



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 20, 2020 – 05:36 PM BST

PDB ID : 4EKW
Title : Crystal structure of the NavAb voltage-gated sodium channel (wild-type, 3.2 Å)
Authors : Payandeh, J.; Gamal El-Din, T.M.; Scheuer, T.; Zheng, N.; Catterall, W.A.
Deposited on : 2012-04-10
Resolution : 3.21 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

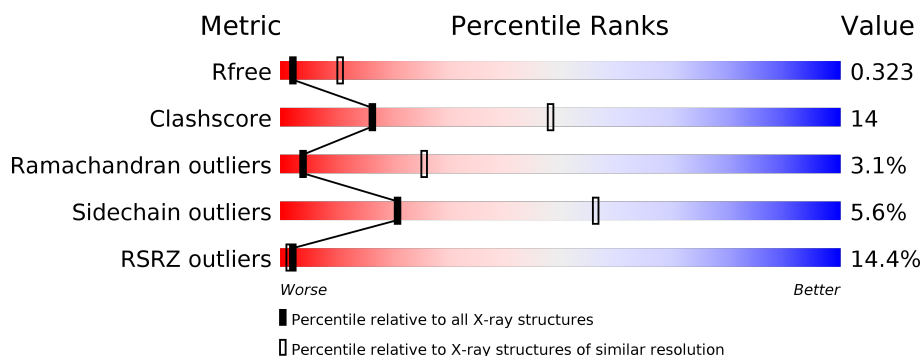
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.21 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	1335 (3.24-3.20)
Clashscore	141614	1460 (3.24-3.20)
Ramachandran outliers	138981	1437 (3.24-3.20)
Sidechain outliers	138945	1436 (3.24-3.20)
RSRZ outliers	127900	1291 (3.24-3.20)

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	285	<div> <div>11%</div> <div> <div>52%</div> <div>23%</div> <div>•</div> <div>23%</div> </div> </div>
1	B	285	<div> <div>11%</div> <div> <div>41%</div> <div>31%</div> <div>•</div> <div>23%</div> </div> </div>
1	C	285	<div> <div>8%</div> <div> <div>49%</div> <div>24%</div> <div>•</div> <div>25%</div> </div> </div>
1	D	285	<div> <div>12%</div> <div> <div>43%</div> <div>31%</div> <div>•</div> <div>24%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7218 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	219	Total	C	N	O	Se	0	0	0
			1799	1228	268	292	11			
1	B	219	Total	C	N	O	Se	0	0	0
			1799	1228	268	292	11			
1	C	213	Total	C	N	O	Se	0	0	0
			1756	1199	262	284	11			
1	D	217	Total	C	N	O	Se	0	0	0
			1784	1219	266	288	11			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	983	MSE	-	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
B	983	MSE	-	INITIATING METHIONINE	UNP A8EVM5
B	984	ASP	-	EXPRESSION TAG	UNP A8EVM5
B	985	TYR	-	EXPRESSION TAG	UNP A8EVM5

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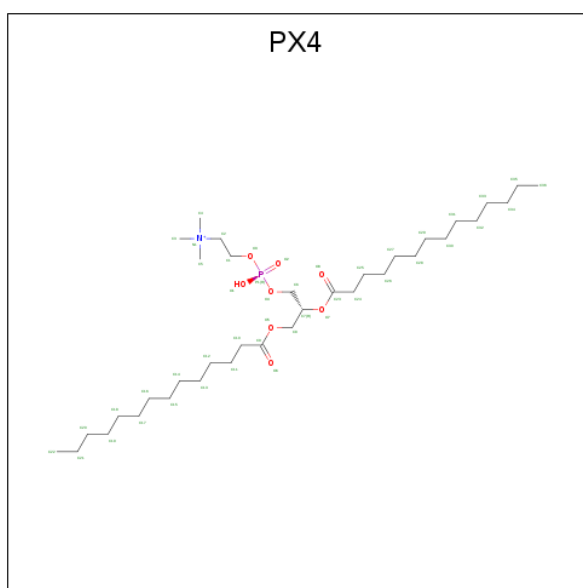
Chain	Residue	Modelled	Actual	Comment	Reference
B	986	LYS	-	EXPRESSION TAG	UNP A8EVM5
B	987	ASP	-	EXPRESSION TAG	UNP A8EVM5
B	988	ASP	-	EXPRESSION TAG	UNP A8EVM5
B	989	ASP	-	EXPRESSION TAG	UNP A8EVM5
B	990	ASP	-	EXPRESSION TAG	UNP A8EVM5
B	991	LYS	-	EXPRESSION TAG	UNP A8EVM5
B	992	GLY	-	EXPRESSION TAG	UNP A8EVM5
B	993	SER	-	EXPRESSION TAG	UNP A8EVM5
B	994	LEU	-	EXPRESSION TAG	UNP A8EVM5
B	995	VAL	-	EXPRESSION TAG	UNP A8EVM5
B	996	PRO	-	EXPRESSION TAG	UNP A8EVM5
B	997	ARG	-	EXPRESSION TAG	UNP A8EVM5
B	998	GLY	-	EXPRESSION TAG	UNP A8EVM5
B	999	SER	-	EXPRESSION TAG	UNP A8EVM5
B	1000	HIS	-	EXPRESSION TAG	UNP A8EVM5
C	983	MSE	-	INITIATING METHIONINE	UNP A8EVM5
C	984	ASP	-	EXPRESSION TAG	UNP A8EVM5
C	985	TYR	-	EXPRESSION TAG	UNP A8EVM5
C	986	LYS	-	EXPRESSION TAG	UNP A8EVM5
C	987	ASP	-	EXPRESSION TAG	UNP A8EVM5
C	988	ASP	-	EXPRESSION TAG	UNP A8EVM5
C	989	ASP	-	EXPRESSION TAG	UNP A8EVM5
C	990	ASP	-	EXPRESSION TAG	UNP A8EVM5
C	991	LYS	-	EXPRESSION TAG	UNP A8EVM5
C	992	GLY	-	EXPRESSION TAG	UNP A8EVM5
C	993	SER	-	EXPRESSION TAG	UNP A8EVM5
C	994	LEU	-	EXPRESSION TAG	UNP A8EVM5
C	995	VAL	-	EXPRESSION TAG	UNP A8EVM5
C	996	PRO	-	EXPRESSION TAG	UNP A8EVM5
C	997	ARG	-	EXPRESSION TAG	UNP A8EVM5
C	998	GLY	-	EXPRESSION TAG	UNP A8EVM5
C	999	SER	-	EXPRESSION TAG	UNP A8EVM5
C	1000	HIS	-	EXPRESSION TAG	UNP A8EVM5
D	983	MSE	-	INITIATING METHIONINE	UNP A8EVM5
D	984	ASP	-	EXPRESSION TAG	UNP A8EVM5
D	985	TYR	-	EXPRESSION TAG	UNP A8EVM5
D	986	LYS	-	EXPRESSION TAG	UNP A8EVM5
D	987	ASP	-	EXPRESSION TAG	UNP A8EVM5
D	988	ASP	-	EXPRESSION TAG	UNP A8EVM5
D	989	ASP	-	EXPRESSION TAG	UNP A8EVM5
D	990	ASP	-	EXPRESSION TAG	UNP A8EVM5
D	991	LYS	-	EXPRESSION TAG	UNP A8EVM5

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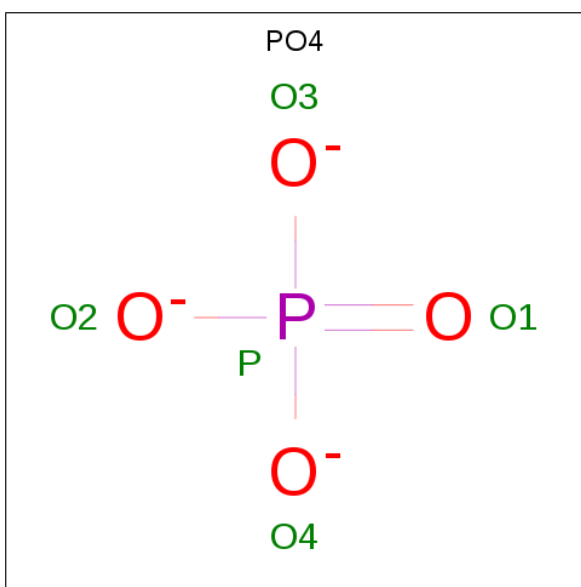
Chain	Residue	Modelled	Actual	Comment	Reference
D	992	GLY	-	EXPRESSION TAG	UNP A8EVM5
D	993	SER	-	EXPRESSION TAG	UNP A8EVM5
D	994	LEU	-	EXPRESSION TAG	UNP A8EVM5
D	995	VAL	-	EXPRESSION TAG	UNP A8EVM5
D	996	PRO	-	EXPRESSION TAG	UNP A8EVM5
D	997	ARG	-	EXPRESSION TAG	UNP A8EVM5
D	998	GLY	-	EXPRESSION TAG	UNP A8EVM5
D	999	SER	-	EXPRESSION TAG	UNP A8EVM5
D	1000	HIS	-	EXPRESSION TAG	UNP A8EVM5

