



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:04 PM UTC

PDB ID : 8K1Y / pdb_00008k1y
Title : Crystal structure of human serum albumin and ruthenium Val complex adduct
Authors : Gong, W.J.; Xie, L.L.; Wang, W.M.; Wang, H.F.
Deposited on : 2023-07-11
Resolution : 2.15 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

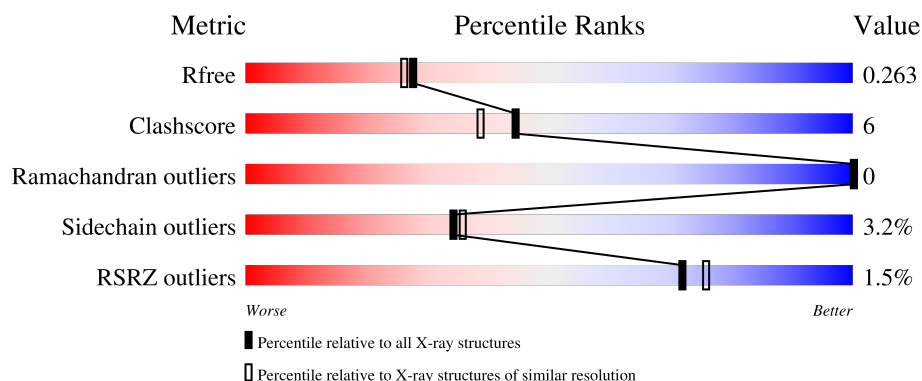
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

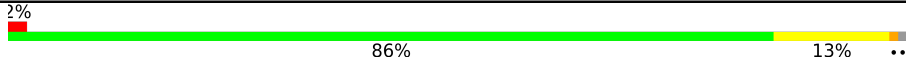
The reported resolution of this entry is 2.15 Å.

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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	585	

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
8	DMS	A	616	-	-	X	-

2 Entry composition [i](#)

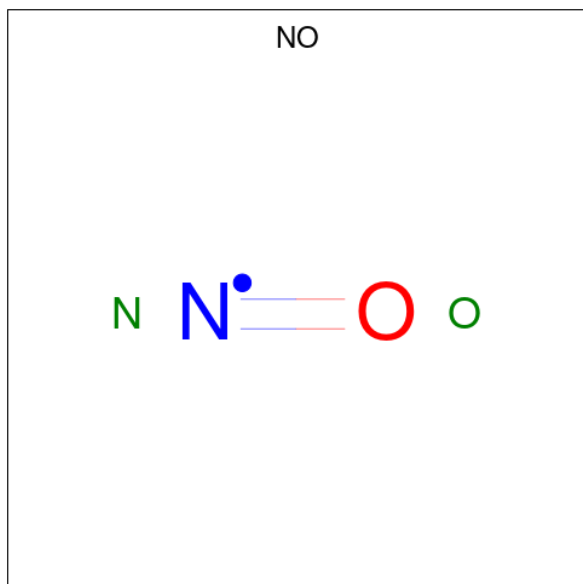
There are 9 unique types of molecules in this entry. The entry contains 4894 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Albumin.

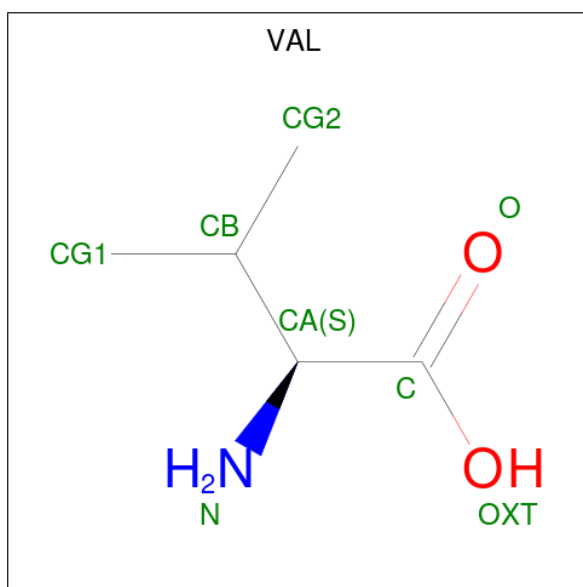
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	581	Total	C	N	O	S	0	0	0
			4606	2909	776	880	41			

- Molecule 2 is NITRIC OXIDE (CCD ID: NO) (formula: NO).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	N	O	0	0
			2	1	1		

- Molecule 3 is VALINE (CCD ID: VAL) (formula: C₅H₁₁NO₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			8	5	1	2		

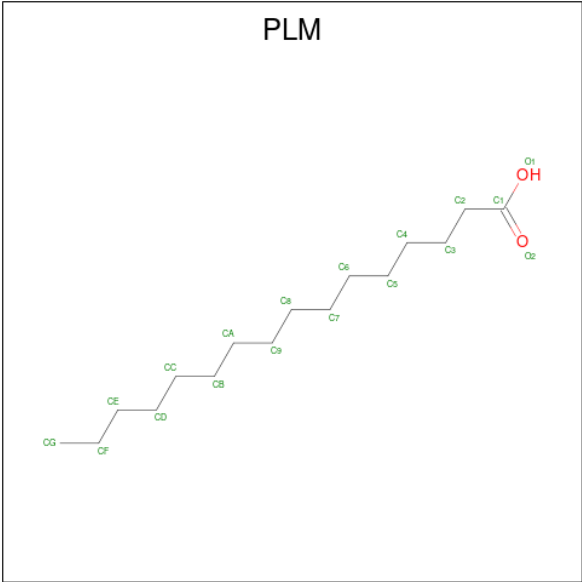
- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		

- Molecule 5 is RUTHENIUM ION (CCD ID: RU) (formula: Ru).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Ru	0	0
			1	1		

- Molecule 6 is PALMITIC ACID (CCD ID: PLM) (formula: C₁₆H₃₂O₂).



