2.22	0.71
4:A:1305:LMT:H4'	4:A:1305:LMT:H5B	1.73	0.68
1:A:1185:ARG:NH2	6:A:1402:HOH:O	2.28	0.67
1:A:1115:GLN:HG2	3:A:1304:1N7:H26	1.78	0.65
1:A:1029:MET:HE2	1:A:1102:ARG:NH1	2.12	0.65
5:A:1312:PX4:H33	5:A:1312:PX4:H61	1.77	0.64
5:A:1307:PX4:H35	5:A:1307:PX4:H65	1.81	0.62
1:A:1107:PHE:O	1:A:1110:VAL:HG22	2.02	0.60
1:A:1188:MET:SD	5:A:1311:PX4:H10	2.42	0.60
1:A:1027:ILE:HG12	5:A:1312:PX4:H63	1.83	0.59
1:A:1155:ARG:O	1:A:1155:ARG:HD2	2.02	0.59
5:A:1313:PX4:H66	5:A:1313:PX4:H33	1.84	0.58
1:A:1046:THR:O	1:A:1050:GLN:OE1	2.21	0.57
1:A:1028:THR:HA	1:A:1031:LEU:HD12	1.87	0.57
1:A:1022:ILE:HD11	1:A:1108:ARG:HG2	1.88	0.56
5:A:1306:PX4:H53	5:A:1307:PX4:H19	1.88	0.56
1:A:1067:HIS:HB3	1:A:1070:SER:HB2	1.89	0.55
1:A:1034:SER:HA	5:A:1312:PX4:H4	1.88	0.55
1:A:1050:GLN:O	1:A:1054:THR:HG23	2.07	0.54
1:A:1049:LYS:O	1:A:1053:ILE:HG13	2.06	0.54
1:A:1028:THR:O	1:A:1032:GLU:HG3	2.08	0.54
1:A:1158:GLU:O	1:A:1158:GLU:HG3	2.08	0.53
4:A:1305:LMT:H4'	4:A:1305:LMT:C5B	2.37	0.52
5:A:1306:PX4:H31	5:A:1307:PX4:H60	1.92	0.51
5:A:1311:PX4:H25	5:A:1311:PX4:H55	1.93	0.51
1:A:1099:ARG:HA	1:A:1102:ARG:HD3	1.93	0.51
1:A:1082:PHE:O	1:A:1086:ILE:HG23	2.12	0.50
5:A:1309:PX4:H71	5:A:1309:PX4:H37	1.93	0.50
1:A:1151:LEU:HD22	5:A:1313:PX4:H57	1.94	0.50
1:A:1060:ILE:HG13	1:A:1061:ILE:N	2.27	0.50
1:A:1125:SER:O	1:A:1128:PRO:HD2	2.12	0.49
1:A:1105:ARG:HG3	1:A:1106:LEU:HD22	1.95	0.49
1:A:1188:MET:HE1	5:A:1311:PX4:H6	1.94	0.49
1:A:1105:ARG:O	1:A:1108:ARG:HB3	2.13	0.49
4:A:1305:LMT:O6'	4:A:1305:LMT:H3B	2.13	0.48
1:A:1083:VAL:HG21	1:A:1105:ARG:CA	2.42	0.48
1:A:1101:LEU:O	1:A:1104:LEU:HD23	2.14	0.47
1:A:1080:ASP:O	1:A:1084:VAL:HG22	2.14	0.47
5:A:1307:PX4:H47	5:A:1308:PX4:H26	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1029:MET:HE2	1:A:1102:ARG:HH11	1.79	0.47
5:A:1313:PX4:H71	5:A:1313:PX4:H37	1.96	0.47
3:A:1304:1N7:H37	3:A:1304:1N7:H56	1.97	0.46
5:A:1311:PX4:H9	5:A:1311:PX4:H1	1.41	0.46
1:A:1143:ILE:O	1:A:1147:MET:HG3	2.16	0.45
1:A:1155:ARG:O	1:A:1157:PRO:HD3	2.17	0.45
1:A:1199:ILE:HB	1:A:1200:PRO:HD3	1.98	0.45
1:A:1079:PHE:HD1	1:A:1111:THR:HG21	1.81	0.45
1:A:1204:VAL:O	1:A:1208:VAL:HG23	2.17	0.44
1:A:1209:MET:HE2	4:A:1305:LMT:O5'	2.17	0.44
1:A:1188:MET:HA	1:A:1191:TYR:O	2.18	0.44
1:A:1202:ILE:O	1:A:1203:PHE:C	2.59	0.44
1:A:1078:LEU:HD11	5:A:1308:PX4:H51	2.00	0.44
5:A:1309:PX4:H1	5:A:1309:PX4:H7	1.64	0.44
1:A:1037:PHE:HD1	1:A:1038:MET:CE	2.31	0.43
1:A:1109:LEU:HD23	1:A:1109:LEU:O	2.19	0.43
1:A:1037:PHE:HD1	1:A:1038:MET:HE1	1.82	0.43
1:A:1140:PHE:CE1	1:A:1144:PHE:CE2	3.07	0.43
5:A:1310:PX4:H21	5:A:1310:PX4:H28	1.67	0.43
1:A:1049:LYS:HE2	1:A:1102:ARG:HH11	1.84	0.42
1:A:1004:ARG:O	1:A:1004:ARG:NH2	2.53	0.42
1:A:1082:PHE:CE1	5:A:1308:PX4:H63	2.54	0.42
1:A:1184:VAL:O	1:A:1188:MET:HG3	2.19	0.42
5:A:1307:PX4:H47	5:A:1308:PX4:H20	2.01	0.41
1:A:1180:SER:O	1:A:1185:ARG:HG3	2.20	0.41
5:A:1312:PX4:H57	5:A:1312:PX4:H29	2.02	0.41
5:A:1313:PX4:H1	5:A:1313:PX4:H10	1.62	0.41
1:A:1181:MET:HG2	1:A:1185:ARG:HD3	2.01	0.41
1:A:1050:GLN:OE1	1:A:1050:GLN:N	2.43	0.41
1:A:1206:THR:HA	4:A:1305:LMT:C6B	2.51	0.41
1:A:1060:ILE:HG21	1:A:1084:VAL:HG21	2.03	0.41
1:A:1079:PHE:CD1	1:A:1111:THR:HG21	2.56	0.41
1:A:1124:ILE:HG13	5:A:1306:PX4:H20	2.03	0.41
5:A:1306:PX4:H10	5:A:1306:PX4:H2	1.68	0.41
1:A:1003:LEU:HG	1:A:1007:ASN:ND2	2.35	0.41
1:A:1026:GLY:CA	1:A:1106:LEU:HD11	2.50	0.41
5:A:1310:PX4:H54	5:A:1310:PX4:H48	1.87	0.41
1:A:1155:ARG:HD2	1:A:1155:ARG:HA	1.60	0.41
1:A:1032:GLU:HA	1:A:1038:MET:HE3	2.02	0.40
5:A:1307:PX4:H30	5:A:1307:PX4:H24	1.93	0.40
5:A:1313:PX4:C9	5:A:1313:PX4:H53	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:1305:LMT:H6E	6:A:1404:HOH:O	2.20	0.40

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)
6:A:1406:HOH:O	6:A:1406:HOH:O[7_555]	1.87	0.33

## 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	227/271 (84%)	223 (98%)	4 (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	208/249 (84%)	208 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

Mol	Chain	Res	Type
1	A	1211	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

### 5.6 Ligand geometry ⓘ

Of 13 ligands modelled in this entry, 3 are monoatomic - leaving 10 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	$\# Z  > 2$
5	PX4	A	1308	-	45,45,45	1.05	2 (4%)	51,53,53	0.96	3 (5%)
4	LMT	A	1305	2	36,36,36	1.16	5 (13%)	47,47,47	1.59	10 (21%)
3	1N7	A	1304	-	44,46,46	0.59	1 (2%)	69,72,72	1.59	11 (15%)
5	PX4	A	1311	-	45,45,45	1.15	2 (4%)	51,53,53	1.53	5 (9%)
5	PX4	A	1313	-	45,45,45	1.20	2 (4%)	51,53,53	1.15	3 (5%)
5	PX4	A	1312	-	45,45,45	1.09	2 (4%)	51,53,53	1.14	5 (9%)
5	PX4	A	1310	-	45,45,45	1.00	2 (4%)	51,53,53	1.12	5 (9%)
5	PX4	A	1306	-	45,45,45	1.04	2 (4%)	51,53,53	1.09	3 (5%)
5	PX4	A	1309	-	45,45,45	1.13	2 (4%)	51,53,53	1.02	3 (5%)
5	PX4	A	1307	-	45,45,45	1.04	2 (4%)	51,53,53	0.90	2 (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	PX4	A	1308	-	-	22/49/49/49	-
4	LMT	A	1305	2	-	13/21/61/61	0/2/2/2
3	1N7	A	1304	-	-	15/27/92/92	0/4/4/4
5	PX4	A	1311	-	-	12/49/49/49	-
5	PX4	A	1313	-	-	23/49/49/49	-
5	PX4	A	1312	-	-	16/49/49/49	-
5	PX4	A	1310	-	-	21/49/49/49	-
5	PX4	A	1306	-	-	23/49/49/49	-
5	PX4	A	1309	-	-	21/49/49/49	-
5	PX4	A	1307	-	-	20/49/49/49	-

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1313	PX4	O5-C9	5.39	1.49	1.33
5	A	1311	PX4	O5-C9	5.10	1.48	1.33
5	A	1309	PX4	O7-C23	4.79	1.47	1.34
5	A	1309	PX4	O5-C9	4.74	1.47	1.33
5	A	1312	PX4	O7-C23	4.71	1.47	1.34
5	A	1313	PX4	O7-C23	4.56	1.47	1.34
5	A	1308	PX4	O7-C23	4.46	1.46	1.34
5	A	1306	PX4	O7-C23	4.45	1.46	1.34
5	A	1307	PX4	O5-C9	4.43	1.46	1.33
5	A	1310	PX4	O5-C9	4.33	1.46	1.33
5	A	1311	PX4	O7-C23	4.33	1.46	1.34
5	A	1307	PX4	O7-C23	4.32	1.46	1.34
5	A	1308	PX4	O5-C9	4.30	1.45	1.33
5	A	1312	PX4	O5-C9	4.27	1.45	1.33
5	A	1306	PX4	O5-C9	4.25	1.45	1.33
5	A	1310	PX4	O7-C23	4.02	1.45	1.34
4	A	1305	LMT	O2'-C2'	-3.21	1.35	1.43
4	A	1305	LMT	C3'-C2'	2.89	1.59	1.52
4	A	1305	LMT	O3'-C3'	-2.40	1.37	1.43
4	A	1305	LMT	C4B-C3B	2.38	1.58	1.52
4	A	1305	LMT	C3B-C2B	2.32	1.58	1.52
3	A	1304	1N7	O8-S1	2.21	1.55	1.47