- Molecule 2 is 1,2-DIMYRISTOYL-SN-GLYCERO-3-PHOSPHOCHOLINE (three-letter code: PX4) (formula: C₃₆H₇₃NO₈P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			14	6	7	1		
2	B	1	Total	C	O	P	0	0
			9	3	5	1		
2	B	1	Total	C	O	P	0	0
			6	1	4	1		
2	C	1	Total	C	O	P	0	0
			20	12	7	1		
2	D	1	Total	C	N	O	P	0
			8	2	1	4	1	

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	D	1	Total	O	P	0	0
			5	4	1		

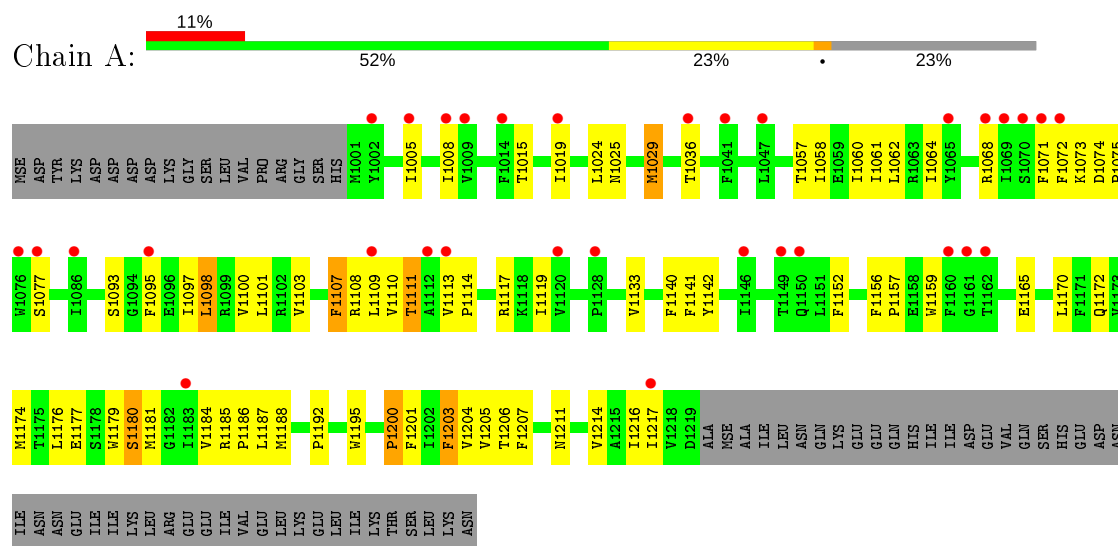
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	O	0	0
			1	1		
4	B	1	Total	O	0	0
			1	1		
4	D	1	Total	O	0	0
			1	1		

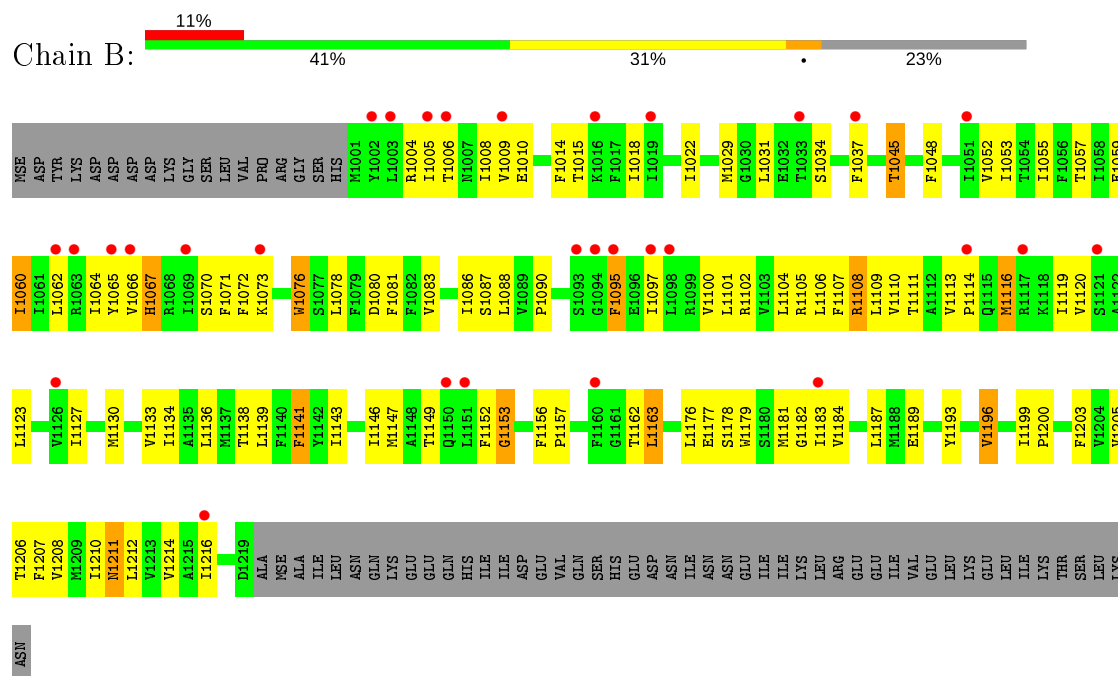
3 Residue-property plots

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.