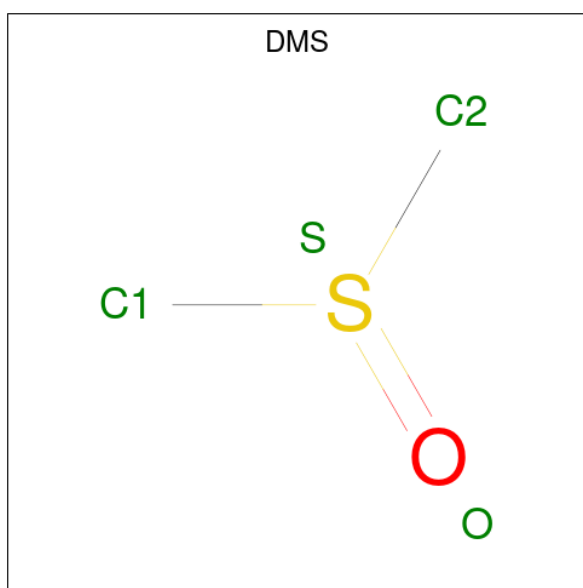
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			18	16	2		
6	A	1	Total	C	O	0	0
			18	16	2		
6	A	1	Total	C	O	0	0
			18	16	2		
6	A	1	Total	C	O	0	0
			18	16	2		
6	A	1	Total	C	O	0	0
			18	16	2		
6	A	1	Total	C	O	0	0
			18	16	2		

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 8 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	O	S	0	0
			4	2	1	1		
8	A	1	Total	C	O	S	0	0
			4	2	1	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	O	S	0	0
			4	2	1	1		
8	A	1	Total	C	O	S	0	0
			4	2	1	1		

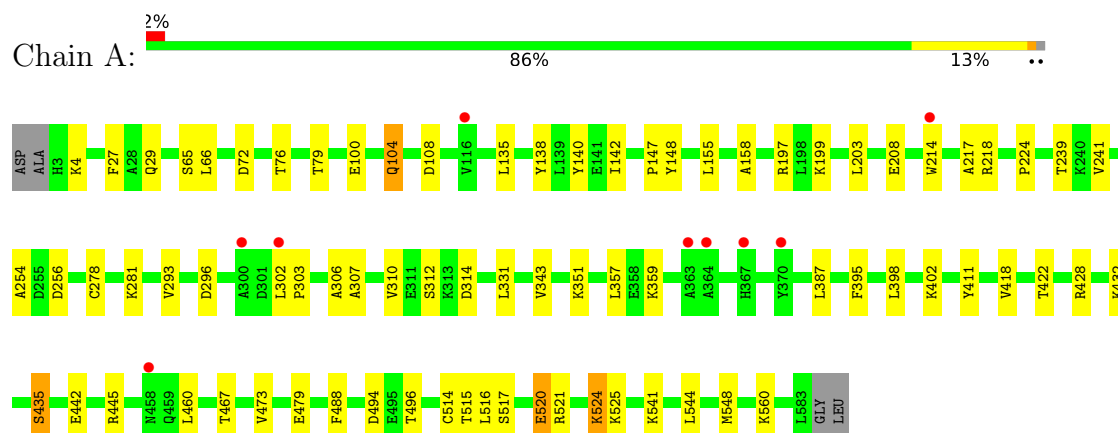
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	140	Total	O	0	0
			140	140		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Albumin



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	185.00Å 38.42Å 95.78Å 90.00° 105.18° 90.00°	Depositor
Resolution (Å)	34.15 – 2.15 34.15 – 2.15	Depositor EDS
% Data completeness (in resolution range)	98.4 (34.15-2.15) 98.4 (34.15-2.15)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 2.16Å)	Xtriage
Refinement program	PHENIX (1.19.2_4158: ???)	Depositor
R, R_{free}	0.225 , 0.269 0.227 , 0.263	Depositor DCC
R_{free} test set	1818 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	44.4	Xtriage
Anisotropy	0.589	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.5	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	0.95	EDS
Total number of atoms	4894	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.83% 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: NO, RU, CL, GOL, DMS, PLM

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	0.39	0/4696	0.60	0/6339

There are no bond length outliers.

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	4606	0	4502	54	0
2	A	2	0	0	0	0
3	A	8	0	8	0	0
4	A	1	0	0	0	0
5	A	1	0	0	0	0
6	A	108	0	186	12	0
7	A	12	0	16	2	0
8	A	16	0	24	7	0
9	A	140	0	0	11	0
All	All	4894	0	4736	57	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 6.

