



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:45 PM UTC

PDB ID : 6P6Y / pdb_00006p6y
Title : Crystal structure of voltage-gated sodium channel NavAb V100C/Q150C disulfide crosslinked mutant in the activated state
Authors : Wisedchaisri, G.; Tonggu, L.; McCord, E.; Gamal El-din, T.M.; Wang, L.; Zheng, N.; Catterall, W.A.
Deposited on : 2019-06-04
Resolution : 2.89 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) 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

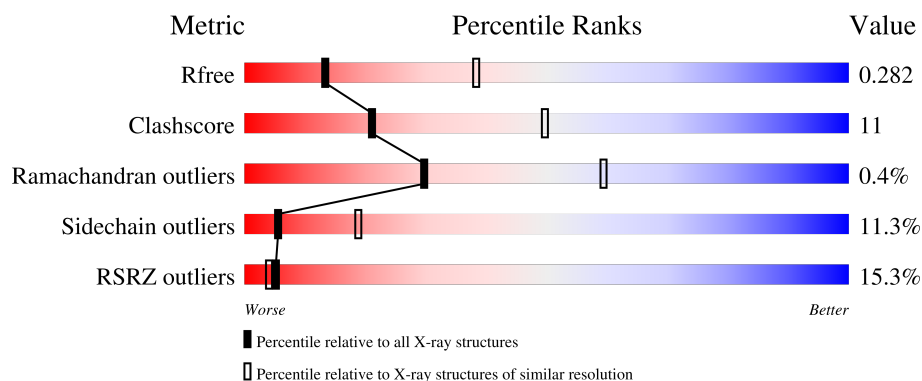
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	257	14% 66% 25% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2201 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.

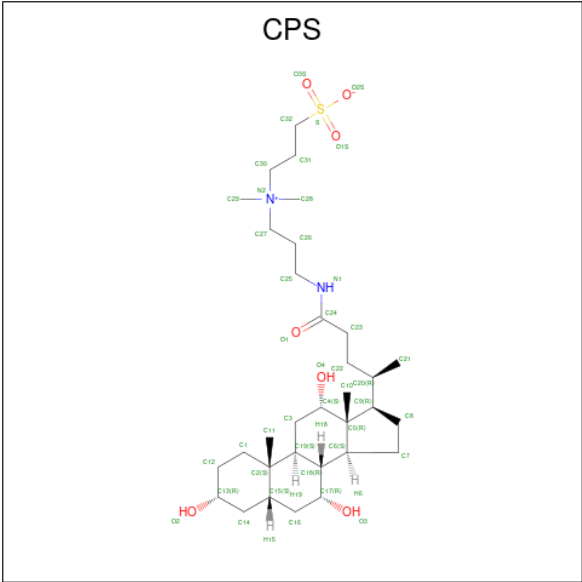
- Molecule 1 is a protein called Ion transport protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	242	Total	C	N	O	S	0	0	0
			1977	1334	301	328	14			

There are 20 discrepancies between the modelled and reference sequences:

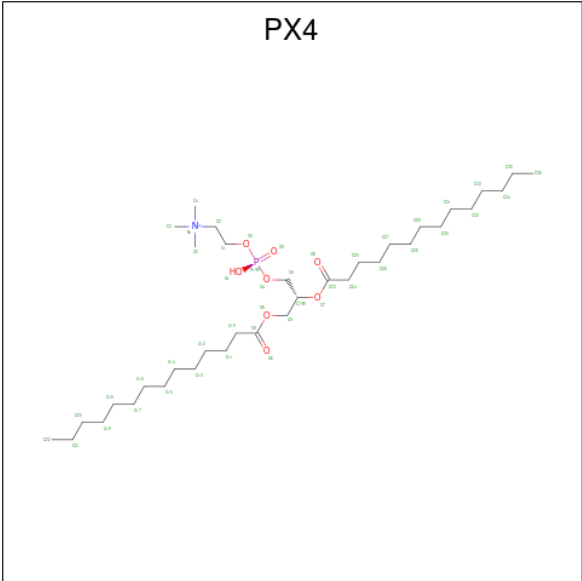
Chain	Residue	Modelled	Actual	Comment	Reference
A	983	MET	-	initiating methionine	UNP A8EVM5
A	984	ASP	-	expression tag	UNP A8EVM5
A	985	TYR	-	expression tag	UNP A8EVM5
A	986	LYS	-	expression tag	UNP A8EVM5
A	987	ASP	-	expression tag	UNP A8EVM5
A	988	ASP	-	expression tag	UNP A8EVM5
A	989	ASP	-	expression tag	UNP A8EVM5
A	990	ASP	-	expression tag	UNP A8EVM5
A	991	LYS	-	expression tag	UNP A8EVM5
A	992	GLY	-	expression tag	UNP A8EVM5
A	993	SER	-	expression tag	UNP A8EVM5
A	994	LEU	-	expression tag	UNP A8EVM5
A	995	VAL	-	expression tag	UNP A8EVM5
A	996	PRO	-	expression tag	UNP A8EVM5
A	997	ARG	-	expression tag	UNP A8EVM5
A	998	GLY	-	expression tag	UNP A8EVM5
A	999	SER	-	expression tag	UNP A8EVM5
A	1000	HIS	-	expression tag	UNP A8EVM5
A	1100	CYS	VAL	engineered mutation	UNP A8EVM5
A	1150	CYS	GLN	engineered mutation	UNP A8EVM5

- Molecule 2 is 3-[(3-CHOLAMIDOPROPYL)DIMETHYLAMMONIO]-1-PROPANESULFONATE (CCD ID: CPS) (formula: C₃₂H₅₈N₂O₇S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			31	26	1	4		
2	A	1	Total	C	N	O	0	0
			29	24	1	4		

- Molecule 3 is 1,2-DIMYRISTOYL-SN-GLYCERO-3-PHOSPHOCHOLINE (CCD ID: PX4) (formula: C₃₆H₇₃NO₈P).