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1304	1N7	C9-C5-C6	-5.40	94.70	100.11
5	A	1311	PX4	C7-O7-C23	-4.80	106.31	117.80
5	A	1311	PX4	O7-C23-C24	4.77	121.80	111.48
3	A	1304	1N7	C7-C6-C18	4.63	124.71	118.36
5	A	1312	PX4	O7-C23-C24	4.57	121.36	111.48
5	A	1309	PX4	O7-C23-C24	4.50	121.22	111.48
3	A	1304	1N7	C28-C31-C32	4.29	117.55	108.44
5	A	1308	PX4	O7-C23-C24	4.28	120.74	111.48
4	A	1305	LMT	C2'-C3'-C4'	4.25	119.32	109.68
5	A	1313	PX4	O7-C23-C24	4.07	120.28	111.48
3	A	1304	1N7	C9-C5-C4	3.98	121.25	117.67
5	A	1306	PX4	O7-C23-C24	3.97	120.06	111.48
4	A	1305	LMT	O5B-C5B-C4B	3.96	116.84	109.70
5	A	1313	PX4	O5-C9-C10	3.68	123.06	111.83
3	A	1304	1N7	C31-C28-N2	3.55	122.87	116.69
5	A	1311	PX4	O5-C9-C10	3.42	122.26	111.83
5	A	1310	PX4	O7-C23-C24	3.38	118.79	111.48
5	A	1311	PX4	C1-C2-N1	-3.37	105.02	115.82
5	A	1306	PX4	O5-C9-C10	3.36	122.08	111.83
3	A	1304	1N7	C6-C18-C17	3.29	116.22	111.85
5	A	1312	PX4	O5-C9-C10	3.28	121.85	111.83
4	A	1305	LMT	C1'-C2'-C3'	3.11	116.56	110.01
5	A	1307	PX4	O7-C23-C24	3.04	118.07	111.48
5	A	1312	PX4	O5-C9-O6	-3.01	116.09	123.63
5	A	1310	PX4	C7-O7-C23	-2.92	110.80	117.80
5	A	1310	PX4	O5-C9-C10	2.89	120.64	111.83
5	A	1307	PX4	O5-C9-C10	2.84	120.50	111.83
4	A	1305	LMT	O1B-C4'-C3'	2.81	114.38	107.23
3	A	1304	1N7	C8-C9-C5	-2.79	100.83	103.54
4	A	1305	LMT	O1B-C1B-C2B	2.73	114.81	108.09
3	A	1304	1N7	C7-C6-C5	-2.66	100.96	103.54
4	A	1305	LMT	O2'-C2'-C1'	-2.64	103.80	110.08
3	A	1304	1N7	C8-C7-C6	-2.56	100.13	105.14
5	A	1312	PX4	C7-O7-C23	-2.45	111.92	117.80
3	A	1304	1N7	C5-C9-C20	2.42	122.41	119.48
4	A	1305	LMT	C1B-O5B-C5B	2.30	118.20	113.72
3	A	1304	1N7	C19-C18-C6	-2.29	106.56	109.75
5	A	1309	PX4	O5-C9-C10	2.21	118.57	111.83
5	A	1306	PX4	C11-C10-C9	-2.15	105.81	113.69
4	A	1305	LMT	O2B-C2B-C3B	2.15	115.44	110.38
5	A	1311	PX4	C4-N1-C3	2.12	114.54	108.98
5	A	1309	PX4	C1-C2-N1	-2.10	109.08	115.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1305	LMT	O3'-C3'-C4'	-2.10	104.57	109.94
5	A	1310	PX4	C8-C7-C6	-2.08	106.93	111.78
5	A	1308	PX4	C8-C7-C6	-2.07	106.96	111.78
4	A	1305	LMT	O3B-C3B-C2B	2.06	115.24	110.38
5	A	1308	PX4	O5-C9-C10	2.06	118.13	111.83
5	A	1310	PX4	O5-C9-O6	-2.06	118.48	123.63
5	A	1312	PX4	C25-C24-C23	-2.04	106.21	113.69
5	A	1313	PX4	O1-P1-O2	2.03	121.88	112.44

There are no chirality outliers.

All (186) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1304	1N7	N2-C28-C31-C32
3	A	1304	1N7	O5-C31-C32-S1
4	A	1305	LMT	C2'-C1'-O1'-C1
5	A	1306	PX4	C6-O4-P1-O1
5	A	1306	PX4	C6-O4-P1-O3
5	A	1306	PX4	O3-C1-C2-N1
5	A	1307	PX4	O3-C1-C2-N1
5	A	1307	PX4	C24-C23-O7-C7
5	A	1308	PX4	C1-O3-P1-O1
5	A	1308	PX4	C1-O3-P1-O2
5	A	1309	PX4	C1-O3-P1-O1
5	A	1309	PX4	C1-O3-P1-O2
5	A	1309	PX4	C1-O3-P1-O4
5	A	1309	PX4	C6-O4-P1-O1
5	A	1309	PX4	C6-O4-P1-O2
5	A	1309	PX4	C6-O4-P1-O3
5	A	1310	PX4	C1-O3-P1-O4
5	A	1310	PX4	C6-O4-P1-O1
5	A	1310	PX4	C6-O4-P1-O2
5	A	1310	PX4	C24-C23-O7-C7
5	A	1312	PX4	C1-O3-P1-O2
5	A	1312	PX4	C6-O4-P1-O1
5	A	1312	PX4	C6-O4-P1-O2
5	A	1312	PX4	C6-O4-P1-O3
5	A	1312	PX4	O3-C1-C2-N1
5	A	1313	PX4	C1-O3-P1-O1
5	A	1313	PX4	C6-O4-P1-O1
5	A	1313	PX4	C6-O4-P1-O2
5	A	1313	PX4	C6-O4-P1-O3

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Mol	Chain	Res	Type	Atoms
5	A	1313	PX4	O3-C1-C2-N1
5	A	1313	PX4	C24-C23-O7-C7
4	A	1305	LMT	O5B-C1B-O1B-C4'
5	A	1307	PX4	O8-C23-O7-C7
5	A	1313	PX4	O8-C23-O7-C7
5	A	1310	PX4	O8-C23-O7-C7
4	A	1305	LMT	O5B-C5B-C6B-O6B
5	A	1313	PX4	C10-C9-O5-C8
5	A	1309	PX4	C24-C23-O7-C7
5	A	1313	PX4	O6-C9-O5-C8
4	A	1305	LMT	O5'-C1'-O1'-C1
4	A	1305	LMT	C4B-C5B-C6B-O6B
5	A	1308	PX4	C1-C2-N1-C5
5	A	1306	PX4	C18-C19-C20-C21
5	A	1312	PX4	C23-C24-C25-C26
3	A	1304	1N7	O1-C24-N1-C25
3	A	1304	1N7	C23-C24-N1-C25
5	A	1310	PX4	O7-C7-C8-O5
5	A	1309	PX4	O8-C23-O7-C7
3	A	1304	1N7	N1-C25-C26-C27
3	A	1304	1N7	C22-C20-C9-C5
3	A	1304	1N7	C21-C20-C9-C8
5	A	1307	PX4	C9-C10-C11-C12
3	A	1304	1N7	C21-C20-C9-C5
5	A	1307	PX4	C23-C24-C25-C26
5	A	1306	PX4	C1-C2-N1-C4
5	A	1306	PX4	C1-C2-N1-C5
5	A	1308	PX4	C1-C2-N1-C3
5	A	1313	PX4	C23-C24-C25-C26
5	A	1313	PX4	O4-C6-C7-O7
5	A	1306	PX4	C1-C2-N1-C3
5	A	1311	PX4	C17-C18-C19-C20
5	A	1306	PX4	C23-C24-C25-C26
5	A	1310	PX4	C28-C29-C30-C31
5	A	1311	PX4	C24-C23-O7-C7
5	A	1307	PX4	C26-C27-C28-C29
5	A	1310	PX4	C9-C10-C11-C12
5	A	1312	PX4	C18-C19-C20-C21
5	A	1308	PX4	C26-C27-C28-C29
5	A	1313	PX4	C31-C32-C33-C34
5	A	1309	PX4	C32-C33-C34-C35
5	A	1307	PX4	C15-C16-C17-C18

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Mol	Chain	Res	Type	Atoms
5	A	1306	PX4	C14-C15-C16-C17
5	A	1306	PX4	C27-C28-C29-C30
5	A	1308	PX4	C16-C17-C18-C19
5	A	1307	PX4	C30-C31-C32-C33
5	A	1306	PX4	C28-C29-C30-C31
5	A	1306	PX4	C13-C14-C15-C16
5	A	1306	PX4	C33-C34-C35-C36
5	A	1311	PX4	O8-C23-O7-C7
5	A	1309	PX4	C13-C14-C15-C16
5	A	1308	PX4	C1-C2-N1-C4
5	A	1308	PX4	C29-C30-C31-C32
5	A	1306	PX4	C25-C26-C27-C28
5	A	1307	PX4	O7-C7-C8-O5
5	A	1306	PX4	C11-C12-C13-C14
5	A	1306	PX4	C12-C13-C14-C15
5	A	1311	PX4	C27-C28-C29-C30
5	A	1313	PX4	O4-C6-C7-C8
4	A	1305	LMT	O5'-C5'-C6'-O6'
5	A	1306	PX4	C6-C7-C8-O5
3	A	1304	1N7	N2-C28-C31-O5
5	A	1309	PX4	C14-C15-C16-C17
5	A	1313	PX4	C29-C30-C31-C32
5	A	1306	PX4	C29-C30-C31-C32
5	A	1307	PX4	O4-C6-C7-O7
5	A	1313	PX4	C26-C27-C28-C29
5	A	1308	PX4	C17-C18-C19-C20
5	A	1312	PX4	C31-C32-C33-C34
3	A	1304	1N7	C20-C22-C23-C24
5	A	1309	PX4	C7-C8-O5-C9
4	A	1305	LMT	C2-C1-O1'-C1'
5	A	1306	PX4	C26-C27-C28-C29
5	A	1308	PX4	C27-C28-C29-C30
5	A	1309	PX4	C17-C18-C19-C20
5	A	1307	PX4	C33-C34-C35-C36
5	A	1307	PX4	C6-C7-C8-O5
5	A	1310	PX4	C6-C7-C8-O5
5	A	1307	PX4	C14-C15-C16-C17
5	A	1306	PX4	O7-C7-C8-O5
5	A	1310	PX4	C18-C19-C20-C21
5	A	1310	PX4	C10-C11-C12-C13
5	A	1309	PX4	C28-C29-C30-C31
5	A	1308	PX4	C11-C12-C13-C14