All (57) 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:293:VAL:HG12	9:A:769:HOH:O	1.72	0.90
8:A:616:DMS:S	9:A:816:HOH:O	2.30	0.89
7:A:612:GOL:H12	9:A:794:HOH:O	1.75	0.85
1:A:293:VAL:CG1	9:A:769:HOH:O	2.26	0.80
1:A:418:VAL:HG21	6:A:608:PLM:HB1	1.71	0.71
1:A:524:LYS:HE2	9:A:823:HOH:O	1.93	0.68
1:A:27:PHE:CD1	9:A:826:HOH:O	2.51	0.63
1:A:254:ALA:HB1	6:A:607:PLM:H52	1.80	0.63
1:A:411:TYR:HB3	6:A:608:PLM:H91	1.81	0.62
7:A:612:GOL:C1	9:A:794:HOH:O	2.40	0.59
1:A:27:PHE:CE1	9:A:826:HOH:O	2.56	0.58
1:A:473:VAL:HG11	6:A:608:PLM:HD1	1.87	0.56
1:A:488:PHE:HB3	6:A:608:PLM:H51	1.87	0.56
1:A:224:PRO:HD2	1:A:296:ASP:HB3	1.87	0.54
1:A:281:LYS:NZ	9:A:706:HOH:O	2.34	0.53
1:A:217:ALA:HB3	1:A:343:VAL:HG13	1.89	0.53
1:A:79:THR:OG1	8:A:614:DMS:O	2.26	0.53
1:A:387:LEU:HB3	6:A:605:PLM:H71	1.90	0.53
1:A:422:THR:HG21	6:A:608:PLM:HG2	1.90	0.53
1:A:27:PHE:CG	9:A:826:HOH:O	2.62	0.52
1:A:428:ARG:O	1:A:432:LYS:HG3	2.11	0.51
1:A:135:LEU:HB3	6:A:610:PLM:HB1	1.93	0.50
1:A:199:LYS:HD3	8:A:613:DMS:H22	1.93	0.50
1:A:521:ARG:HG2	1:A:525:LYS:HE2	1.94	0.48
1:A:541:LYS:HA	1:A:541:LYS:HD2	1.68	0.48
1:A:138:TYR:O	1:A:142:ILE:HG12	2.13	0.48
1:A:108:ASP:HB2	1:A:148:TYR:HE2	1.80	0.47
1:A:516:LEU:HB3	1:A:520:GLU:HB2	1.95	0.47
1:A:307:ALA:O	1:A:312:SER:HB3	2.15	0.47
1:A:241:VAL:HG22	1:A:256:ASP:HB3	1.96	0.46
1:A:494:ASP:OD1	1:A:496:THR:HG22	2.16	0.46
1:A:65:SER:HA	8:A:616:DMS:H13	1.96	0.46
1:A:66:LEU:H	8:A:616:DMS:C1	2.29	0.46
1:A:306:ALA:HA	1:A:310:VAL:HG22	1.97	0.46
1:A:442:GLU:HA	1:A:445:ARG:HD2	1.99	0.45
1:A:218:ARG:HA	1:A:218:ARG:HD2	1.73	0.45
1:A:214:TRP:CE2	8:A:613:DMS:H21	2.51	0.45
1:A:155:LEU:HD11	6:A:610:PLM:HG3	1.99	0.44
1:A:158:ALA:HB1	6:A:610:PLM:HA1	1.98	0.44
1:A:351:LYS:HG3	6:A:609:PLM:H41	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:PHE:CZ	9:A:826:HOH:O	2.71	0.43
1:A:108:ASP:HB2	1:A:148:TYR:CE2	2.53	0.43
1:A:460:LEU:HD21	6:A:608:PLM:HC1	2.00	0.43
1:A:100:GLU:O	1:A:104:GLN:HG2	2.18	0.43
1:A:398:LEU:HB3	1:A:402:LYS:HB2	2.01	0.42
1:A:560:LYS:HB2	1:A:560:LYS:HE2	1.62	0.42
1:A:544:LEU:HB3	1:A:548:MET:HE2	2.01	0.42
1:A:66:LEU:H	8:A:616:DMS:H13	1.85	0.42
1:A:72:ASP:O	1:A:76:THR:HG23	2.20	0.42
1:A:104:GLN:HG2	1:A:104:GLN:H	1.64	0.41
1:A:278:CYS:HA	1:A:281:LYS:HG3	2.01	0.41
1:A:197:ARG:HA	1:A:197:ARG:HD2	1.75	0.41
1:A:331:LEU:HD12	1:A:331:LEU:HA	1.91	0.41
1:A:208:GLU:HG3	1:A:239:THR:HG21	2.03	0.41
1:A:302:LEU:HA	1:A:303:PRO:HD3	1.88	0.41
1:A:29:GLN:HG2	1:A:147:PRO:HA	2.03	0.40
1:A:395:PHE:CZ	1:A:435:SER:HA	2.56	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	579/585 (99%)	568 (98%)	11 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	503/511 (98%)	488 (97%)	15 (3%)	36	38

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LYS
1	A	104	GLN
1	A	140	TYR
1	A	203	LEU
1	A	314	ASP
1	A	357	LEU
1	A	359	LYS
1	A	435	SER
1	A	467	THR
1	A	479	GLU
1	A	514	CYS
1	A	515	THR
1	A	517	SER
1	A	520	GLU
1	A	524	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	196	GLN
1	A	247	HIS
1	A	288	HIS
1	A	397	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 2 are monoatomic - leaving 14 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	$\# Z > 2$
6	PLM	A	610	-	17,17,17	1.03	1 (5%)	17,17,17	0.73	2 (11%)
6	PLM	A	608	-	17,17,17	0.94	1 (5%)	17,17,17	0.73	2 (11%)
6	PLM	A	609	-	17,17,17	0.96	1 (5%)	17,17,17	0.73	1 (5%)
7	GOL	A	612	-	5,5,5	0.93	0	5,5,5	1.07	0
2	NO	A	601	5	0,1,1	-	-	-		
6	PLM	A	605	-	17,17,17	0.90	1 (5%)	17,17,17	0.74	2 (11%)
8	DMS	A	615	-	3,3,3	0.66	0	3,3,3	0.53	0
6	PLM	A	607	-	17,17,17	0.95	1 (5%)	17,17,17	0.72	1 (5%)
6	PLM	A	606	-	17,17,17	0.91	1 (5%)	17,17,17	0.73	2 (11%)
8	DMS	A	614	-	3,3,3	0.67	0	3,3,3	0.54	0
7	GOL	A	611	-	5,5,5	0.91	0	5,5,5	1.08	0
3	VAL	A	602	5	6,7,7	2.08	2 (33%)	8,9,9	2.73	1 (12%)
8	DMS	A	616	-	3,3,3	0.65	0	3,3,3	0.50	0
8	DMS	A	613	-	3,3,3	0.67	0	3,3,3	0.54	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
6	PLM	A	610	-	-	4/15/15/15	-
6	PLM	A	608	-	-	9/15/15/15	-
6	PLM	A	609	-	-	9/15/15/15	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	GOL	A	612	-	-	2/4/4/4	-
6	PLM	A	605	-	-	7/15/15/15	-
6	PLM	A	607	-	-	9/15/15/15	-
6	PLM	A	606	-	-	5/15/15/15	-
7	GOL	A	611	-	-	2/4/4/4	-
3	VAL	A	602	5	-	5/8/8/8	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	VAL	O-C	4.58	1.35	1.22
6	A	610	PLM	C2-C1	3.25	1.58	1.50
6	A	609	PLM	C2-C1	3.08	1.57	1.50
6	A	607	PLM	C2-C1	3.06	1.57	1.50
6	A	608	PLM	C2-C1	3.00	1.57	1.50
6	A	606	PLM	C2-C1	2.81	1.57	1.50
6	A	605	PLM	C2-C1	2.78	1.57	1.50
3	A	602	VAL	OXT-C	-2.12	1.23	1.30

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	VAL	OXT-C-O	-7.07	108.03	124.08
6	A	605	PLM	O1-C1-O2	2.14	128.84	123.33
6	A	607	PLM	O1-C1-O2	2.14	128.84	123.33
6	A	606	PLM	O1-C1-O2	2.14	128.83	123.33
6	A	609	PLM	O1-C1-O2	2.12	128.80	123.33
6	A	608	PLM	O1-C1-O2	2.09	128.71	123.33
6	A	610	PLM	O1-C1-O2	2.08	128.68	123.33
6	A	608	PLM	O2-C1-C2	-2.06	116.56	123.09
6	A	606	PLM	O2-C1-C2	-2.05	116.59	123.09
6	A	605	PLM	O2-C1-C2	-2.03	116.64	123.09
6	A	610	PLM	O2-C1-C2	-2.01	116.71	123.09

There are no chirality outliers.