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

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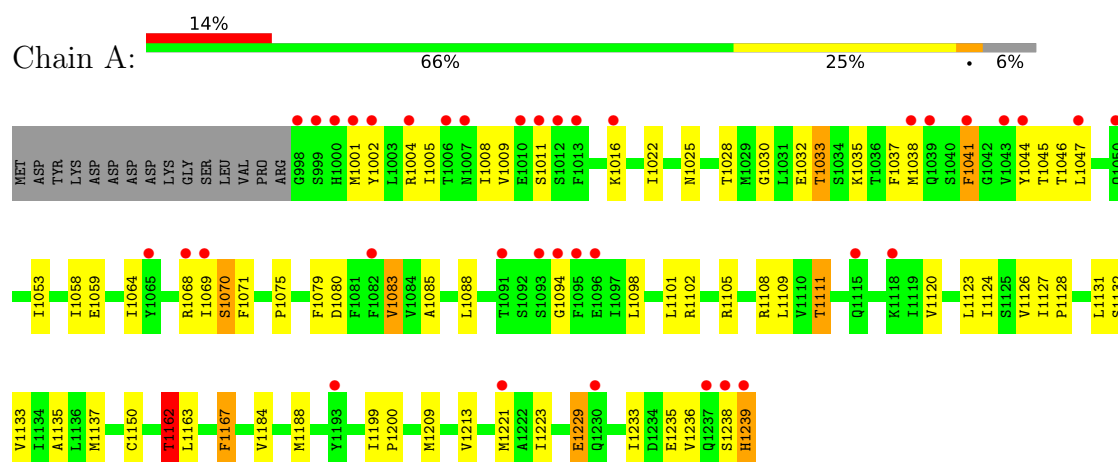
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	21	1	8	1		
3	A	1	Total	C	N	O	P	0	0
			42	32	1	8	1		
3	A	1	Total	C	N	O	P	0	0
			32	23	1	7	1		
3	A	1	Total	C				0	0
			13	13					

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: Ion transport protein



4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	125.83Å 125.83Å 189.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.53 – 2.89 44.53 – 2.89	Depositor EDS
% Data completeness (in resolution range)	97.6 (44.53-2.89) 97.6 (44.53-2.89)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.214 , 0.273 0.225 , 0.282	Depositor DCC
R_{free} test set	923 reflections (5.31%)	wwPDB-VP
Wilson B-factor (Å ²)	76.2	Xtriage
Anisotropy	0.004	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 72.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2201	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% 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: PX4, CPS

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	1.01	0/2031	1.54	5/2760 (0.2%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1162	THR	CA-CB-OG1	-6.42	99.98	109.60
1	A	1070	SER	CA-C-N	5.20	127.20	120.44
1	A	1070	SER	C-N-CA	5.20	127.20	120.44
1	A	1044	TYR	CA-C-N	5.10	127.11	120.28
1	A	1044	TYR	C-N-CA	5.10	127.11	120.28

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	1977	0	2039	44	0
2	A	60	0	81	2	0
3	A	164	0	239	10	0
All	All	2201	0	2359	50	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 11.