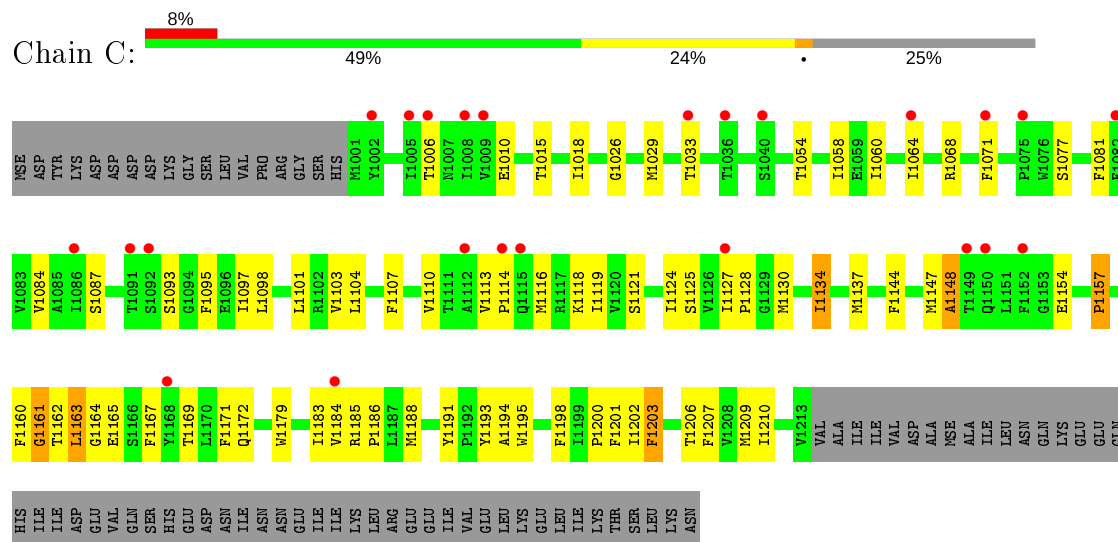
• Molecule 1: Ion transport protein



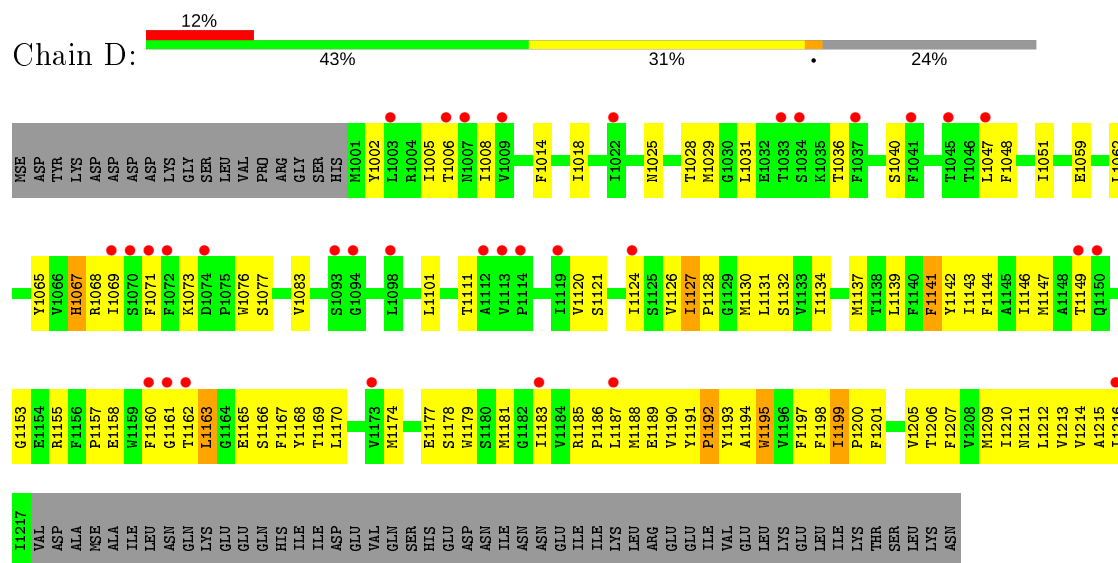
• Molecule 1: Ion transport protein



- Molecule 1: Ion transport protein



- Molecule 1: Ion transport protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 42	Depositor
Cell constants a, b, c, α , β , γ	125.86Å 125.86Å 192.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.21 48.60 – 3.21	Depositor EDS
% Data completeness (in resolution range)	96.4 (50.00-3.21) 96.6 (48.60-3.21)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.21 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.308 , 0.322 0.315 , 0.323	Depositor DCC
R_{free} test set	2401 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	99.5	Xtriage
Anisotropy	0.358	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 62.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.448 for h,-k,-l	Xtriage
Reported twinning fraction	0.496 for H, K, L 0.504 for K, H, -L	Depositor
Outliers	0 of 47387 reflections	Xtriage
F_o, F_c correlation	0.81	EDS
Total number of atoms	7218	wwPDB-VP
Average B, all atoms (Å ²)	102.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 25.44 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.1712e-03.*

¹ Intensities estimated from amplitudes.

² 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: PO4, PX4

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	0.56	0/1840	0.63	0/2487
1	B	0.57	1/1840 (0.1%)	0.62	0/2487
1	C	0.54	0/1797	0.59	0/2427
1	D	0.55	1/1825 (0.1%)	0.60	0/2466
All	All	0.56	2/7302 (0.0%)	0.61	0/9867

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	1195	TRP	CD2-CE2	5.07	1.47	1.41
1	B	1076	TRP	CD2-CE2	5.05	1.47	1.41

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1799	0	1875	47	1
1	B	1799	0	1875	68	1
1	C	1756	0	1826	47	0
1	D	1784	0	1862	69	0
2	A	14	0	5	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	15	0	4	0	0
2	C	20	0	17	0	0
2	D	8	0	4	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	5	0	0	0	0
3	D	5	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	D	1	0	0	0	0
All	All	7218	0	7468	210	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 14.