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Mol	Chain	Res	Type	Atoms
5	A	1313	PX4	C11-C12-C13-C14
5	A	1308	PX4	O4-C6-C7-C8
5	A	1308	PX4	C31-C32-C33-C34
5	A	1313	PX4	C18-C19-C20-C21
5	A	1308	PX4	O4-C6-C7-O7
5	A	1312	PX4	C6-C7-C8-O5
5	A	1313	PX4	C2-C1-O3-P1
5	A	1309	PX4	O7-C7-C8-O5
4	A	1305	LMT	C11-C10-C9-C8
5	A	1308	PX4	O3-C1-C2-N1
5	A	1309	PX4	O3-C1-C2-N1
5	A	1311	PX4	C26-C27-C28-C29
5	A	1312	PX4	C17-C18-C19-C20
5	A	1307	PX4	C1-C2-N1-C5
5	A	1309	PX4	C27-C28-C29-C30
5	A	1310	PX4	C24-C25-C26-C27
5	A	1308	PX4	C1-O3-P1-O4
5	A	1310	PX4	C1-O3-P1-O2
5	A	1310	PX4	C6-O4-P1-O3
5	A	1311	PX4	C6-O4-P1-O2
5	A	1312	PX4	C1-O3-P1-O1
5	A	1312	PX4	C1-O3-P1-O4
5	A	1308	PX4	C11-C10-C9-O5
5	A	1309	PX4	C15-C16-C17-C18
5	A	1311	PX4	C19-C20-C21-C22
4	A	1305	LMT	C4-C5-C6-C7
5	A	1307	PX4	C1-C2-N1-C4
5	A	1311	PX4	C25-C26-C27-C28
5	A	1309	PX4	C18-C19-C20-C21
5	A	1313	PX4	O7-C23-C24-C25
5	A	1308	PX4	C10-C9-O5-C8
5	A	1311	PX4	C13-C14-C15-C16
5	A	1313	PX4	C10-C11-C12-C13
3	A	1304	1N7	C22-C23-C24-O1
5	A	1311	PX4	C24-C25-C26-C27
5	A	1312	PX4	C29-C30-C31-C32
5	A	1308	PX4	O6-C9-O5-C8
3	A	1304	1N7	C22-C20-C9-C8
5	A	1313	PX4	C15-C16-C17-C18
5	A	1312	PX4	C19-C20-C21-C22
5	A	1312	PX4	C13-C14-C15-C16
5	A	1308	PX4	C30-C31-C32-C33

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Mol	Chain	Res	Type	Atoms
5	A	1307	PX4	C31-C32-C33-C34
4	A	1305	LMT	C6-C7-C8-C9
4	A	1305	LMT	C3'-C4'-O1B-C1B
3	A	1304	1N7	C22-C23-C24-N1
5	A	1307	PX4	C1-C2-N1-C3
5	A	1313	PX4	C13-C14-C15-C16
5	A	1306	PX4	C15-C16-C17-C18
5	A	1311	PX4	C28-C29-C30-C31
5	A	1307	PX4	O4-C6-C7-C8
5	A	1308	PX4	C24-C25-C26-C27
5	A	1307	PX4	C11-C12-C13-C14
5	A	1307	PX4	C16-C17-C18-C19
5	A	1310	PX4	C11-C12-C13-C14
3	A	1304	1N7	C26-C25-N1-C24
5	A	1309	PX4	C6-C7-C8-O5
5	A	1310	PX4	C25-C26-C27-C28
5	A	1311	PX4	C31-C32-C33-C34
5	A	1308	PX4	O7-C7-C8-O5
5	A	1313	PX4	C1-C2-N1-C5
5	A	1309	PX4	C10-C11-C12-C13
5	A	1310	PX4	O3-C1-C2-N1
4	A	1305	LMT	C9-C10-C11-C12
5	A	1310	PX4	C23-C24-C25-C26
5	A	1306	PX4	O7-C23-C24-C25
5	A	1310	PX4	C11-C10-C9-O5
3	A	1304	1N7	C25-C26-C27-N2
5	A	1306	PX4	O8-C23-C24-C25
4	A	1305	LMT	C7-C8-C9-C10
5	A	1310	PX4	C11-C10-C9-O6
5	A	1312	PX4	C24-C25-C26-C27
5	A	1310	PX4	C31-C32-C33-C34

There are no ring outliers.

10 monomers are involved in 36 short contacts:

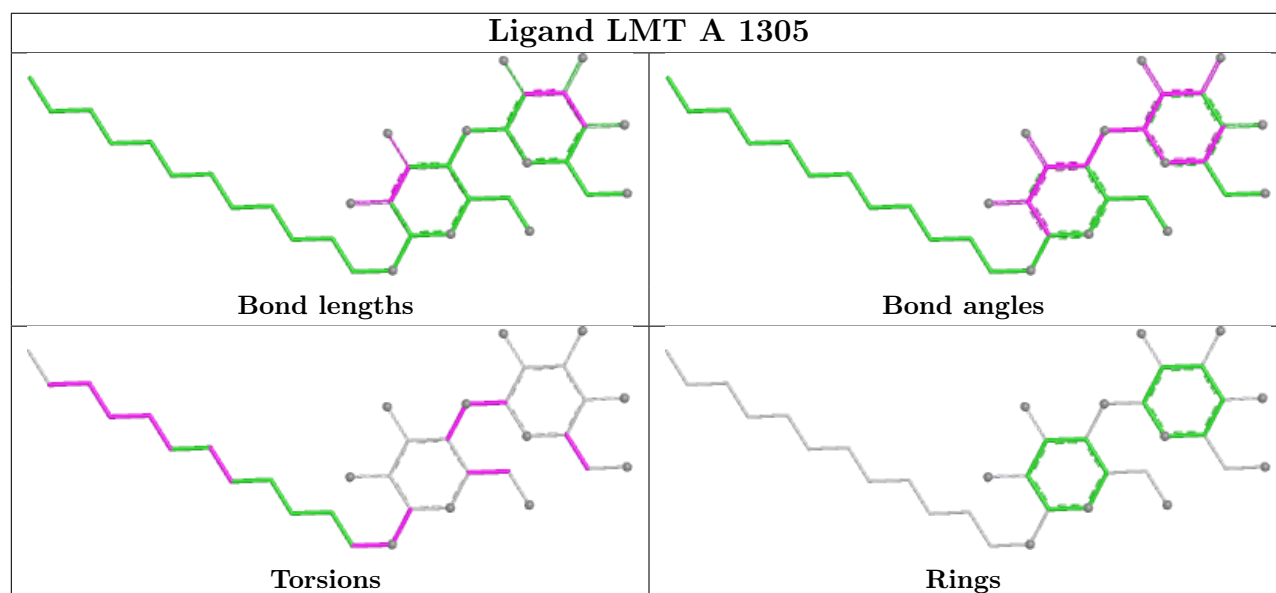
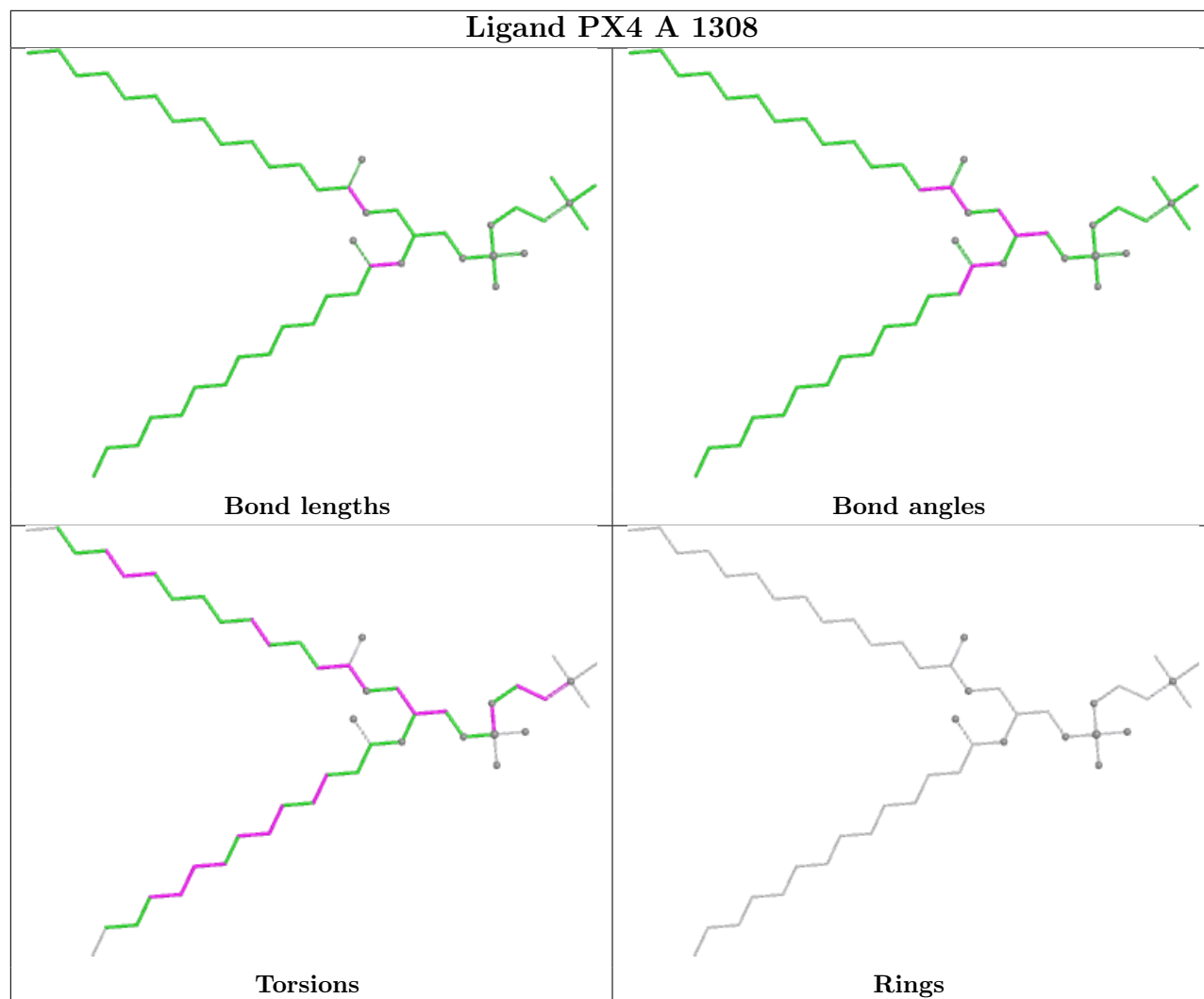
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1308	PX4	4	0
4	A	1305	LMT	7	0
3	A	1304	1N7	2	0
5	A	1311	PX4	4	0
5	A	1313	PX4	5	0
5	A	1312	PX4	4	0