All (50) 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)
3:A:1305:PX4:O2	3:A:1305:PX4:H13	1.58	1.03
1:A:1008:ILE:O	1:A:1011:SER:HB3	1.63	0.97
1:A:1080:ASP:OD1	1:A:1111:THR:HG21	1.81	0.80
1:A:1162:THR:HG23	3:A:1303:PX4:H18	1.63	0.78
1:A:1239:HIS:ND1	1:A:1239:HIS:C	2.49	0.69
2:A:1301:CPS:H10B	2:A:1301:CPS:H21A	1.81	0.63
3:A:1305:PX4:O2	3:A:1305:PX4:C5	2.43	0.61
1:A:1032:GLU:HA	1:A:1038:MET:HE3	1.84	0.60
1:A:1008:ILE:O	1:A:1011:SER:CB	2.44	0.60
1:A:1030:GLY:O	1:A:1033:THR:HB	2.03	0.58
1:A:1083:VAL:HG13	1:A:1105:ARG:HA	1.86	0.57
1:A:1032:GLU:HG2	1:A:1045:THR:HG21	1.86	0.57
1:A:1032:GLU:CG	1:A:1045:THR:HG21	2.35	0.56
1:A:1108:ARG:O	1:A:1111:THR:HG22	2.06	0.56
1:A:1022:ILE:HG21	1:A:1109:LEU:HB2	1.89	0.53
1:A:1001:MET:O	1:A:1004:ARG:HB2	2.12	0.50
1:A:1005:ILE:O	1:A:1008:ILE:HB	2.14	0.48
1:A:1101:LEU:O	1:A:1102:ARG:C	2.56	0.48
1:A:1209:MET:HE3	3:A:1305:PX4:H45	1.96	0.47
1:A:1199:ILE:HB	1:A:1200:PRO:HD3	1.97	0.46
1:A:1127:ILE:N	1:A:1128:PRO:CD	2.79	0.46
3:A:1306:PX4:O2	3:A:1306:PX4:O8	2.33	0.46
1:A:1022:ILE:HD12	1:A:1108:ARG:HB2	1.97	0.46
1:A:1079:PHE:CE1	1:A:1083:VAL:HG21	2.51	0.46
1:A:1085:ALA:O	1:A:1088:LEU:HB2	2.16	0.45
1:A:1064:ILE:O	1:A:1068:ARG:N	2.48	0.45
1:A:1002:TYR:CD1	1:A:1002:TYR:C	2.94	0.45
1:A:1002:TYR:O	1:A:1005:ILE:HG12	2.16	0.44
1:A:1223:ILE:HG21	2:A:1302:CPS:H20	1.99	0.44
1:A:1075:PRO:HB2	3:A:1306:PX4:H46	1.98	0.44
1:A:1080:ASP:CG	1:A:1111:THR:HG21	2.41	0.44
1:A:1132:SER:O	1:A:1135:ALA:HB3	2.18	0.44
1:A:1041:PHE:CD1	1:A:1041:PHE:N	2.85	0.44
3:A:1304:PX4:O2	3:A:1304:PX4:H16	2.18	0.43
1:A:1041:PHE:N	1:A:1041:PHE:HD1	2.16	0.43
1:A:1184:VAL:O	1:A:1188:MET:HG3	2.18	0.43
1:A:1120:VAL:HG12	1:A:1124:ILE:HD12	2.00	0.43
1:A:1037:PHE:C	1:A:1037:PHE:CD1	2.96	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1162:THR:HG23	3:A:1303:PX4:C8	2.42	0.43
1:A:1079:PHE:CZ	1:A:1083:VAL:HG21	2.54	0.42
1:A:1123:LEU:O	1:A:1126:VAL:HG23	2.19	0.42
1:A:1209:MET:O	1:A:1213:VAL:HG23	2.20	0.42
1:A:1167:PHE:HB3	3:A:1305:PX4:O6	2.19	0.41
1:A:1058:ILE:O	1:A:1059:GLU:C	2.64	0.41
1:A:1025:ASN:OD1	1:A:1105:ARG:NH1	2.41	0.41
1:A:1229:GLU:O	1:A:1233:ILE:HB	2.20	0.41
3:A:1303:PX4:C13	3:A:1305:PX4:O8	2.68	0.41
1:A:1035:LYS:HA	1:A:1038:MET:HB2	2.02	0.41
1:A:1071:PHE:CD1	1:A:1071:PHE:C	2.97	0.41
1:A:1047:LEU:HD13	1:A:1047:LEU:HA	1.95	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	240/257 (93%)	209 (87%)	30 (12%)	1 (0%)	30 59

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1094	GLY

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	222/236 (94%)	197 (89%)	25 (11%)	5 19

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1009	VAL
1	A	1016	LYS
1	A	1028	THR
1	A	1033	THR
1	A	1041	PHE
1	A	1046	THR
1	A	1053	ILE
1	A	1069	ILE
1	A	1070	SER
1	A	1083	VAL
1	A	1098	LEU
1	A	1111	THR
1	A	1131	LEU
1	A	1133	VAL
1	A	1137	MET
1	A	1150	CYS
1	A	1162	THR
1	A	1163	LEU
1	A	1167	PHE
1	A	1221	MET
1	A	1229	GLU
1	A	1235	GLU
1	A	1236	VAL
1	A	1238	SER
1	A	1239	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1039	GLN
1	A	1237	GLN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 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	$\# Z > 2$
3	PX4	A	1306	-	31,31,45	0.49	0	36,38,53	0.53	0
2	CPS	A	1301	-	34,34,45	0.56	0	53,53,70	0.88	1 (1%)
3	PX4	A	1307	-	12,12,45	0.27	0	11,11,53	0.23	0
3	PX4	A	1303	-	45,45,45	0.37	0	51,53,53	0.72	2 (3%)
2	CPS	A	1302	-	32,32,45	0.47	0	51,51,70	0.75	0
3	PX4	A	1304	-	30,30,45	0.50	0	36,38,53	0.58	0
3	PX4	A	1305	-	41,41,45	0.39	0	47,49,53	0.96	1 (2%)

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
3	PX4	A	1306	-	-	17/34/34/49	-
2	CPS	A	1301	-	-	3/12/77/90	0/4/4/4
3	PX4	A	1307	-	-	6/10/10/49	-
3	PX4	A	1303	-	-	20/49/49/49	-
2	CPS	A	1302	-	-	0/9/74/90	0/4/4/4
3	PX4	A	1304	-	-	8/34/34/49	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PX4	A	1305	-	-	28/45/45/49	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1305	PX4	O7-C23-C24	4.29	120.76	111.48
3	A	1303	PX4	O7-C23-C24	3.36	118.75	111.48
3	A	1303	PX4	O7-C7-C6	2.70	118.03	108.34
2	A	1301	CPS	C6-C5-C4	-2.23	105.38	107.42

There are no chirality outliers.