All (210) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1130:MSE:HE3	1:B:1212:LEU:HD11	1.36	1.07
1:A:1141:PHE:HZ	1:A:1174:MSE:SE	2.11	0.83
1:D:1131:LEU:HA	1:D:1134:ILE:HG22	1.62	0.82
1:C:1147:MSE:HA	1:C:1147:MSE:HE2	1.58	0.82
1:A:1174:MSE:HG3	1:A:1205:VAL:HG11	1.65	0.79
1:A:1107:PHE:CD1	1:B:1143:ILE:HD13	2.19	0.78
1:D:1174:MSE:HG3	1:D:1205:VAL:CG1	2.14	0.76
1:D:1174:MSE:HG3	1:D:1205:VAL:HG13	1.69	0.74
1:D:1183:ILE:O	1:D:1187:LEU:HB2	1.89	0.73
1:B:1018:ILE:HD11	1:B:1108:ARG:NH1	2.04	0.73
1:C:1103:VAL:HG21	1:D:1147:MSE:HG2	1.69	0.72
1:A:1058:ILE:O	1:A:1062:LEU:HG	1.90	0.72
1:A:1141:PHE:CZ	1:A:1174:MSE:SE	2.91	0.72
1:B:1018:ILE:HD11	1:B:1108:ARG:HH12	1.55	0.71
1:C:1209:MSE:HA	1:C:1209:MSE:HE2	1.71	0.71
1:D:1170:LEU:HB3	1:D:1201:PHE:CZ	2.26	0.70
1:D:1181:MSE:HA	1:D:1185:ARG:HG3	1.74	0.69
1:C:1200:PRO:HA	1:C:1203:PHE:HB2	1.74	0.69
1:D:1211:ASN:HA	1:D:1214:VAL:HG23	1.76	0.68
1:B:1130:MSE:CE	1:B:1212:LEU:HD11	2.22	0.67
1:A:1015:THR:O	1:A:1019:ILE:HG12	1.93	0.67
1:B:1060:ILE:O	1:B:1064:ILE:HG12	1.94	0.67
1:B:1076:TRP:HB3	1:B:1111:THR:HG23	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1130:MSE:HE3	1:B:1212:LEU:CD1	2.21	0.66
1:D:1206:THR:O	1:D:1210:ILE:HG13	1.96	0.65
1:A:1177:GLU:OE1	1:B:1178:SER:HA	1.96	0.65
1:D:1028:THR:HA	1:D:1031:LEU:HD12	1.78	0.65
1:A:1119:ILE:HD13	1:B:1136:LEU:HB2	1.78	0.64
1:A:1185:ARG:HB2	1:A:1186:PRO:HD3	1.79	0.64
1:B:1130:MSE:HE2	1:B:1134:ILE:HD11	1.80	0.63
1:B:1005:ILE:HA	1:B:1008:ILE:HD12	1.79	0.63
1:C:1185:ARG:HB2	1:C:1186:PRO:HD3	1.80	0.63
1:B:1119:ILE:O	1:B:1123:LEU:HG	1.98	0.62
1:D:1144:PHE:CD2	1:D:1201:PHE:HD1	2.17	0.61
1:B:1052:VAL:O	1:B:1055:ILE:HG22	2.01	0.61
1:C:1147:MSE:CE	1:C:1147:MSE:HA	2.30	0.60
1:C:1144:PHE:CE2	1:C:1201:PHE:HB2	2.36	0.60
1:A:1152:PHE:HB3	1:A:1187:LEU:HD11	1.84	0.60
1:D:1137:MSE:HA	1:D:1137:MSE:HE2	1.84	0.59
1:B:1022:ILE:HG21	1:B:1109:LEU:HB2	1.85	0.59
1:B:1182:GLY:O	1:B:1183:ILE:HD13	2.03	0.59
1:D:1120:VAL:O	1:D:1124:ILE:HG13	2.03	0.58
1:D:1191:TYR:O	1:D:1193:TYR:N	2.37	0.58
1:B:1153:GLY:HA2	1:B:1156:PHE:O	2.04	0.58
1:C:1103:VAL:HG21	1:D:1147:MSE:CG	2.33	0.58
1:B:1101:LEU:HB3	1:B:1104:LEU:HD12	1.84	0.58
1:D:1185:ARG:HB2	1:D:1186:PRO:HD3	1.86	0.57
1:D:1008:ILE:HG22	1:D:1014:PHE:HD1	1.68	0.57
1:B:1176:LEU:HA	1:B:1179:TRP:HB3	1.86	0.57
1:A:1214:VAL:HA	1:A:1217:ILE:HD12	1.86	0.57
1:B:1183:ILE:HG22	1:B:1184:VAL:N	2.20	0.56
1:D:1177:GLU:OE2	1:D:1178:SER:OG	2.20	0.56
1:B:1053:ILE:HD11	1:B:1088:LEU:HD23	1.86	0.56
1:A:1184:VAL:O	1:A:1188:MSE:HG3	2.06	0.56
1:A:1107:PHE:CG	1:B:1143:ILE:HD13	2.41	0.56
1:B:1133:VAL:HG11	1:B:1212:LEU:HB2	1.88	0.55
1:C:1015:THR:HA	1:C:1018:ILE:HD12	1.89	0.55
1:A:1180:SER:OG	1:A:1181:MSE:N	2.38	0.55
1:C:1033:THR:HG23	1:D:1149:THR:HG21	1.89	0.55
1:C:1157:PRO:HA	1:C:1161:GLY:HA3	1.89	0.55
1:B:1127:ILE:HA	1:B:1130:MSE:HB2	1.88	0.55
1:D:1210:ILE:O	1:D:1213:VAL:HB	2.06	0.55
1:C:1188:MSE:HA	1:C:1191:TYR:O	2.06	0.54
1:C:1193:TYR:O	1:C:1195:TRP:N	2.39	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1130:MSE:HG3	1:B:1216:ILE:HD11	1.90	0.54
1:D:1067:HIS:ND1	1:D:1067:HIS:N	2.56	0.54
1:D:1191:TYR:O	1:D:1194:ALA:N	2.40	0.54
1:C:1119:ILE:HG13	1:D:1132:SER:HB3	1.90	0.53
1:D:1124:ILE:HG23	1:D:1127:ILE:HD12	1.90	0.53
1:D:1155:ARG:O	1:D:1157:PRO:HD3	2.07	0.53
1:A:1100:VAL:O	1:A:1103:VAL:HG22	2.09	0.53
1:D:1207:PHE:HA	1:D:1210:ILE:HD12	1.90	0.53
1:B:1083:VAL:HA	1:B:1086:ILE:HD12	1.90	0.53
1:D:1008:ILE:HG22	1:D:1014:PHE:CD1	2.44	0.53
1:D:1146:ILE:HG12	1:D:1163:LEU:HD21	1.90	0.53
1:D:1005:ILE:HG23	1:D:1006:THR:H	1.74	0.53
1:C:1064:ILE:HD11	1:C:1081:PHE:HE1	1.73	0.52
1:D:1005:ILE:HD11	1:D:1062:LEU:HA	1.91	0.52
1:B:1087:SER:HA	1:B:1102:ARG:HG2	1.91	0.52
1:C:1107:PHE:CE1	1:D:1143:ILE:HD13	2.45	0.52
1:C:1127:ILE:HB	1:C:1128:PRO:HD3	1.92	0.52
1:A:1172:GLN:HE22	1:B:1181:MSE:SE	2.43	0.51
1:D:1071:PHE:CE1	1:D:1077:SER:HB3	2.46	0.51
1:D:1025:ASN:HA	1:D:1028:THR:HG22	1.93	0.51
1:C:1179:TRP:O	1:C:1184:VAL:HG23	2.11	0.51
1:D:1160:PHE:CZ	1:D:1198:PHE:HZ	2.29	0.51
1:C:1121:SER:HA	1:C:1124:ILE:HD12	1.93	0.51
1:D:1025:ASN:O	1:D:1029:MSE:HB2	2.11	0.51
1:A:1176:LEU:O	1:A:1179:TRP:HD1	1.92	0.50
1:D:1065:TYR:HE2	1:D:1068:ARG:HH21	1.59	0.50
1:B:1193:TYR:O	1:B:1196:VAL:HG23	2.11	0.50
1:B:1004:ARG:O	1:B:1008:ILE:HG13	2.12	0.50
1:D:1069:ILE:O	1:D:1073:LYS:HG3	2.11	0.50
1:A:1097:ILE:O	1:A:1101:LEU:HD12	2.11	0.50
1:D:1201:PHE:O	1:D:1205:VAL:HB	2.12	0.50
1:D:1201:PHE:CE2	1:D:1205:VAL:HG21	2.47	0.50
1:A:1103:VAL:HG23	1:B:1147:MSE:HG2	1.94	0.50
1:D:1160:PHE:CZ	1:D:1198:PHE:CZ	2.99	0.50
1:B:1067:HIS:HB2	1:B:1070:SER:HB2	1.93	0.49
1:B:1207:PHE:O	1:B:1211:ASN:ND2	2.45	0.49
1:B:1045:THR:HA	1:B:1048:PHE:HB3	1.94	0.49
1:D:1160:PHE:CZ	1:D:1169:THR:HG21	2.47	0.49
1:C:1033:THR:OG1	1:D:1163:LEU:HD23	2.12	0.49
1:D:1048:PHE:HA	1:D:1051:ILE:HD12	1.93	0.49
1:A:1156:PHE:CZ	1:D:1158:GLU:OE2	2.65	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1140:PHE:CZ	1:A:1201:PHE:HA	2.47	0.49
1:D:1076:TRP:HB3	1:D:1111:THR:HG23	1.94	0.49
1:C:1130:MSE:O	1:C:1134:ILE:HG13	2.12	0.49
1:A:1109:LEU:HG	1:B:1139:LEU:HD21	1.95	0.48
1:A:1114:PRO:HA	1:A:1117:ARG:HB2	1.94	0.48
1:C:1209:MSE:HA	1:C:1209:MSE:CE	2.41	0.48
1:B:1207:PHE:HA	1:B:1210:ILE:HD12	1.94	0.48
1:B:1106:LEU:C	1:B:1108:ARG:H	2.17	0.48
1:A:1098:LEU:HA	1:A:1101:LEU:HD12	1.96	0.47
1:B:1206:THR:HG22	1:B:1210:ILE:HD11	1.96	0.47
1:D:1174:MSE:HE3	1:D:1209:MSE:HE3	1.95	0.47
1:B:1162:THR:HG22	1:B:1163:LEU:N	2.30	0.47
1:D:1174:MSE:HG3	1:D:1205:VAL:HG11	1.94	0.47
1:A:1179:TRP:O	1:A:1184:VAL:HG23	2.15	0.47
1:A:1057:THR:O	1:A:1060:ILE:HG22	2.14	0.47
1:B:1006:THR:HG23	1:B:1066:VAL:HG13	1.97	0.47
1:C:1071:PHE:CE2	1:C:1077:SER:HB3	2.50	0.47
1:B:1208:VAL:HG12	1:B:1208:VAL:O	2.14	0.47
1:A:1140:PHE:HZ	1:A:1201:PHE:HA	1.80	0.46
1:B:1029:MSE:C	1:B:1031:LEU:H	2.17	0.46
1:B:1199:ILE:HB	1:B:1200:PRO:HD3	1.96	0.46
1:B:1146:ILE:O	1:B:1149:THR:N	2.47	0.46
1:C:1162:THR:O	1:C:1165:GLU:N	2.47	0.46
1:D:1158:GLU:HA	1:D:1158:GLU:OE1	2.15	0.46
1:D:1179:TRP:O	1:D:1183:ILE:HB	2.15	0.46
1:B:1059:GLU:HA	1:B:1062:LEU:HD12	1.97	0.46
1:B:1138:THR:HA	1:B:1141:PHE:HB2	1.98	0.46
1:C:1207:PHE:O	1:C:1210:ILE:HB	2.15	0.46
1:C:1162:THR:O	1:C:1163:LEU:C	2.54	0.46
1:A:1110:VAL:HG22	1:B:1139:LEU:HD23	1.98	0.46
1:A:1005:ILE:HA	1:A:1008:ILE:HD12	1.98	0.46
1:D:1191:TYR:O	1:D:1192:PRO:C	2.54	0.46
1:D:1008:ILE:O	1:D:1014:PHE:HB2	2.16	0.46
1:B:1110:VAL:HG11	1:B:1120:VAL:HG21	1.97	0.45
1:C:1098:LEU:HA	1:C:1101:LEU:HD12	1.98	0.45
1:D:1130:MSE:HE3	1:D:1215:ALA:HB3	1.98	0.45
1:A:1103:VAL:CG2	1:B:1147:MSE:HG2	2.45	0.45
1:C:1104:LEU:HD23	1:C:1107:PHE:CD1	2.51	0.45
1:B:1029:MSE:HE1	1:B:1105:ARG:HH21	1.81	0.45
1:A:1071:PHE:CE1	1:A:1077:SER:HB3	2.52	0.45
1:A:1188:MSE:CE	1:A:1195:TRP:HD1	2.29	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1203:PHE:O	1:A:1204:VAL:C	2.55	0.45
1:D:1128:PRO:HA	1:D:1131:LEU:HD12	1.98	0.45
1:C:1097:ILE:H	1:C:1097:ILE:HG13	1.61	0.45
1:A:1200:PRO:O	1:A:1204:VAL:N	2.41	0.45
1:C:1118:LYS:O	1:C:1121:SER:HB2	2.17	0.45
1:C:1148:ALA:HB1	1:C:1160:PHE:HD2	1.81	0.45
1:D:1126:VAL:HG11	1:D:1216:ILE:HD13	1.97	0.45
1:C:1084:VAL:O	1:C:1087:SER:HB3	2.17	0.44
1:A:1074:ASP:HA	1:A:1075:PRO:HD3	1.88	0.44
1:B:1113:VAL:HA	1:B:1114:PRO:HD3	1.77	0.44
1:A:1174:MSE:CG	1:A:1205:VAL:HG11	2.43	0.44
1:B:1100:VAL:C	1:B:1102:ARG:H	2.19	0.44
1:B:1203:PHE:O	1:B:1207:PHE:HB2	2.17	0.44
1:B:1057:THR:O	1:B:1060:ILE:HG22	2.17	0.44
1:A:1159:TRP:O	1:A:1165:GLU:O	2.35	0.44
1:D:1181:MSE:O	1:D:1186:PRO:HD3	2.17	0.44
1:C:1162:THR:O	1:C:1164:GLY:N	2.50	0.44
1:B:1178:SER:OG	1:B:1178:SER:O	2.34	0.44
1:C:1113:VAL:HA	1:C:1114:PRO:HD3	1.80	0.44
1:C:1026:GLY:O	1:D:1142:TYR:OH	2.35	0.43
1:D:1139:LEU:O	1:D:1143:ILE:HD12	2.18	0.43
1:B:1095:PHE:HB2	1:B:1097:ILE:HG13	2.01	0.43
1:A:1157:PRO:HG2	1:D:1155:ARG:O	2.17	0.43
1:B:1108:ARG:HH11	1:B:1108:ARG:HB3	1.82	0.43
1:B:1139:LEU:O	1:B:1143:ILE:HD12	2.18	0.43
1:D:1211:ASN:HA	1:D:1214:VAL:CG2	2.47	0.43
1:B:1005:ILE:O	1:B:1009:VAL:HG23	2.18	0.43
1:C:1193:TYR:C	1:C:1195:TRP:H	2.22	0.43
1:D:1166:SER:C	1:D:1168:TYR:H	2.22	0.43
1:A:1113:VAL:HA	1:A:1114:PRO:HD3	1.92	0.43
1:D:1188:MSE:C	1:D:1190:VAL:H	2.22	0.42
1:D:1199:ILE:HB	1:D:1200:PRO:HD3	2.02	0.42
1:C:1198:PHE:O	1:C:1202:ILE:HG13	2.18	0.42
1:A:1203:PHE:O	1:A:1206:THR:N	2.51	0.42
1:D:1130:MSE:O	1:D:1134:ILE:N	2.42	0.42
1:A:1071:PHE:C	1:A:1073:LYS:H	2.22	0.42
1:C:1107:PHE:CE1	1:D:1143:ILE:HG21	2.54	0.42
1:B:1080:ASP:OD2	1:B:1111:THR:HG21	2.19	0.42
1:A:1207:PHE:O	1:A:1211:ASN:HB2	2.19	0.42
1:B:1034:SER:HB3	1:B:1037:PHE:HB2	2.02	0.42
1:B:1072:PHE:HD2	1:B:1078:LEU:HD21	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1083:VAL:HG12	1:B:1105:ARG:HA	2.01	0.42
1:C:1179:TRP:O	1:C:1183:ILE:HB	2.19	0.42
1:C:1207:PHE:HA	1:C:1210:ILE:HD12	2.02	0.42
1:B:1189:GLU:HB3	1:D:1189:GLU:OE1	2.20	0.41
1:C:1137:MSE:HE2	1:C:1137:MSE:HA	2.02	0.41
1:B:1014:PHE:HZ	1:B:1062:LEU:HD13	1.84	0.41
1:A:1108:ARG:HA	1:A:1111:THR:OG1	2.21	0.41
1:D:1162:THR:OG1	1:D:1165:GLU:OE2	2.38	0.41
1:A:1061:ILE:HA	1:A:1064:ILE:HD12	2.01	0.41
1:A:1216:ILE:HG22	1:B:1214:VAL:HG21	2.02	0.41
1:B:1116:MSE:H	1:B:1116:MSE:HG2	1.61	0.41
1:C:1054:THR:O	1:C:1058:ILE:HG12	2.20	0.41
1:C:1206:THR:HA	1:C:1209:MSE:HB2	2.02	0.41
1:D:1197:PHE:O	1:D:1200:PRO:HD2	2.21	0.41
1:A:1170:LEU:HD22	1:A:1201:PHE:CZ	2.56	0.41
1:C:1169:THR:O	1:C:1172:GLN:HB3	2.21	0.41
1:B:1071:PHE:C	1:B:1073:LYS:H	2.24	0.41
1:A:1029:MSE:SE	1:A:1103:VAL:HG12	2.71	0.41
1:C:1110:VAL:HG22	1:C:1116:MSE:HG2	2.02	0.40
1:C:1160:PHE:O	1:C:1162:THR:N	2.53	0.40
1:D:1195:TRP:O	1:D:1199:ILE:HG12	2.20	0.40
1:D:1018:ILE:HG13	1:D:1059:GLU:OE2	2.21	0.40
1:C:1006:THR:O	1:C:1010:GLU:HG2	2.21	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)
1:A:1180:SER:OG	1:B:1177:GLU:OE2[2_565]	2.09	0.11