All (52) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	602	VAL	N-CA-CB-CG2
7	A	612	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
6	A	608	PLM	C1-C2-C3-C4
6	A	609	PLM	C1-C2-C3-C4
6	A	606	PLM	C1-C2-C3-C4
7	A	611	GOL	O1-C1-C2-C3
6	A	609	PLM	C7-C8-C9-CA
6	A	607	PLM	C4-C5-C6-C7
7	A	612	GOL	O1-C1-C2-O2
6	A	608	PLM	C7-C8-C9-CA
6	A	608	PLM	C9-CA-CB-CC
6	A	609	PLM	C2-C3-C4-C5
6	A	609	PLM	C9-CA-CB-CC
6	A	607	PLM	C5-C6-C7-C8
6	A	607	PLM	C6-C7-C8-C9
6	A	610	PLM	CC-CD-CE-CF
3	A	602	VAL	O-C-CA-CB
3	A	602	VAL	OXT-C-CA-CB
6	A	610	PLM	CB-CC-CD-CE
6	A	608	PLM	C3-C4-C5-C6
6	A	605	PLM	CB-CC-CD-CE
6	A	607	PLM	C7-C8-C9-CA
6	A	608	PLM	C6-C7-C8-C9
6	A	609	PLM	CA-CB-CC-CD
6	A	607	PLM	CC-CD-CE-CF
6	A	607	PLM	C3-C4-C5-C6
6	A	606	PLM	C3-C4-C5-C6
6	A	607	PLM	C9-CA-CB-CC
6	A	606	PLM	C2-C3-C4-C5
6	A	605	PLM	CA-CB-CC-CD
6	A	608	PLM	C5-C6-C7-C8
7	A	611	GOL	O1-C1-C2-O2
6	A	605	PLM	C9-CA-CB-CC
6	A	608	PLM	C2-C3-C4-C5
3	A	602	VAL	OXT-C-CA-N
6	A	605	PLM	C8-C9-CA-CB
6	A	607	PLM	CD-CE-CF-CG
6	A	609	PLM	O1-C1-C2-C3
6	A	609	PLM	O2-C1-C2-C3
6	A	606	PLM	O1-C1-C2-C3
6	A	606	PLM	O2-C1-C2-C3
6	A	608	PLM	CB-CC-CD-CE
6	A	607	PLM	C2-C3-C4-C5
6	A	610	PLM	O2-C1-C2-C3

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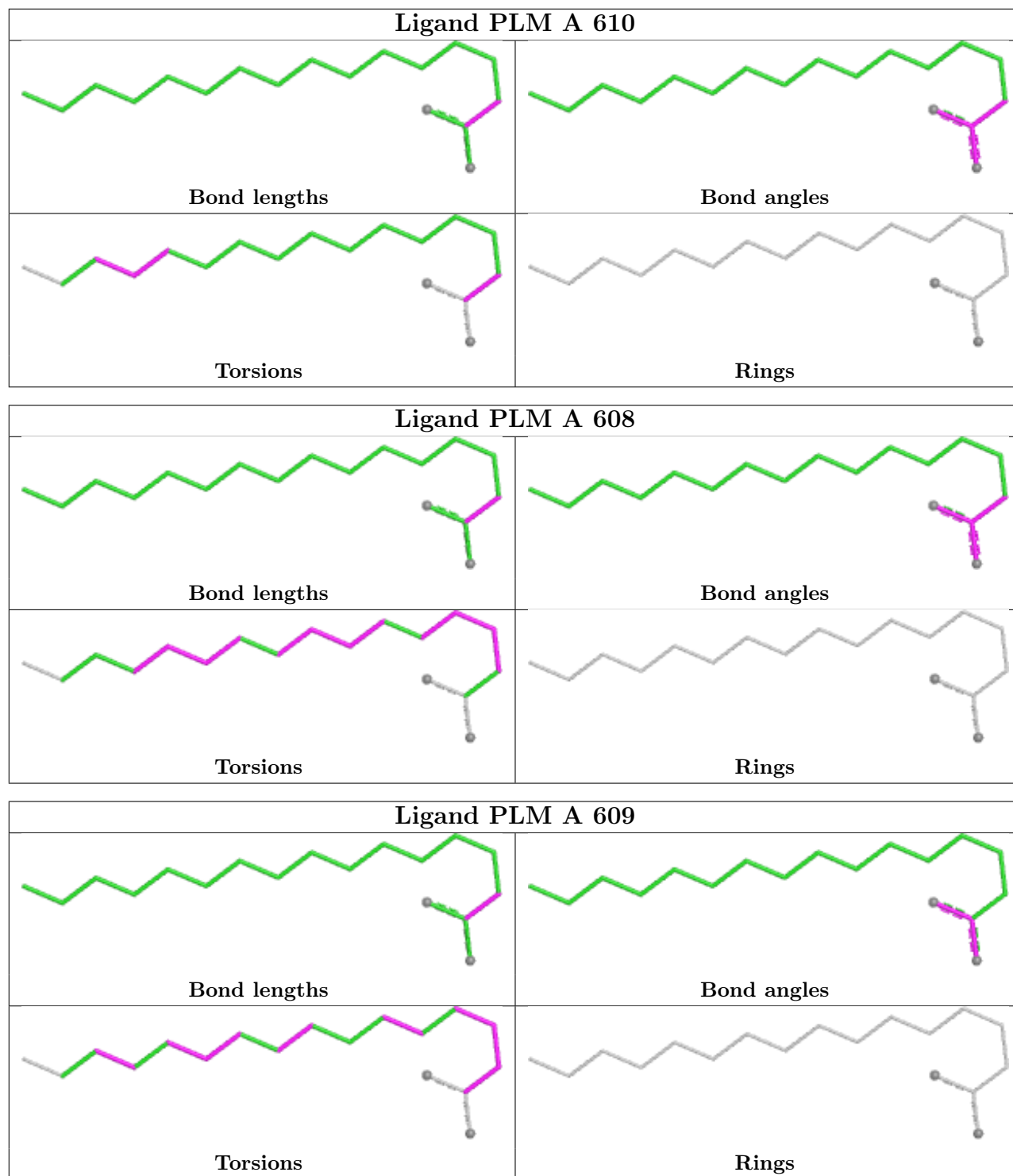
Mol	Chain	Res	Type	Atoms
6	A	609	PLM	C4-C5-C6-C7
6	A	610	PLM	O1-C1-C2-C3
6	A	608	PLM	CA-CB-CC-CD
6	A	605	PLM	C7-C8-C9-CA
6	A	609	PLM	CC-CD-CE-CF
6	A	605	PLM	O1-C1-C2-C3
6	A	605	PLM	O2-C1-C2-C3
3	A	602	VAL	C-CA-CB-CG1

