



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:45 PM UTC

PDB ID : 4RW3 / pdb_00004rw3
Title : Structural insights into substrate binding of brown spider venom class II phospholipases D
Authors : Coronado, M.A.; Ullah, A.; da Silva, L.S.; Chaves-Moreira, D.; Vuitika, L.; Chaim, O.M.; Veiga, S.S.; Chahine, J.; Murakami, M.T.; Arni, R.K.
Deposited on : 2014-12-01
Resolution : 1.72 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

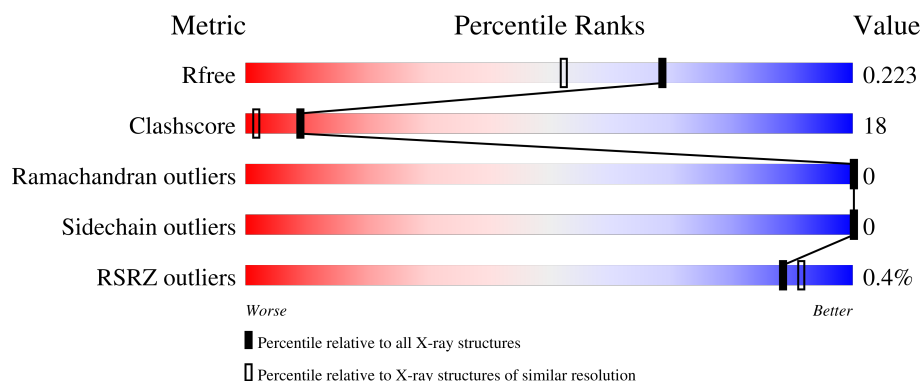
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.72 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	1039 (1.72-1.72)
Clashscore	190562	1049 (1.72-1.72)
Ramachandran outliers	187476	1041 (1.72-1.72)
Sidechain outliers	187428	1041 (1.72-1.72)
RSRZ outliers	180081	1039 (1.72-1.72)

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	302	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
10	TDA	A	313	-	-	X	-
4	IPD	A	304	-	-	X	-
6	OCA	A	306	-	-	X	-
8	DKA	A	312	-	-	X	-
9	PLM	A	309	-	-	X	-
9	PLM	A	311	-	-	X	-

2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 2558 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 Phospholipase D LiSicTox-alphaIA1bii.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	279	Total	C	N	O	S	2	7	0
			2241	1412	395	426	8			

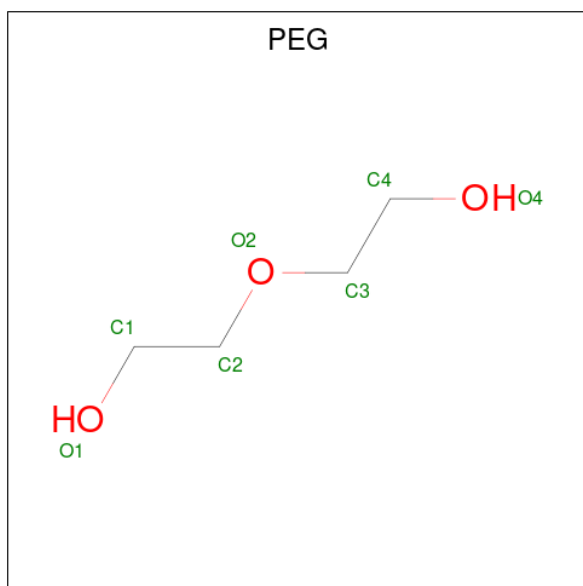
There are 23 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	expression tag	UNP P0CE82
A	-20	GLY	-	expression tag	UNP P0CE82
A	-19	SER	-	expression tag	UNP P0CE82
A	-18	SER	-	expression tag	UNP P0CE82
A	-17	HIS	-	expression tag	UNP P0CE82
A	-16	HIS	-	expression tag	UNP P0CE82
A	-15	HIS	-	expression tag	UNP P0CE82
A	-14	HIS	-	expression tag	UNP P0CE82
A	-13	HIS	-	expression tag	UNP P0CE82
A	-12	HIS	-	expression tag	UNP P0CE82
A	-11	SER	-	expression tag	UNP P0CE82
A	-10	SER	-	expression tag	UNP P0CE82
A	-9	GLY	-	expression tag	UNP P0CE82
A	-8	LEU	-	expression tag	UNP P0CE82
A	-7	VAL	-	expression tag	UNP P0CE82
A	-6	PRO	-	expression tag	UNP P0CE82
A	-5	ARG	-	expression tag	UNP P0CE82
A	-4	GLY	-	expression tag	UNP P0CE82
A	-3	SER	-	expression tag	UNP P0CE82
A	-2	HIS	-	expression tag	UNP P0CE82
A	-1	MET	-	expression tag	UNP P0CE82
A	0	LEU	-	expression tag	UNP P0CE82
A	1	GLU	-	expression tag	UNP P0CE82

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

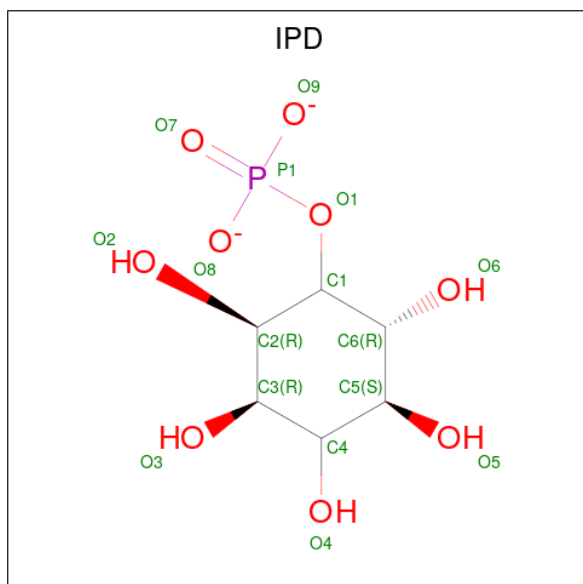
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



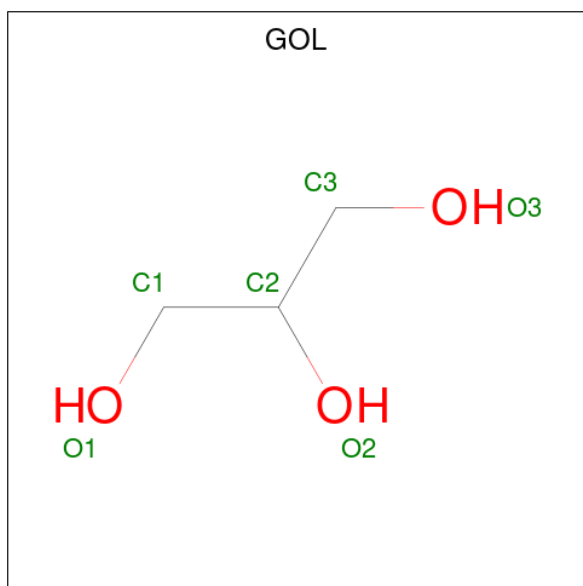
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		
3	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is D-MYO-INOSITOL-1-PHOSPHATE (CCD ID: IPD) (formula: $C_6H_{11}O_9P$).