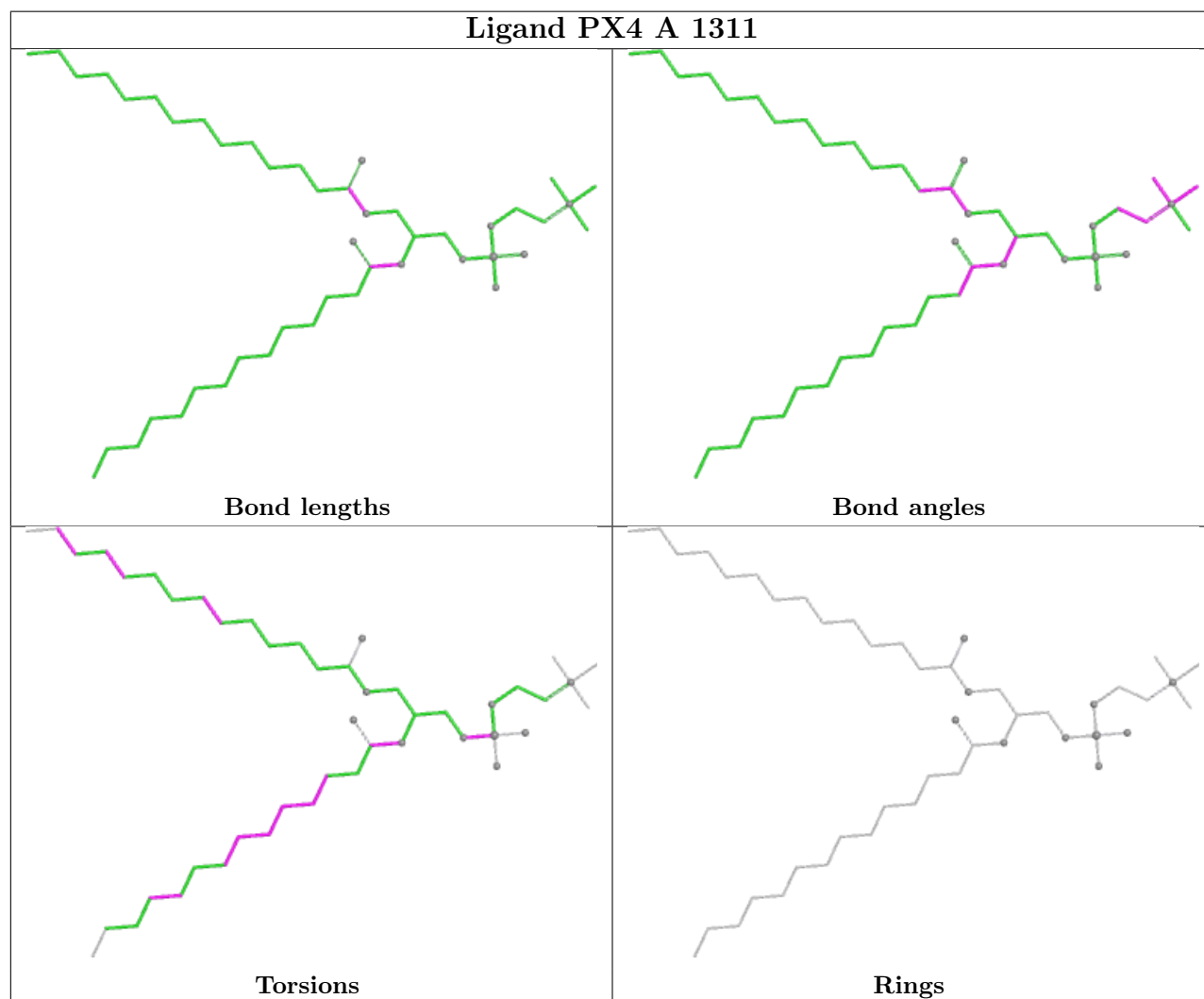
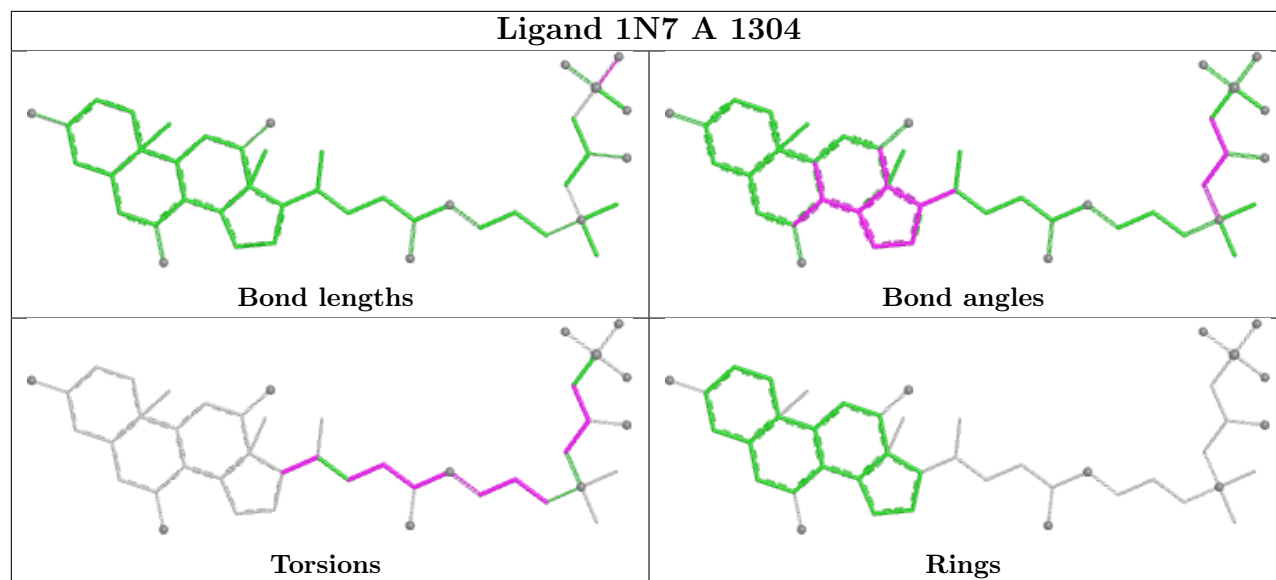
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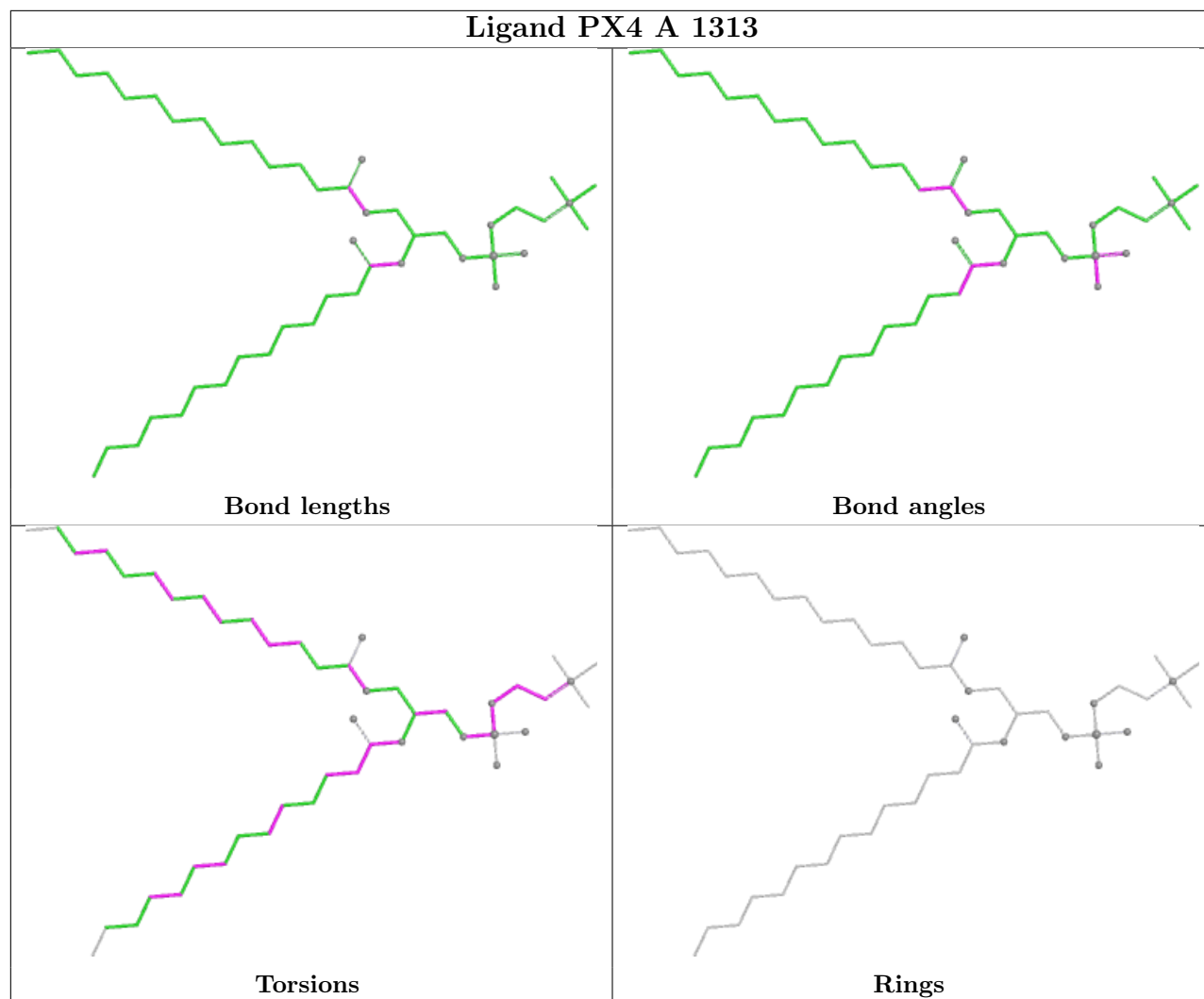
*Continued from previous page...*

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1310	PX4	2	0
5	A	1306	PX4	4	0
5	A	1309	PX4	2	0
5	A	1307	PX4	6	0

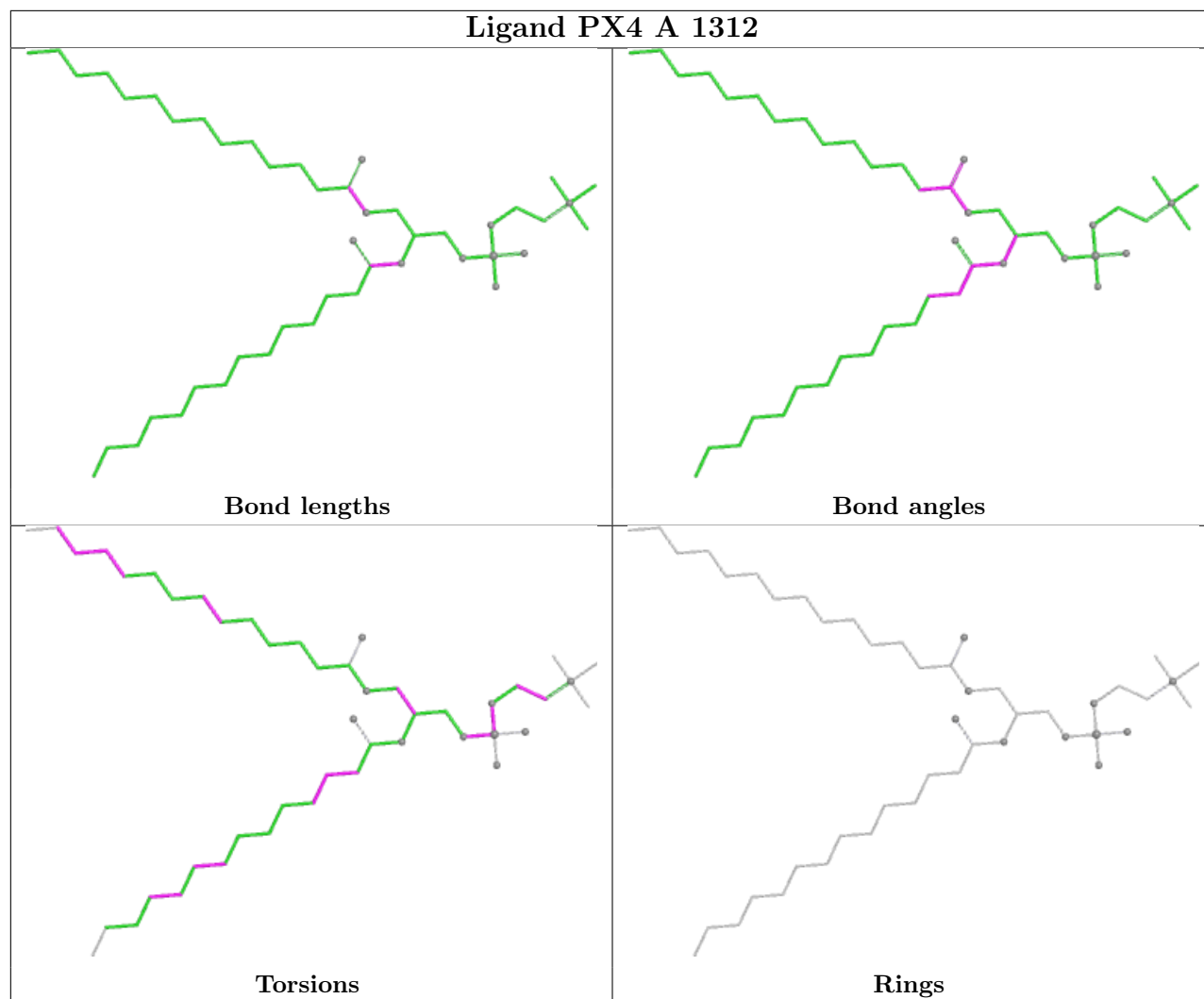
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