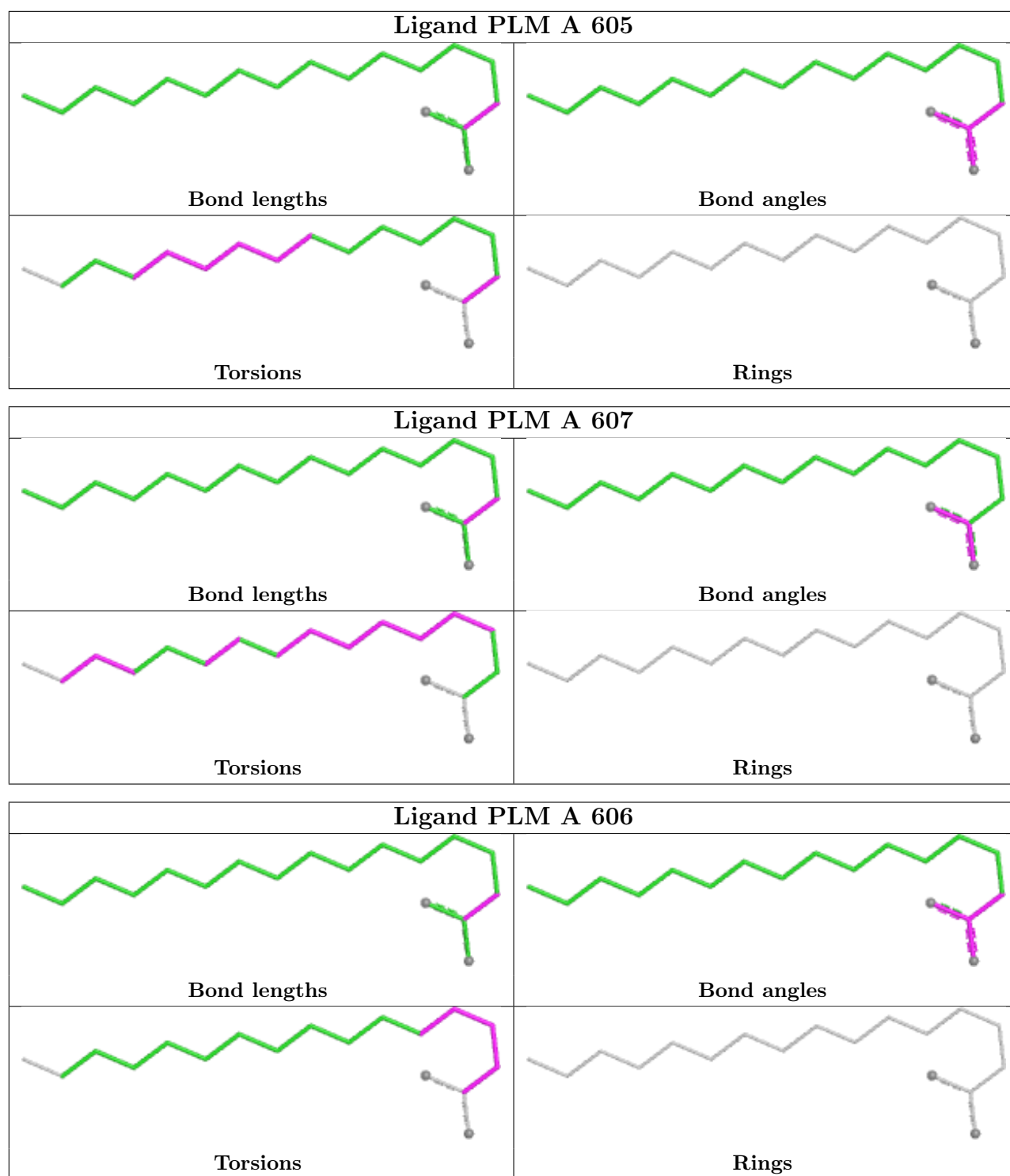
There are no ring outliers.