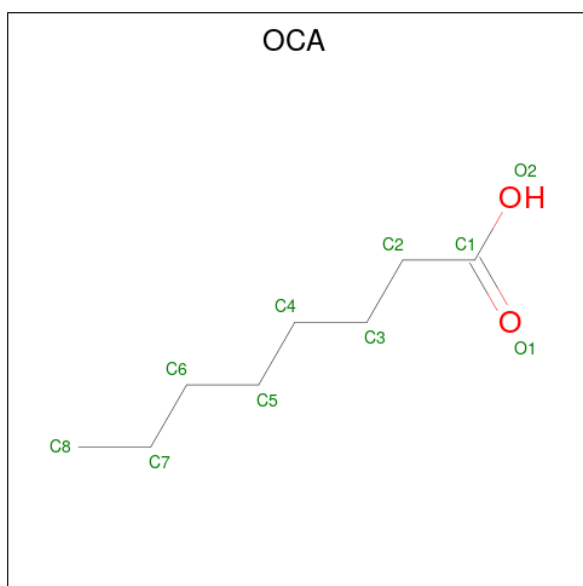
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	O	P	0	0
			16	6	9	1		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



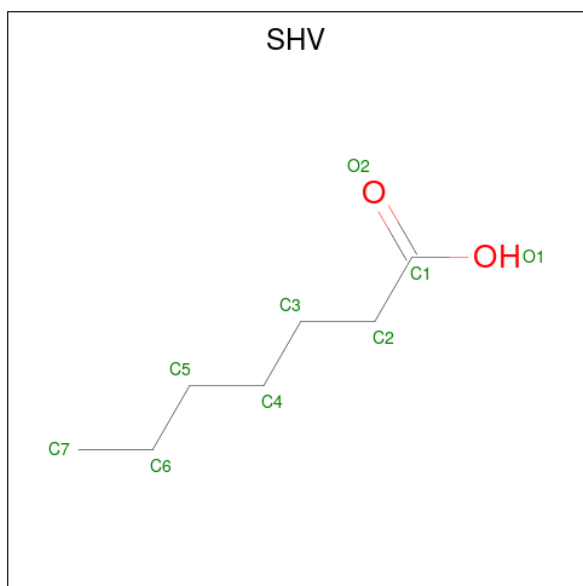
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is OCTANOIC ACID (CAPRYLIC ACID) (CCD ID: OCA) (formula: $C_8H_{16}O_2$).



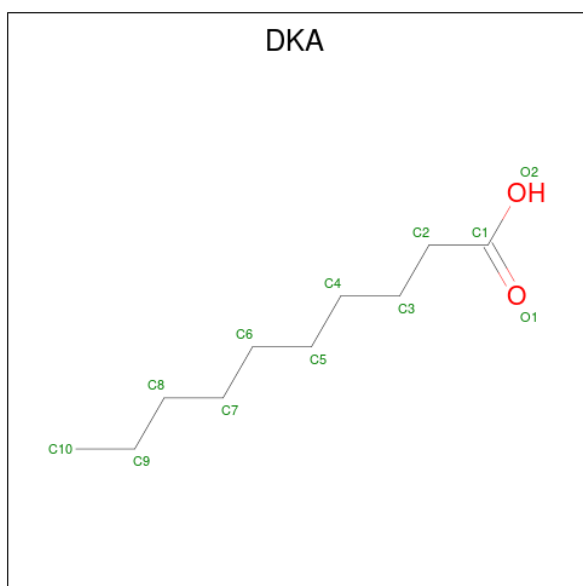
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			10	8	2		

- Molecule 7 is HEPTANOIC ACID (CCD ID: SHV) (formula: $C_7H_{14}O_2$).



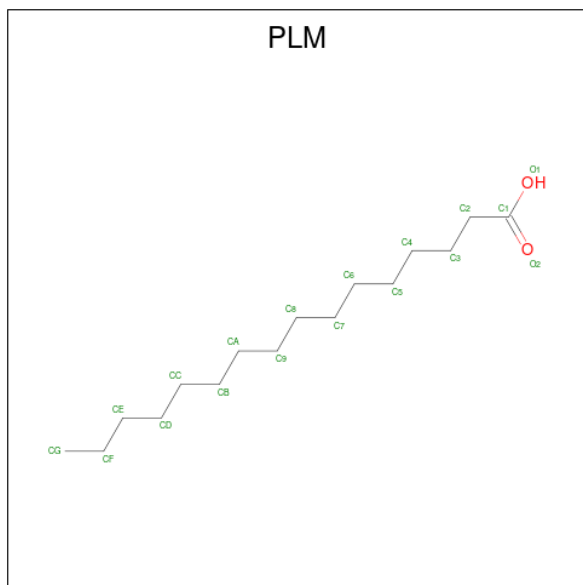
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			9	7	2		

- Molecule 8 is DECANOIC ACID (CCD ID: DKA) (formula: $C_{10}H_{20}O_2$).



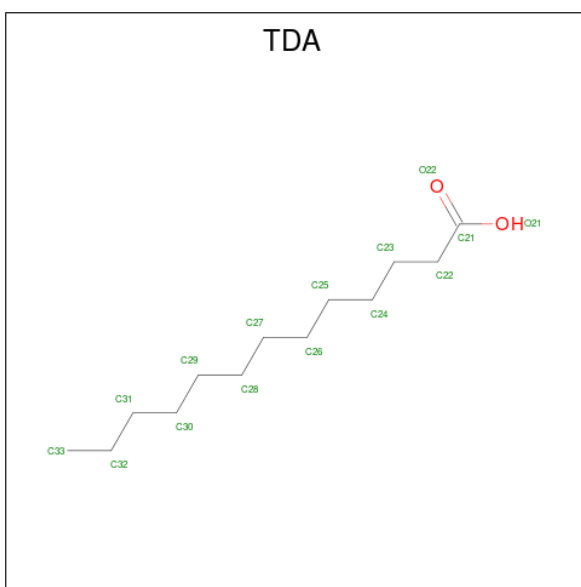
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			12	10	2		
8	A	1	Total	C	O	0	0
			12	10	2		