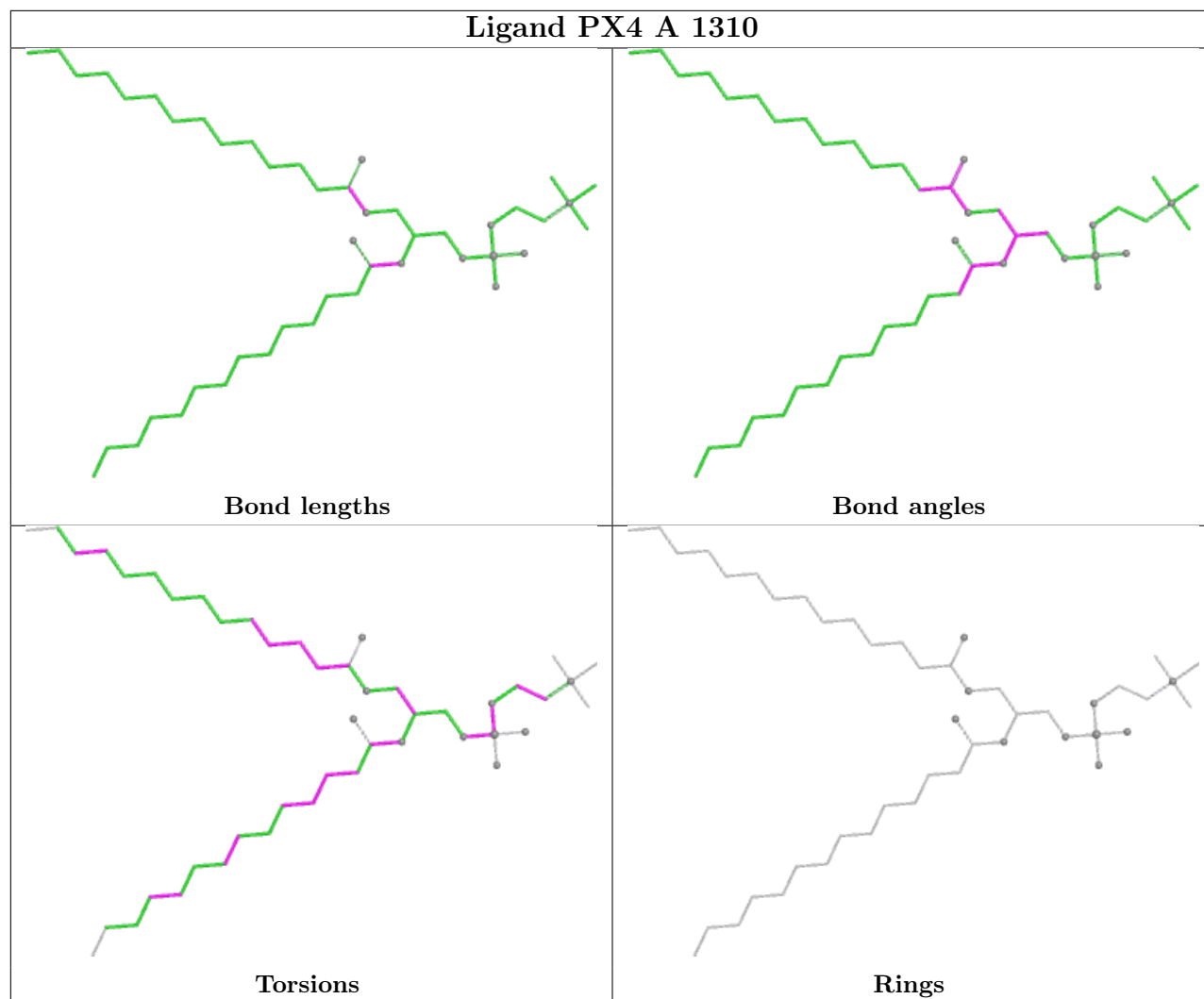


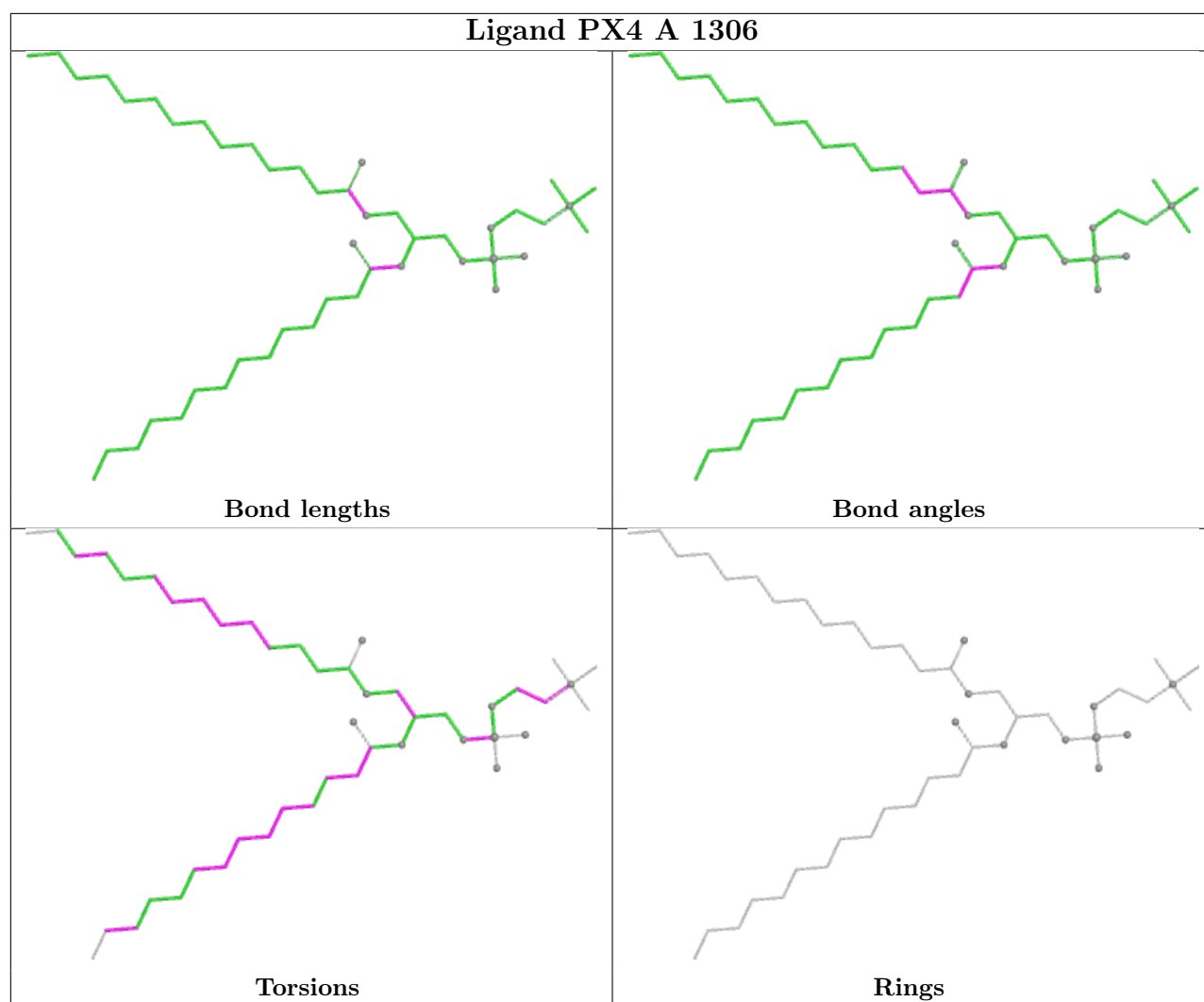


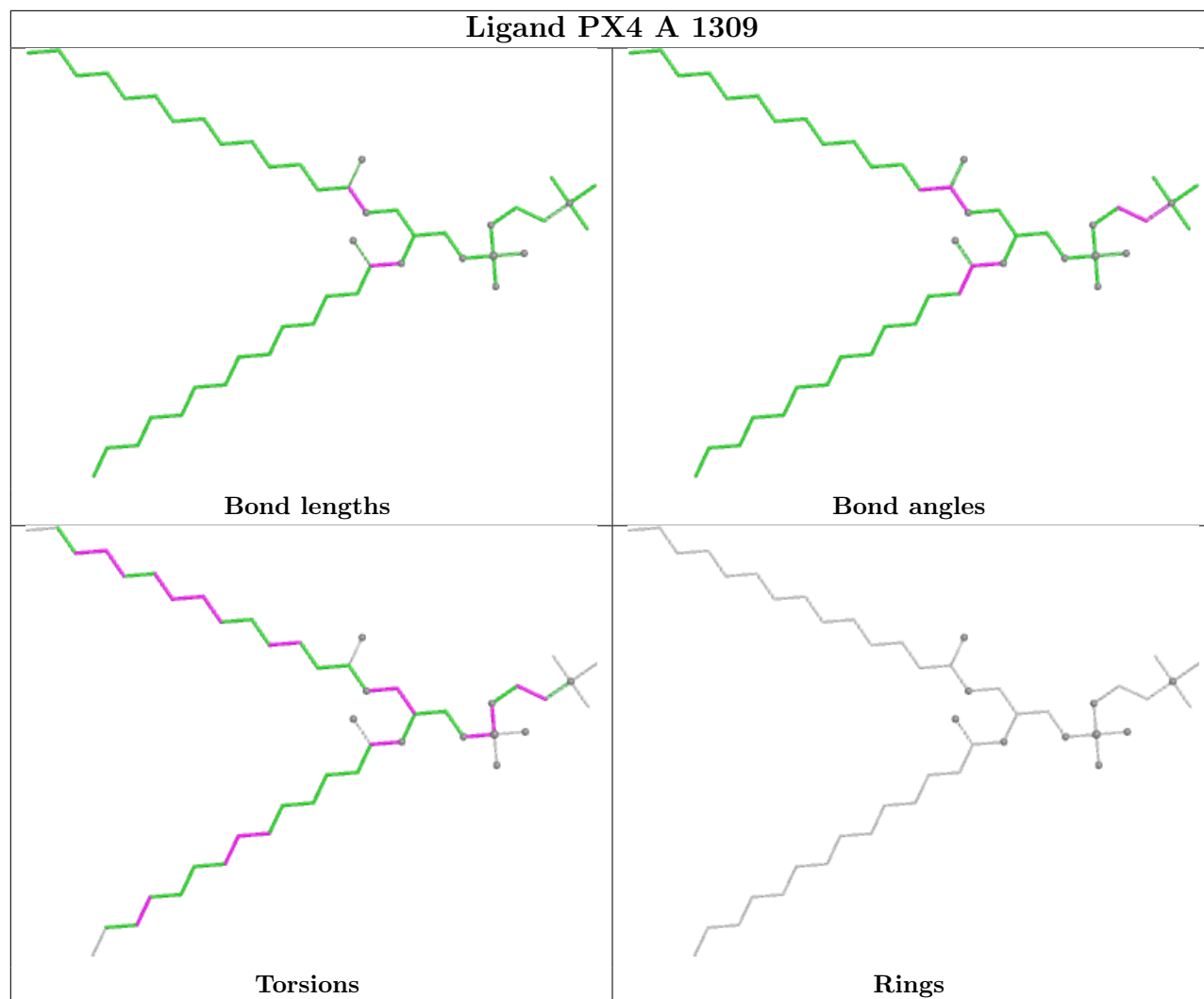


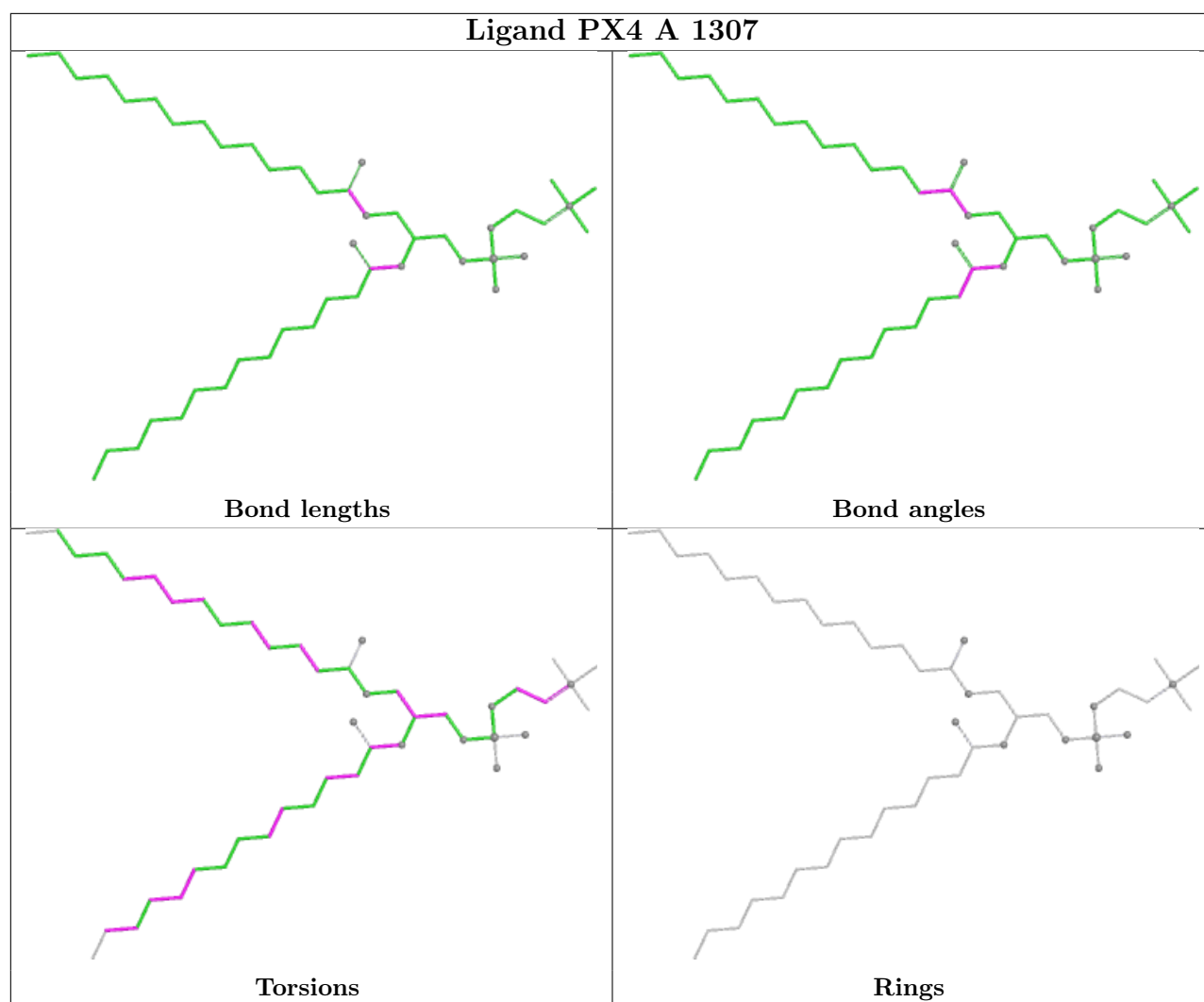












## 5.7 Other polymers [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	229/271 (84%)	2.25	75 (32%) <b>1</b> <b>1</b>	23, 76, 129, 148	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1094	GLY	22.9
1	A	1093	SER	21.3
1	A	1091	THR	21.2
1	A	1092	SER	19.6
1	A	1095	PHE	19.6
1	A	1096	GLU	18.0
1	A	999	GLY	17.3
1	A	1000	SER	16.1
1	A	1002	TYR	15.3
1	A	1068	ARG	13.0
1	A	1004	ARG	9.8
1	A	1001	MET	9.0
1	A	1065	TYR	8.3
1	A	1089	VAL	7.8
1	A	1090	PRO	7.8
1	A	1003	LEU	7.7
1	A	1012	SER	7.4
1	A	1007	ASN	7.2
1	A	1013	PHE	6.1
1	A	1008	ILE	5.9
1	A	1097	ILE	5.8
1	A	1098	LEU	5.3
1	A	1217	ILE	5.3
1	A	1218	VAL	5.0
1	A	1006	THR	4.7
1	A	1211	ASN	4.5
1	A	1099	ARG	4.4

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Mol	Chain	Res	Type	RSRZ
1	A	1009	VAL	4.3
1	A	1061	ILE	4.1
1	A	1081	PHE	4.0
1	A	1223	ILE	3.9
1	A	1005	ILE	3.7
1	A	1078	LEU	3.6
1	A	1222	ALA	3.6
1	A	1227	LYS	3.5
1	A	1216	ILE	3.3
1	A	1150	GLN	3.2
1	A	1011	SER	3.1
1	A	1067	HIS	3.0
1	A	1075	PRO	3.0
1	A	1054	THR	2.9
1	A	1060	ILE	2.9
1	A	1213	VAL	2.9
1	A	1088	LEU	2.8
1	A	1036	THR	2.7
1	A	1214	VAL	2.7
1	A	1052	VAL	2.7
1	A	1105	ARG	2.6
1	A	1058	ILE	2.6
1	A	1015	THR	2.6
1	A	1010	GLU	2.6
1	A	1049	LYS	2.5
1	A	1101	LEU	2.5