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	217/285 (76%)	182 (84%)	30 (14%)	5 (2%)	6	33
1	B	217/285 (76%)	180 (83%)	29 (13%)	8 (4%)	3	21
1	C	211/285 (74%)	185 (88%)	17 (8%)	9 (4%)	2	19
1	D	215/285 (75%)	185 (86%)	25 (12%)	5 (2%)	6	33
All	All	860/1140 (75%)	732 (85%)	101 (12%)	27 (3%)	4	26

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1068	ARG
1	B	1153	GLY
1	C	1125	SER
1	C	1148	ALA
1	C	1161	GLY
1	C	1194	ALA
1	D	1153	GLY
1	D	1161	GLY
1	A	1093	SER
1	A	1180	SER
1	B	1157	PRO
1	C	1068	ARG
1	D	1141	PHE
1	B	1081	PHE
1	B	1107	PHE
1	C	1163	LEU
1	D	1192	PRO
1	B	1152	PHE
1	C	1167	PHE
1	A	1072	PHE
1	B	1065	TYR
1	C	1093	SER
1	D	1199	ILE
1	A	1200	PRO
1	B	1205	VAL
1	C	1157	PRO
1	B	1090	PRO

5.3.2 Protein sidechains ⓘ

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	202/251 (80%)	190 (94%)	12 (6%)	19	54
1	B	202/251 (80%)	189 (94%)	13 (6%)	17	51
1	C	197/251 (78%)	190 (96%)	7 (4%)	35	68
1	D	200/251 (80%)	187 (94%)	13 (6%)	17	50
All	All	801/1004 (80%)	756 (94%)	45 (6%)	21	56

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1024	LEU
1	A	1025	ASN
1	A	1029	MSE
1	A	1036	THR
1	A	1095	PHE
1	A	1098	LEU
1	A	1107	PHE
1	A	1111	THR
1	A	1133	VAL
1	A	1142	TYR
1	A	1192	PRO
1	A	1203	PHE
1	B	1010	GLU
1	B	1015	THR
1	B	1045	THR
1	B	1060	ILE
1	B	1067	HIS
1	B	1095	PHE
1	B	1108	ARG
1	B	1116	MSE
1	B	1141	PHE
1	B	1163	LEU
1	B	1187	LEU
1	B	1196	VAL
1	B	1211	ASN
1	C	1029	MSE
1	C	1060	ILE
1	C	1095	PHE

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Mol	Chain	Res	Type
1	C	1134	ILE
1	C	1154	GLU
1	C	1171	PHE
1	C	1203	PHE
1	D	1002	TYR
1	D	1036	THR
1	D	1040	SER
1	D	1047	LEU
1	D	1067	HIS
1	D	1083	VAL
1	D	1101	LEU
1	D	1121	SER
1	D	1127	ILE
1	D	1141	PHE
1	D	1163	LEU
1	D	1167	PHE
1	D	1212	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1172	GLN
1	A	1211	ASN
1	B	1025	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 monosaccharides in this entry.