- Molecule 9 is PALMITIC ACID (CCD ID: PLM) (formula: $C_{16}H_{32}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			18	16	2		
9	A	1	Total	C	O	0	0
			18	16	2		
9	A	1	Total	C	O	0	0
			18	16	2		

- Molecule 10 is N-TRIDEcanoic ACID (CCD ID: TDA) (formula: $C_{13}H_{26}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			15	13	2		

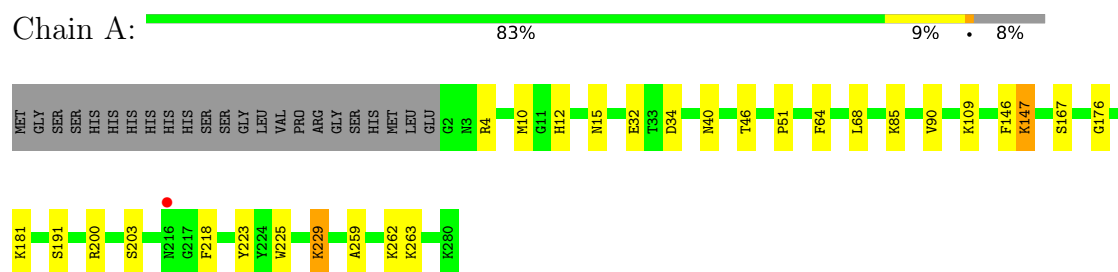
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	168	Total	O	0	0
			168	168		

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: Phospholipase D LiSicTox-alphaIA1bii



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.81Å 49.30Å 56.30Å 90.00° 105.83° 90.00°	Depositor
Resolution (Å)	30.00 – 1.72 30.00 – 1.72	Depositor EDS
% Data completeness (in resolution range)	99.8 (30.00-1.72) 99.8 (30.00-1.72)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 1.72Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.183 , 0.223 0.185 , 0.223	Depositor DCC
R_{free} test set	2808 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å ²)	17.9	Xtriage
Anisotropy	0.074	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 35.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2558	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.23% 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: SHV, MG, OCA, IPD, TDA, PEG, DKA, PLM, GOL

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	0.70	4/2306 (0.2%)	0.77	0/3122

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	146	PHE	C-N	-10.67	1.20	1.33
1	A	147	LYS	C-N	-7.26	1.24	1.34
1	A	85	LYS	C-O	-5.19	1.17	1.24
1	A	229	LYS	C-N	-5.09	1.27	1.33

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	2241	0	2180	81	0
2	A	1	0	0	0	0
3	A	14	0	20	1	0
4	A	16	0	8	10	0
5	A	6	0	8	0	0
6	A	10	0	15	10	0
7	A	9	0	13	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	24	0	38	16	0
9	A	54	0	93	27	0
10	A	15	0	25	13	0
11	A	168	0	0	4	0
All	All	2558	0	2400	88	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 18.

