



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 28, 2021 – 04:23 PM JST

PDB ID : 7CHG  
Title : The structure of human pregnane X receptor in complex with an SRC-1 coactivator peptide and a limonoid compound  
Authors : Xia, Y.; Yao, D.; Huang, C.; Cao, Y.  
Deposited on : 2020-07-05  
Resolution : 1.93 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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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 : 2.20  
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) : 2.20

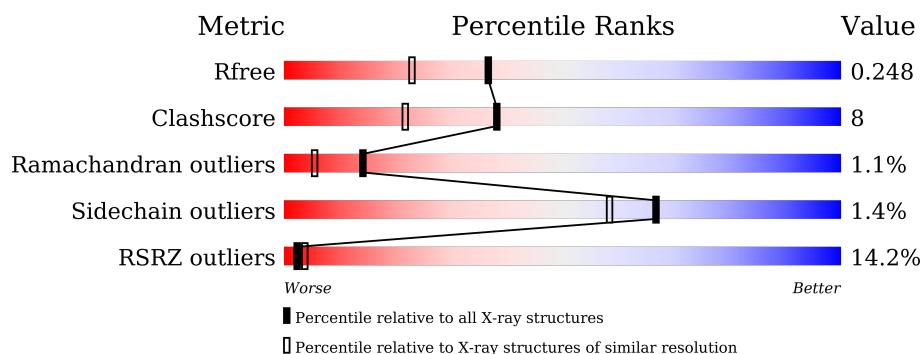
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.93 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	4310 (1.96-1.92)
Clashscore	141614	1023 (1.94-1.94)
Ramachandran outliers	138981	1007 (1.94-1.94)
Sidechain outliers	138945	1007 (1.94-1.94)
RSRZ outliers	127900	4250 (1.96-1.92)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	337	<div> <div>12%</div> <div> <div></div> <div>71%</div> <div>10%</div> <div>..</div> <div>16%</div> </div> </div>
1	B	337	<div> <div>12%</div> <div> <div></div> <div>72%</div> <div>11%</div> <div>.</div> <div>16%</div> </div> </div>

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
2	G3L	A	601	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4765 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.

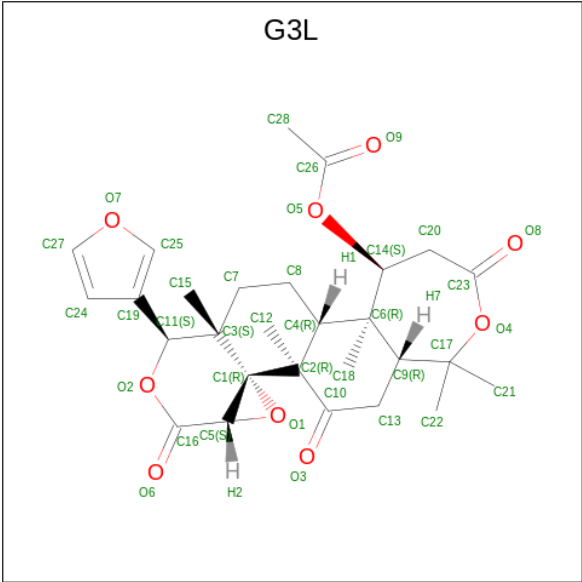
- Molecule 1 is a protein called Nuclear receptor subfamily 1 group I member 2, Nuclear receptor coactivator 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	282	Total	C	N	O	S	0	2	0
			2269	1463	386	401	19			
1	B	283	Total	C	N	O	S	0	1	0
			2293	1476	397	402	18			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	129	MET	-	initiating methionine	UNP O75469
A	435	SER	-	linker	UNP O75469
A	436	SER	-	linker	UNP O75469
A	437	SER	-	linker	UNP O75469
A	438	GLY	-	linker	UNP O75469
A	439	GLY	-	linker	UNP O75469
A	440	THR	-	linker	UNP O75469
B	129	MET	-	initiating methionine	UNP O75469
B	435	SER	-	linker	UNP O75469
B	436	SER	-	linker	UNP O75469
B	437	SER	-	linker	UNP O75469
B	438	GLY	-	linker	UNP O75469
B	439	GLY	-	linker	UNP O75469
B	440	THR	-	linker	UNP O75469