All (82) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1303	PX4	C1-O3-P1-O1
3	A	1303	PX4	C1-O3-P1-O2
3	A	1303	PX4	C1-O3-P1-O4
3	A	1303	PX4	C6-O4-P1-O3
3	A	1304	PX4	C1-O3-P1-O4
3	A	1304	PX4	O3-C1-C2-N1
3	A	1304	PX4	C7-C6-O4-P1
3	A	1304	PX4	O7-C7-C8-O5
3	A	1305	PX4	C1-O3-P1-O1
3	A	1305	PX4	C1-O3-P1-O4
3	A	1305	PX4	C2-C1-O3-P1
3	A	1305	PX4	O3-C1-C2-N1
3	A	1305	PX4	C24-C23-O7-C7
3	A	1306	PX4	C1-O3-P1-O1
3	A	1306	PX4	C1-O3-P1-O4
3	A	1306	PX4	C6-O4-P1-O1
3	A	1306	PX4	C6-O4-P1-O3
3	A	1306	PX4	C24-C23-O7-C7
3	A	1305	PX4	O8-C23-O7-C7
3	A	1303	PX4	O6-C9-O5-C8
3	A	1306	PX4	O8-C23-O7-C7
3	A	1303	PX4	C10-C9-O5-C8
3	A	1303	PX4	C24-C23-O7-C7
2	A	1301	CPS	C21-C20-C22-C23
3	A	1305	PX4	C1-C2-N1-C5
3	A	1303	PX4	O4-C6-C7-O7

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Mol	Chain	Res	Type	Atoms
3	A	1304	PX4	C23-C24-C25-C26
3	A	1303	PX4	O8-C23-O7-C7
3	A	1306	PX4	C23-C24-C25-C26
3	A	1305	PX4	C1-C2-N1-C4
3	A	1305	PX4	C1-C2-N1-C3
2	A	1301	CPS	C20-C22-C23-C24
3	A	1305	PX4	C11-C12-C13-C14
3	A	1307	PX4	C25-C26-C27-C28
3	A	1305	PX4	C26-C27-C28-C29
3	A	1307	PX4	C26-C27-C28-C29
3	A	1306	PX4	C26-C27-C28-C29
3	A	1303	PX4	C11-C12-C13-C14
3	A	1305	PX4	C18-C19-C20-C21
3	A	1305	PX4	C16-C17-C18-C19
3	A	1306	PX4	C27-C28-C29-C30
3	A	1305	PX4	C17-C18-C19-C20
3	A	1305	PX4	C25-C26-C27-C28
3	A	1303	PX4	C16-C17-C18-C19
3	A	1307	PX4	C28-C29-C30-C31
3	A	1303	PX4	O4-C6-C7-C8
3	A	1305	PX4	C27-C28-C29-C30
3	A	1304	PX4	C6-C7-C8-O5
3	A	1307	PX4	C29-C30-C31-C32
3	A	1306	PX4	C31-C32-C33-C34
3	A	1306	PX4	C29-C30-C31-C32
3	A	1303	PX4	C28-C29-C30-C31
3	A	1306	PX4	C32-C33-C34-C35
3	A	1303	PX4	C15-C16-C17-C18
3	A	1303	PX4	C33-C34-C35-C36
3	A	1306	PX4	C33-C34-C35-C36
3	A	1306	PX4	C30-C31-C32-C33
3	A	1305	PX4	C10-C11-C12-C13
3	A	1303	PX4	C25-C26-C27-C28
3	A	1306	PX4	C2-C1-O3-P1
3	A	1303	PX4	C23-C24-C25-C26
3	A	1303	PX4	O7-C7-C8-O5
3	A	1306	PX4	O4-C6-C7-C8
3	A	1307	PX4	C31-C32-C33-C34
3	A	1305	PX4	C6-C7-C8-O5
3	A	1304	PX4	C25-C26-C27-C28
3	A	1304	PX4	C1-O3-P1-O2
3	A	1305	PX4	C6-O4-P1-O2

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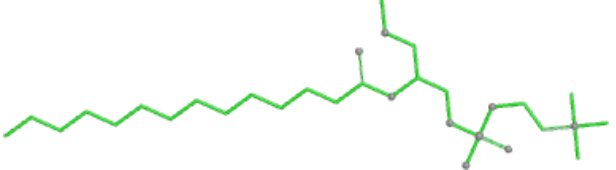
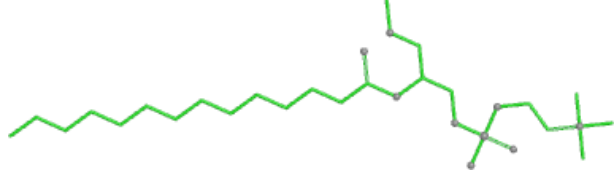
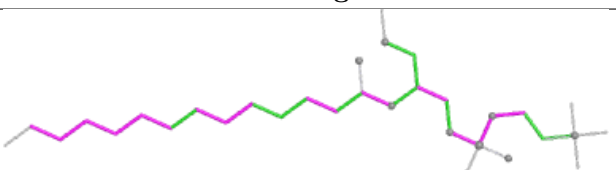
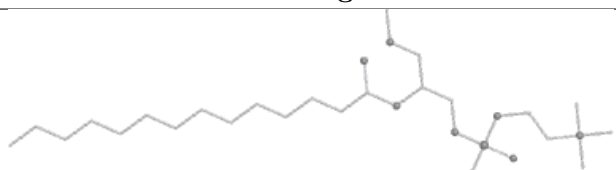
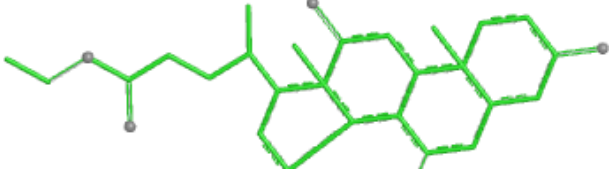
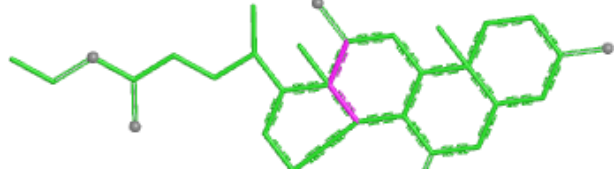
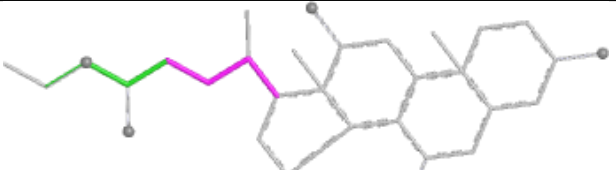
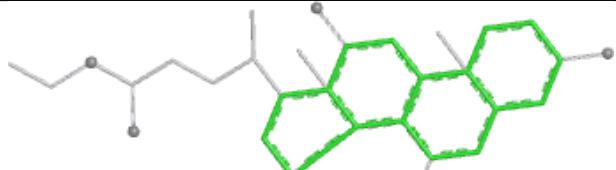
Mol	Chain	Res	Type	Atoms
3	A	1305	PX4	C6-O4-P1-O3
3	A	1306	PX4	C1-O3-P1-O2
3	A	1305	PX4	C8-C7-O7-C23
3	A	1305	PX4	C19-C20-C21-C22
3	A	1305	PX4	O7-C23-C24-C25
2	A	1301	CPS	C21-C20-C9-C5
3	A	1305	PX4	C24-C25-C26-C27
3	A	1305	PX4	C13-C14-C15-C16
3	A	1305	PX4	C6-C7-O7-C23
3	A	1305	PX4	C14-C15-C16-C17
3	A	1305	PX4	C7-C6-O4-P1
3	A	1307	PX4	C24-C25-C26-C27
3	A	1303	PX4	C30-C31-C32-C33
3	A	1303	PX4	C6-C7-C8-O5