All (88) 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:229:LYS:HE3	9:A:309:PLM:CD	1.62	1.29
1:A:181:LYS:NZ	10:A:313:TDA:H281	1.50	1.24
1:A:229:LYS:NZ	9:A:309:PLM:HD1	1.56	1.19
1:A:147:LYS:HE2	8:A:312:DKA:H61	1.31	1.13
1:A:181:LYS:HZ1	10:A:313:TDA:C28	1.62	1.12
1:A:229:LYS:CE	9:A:309:PLM:HD2	1.83	1.07
1:A:229:LYS:CE	9:A:309:PLM:CD	2.31	1.07
1:A:229:LYS:HZ1	9:A:309:PLM:HD1	1.12	1.07
1:A:147:LYS:HD2	8:A:312:DKA:H81	1.37	1.06
1:A:181:LYS:NZ	10:A:313:TDA:H251	1.71	1.06
1:A:181:LYS:HZ2	10:A:313:TDA:H251	0.93	1.05
1:A:109:LYS:HZ1	9:A:311:PLM:H82	1.24	1.00
1:A:147:LYS:CE	8:A:312:DKA:H61	1.90	1.00
1:A:181:LYS:HZ1	10:A:313:TDA:H281	0.85	0.99
1:A:263:LYS:CE	6:A:306:OCA:H51	1.94	0.98
1:A:229:LYS:HE3	9:A:309:PLM:HD2	0.93	0.93
1:A:229:LYS:CE	9:A:309:PLM:HD1	1.99	0.92
1:A:181:LYS:NZ	10:A:313:TDA:H312	1.90	0.85
1:A:109:LYS:NZ	9:A:311:PLM:H82	1.92	0.83
1:A:229:LYS:HZ2	9:A:309:PLM:HG2	1.44	0.83
1:A:147:LYS:HE3	8:A:312:DKA:H31	1.63	0.81
1:A:263:LYS:NZ	6:A:306:OCA:H51	1.95	0.80
1:A:181:LYS:HZ2	10:A:313:TDA:C25	1.86	0.80
1:A:147:LYS:HE3	8:A:312:DKA:C3	2.13	0.79
1:A:229:LYS:NZ	9:A:309:PLM:HG2	2.02	0.75
1:A:263:LYS:HE3	6:A:306:OCA:H51	1.67	0.75
1:A:218:PHE:HB2	7:A:307:SHV:H51	1.69	0.74
1:A:109:LYS:HE2	9:A:311:PLM:H61	1.71	0.73
1:A:12:HIS:HE2	4:A:304:IPD:H5	1.52	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:A:311:PLM:HG2	9:A:311:PLM:O1	1.90	0.71
1:A:147:LYS:HE2	8:A:312:DKA:C6	2.14	0.71
1:A:225:TRP:CD2	4:A:304:IPD:O8	2.45	0.69
1:A:229:LYS:HZ1	9:A:309:PLM:CD	1.99	0.68
1:A:147:LYS:CE	8:A:312:DKA:H32	2.23	0.68
1:A:225:TRP:CG	4:A:304:IPD:O8	2.46	0.68
1:A:181:LYS:HZ3	10:A:313:TDA:H312	1.57	0.67
1:A:259:ALA:HB1	6:A:306:OCA:H81	1.77	0.66
1:A:109:LYS:CE	9:A:311:PLM:H61	2.26	0.65
1:A:229:LYS:HZ2	9:A:309:PLM:CG	2.10	0.64
6:A:306:OCA:H72	11:A:422:HOH:O	1.98	0.63
1:A:203[B]:SER:OG	3:A:302:PEG:H22	1.99	0.62
1:A:263:LYS:HZ1	6:A:306:OCA:H21	1.62	0.62
1:A:181:LYS:NZ	10:A:313:TDA:C28	2.38	0.61
1:A:51:PRO:HB2	8:A:308:DKA:H22	1.82	0.61
1:A:262:LYS:HZ3	9:A:310:PLM:HC2	1.66	0.59
1:A:147:LYS:CE	8:A:312:DKA:C3	2.79	0.59
1:A:181:LYS:HZ1	10:A:313:TDA:H312	1.65	0.59
1:A:109:LYS:HZ2	9:A:311:PLM:HC1	1.66	0.58
1:A:263:LYS:NZ	6:A:306:OCA:C5	2.68	0.57
1:A:147:LYS:CD	8:A:312:DKA:H81	2.26	0.56
1:A:147:LYS:HD2	8:A:312:DKA:C8	2.26	0.56
1:A:15:ASN:HB3	1:A:46:THR:O	2.06	0.55
1:A:4[A]:ARG:HD3	11:A:402:HOH:O	2.06	0.55
1:A:181:LYS:NZ	10:A:313:TDA:C31	2.68	0.54
1:A:147:LYS:HE3	8:A:312:DKA:H61	1.86	0.53
1:A:167:SER:HB2	4:A:304:IPD:O9	2.08	0.53
1:A:181:LYS:HZ1	10:A:313:TDA:C31	2.21	0.53
1:A:40:ASN:HB3	9:A:311:PLM:H91	1.91	0.53
1:A:191:SER:HB3	1:A:223:TYR:CZ	2.43	0.52
1:A:223:TYR:OH	4:A:304:IPD:O7	2.15	0.52
1:A:51:PRO:CB	8:A:308:DKA:H22	2.40	0.52
1:A:12:HIS:NE2	4:A:304:IPD:H5	2.23	0.50
1:A:34:ASP:O	1:A:46:THR:HA	2.12	0.49
9:A:311:PLM:O1	9:A:311:PLM:CG	2.60	0.49
10:A:313:TDA:H281	10:A:313:TDA:H251	1.61	0.48
1:A:259:ALA:CB	6:A:306:OCA:H81	2.44	0.48
1:A:263:LYS:HE2	6:A:306:OCA:H72	1.96	0.48
8:A:312:DKA:C10	11:A:469:HOH:O	2.62	0.48
1:A:225:TRP:CD1	4:A:304:IPD:O8	2.67	0.47
1:A:263:LYS:HZ1	6:A:306:OCA:H51	1.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:SER:HA	1:A:191:SER:OG	2.14	0.47
9:A:309:PLM:HD1	9:A:309:PLM:HG2	1.73	0.47
1:A:109:LYS:HZ1	9:A:311:PLM:H61	1.79	0.46
8:A:312:DKA:H101	11:A:469:HOH:O	2.15	0.46
1:A:10:MET:CE	1:A:90:VAL:HG21	2.45	0.46
1:A:225:TRP:CZ3	4:A:304:IPD:H2	2.51	0.45
1:A:225:TRP:CE2	4:A:304:IPD:O8	2.69	0.45
1:A:262:LYS:HZ3	9:A:310:PLM:H81	1.82	0.45
1:A:109:LYS:HZ1	9:A:311:PLM:C8	2.12	0.44
1:A:147:LYS:HE3	8:A:312:DKA:H32	1.86	0.44
1:A:32:GLU:OE2	4:A:304:IPD:O6	2.29	0.44
1:A:109:LYS:NZ	9:A:311:PLM:H61	2.33	0.44
1:A:4[B]:ARG:NE	1:A:262:LYS:O	2.52	0.43
1:A:64:PHE:CE2	1:A:68:LEU:HD11	2.53	0.42
1:A:262:LYS:NZ	9:A:310:PLM:H81	2.34	0.42
1:A:200:ARG:HH11	1:A:200:ARG:HG3	1.84	0.42
1:A:176:GLY:O	7:A:307:SHV:H22	2.21	0.41
1:A:229:LYS:HZ3	9:A:309:PLM:C6	2.34	0.41

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	284/302 (94%)	282 (99%)	2 (1%)	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	241/254 (95%)	241 (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	116	GLN

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, 1 is monoatomic - leaving 12 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	GOL	A	305	-	5,5,5	0.27	0	5,5,5	0.40	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	IPD	A	304	2	16,16,16	4.41	1 (6%)	25,25,25	3.87	6 (24%)
9	PLM	A	311	-	17,17,17	0.60	1 (5%)	17,17,17	1.31	1 (5%)
9	PLM	A	310	-	17,17,17	0.49	0	17,17,17	0.86	0
7	SHV	A	307	-	8,8,8	0.73	0	8,8,8	0.86	0
8	DKA	A	308	-	11,11,11	0.65	0	11,11,11	0.76	0
10	TDA	A	313	-	14,14,14	0.65	1 (7%)	14,14,14	0.99	1 (7%)
6	OCA	A	306	-	9,9,9	0.73	0	9,9,9	0.94	0
8	DKA	A	312	-	11,11,11	0.90	1 (9%)	11,11,11	1.47	1 (9%)
9	PLM	A	309	-	17,17,17	0.68	1 (5%)	17,17,17	0.91	0
3	PEG	A	303	-	6,6,6	0.48	0	5,5,5	0.23	0
3	PEG	A	302	-	6,6,6	0.44	0	5,5,5	0.48	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
5	GOL	A	305	-	-	4/4/4/4	-
4	IPD	A	304	2	-	0/5/29/29	0/1/1/1
9	PLM	A	311	-	-	12/15/15/15	-
9	PLM	A	310	-	-	9/15/15/15	-
7	SHV	A	307	-	-	3/6/6/6	-
8	DKA	A	308	-	-	3/9/9/9	-
10	TDA	A	313	-	-	9/12/12/12	-
6	OCA	A	306	-	-	5/7/7/7	-
8	DKA	A	312	-	-	6/9/9/9	-
9	PLM	A	309	-	-	13/15/15/15	-
3	PEG	A	303	-	-	4/4/4/4	-
3	PEG	A	302	-	-	1/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	304	IPD	O4-C4	-17.50	0.99	1.43
8	A	312	DKA	O2-C1	-2.51	1.22	1.30
9	A	309	PLM	O1-C1	-2.43	1.22	1.30
10	A	313	TDA	O21-C21	-2.36	1.23	1.30
9	A	311	PLM	O1-C1	-2.12	1.23	1.30

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	304	IPD	O4-C4-C3	14.98	145.70	110.38
4	A	304	IPD	O6-C6-C1	-8.75	87.51	109.94
4	A	304	IPD	O6-C6-C5	5.61	123.60	110.38
8	A	312	DKA	C3-C2-C1	-4.44	102.92	114.51
9	A	311	PLM	C3-C2-C1	-4.24	103.43	114.51
4	A	304	IPD	O4-C4-C5	-3.76	101.52	110.38
4	A	304	IPD	C3-C2-C1	2.35	115.02	109.68
4	A	304	IPD	C6-C5-C4	2.34	114.94	110.83
10	A	313	TDA	O21-C21-C22	2.05	120.48	114.00

There are no chirality outliers.