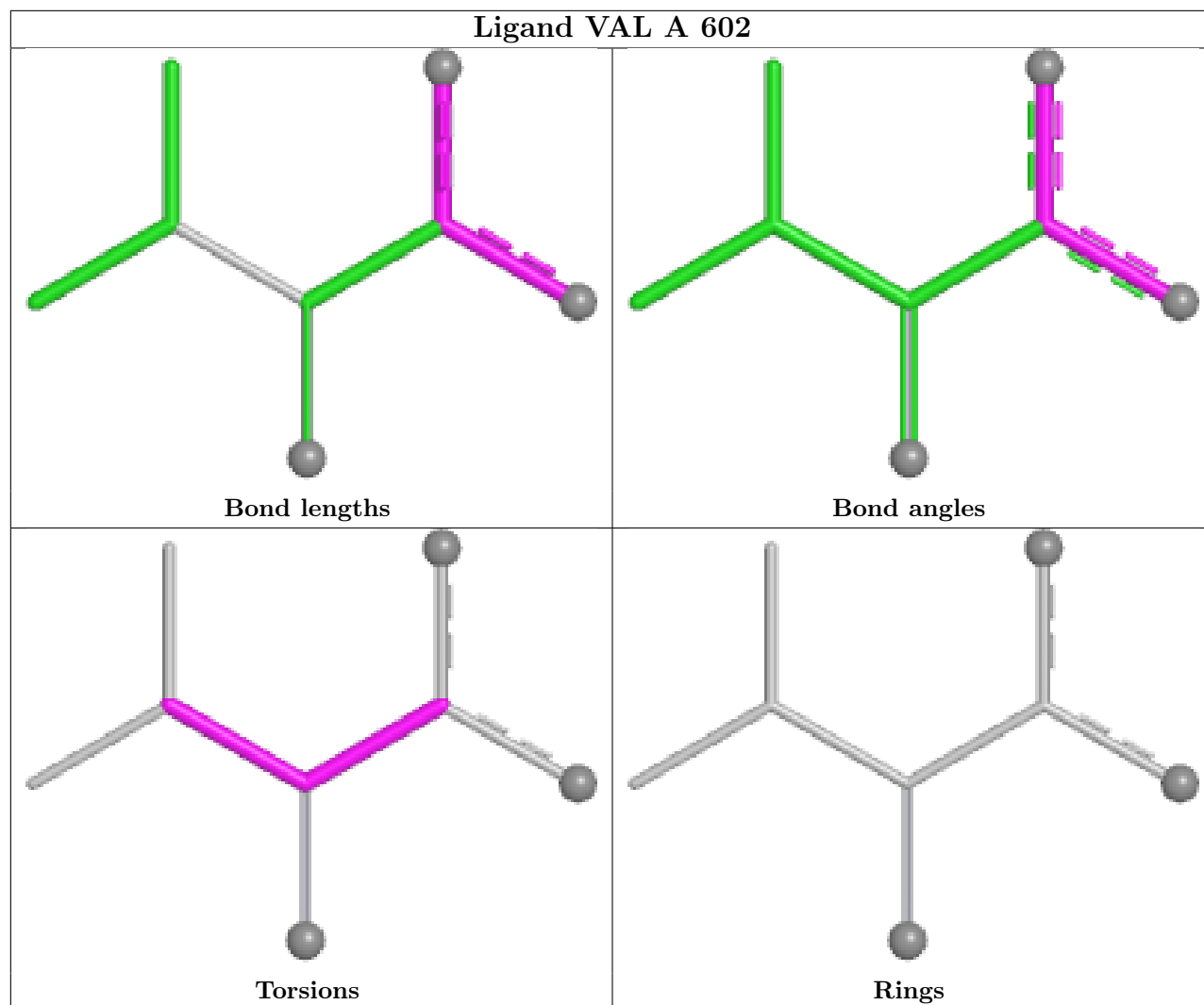
9 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	610	PLM	3	0
6	A	608	PLM	6	0
6	A	609	PLM	1	0
7	A	612	GOL	2	0
6	A	605	PLM	1	0
6	A	607	PLM	1	0
8	A	614	DMS	1	0
8	A	616	DMS	4	0
8	A	613	DMS	2	0

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







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	581/585 (99%)	0.34	9 (1%) 72 76	29, 50, 68, 84	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	370	TYR	2.9
1	A	302	LEU	2.6
1	A	458	ASN	2.5
1	A	364	ALA	2.4
1	A	214	TRP	2.3
1	A	116	VAL	2.2
1	A	300	ALA	2.2
1	A	363	ALA	2.2
1	A	367	HIS	2.1

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](#)

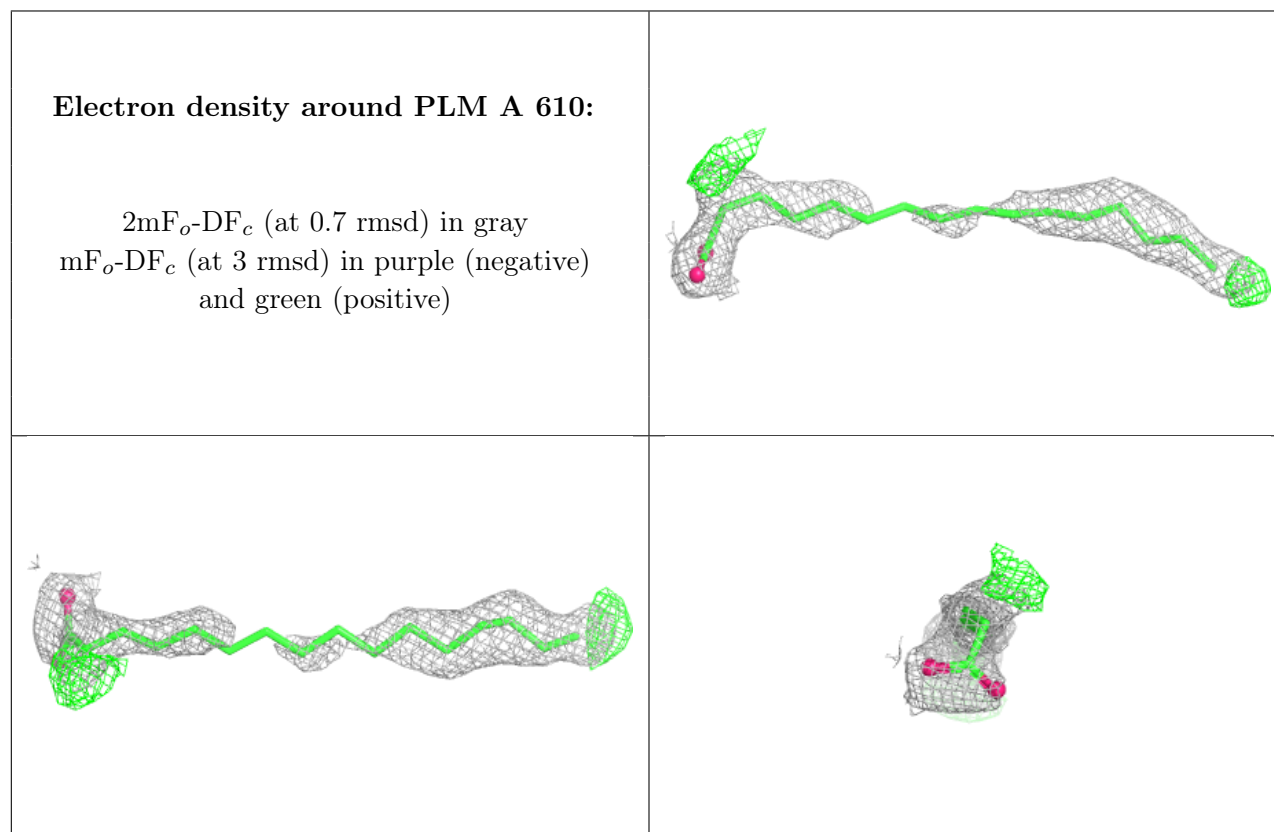
There are no oligosaccharides in this entry.