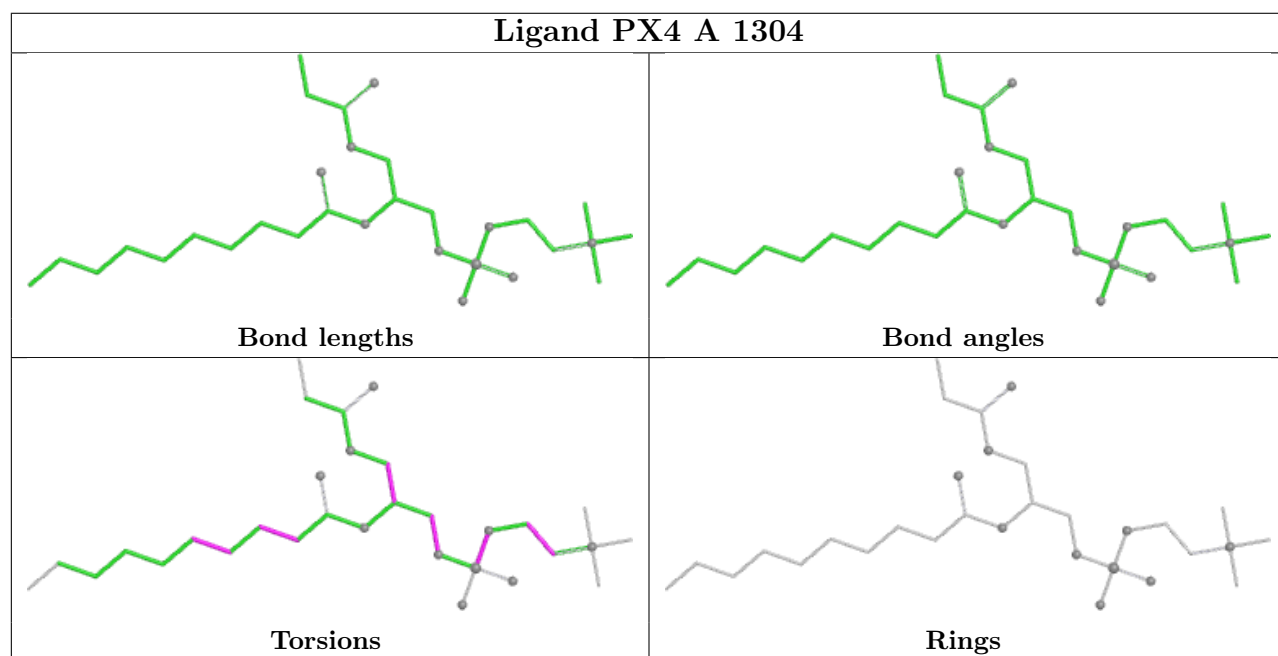
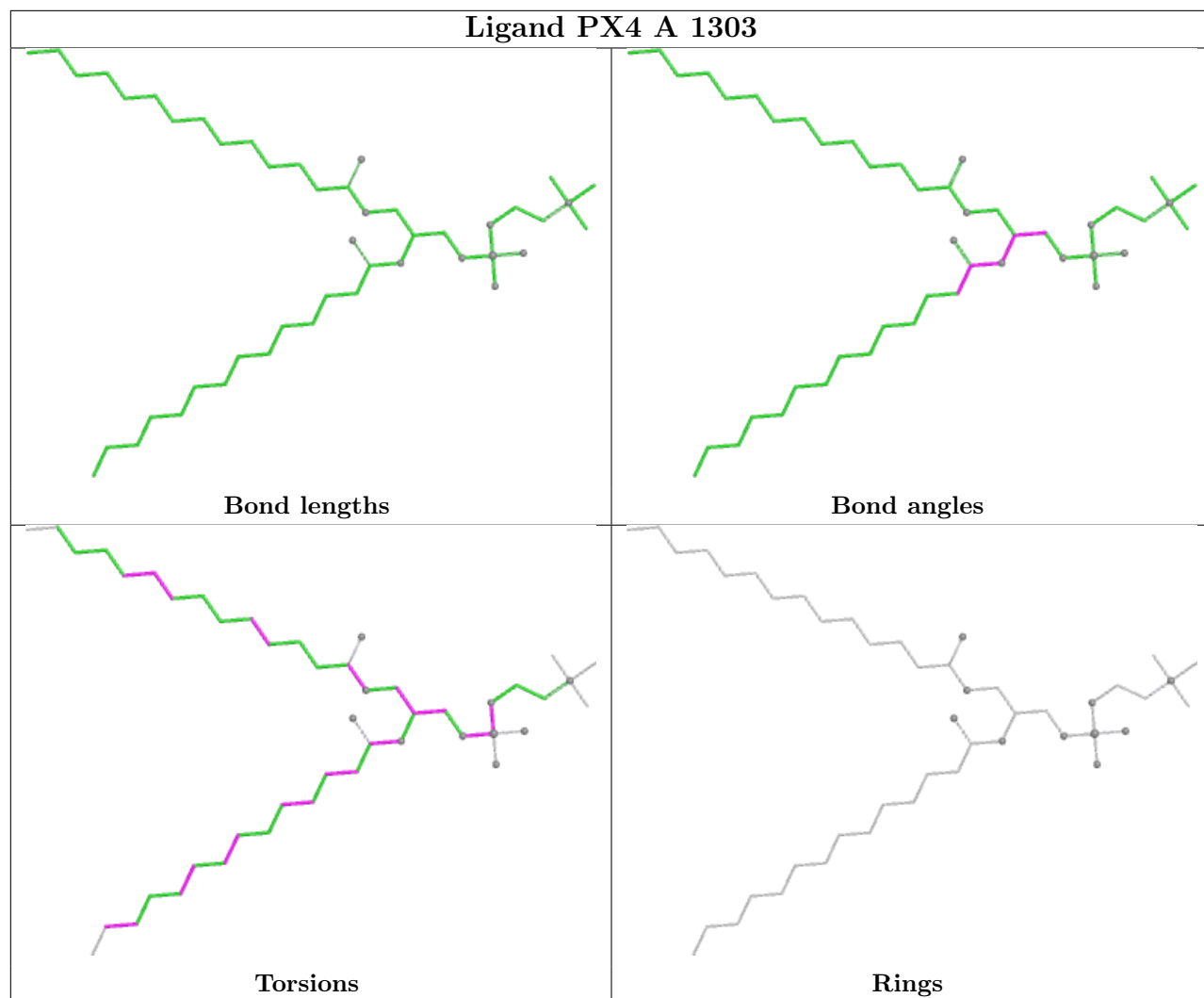
There are no ring outliers.