All (69) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	305	GOL	O1-C1-C2-C3
5	A	305	GOL	C1-C2-C3-O3
3	A	303	PEG	O1-C1-C2-O2
8	A	312	DKA	C3-C4-C5-C6
10	A	313	TDA	C25-C26-C27-C28
9	A	309	PLM	C1-C2-C3-C4
9	A	311	PLM	C5-C6-C7-C8
5	A	305	GOL	O1-C1-C2-O2
5	A	305	GOL	O2-C2-C3-O3
9	A	310	PLM	C8-C9-CA-CB
10	A	313	TDA	C26-C27-C28-C29
8	A	312	DKA	C6-C7-C8-C9
10	A	313	TDA	C21-C22-C23-C24
6	A	306	OCA	C4-C5-C6-C7
9	A	311	PLM	C8-C9-CA-CB
6	A	306	OCA	C2-C3-C4-C5
9	A	311	PLM	C3-C4-C5-C6
9	A	309	PLM	C8-C9-CA-CB
8	A	308	DKA	C3-C4-C5-C6
9	A	311	PLM	C6-C7-C8-C9
3	A	302	PEG	O1-C1-C2-O2
9	A	309	PLM	CC-CD-CE-CF
9	A	309	PLM	C3-C4-C5-C6
9	A	309	PLM	C6-C7-C8-C9
9	A	310	PLM	CA-CB-CC-CD
10	A	313	TDA	C22-C23-C24-C25
9	A	309	PLM	C4-C5-C6-C7

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Mol	Chain	Res	Type	Atoms
7	A	307	SHV	C2-C3-C4-C5
9	A	309	PLM	C2-C3-C4-C5
9	A	309	PLM	CA-CB-CC-CD
10	A	313	TDA	C24-C25-C26-C27
10	A	313	TDA	C30-C31-C32-C33
9	A	309	PLM	CB-CC-CD-CE
9	A	310	PLM	C2-C3-C4-C5
9	A	310	PLM	CB-CC-CD-CE
8	A	312	DKA	C4-C5-C6-C7
9	A	311	PLM	C9-CA-CB-CC
8	A	312	DKA	C1-C2-C3-C4
9	A	310	PLM	C3-C4-C5-C6
9	A	309	PLM	CD-CE-CF-CG
3	A	303	PEG	C4-C3-O2-C2
9	A	310	PLM	C4-C5-C6-C7
3	A	303	PEG	O2-C3-C4-O4
9	A	311	PLM	CC-CD-CE-CF
9	A	311	PLM	CA-CB-CC-CD
9	A	309	PLM	C5-C6-C7-C8
3	A	303	PEG	C1-C2-O2-C3
9	A	310	PLM	C5-C6-C7-C8
7	A	307	SHV	O1-C1-C2-C3
10	A	313	TDA	O21-C21-C22-C23
7	A	307	SHV	O2-C1-C2-C3
6	A	306	OCA	O1-C1-C2-C3
6	A	306	OCA	O2-C1-C2-C3
9	A	310	PLM	O1-C1-C2-C3
9	A	310	PLM	O2-C1-C2-C3
8	A	308	DKA	O2-C1-C2-C3
9	A	309	PLM	O1-C1-C2-C3
8	A	308	DKA	O1-C1-C2-C3
8	A	312	DKA	O2-C1-C2-C3
8	A	312	DKA	O1-C1-C2-C3
9	A	309	PLM	O2-C1-C2-C3
10	A	313	TDA	O22-C21-C22-C23
9	A	311	PLM	CB-CC-CD-CE
9	A	311	PLM	O1-C1-C2-C3
6	A	306	OCA	C3-C4-C5-C6
9	A	311	PLM	O2-C1-C2-C3
9	A	311	PLM	C4-C5-C6-C7
9	A	311	PLM	C7-C8-C9-CA
10	A	313	TDA	C29-C30-C31-C32

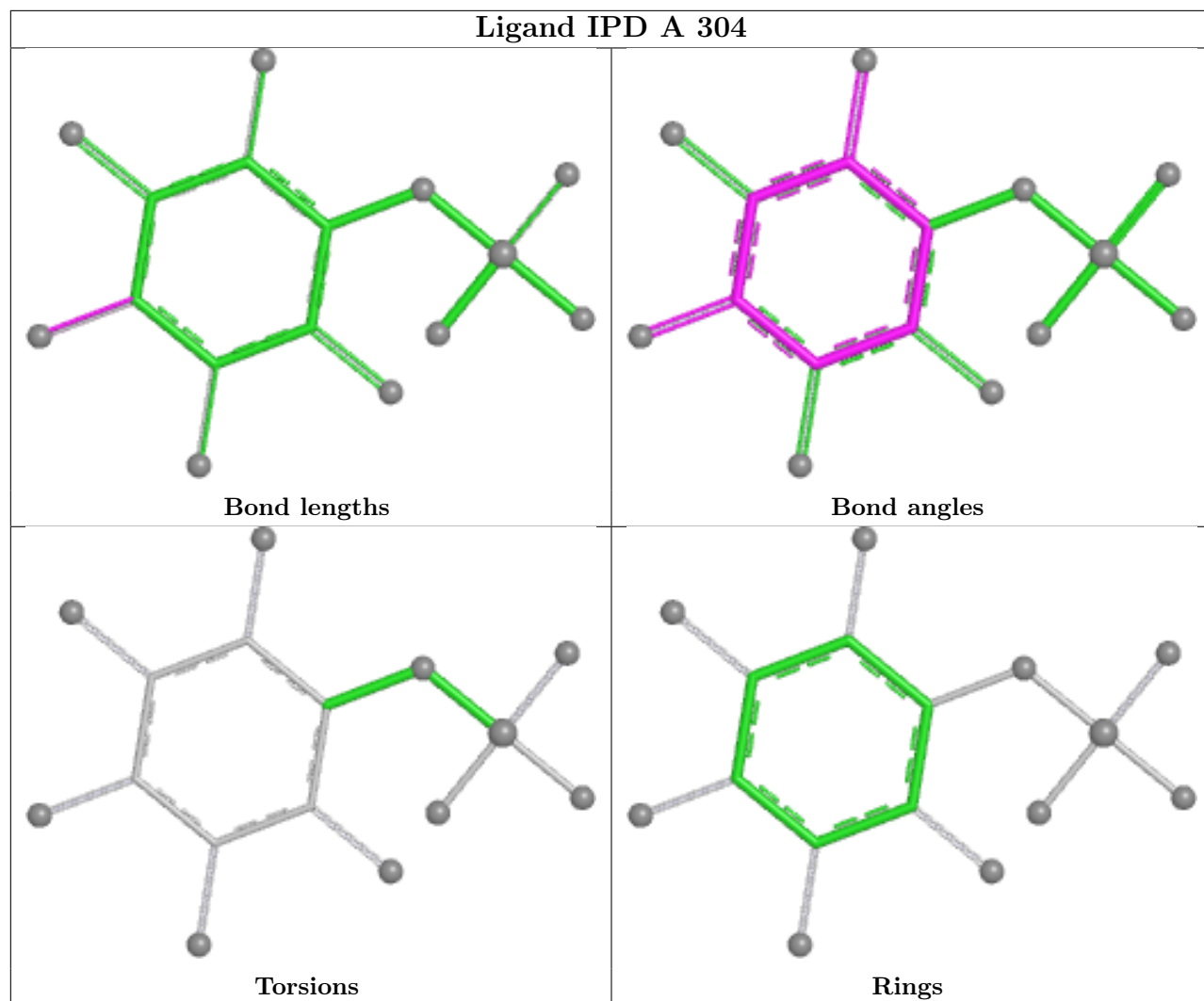
There are no ring outliers.