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.

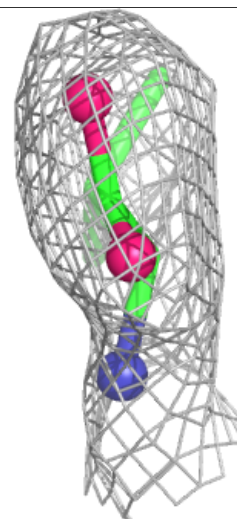
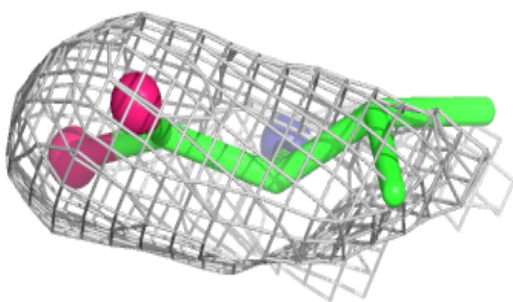
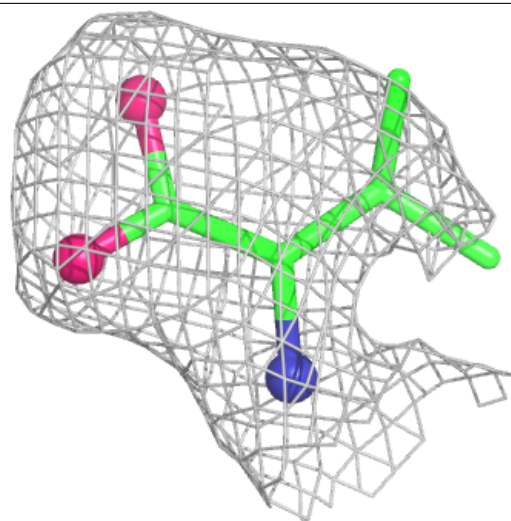
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NO	A	601	2/2	0.74	0.09	76,76,76,77	0
8	DMS	A	616	4/4	0.76	0.17	48,51,67,85	0
6	PLM	A	610	18/18	0.79	0.15	45,58,73,73	0
7	GOL	A	612	6/6	0.81	0.16	51,55,65,67	0
8	DMS	A	615	4/4	0.81	0.17	53,58,71,78	0
3	VAL	A	602	8/8	0.81	0.15	64,69,73,75	0
7	GOL	A	611	6/6	0.83	0.11	58,61,63,63	0
8	DMS	A	614	4/4	0.83	0.18	74,78,78,90	0
8	DMS	A	613	4/4	0.86	0.19	61,65,78,83	0
4	CL	A	603	1/1	0.86	0.10	97,97,97,97	0
6	PLM	A	608	18/18	0.87	0.16	42,54,63,65	0
6	PLM	A	609	18/18	0.88	0.13	40,51,63,70	0
6	PLM	A	607	18/18	0.92	0.11	39,46,53,54	0
6	PLM	A	605	18/18	0.93	0.12	35,48,56,59	0
6	PLM	A	606	18/18	0.94	0.09	35,43,50,52	0
5	RU	A	604	1/1	0.97	0.06	77,77,77,77	1

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



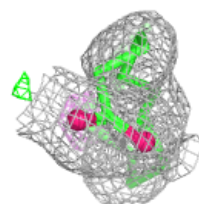
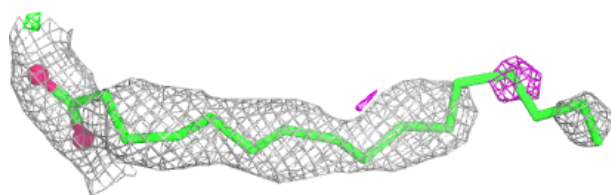
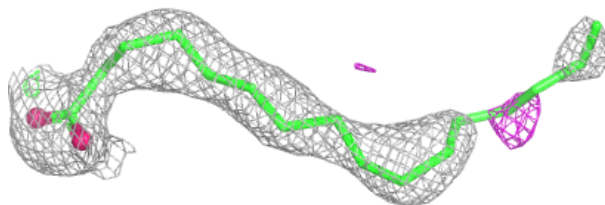
Electron density around VAL A 602:

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

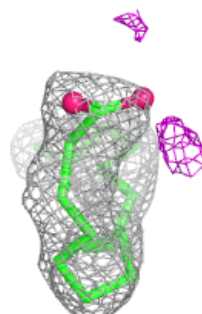
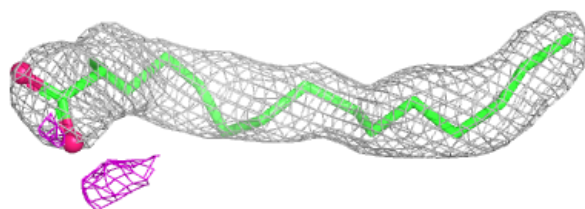
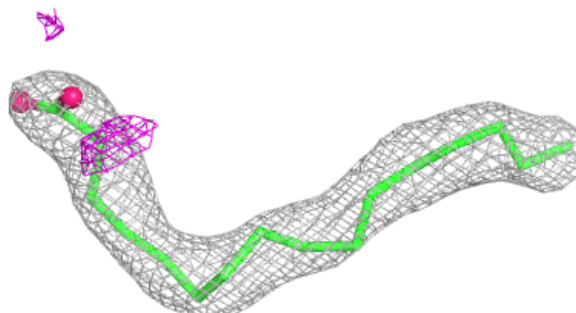