1	A	1069	ILE	2.5
1	A	1055	ILE	2.5
1	A	1137	MET	2.5
1	A	1224	LEU	2.5
1	A	1219	ASP	2.4
1	A	1056	PHE	2.4
1	A	1053	ILE	2.4
1	A	1051	ILE	2.4
1	A	1070	SER	2.4
1	A	1064	ILE	2.4
1	A	1073	LYS	2.3
1	A	1057	THR	2.3
1	A	1142	TYR	2.3
1	A	1043	VAL	2.3
1	A	1072	PHE	2.2
1	A	1045	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	1135	ALA	2.2
1	A	1047	LEU	2.1
1	A	1212	LEU	2.1
1	A	1084	VAL	2.1
1	A	1034	SER	2.0
1	A	1102	ARG	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 oligosaccharides in this entry.

## 6.4 Ligands [i](#)

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

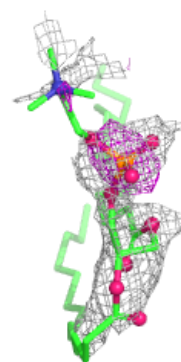
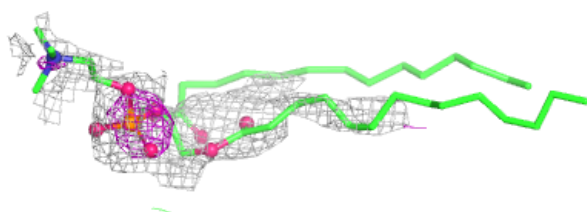
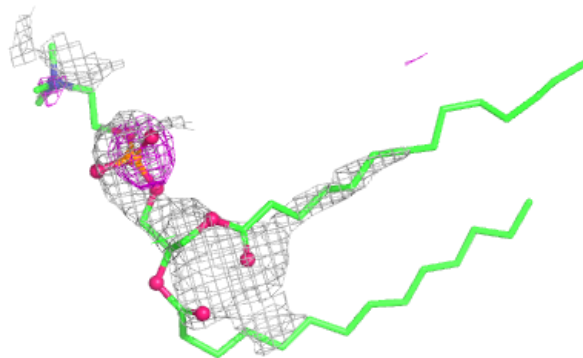
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	PX4	A	1309	46/46	0.64	0.31	60,86,109,113	0
3	1N7	A	1304	43/43	0.65	0.26	70,82,133,139	0
5	PX4	A	1308	46/46	0.68	0.28	84,105,130,144	0
4	LMT	A	1305	35/35	0.69	0.24	54,65,78,80	0
5	PX4	A	1310	46/46	0.70	0.30	79,100,136,141	0
5	PX4	A	1307	46/46	0.71	0.30	81,99,121,135	0
5	PX4	A	1313	46/46	0.71	0.24	40,61,88,102	0
5	PX4	A	1312	46/46	0.78	0.23	47,69,90,95	0
5	PX4	A	1306	46/46	0.80	0.24	67,83,104,121	0
2	CA	A	1301	1/1	0.84	0.08	46,46,46,46	1
5	PX4	A	1311	46/46	0.85	0.21	38,61,74,80	0
2	CA	A	1303	1/1	0.95	0.13	43,43,43,43	0
2	CA	A	1302	1/1	0.98	0.11	76,76,76,76	1

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.