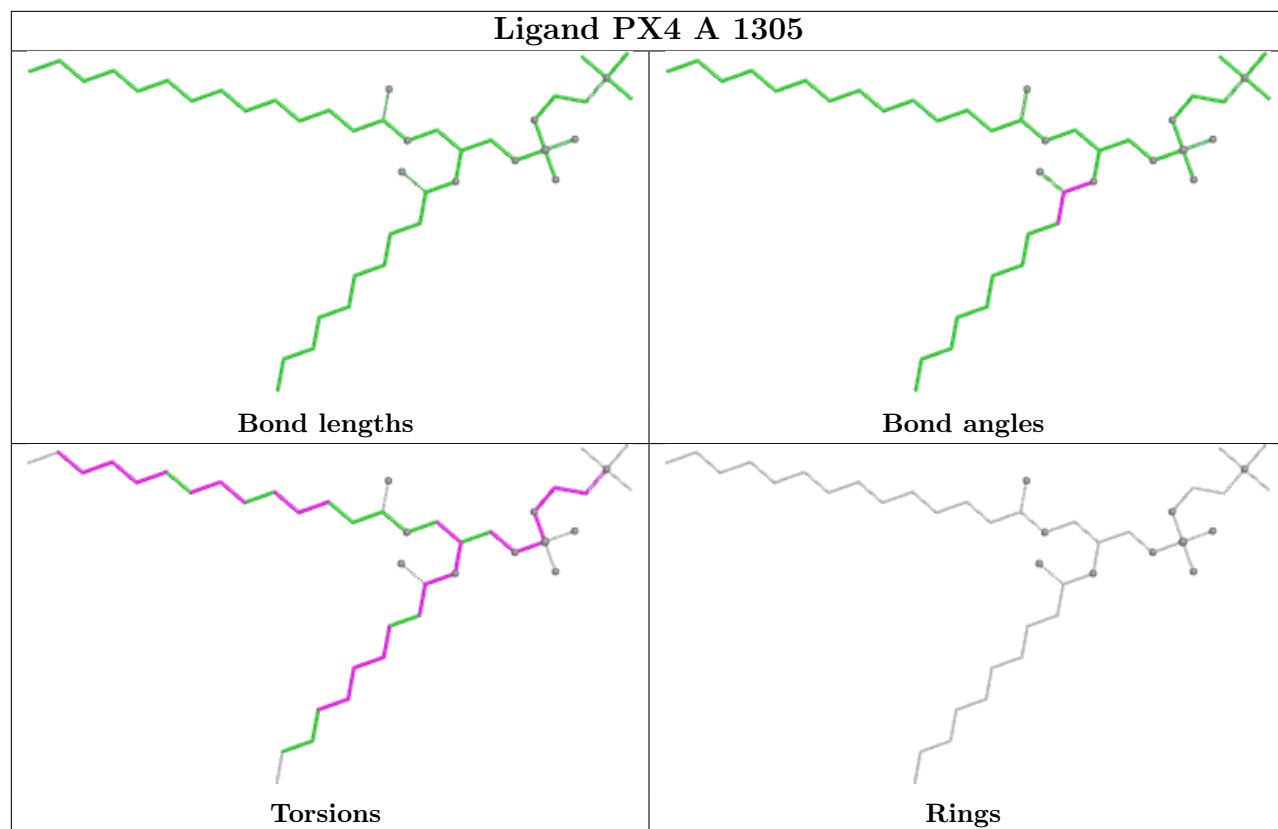
6 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1306	PX4	2	0
2	A	1301	CPS	1	0
3	A	1303	PX4	3	0
2	A	1302	CPS	1	0
3	A	1304	PX4	1	0
3	A	1305	PX4	5	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand PX4 A 1306	
 Bond lengths	 Bond angles
 Torsions	 Rings
Ligand CPS A 1301	
 Bond lengths	 Bond angles
 Torsions	 Rings
Ligand PX4 A 1307	
 Bond lengths	 Bond angles
 Torsions	 Rings