5.6 Ligand geometry

9 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	PO4	C	1302	-	4,4,4	0.95	0	6,6,6	0.62	0
3	PO4	A	1302	-	4,4,4	0.77	0	6,6,6	0.59	0
3	PO4	D	1302	-	4,4,4	0.73	0	6,6,6	0.55	0
3	PO4	B	1303	-	4,4,4	0.75	0	6,6,6	0.64	0
2	PX4	B	1302	-	5,5,45	0.76	0	7,7,53	1.04	0
2	PX4	D	1301	-	7,7,45	0.72	0	9,9,53	0.81	0
2	PX4	A	1301	-	13,13,45	1.22	1 (7%)	13,16,53	0.88	1 (7%)
2	PX4	C	1301	-	19,19,45	1.13	1 (5%)	21,23,53	0.89	1 (4%)
2	PX4	B	1301	-	8,8,45	0.51	0	10,10,53	1.03	0

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
2	PX4	B	1302	-	-	0/1/3/49	-
2	PX4	C	1301	-	-	9/21/21/49	-
2	PX4	B	1301	-	-	2/6/6/49	-
2	PX4	D	1301	-	-	0/5/5/49	-
2	PX4	A	1301	-	-	5/15/15/49	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1301	PX4	O5-C9	4.29	1.45	1.33
2	A	1301	PX4	O5-C9	3.60	1.47	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1301	PX4	O5-C9-C10	2.19	118.78	111.91
2	A	1301	PX4	O5-C8-C7	2.01	114.28	108.43

There are no chirality outliers.

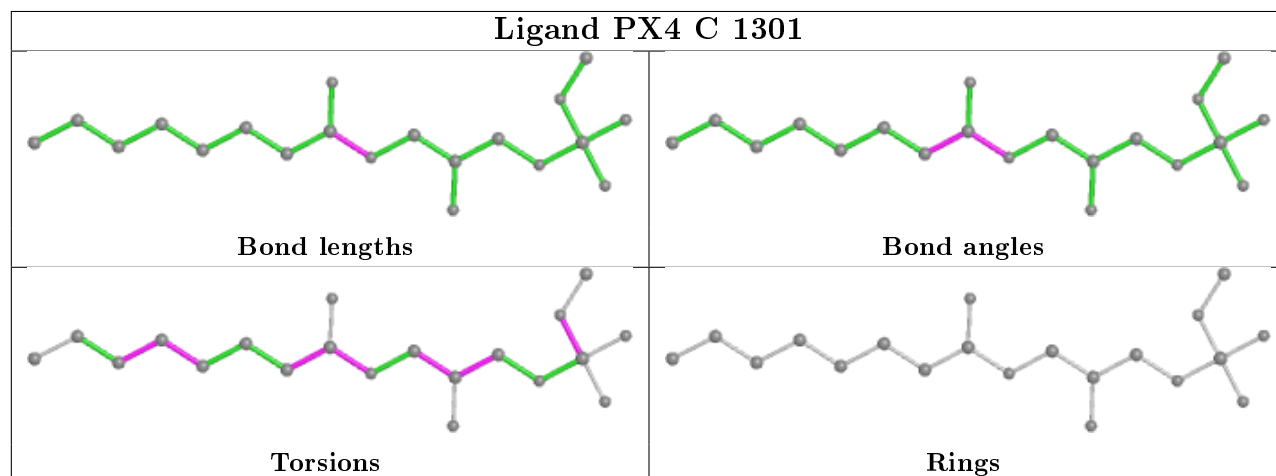
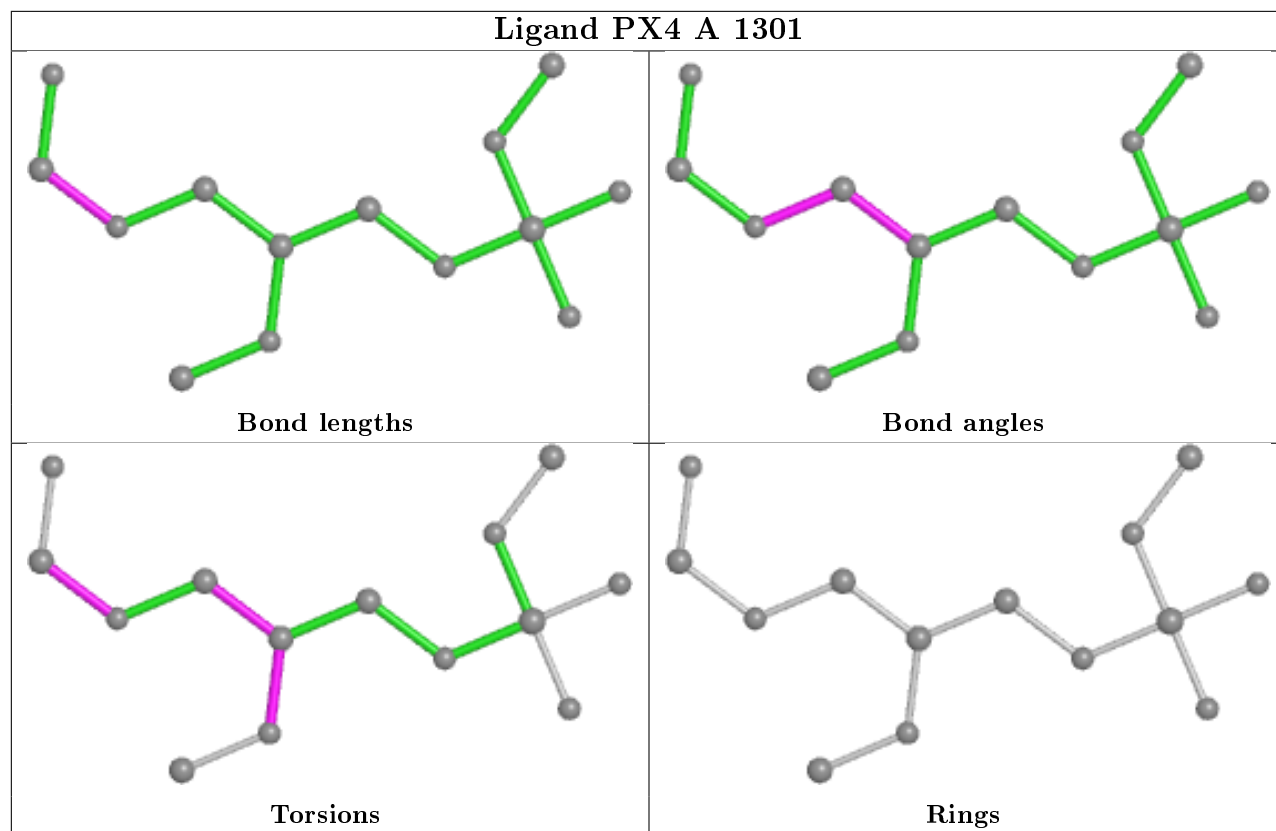
All (16) torsion outliers are listed below:

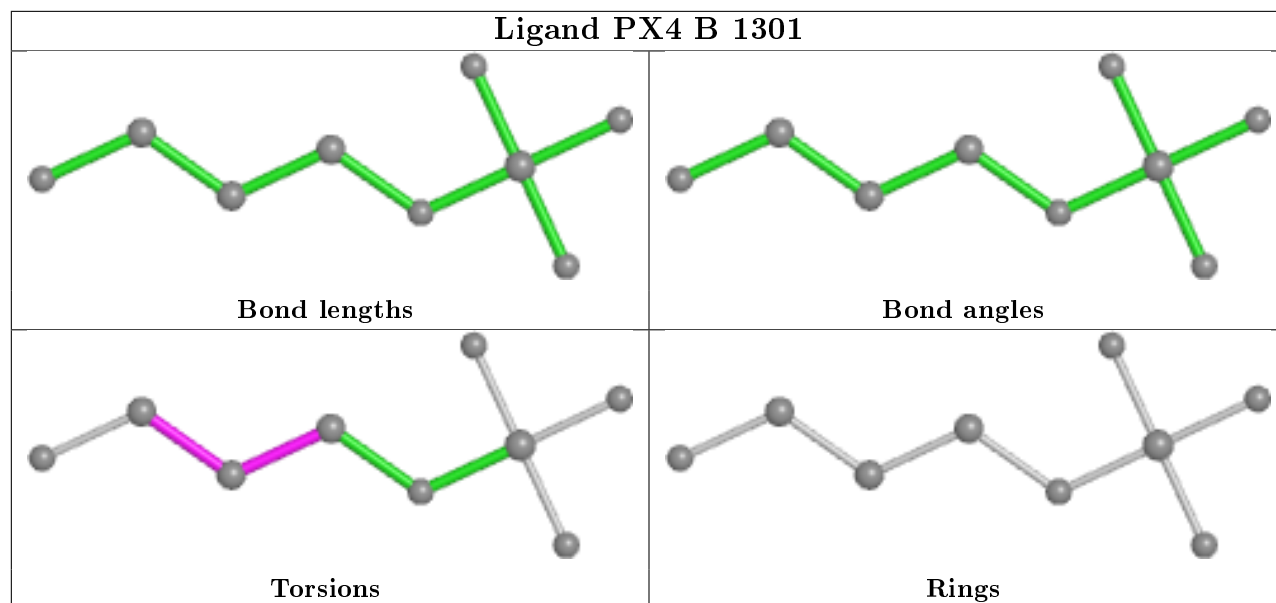
Mol	Chain	Res	Type	Atoms
2	A	1301	PX4	O6-C9-O5-C8
2	B	1301	PX4	O4-C6-C7-C8
2	B	1301	PX4	C6-C7-C8-O5
2	C	1301	PX4	O4-C6-C7-O7
2	C	1301	PX4	O4-C6-C7-C8
2	A	1301	PX4	C6-C7-C8-O5
2	C	1301	PX4	C10-C9-O5-C8
2	C	1301	PX4	C1-O3-P1-O2
2	C	1301	PX4	C12-C13-C14-C15
2	C	1301	PX4	O6-C9-O5-C8
2	A	1301	PX4	O7-C7-C8-O5
2	C	1301	PX4	C11-C12-C13-C14
2	A	1301	PX4	C6-C7-O7-C23
2	A	1301	PX4	C8-C7-O7-C23
2	C	1301	PX4	C11-C10-C9-O5
2	C	1301	PX4	O7-C7-C8-O5

There are no ring outliers.

No monomer is involved in short contacts.

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	208/285 (72%)	1.00	32 (15%) 2 1	67, 101, 119, 131	0
1	B	208/285 (72%)	0.96	30 (14%) 2 1	71, 106, 126, 145	0
1	C	202/285 (70%)	0.92	24 (11%) 4 3	67, 117, 132, 147	0
1	D	206/285 (72%)	1.11	33 (16%) 1 1	65, 105, 126, 135	0
All	All	824/1140 (72%)	1.00	119 (14%) 2 1	65, 106, 129, 147	0

All (119) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	1007	ASN	6.8
1	B	1006	THR	6.4
1	A	1009	VAL	6.1
1	D	1033	THR	6.1
1	D	1009	VAL	6.0
1	B	1073	LYS	5.8
1	D	1094	GLY	5.6
1	D	1183	ILE	5.2
1	A	1014	PHE	5.2
1	D	1162	THR	5.0
1	A	1005	ILE	4.8
1	D	1006	THR	4.8
1	C	1115	GLN	4.7
1	D	1150	GLN	4.6
1	C	1040	SER	4.5
1	A	1036	THR	4.4
1	B	1093	SER	4.2
1	A	1002	TYR	4.2
1	C	1092	SER	4.0
1	B	1160	PHE	4.0
1	C	1114	PRO	4.0

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Mol	Chain	Res	Type	RSRZ
1	C	1009	VAL	3.9
1	C	1071	PHE	3.9
1	B	1098	LEU	3.8
1	A	1065	TYR	3.8
1	C	1086	ILE	3.8
1	C	1005	ILE	3.7
1	D	1160	PHE	3.6
1	A	1183	ILE	3.5
1	D	1093	SER	3.5
1	B	1019	ILE	3.4
1	B	1150	GLN	3.4
1	B	1062	LEU	3.4
1	B	1094	GLY	3.3
1	D	1161	GLY	3.3
1	A	1076	TRP	3.3
1	B	1095	PHE	3.2
1	D	1045	THR	3.2
1	B	1002	TYR	3.2
1	B	1066	VAL	3.2
1	A	1149	THR	3.2
1	B	1037	PHE	3.1
1	D	1216	ILE	3.1
1	D	1037	PHE	3.1
1	C	1036	THR	3.1
1	A	1128	PRO	3.1
1	B	1126	VAL	3.0
1	B	1069	ILE	3.0
1	C	1064	ILE	2.9
1	B	1183	ILE	2.9
1	D	1098	LEU	2.9
1	C	1002	TYR	2.8
1	C	1008	ILE	2.8
1	C	1168	TYR	2.8
1	A	1068	ARG	2.8
1	A	1047	LEU	2.8
1	A	1071	PHE	2.8
1	B	1117	ARG	2.8
1	C	1149	THR	2.8
1	D	1069	ILE	2.8
1	A	1217	ILE	2.7
1	A	1112	ALA	2.7
1	B	1005	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	1151	LEU	2.7
1	B	1121	SER	2.7
1	C	1091	THR	2.6
1	A	1070	SER	2.6
1	C	1150	GLN	2.6
1	A	1069	ILE	2.6
1	A	1120	VAL	2.6
1	D	1072	PHE	2.6
1	C	1112	ALA	2.6
1	D	1113	VAL	2.6
1	B	1016	LYS	2.5
1	C	1033	THR	2.5
1	D	1173	VAL	2.5
1	B	1216	ILE	2.5
1	B	1114	PRO	2.5
1	D	1114	PRO	2.4
1	A	1008	ILE	2.4
1	A	1113	VAL	2.4
1	B	1097	ILE	2.4
1	D	1034	SER	2.4
1	D	1119	ILE	2.4
1	D	1047	LEU	2.4
1	D	1070	SER	2.4
1	B	1063	ARG	2.4
1	D	1074	ASP	2.3
1	B	1033	THR	2.3
1	C	1127	ILE	2.3
1	D	1187	LEU	2.3
1	B	1051	ILE	2.3
1	C	1184	VAL	2.3
1	B	1003	LEU	2.3
1	C	1006	THR	2.2
1	C	1075	PRO	2.2
1	D	1149	THR	2.2
1	D	1071	PHE	2.2
1	A	1160	PHE	2.2
1	C	1082	PHE	2.2
1	A	1019	ILE	2.2
1	B	1065	TYR	2.2
1	A	1161	GLY	2.1
1	A	1072	PHE	2.1
1	D	1112	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	1086	ILE	2.1
1	A	1146	ILE	2.1
1	A	1077	SER	2.1
1	A	1150	GLN	2.1
1	C	1152	PHE	2.1
1	A	1095	PHE	2.1
1	A	1162	THR	2.1
1	A	1109	LEU	2.0
1	D	1041	PHE	2.0
1	D	1022	ILE	2.0
1	B	1009	VAL	2.0
1	D	1003	LEU	2.0
1	D	1124	ILE	2.0
1	A	1041	PHE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