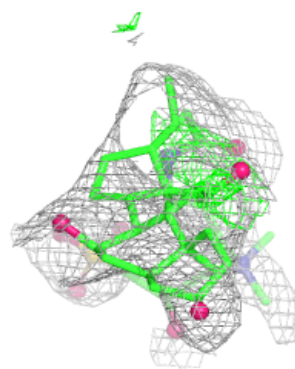
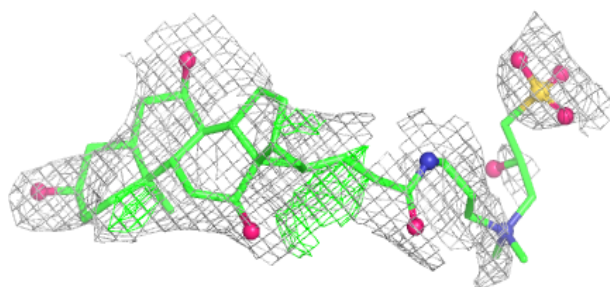
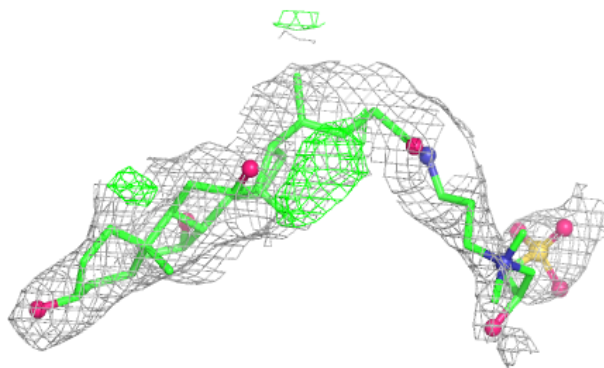


**Electron density around PX4 A 1309:**

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

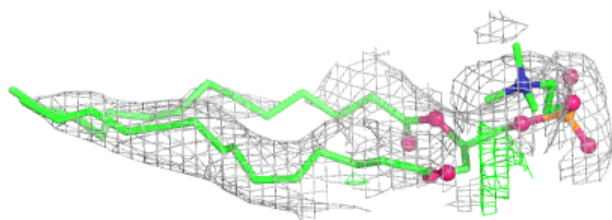
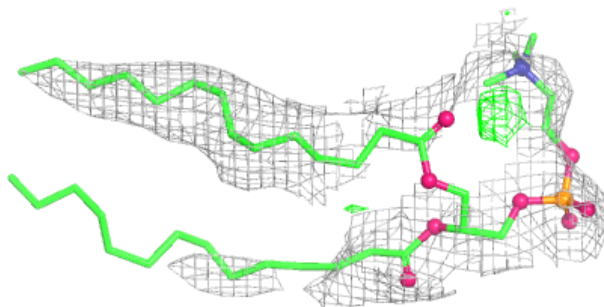
**Electron density around 1N7 A 1304:**

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

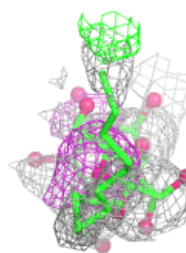
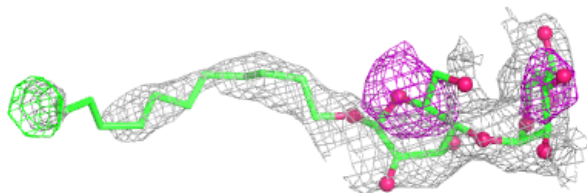
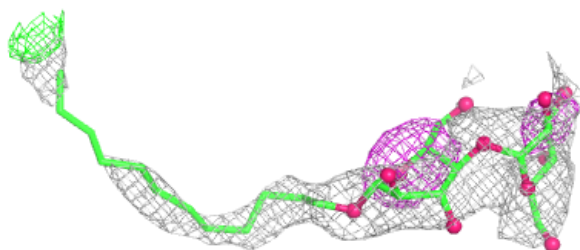


**Electron density around PX4 A 1308:**

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

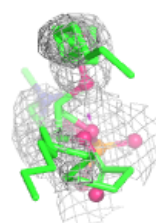
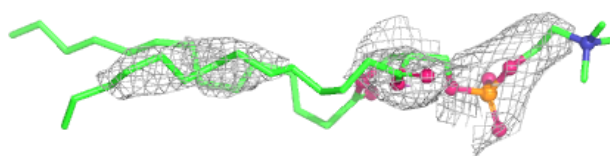
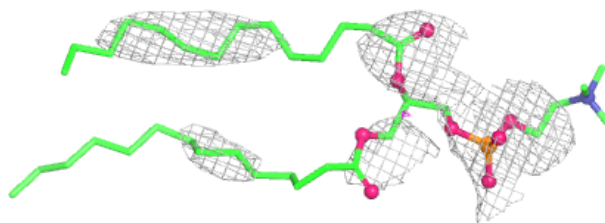
**Electron density around LMT A 1305:**

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

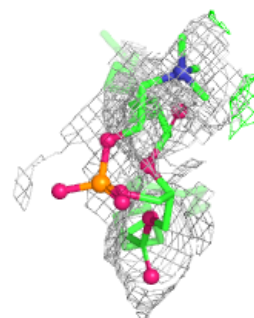
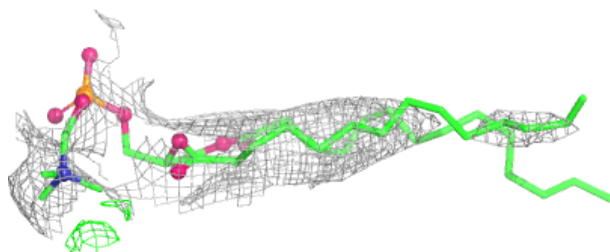
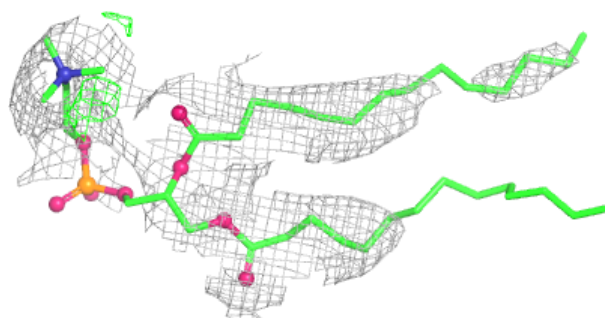


**Electron density around PX4 A 1310:**

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

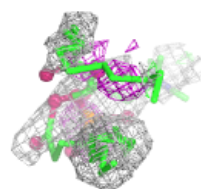
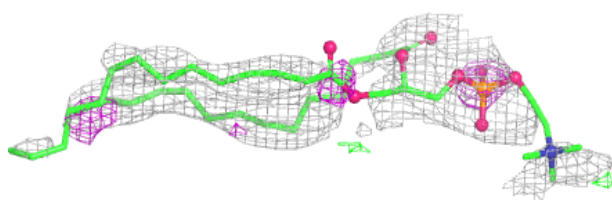
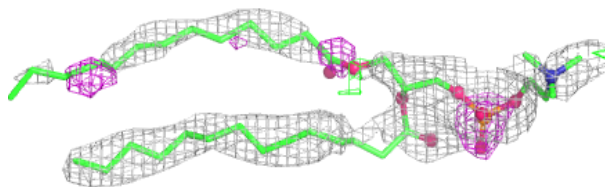
**Electron density around PX4 A 1307:**

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

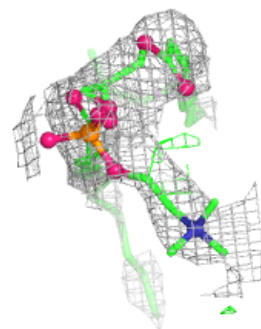
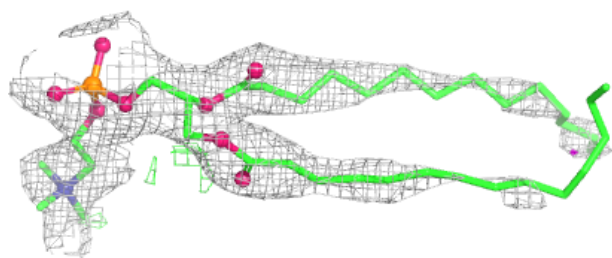
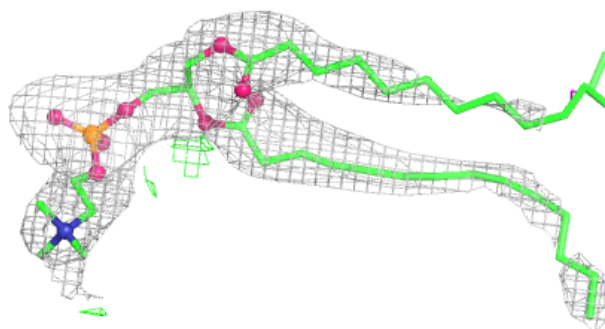


**Electron density around PX4 A 1313:**

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

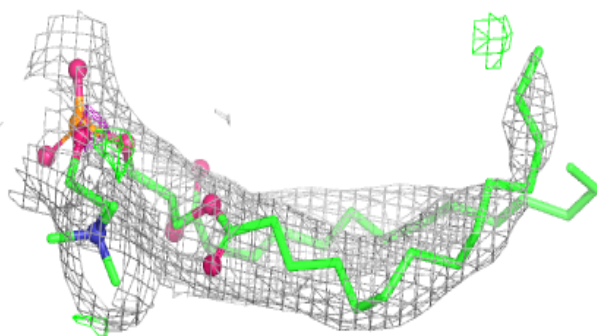
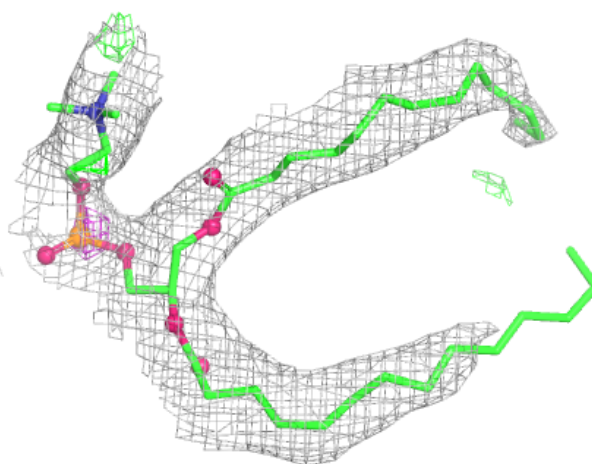
**Electron density around PX4 A 1312:**