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	242/257 (94%)	0.67	37 (15%) 5 4	38, 75, 137, 180	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1095	PHE	6.0
1	A	1238	SER	5.5
1	A	1013	PHE	5.5
1	A	998	GLY	4.9
1	A	1065	TYR	4.6
1	A	1012	SER	4.2
1	A	1007	ASN	4.2
1	A	1002	TYR	3.9
1	A	1039	GLN	3.7
1	A	1239	HIS	3.5
1	A	1237	GLN	3.1
1	A	1094	GLY	3.0
1	A	1044	TYR	3.0
1	A	1096	GLU	3.0
1	A	1068	ARG	2.9
1	A	1010	GLU	2.9
1	A	1069	ILE	2.9
1	A	1016	LYS	2.7
1	A	1011	SER	2.7
1	A	1000	HIS	2.7
1	A	1118	LYS	2.6
1	A	1041	PHE	2.6
1	A	1093	SER	2.6
1	A	1115	GLN	2.4
1	A	1047	LEU	2.4
1	A	1004	ARG	2.3
1	A	999	SER	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	1038	MET	2.3
1	A	1082	PHE	2.3
1	A	1193	TYR	2.2
1	A	1001	MET	2.2
1	A	1050	GLN	2.2
1	A	1043	VAL	2.1
1	A	1221	MET	2.1
1	A	1006	THR	2.1
1	A	1091	THR	2.0
1	A	1230	GLN	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](#)

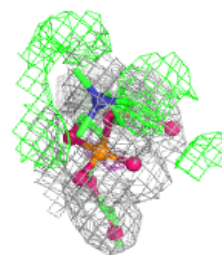
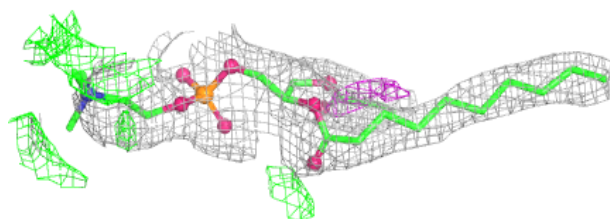
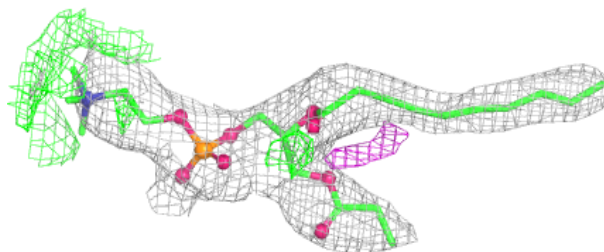
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	PX4	A	1304	31/46	0.78	0.30	64,95,153,157	0
3	PX4	A	1306	32/46	0.87	0.24	88,116,162,181	0
3	PX4	A	1303	46/46	0.89	0.20	69,83,127,131	0
2	CPS	A	1301	31/42	0.90	0.14	76,86,100,103	0
3	PX4	A	1305	42/46	0.94	0.17	64,86,109,118	0
3	PX4	A	1307	13/46	0.94	0.15	60,68,81,81	0
2	CPS	A	1302	29/42	0.95	0.10	66,88,106,121	0

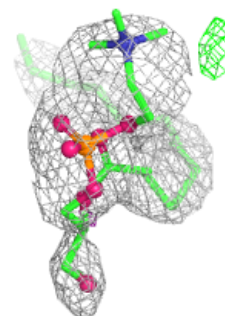
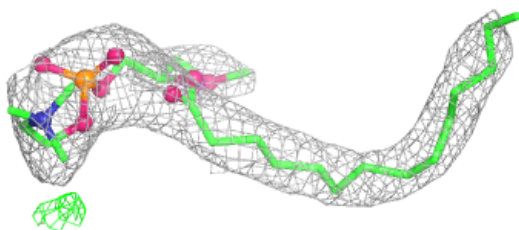
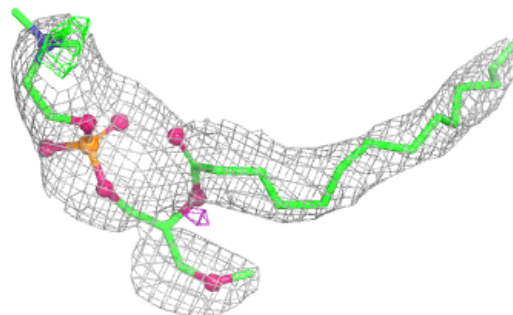
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around PX4 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)