Electron density around PLM A 608:

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

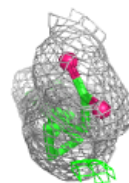
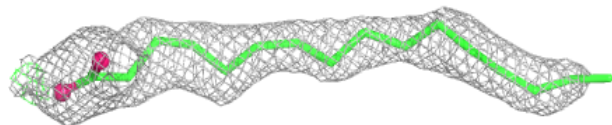
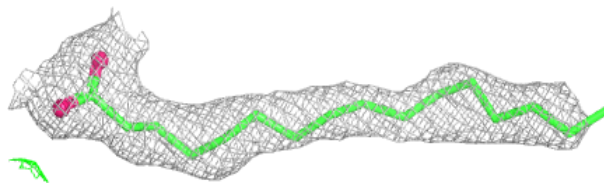
**Electron density around PLM A 609:**

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

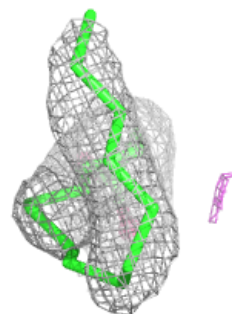
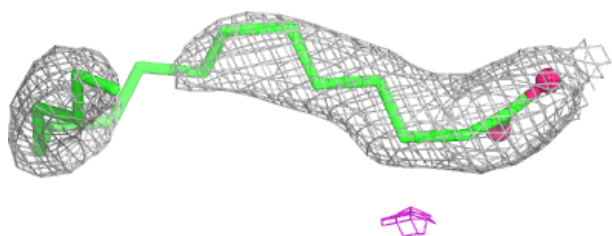
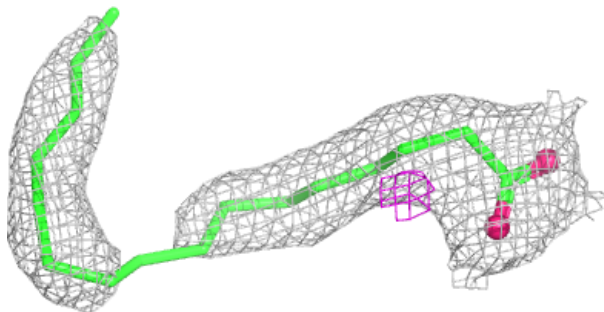


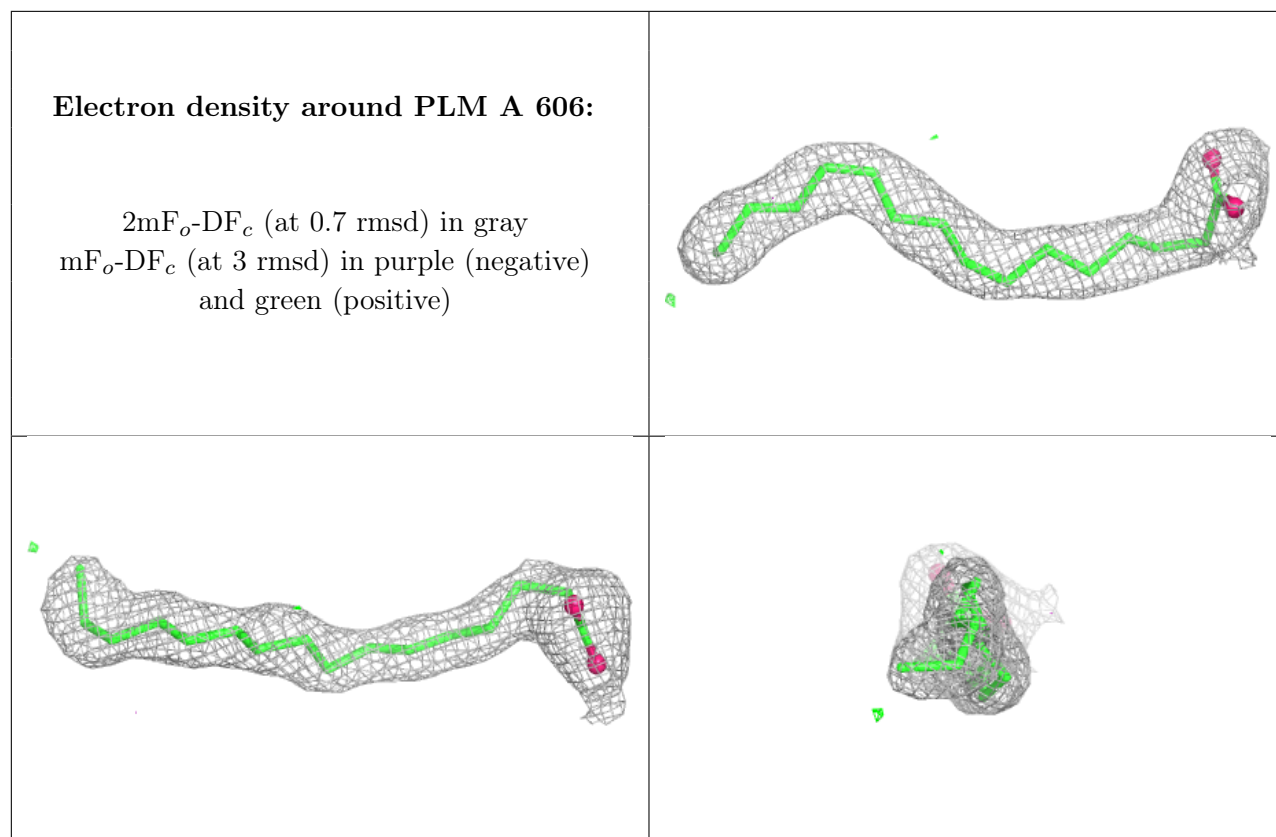
Electron density around PLM A 607:

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

**Electron density around PLM A 605:**

$2mF_o-DF_c$ (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.