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



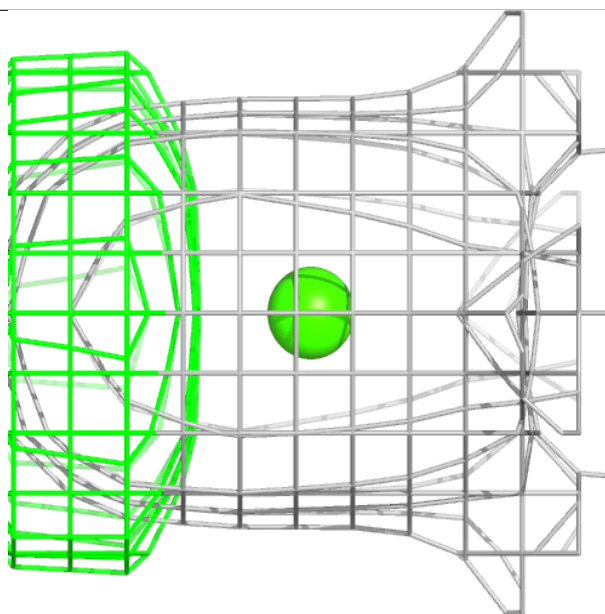
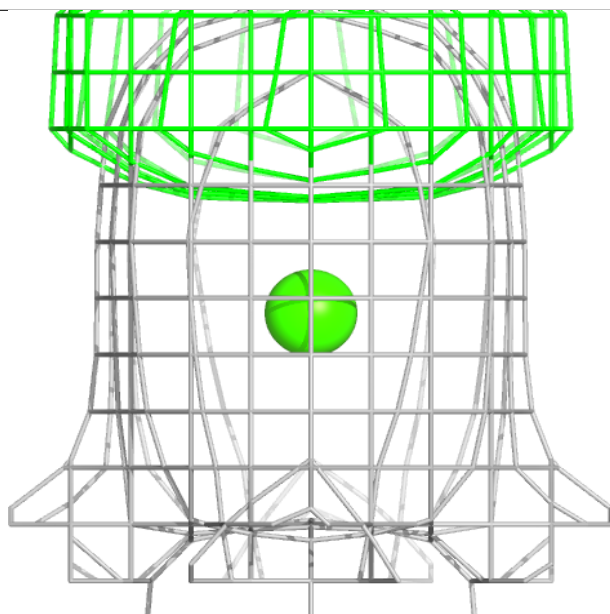
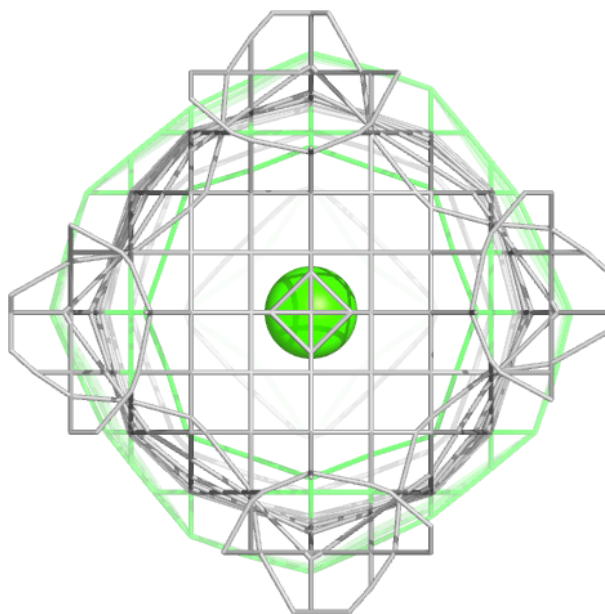
**Electron density around PX4 A 1306:**

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



**Electron density around CA A 1301:**

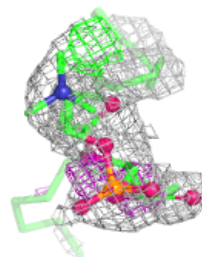
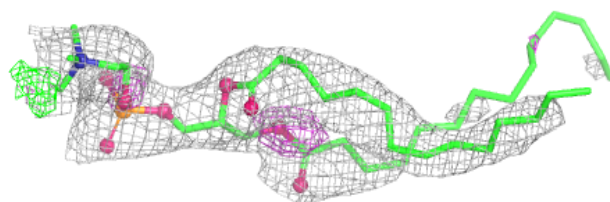
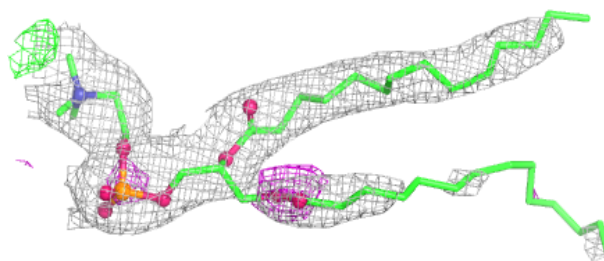
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





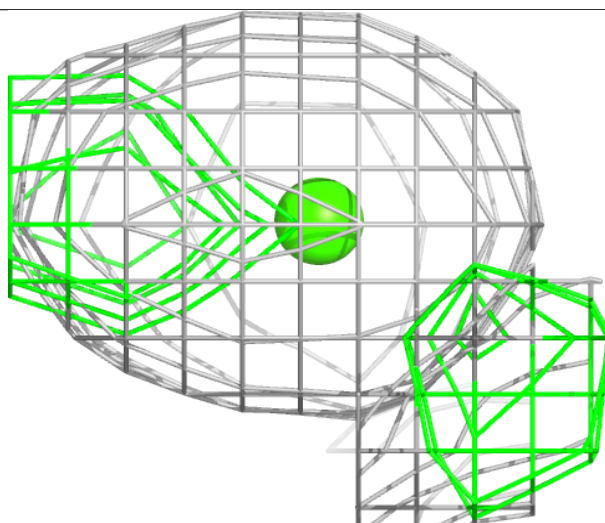
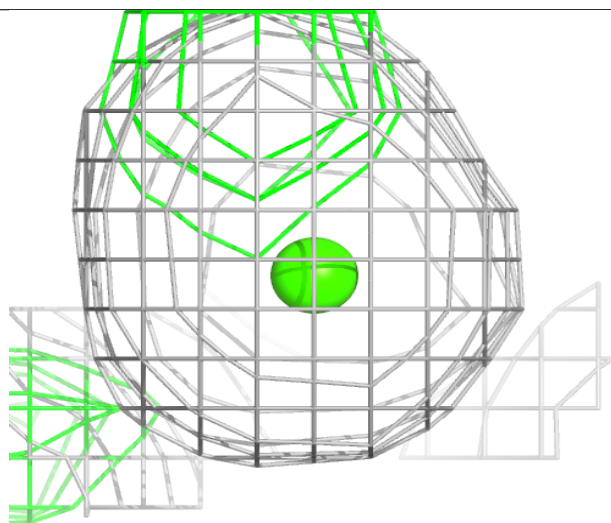
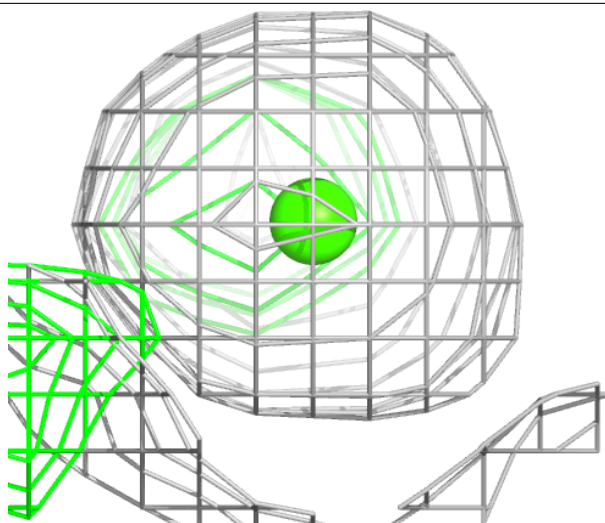
**Electron density around PX4 A 1311:**

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



**Electron density around CA A 1303:**

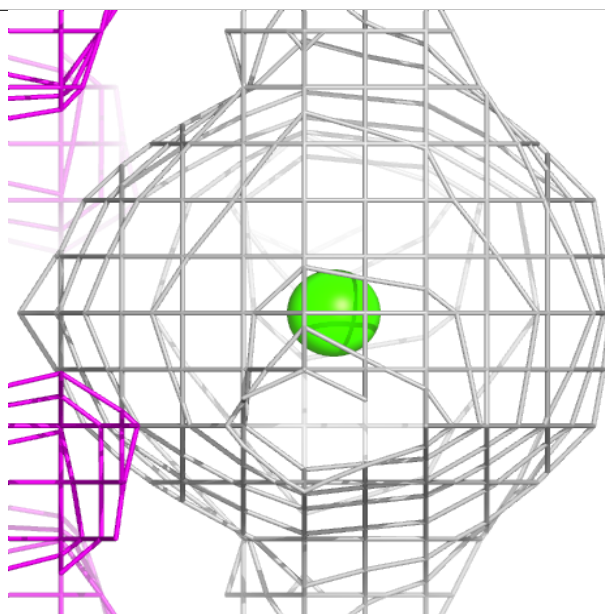
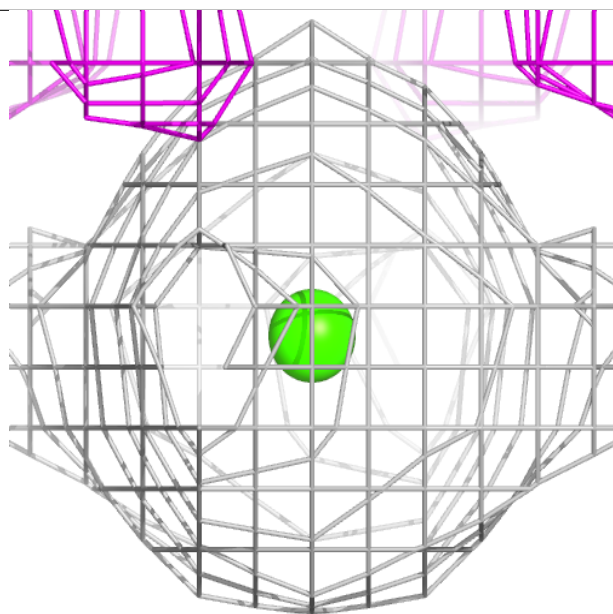
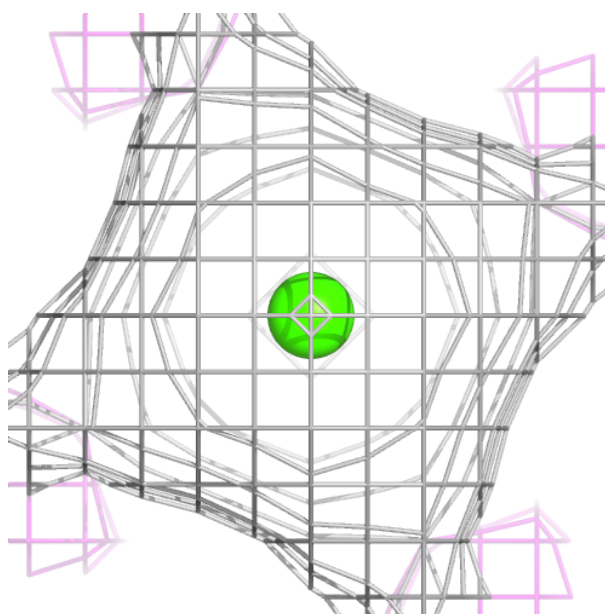
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around CA A 1302:**

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



## 6.5 Other polymers ⓘ

There are no such residues in this entry.