10 monomers are involved in 79 short contacts:

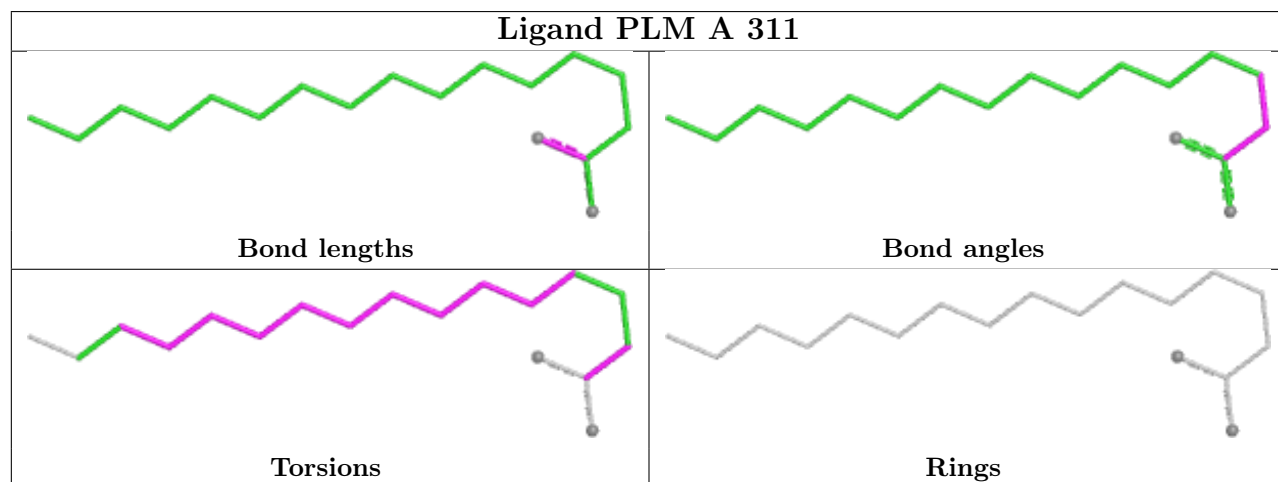
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	304	IPD	10	0
9	A	311	PLM	11	0
9	A	310	PLM	3	0
7	A	307	SHV	2	0
8	A	308	DKA	2	0
10	A	313	TDA	13	0
6	A	306	OCA	10	0
8	A	312	DKA	14	0
9	A	309	PLM	13	0
3	A	302	PEG	1	0

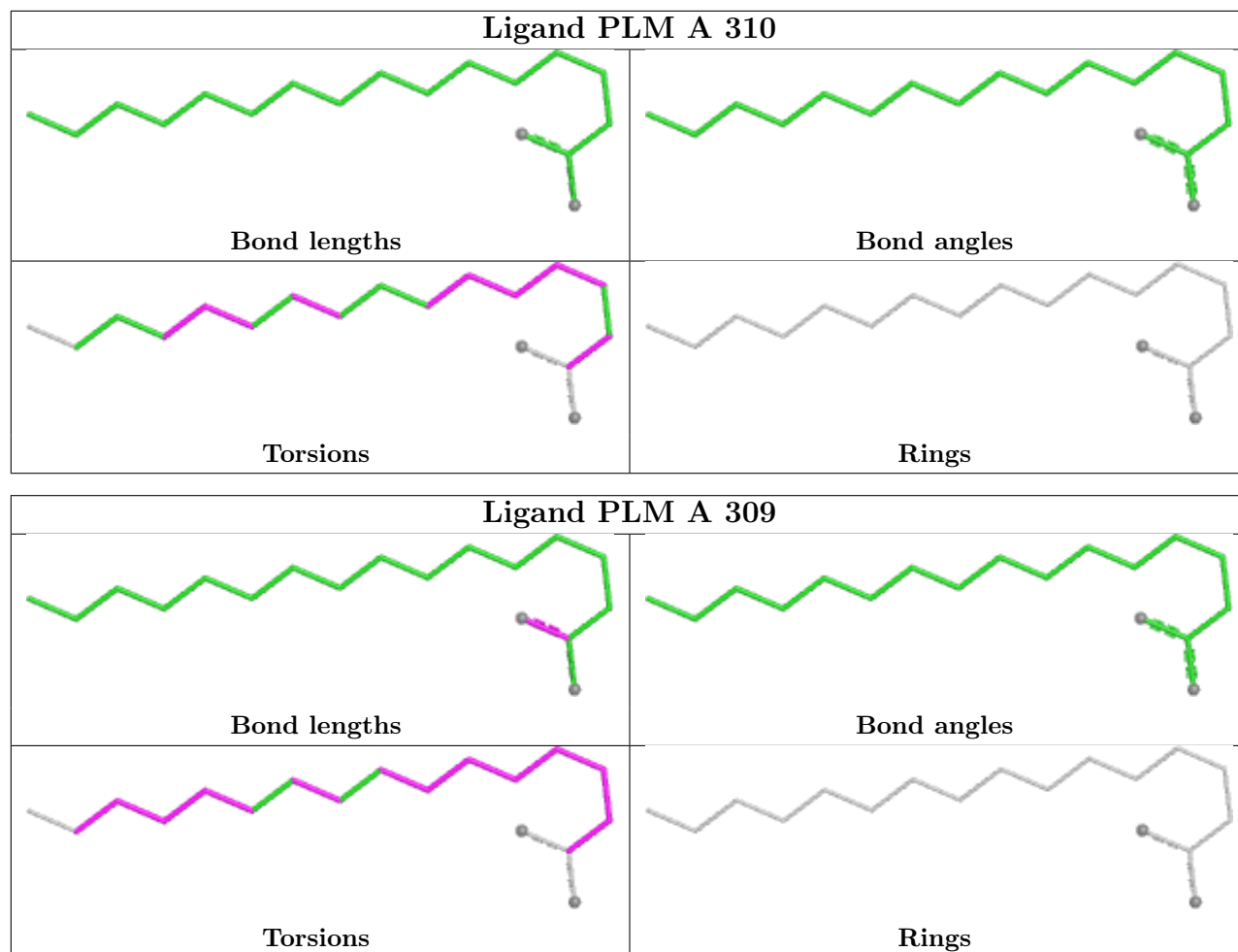
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