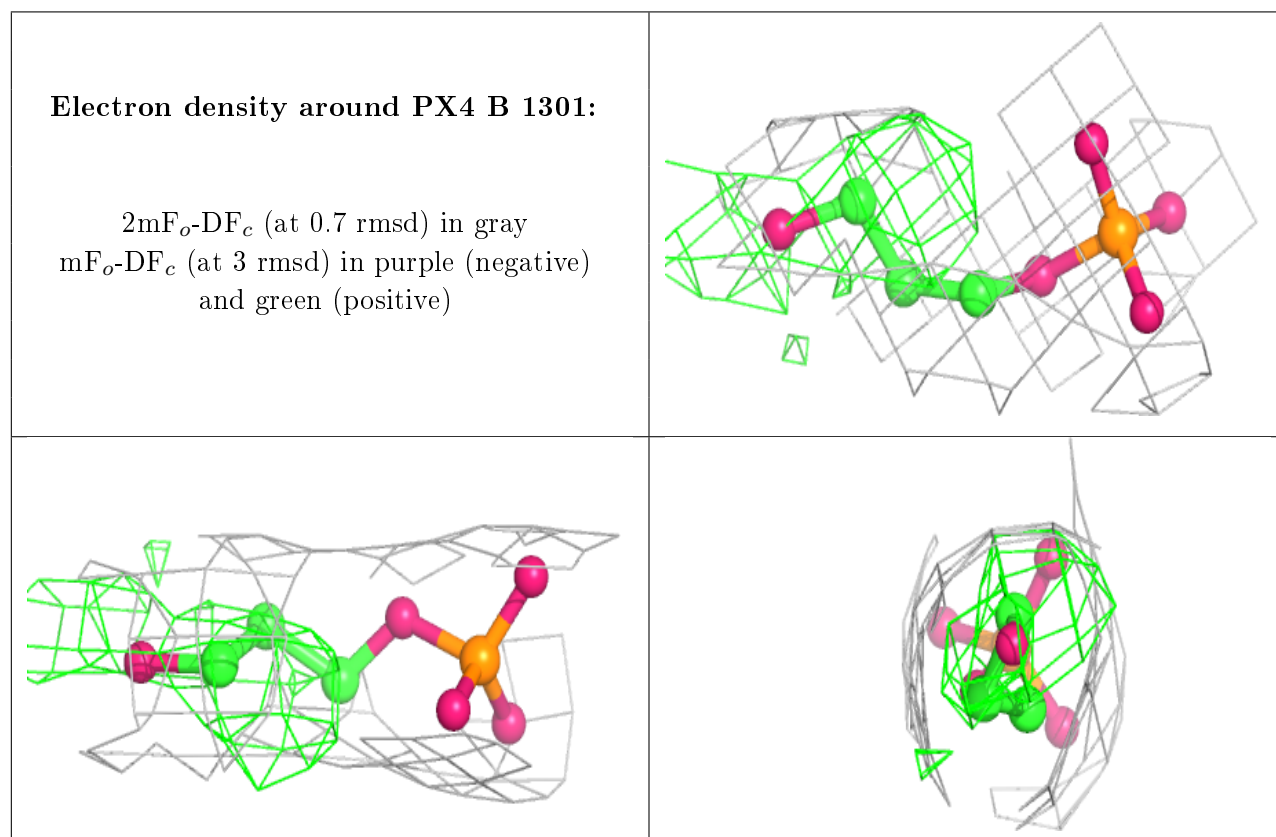
There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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

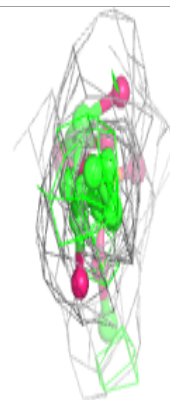
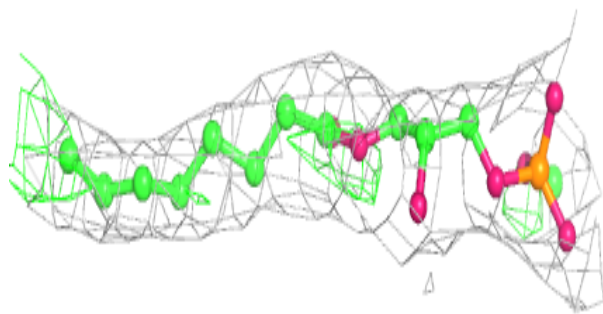
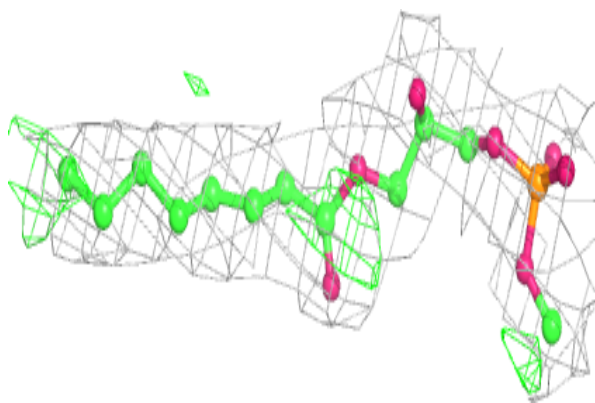
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	PX4	D	1301	8/46	0.83	0.27	82,101,119,125	0
2	PX4	B	1301	9/46	0.83	0.29	91,99,111,115	0
3	PO4	D	1302	5/5	0.84	0.23	95,99,105,113	0
2	PX4	C	1301	20/46	0.87	0.30	73,86,97,107	0
3	PO4	B	1303	5/5	0.88	0.20	107,107,118,123	0
2	PX4	B	1302	6/46	0.90	0.23	102,104,111,114	0
3	PO4	C	1302	5/5	0.90	0.27	106,107,118,120	0
2	PX4	A	1301	14/46	0.92	0.21	65,75,94,96	0
3	PO4	A	1302	5/5	0.94	0.24	87,95,98,102	0

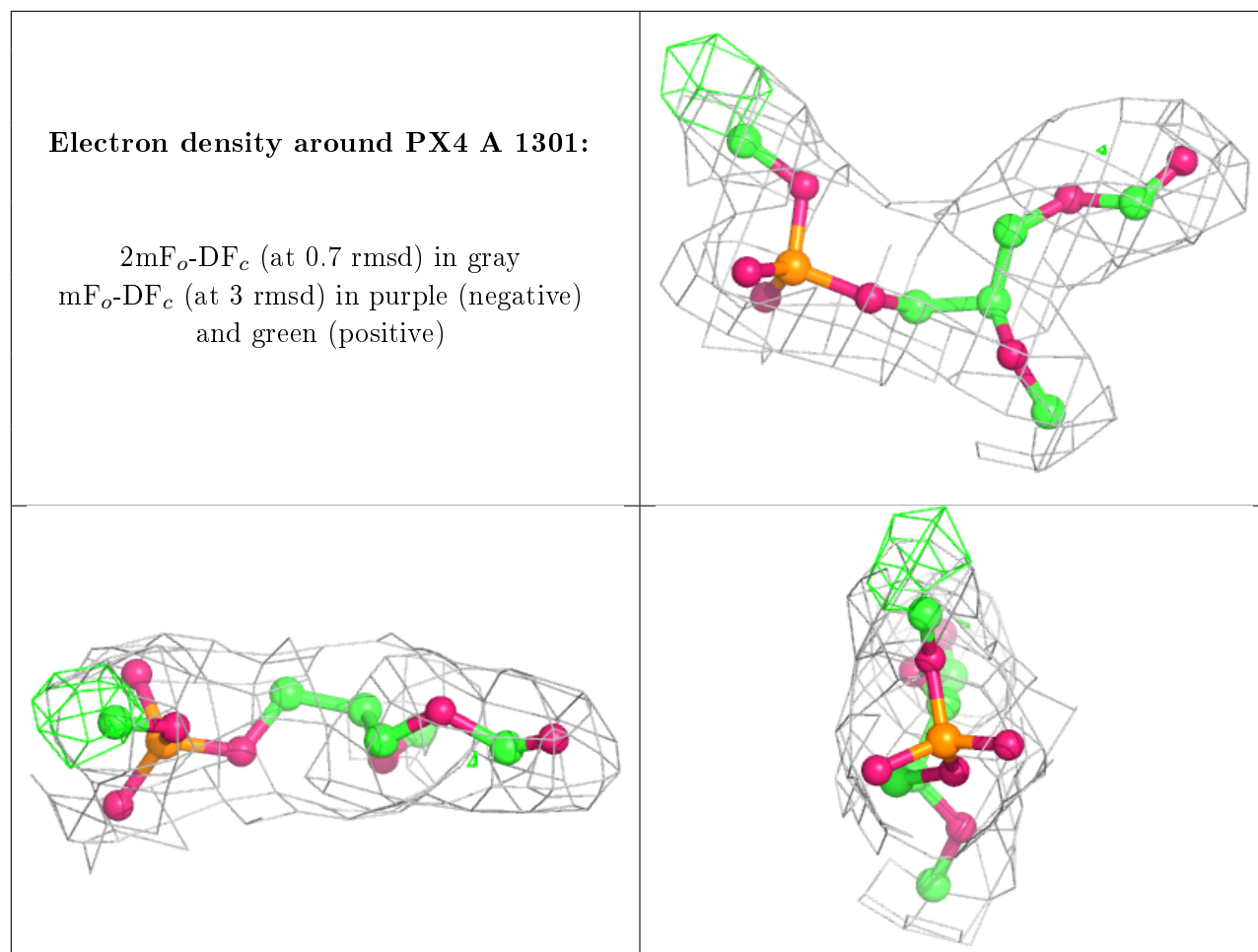
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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



Electron density around PX4 C 1301:

$2mF_o-DF_c$ (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.