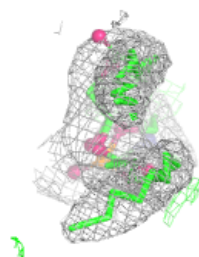
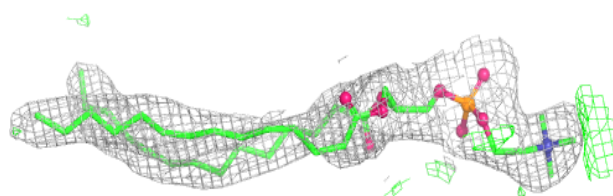
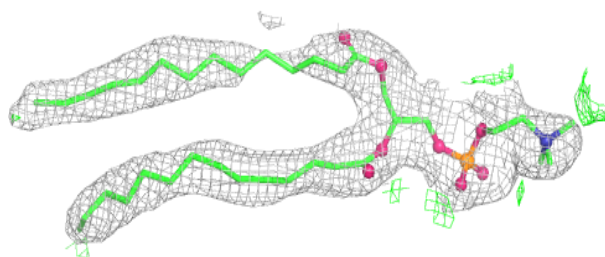
**Electron density around PX4 A 1306:**

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

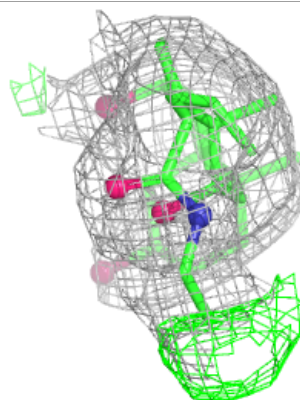
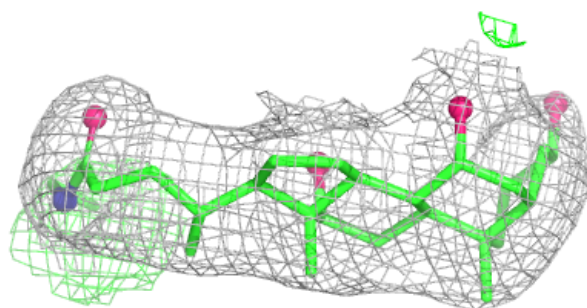
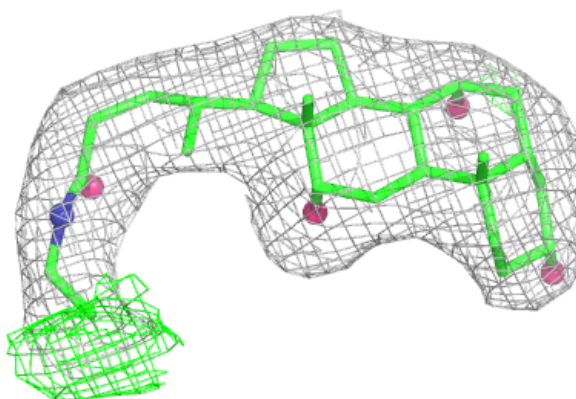


Electron density around PX4 A 1303:

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

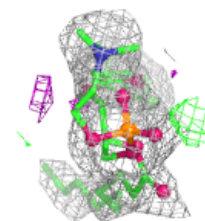
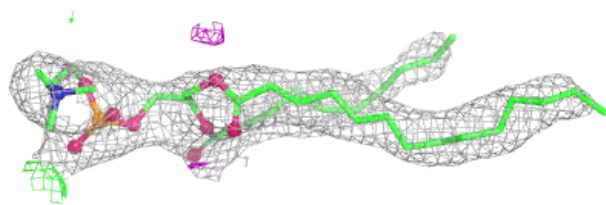
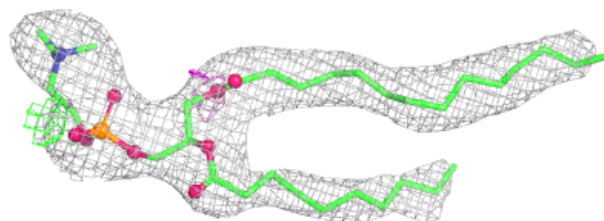
**Electron density around CPS A 1301:**

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and green (positive)

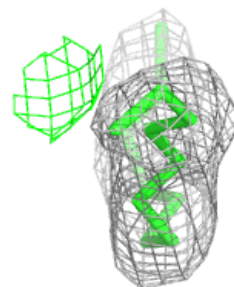
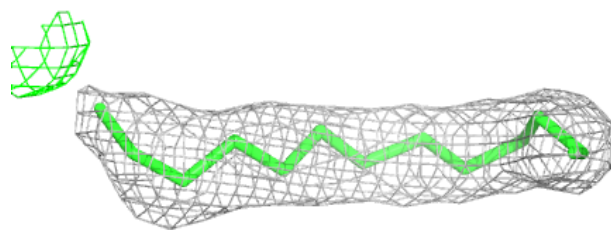
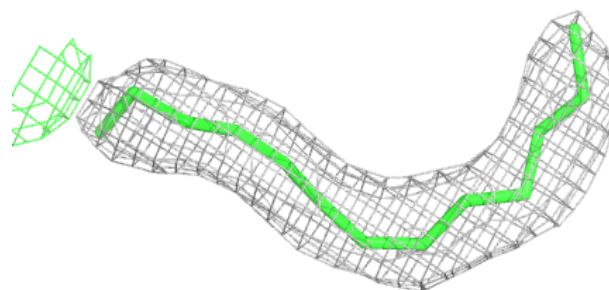


Electron density around PX4 A 1305:

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

**Electron density around PX4 A 1307:**

$2mF_o-DF_c$ (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.