Ligand IPD A 304



Ligand PLM A 311





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	279/302 (92%)	-0.24	1 (0%) 88 91	7, 16, 25, 49	13 (4%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	216	ASN	2.2

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(Å ²)	Q<0.9
4	IPD	A	304	16/16	0.57	0.26	19,24,40,49	8
10	TDA	A	313	15/15	0.65	0.18	41,46,51,51	0
8	DKA	A	312	12/12	0.66	0.16	50,55,61,62	0
8	DKA	A	308	12/12	0.72	0.15	43,45,50,50	0
9	PLM	A	309	18/18	0.75	0.15	36,44,48,48	0
3	PEG	A	302	7/7	0.75	0.16	45,46,48,49	0

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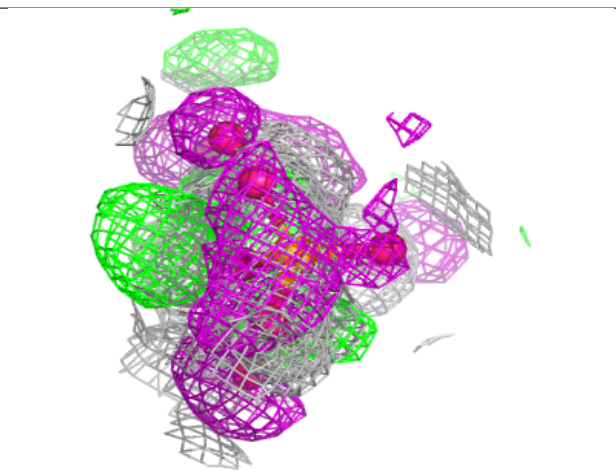
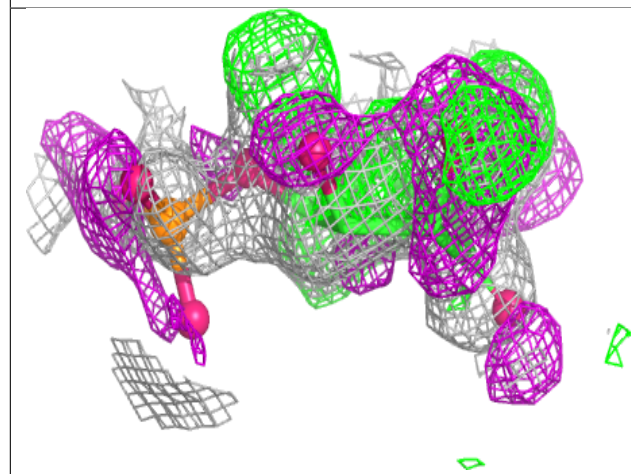
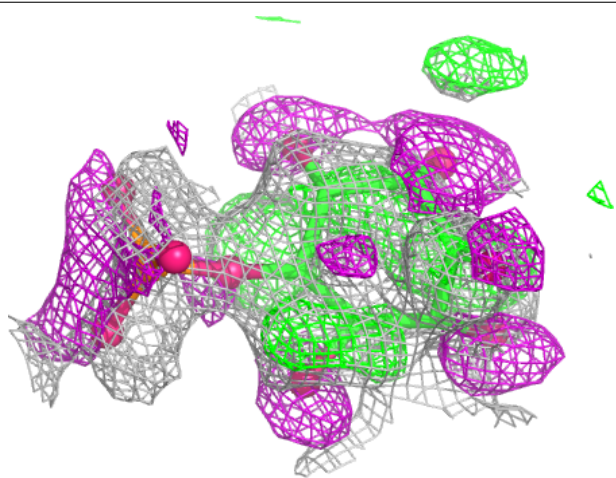
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	PEG	A	303	7/7	0.76	0.13	47,48,48,49	0
9	PLM	A	310	18/18	0.78	0.15	38,44,52,52	0
6	OCA	A	306	10/10	0.78	0.15	47,51,55,56	0
7	SHV	A	307	9/9	0.80	0.15	45,47,51,52	0
9	PLM	A	311	18/18	0.85	0.13	21,25,34,34	0
5	GOL	A	305	6/6	0.86	0.10	40,42,42,43	0
2	MG	A	301	1/1	0.97	0.03	12,12,12,12	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.

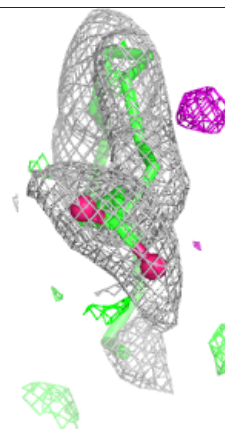
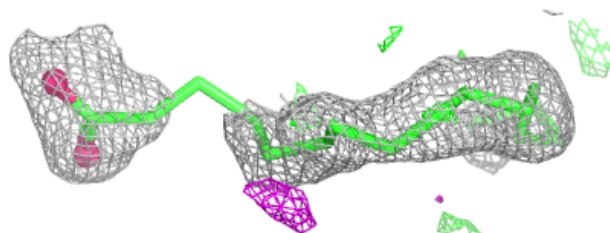
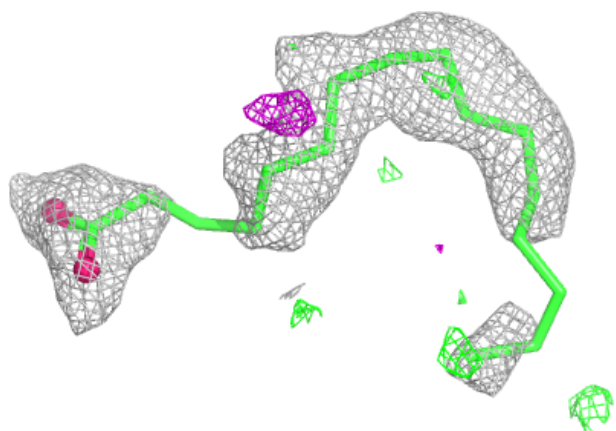
Electron density around IPD A 304:

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



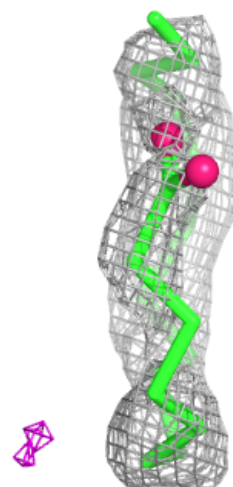
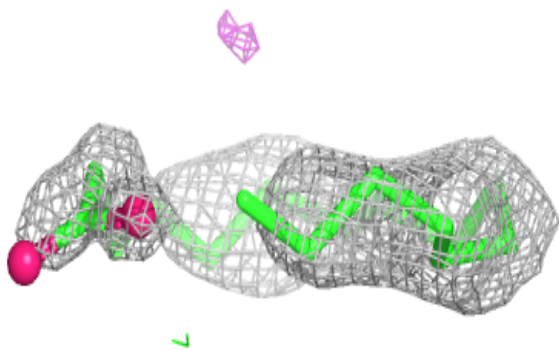
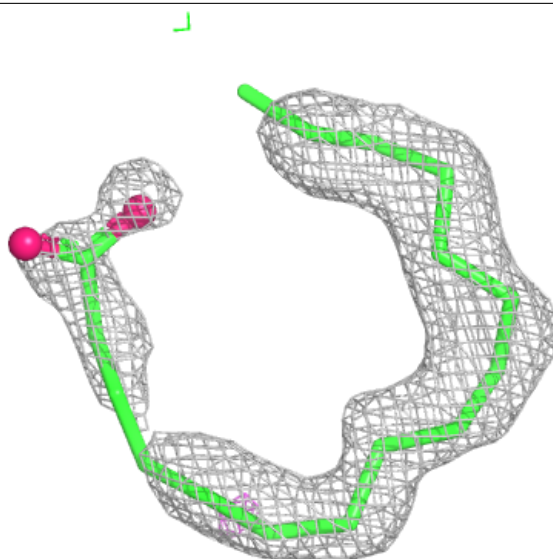
Electron density around PLM A 309:

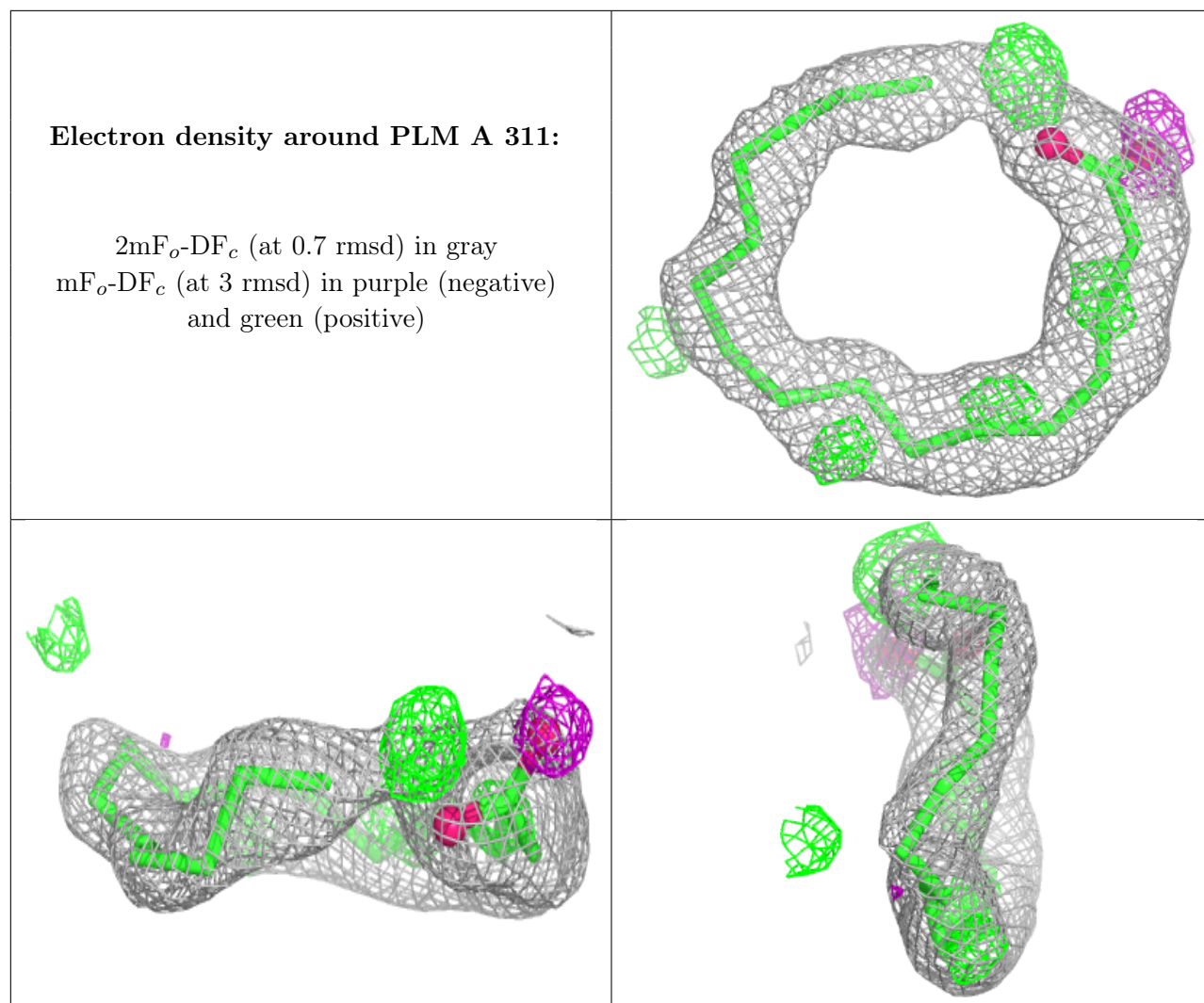
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around PLM A 310:

$2mF_o-DF_c$ (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.