- Molecule 2 is Nomilin (three-letter code: G3L) (formula: C<sub>28</sub>H<sub>34</sub>O<sub>9</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	C O	0	0
			37	28 9		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	86	Total	O	0	0
			86	86		
3	B	80	Total	O	0	0
			80	80		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.93Å 90.11Å 106.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	26.59 – 1.93 26.59 – 1.93	Depositor EDS
% Data completeness (in resolution range)	98.8 (26.59-1.93) 98.8 (26.59-1.93)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.08 (at 1.93Å)	Xtriage
Refinement program	PHENIX v1.13	Depositor
R, $R_{free}$	0.212 , 0.250 0.214 , 0.248	Depositor DCC
$R_{free}$ test set	1994 reflections (3.27%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	43.3	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 52.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4765	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: G3L

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	0.43	0/2324	0.61	3/3133 (0.1%)
1	B	0.41	0/2343	0.56	1/3157 (0.0%)
All	All	0.42	0/4667	0.59	4/6290 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	2
All	All	0	4

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	145	GLU	O-C-N	-5.78	113.45	122.70
1	B	308	LEU	C-N-CA	5.74	136.05	121.70
1	A	383	GLN	C-N-CD	5.55	140.06	128.40
1	A	447	SER	O-C-N	-5.03	114.65	122.70

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	145	GLU	Mainchain
1	A	447	SER	Mainchain

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Mol	Chain	Res	Type	Group
1	B	195	GLU	Mainchain
1	B	208	SER	Mainchain

## 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2269	0	2255	48	0
1	B	2293	0	2284	21	0
2	A	37	0	0	16	0
3	A	86	0	0	0	0
3	B	80	0	0	1	0
All	All	4765	0	4539	69	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 8.

All (69) 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:209:LEU:O	2:A:601:G3L:O9	1.72	1.07
1:A:209:LEU:HA	2:A:601:G3L:C28	1.90	1.02
1:A:319:LEU:H	1:A:319:LEU:HD23	1.27	0.97
1:A:193:ARG:NH1	1:A:193:ARG:HB3	1.82	0.92
1:A:193:ARG:HB3	1:A:193:ARG:HH11	1.35	0.91
1:A:411:LEU:HD12	2:A:601:G3L:C24	2.03	0.87
1:A:206:LEU:HD12	1:A:236:ILE:HD12	1.56	0.85
1:A:285:GLN:NE2	2:A:601:G3L:O3	2.10	0.84
1:A:209:LEU:CA	2:A:601:G3L:C28	2.57	0.82
1:A:407:HIS:HE1	2:A:601:G3L:O2	1.64	0.80
1:A:193:ARG:HH11	1:A:193:ARG:CB	1.97	0.76
1:A:407:HIS:CE1	2:A:601:G3L:O2	2.39	0.75
1:B:205:ASP:O	1:B:208:SER:OG	2.07	0.72
1:A:209:LEU:O	2:A:601:G3L:C26	2.41	0.69
1:A:209:LEU:HD23	2:A:601:G3L:C28	2.24	0.68
1:B:203:ARG:HE	1:B:232:GLY:HA2	1.59	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:165:THR:HG1	1:B:167:SER:HG	1.41	0.66
1:A:230:ASP:OD1	1:A:231:SER:N	2.28	0.65
1:B:254:ILE:HD11	1:B:287:ARG:HD2	1.80	0.64
1:A:415:GLN:HG2	1:A:431:ILE:HD13	1.80	0.63
1:A:285:GLN:HE22	2:A:601:G3L:C10	2.12	0.63
1:A:319:LEU:H	1:A:319:LEU:CD2	2.03	0.63
1:B:358:GLN:HB3	1:B:361:VAL:HG12	1.83	0.61
1:A:206:LEU:CD1	1:A:236:ILE:HD12	2.30	0.60
1:A:195:GLU:OE1	1:A:199:TRP:HZ3	1.85	0.59
1:B:360:ARG:O	1:B:364:GLN:HG3	2.03	0.58
1:B:352:ASP:OD1	1:B:401:ARG:NH2	2.37	0.57
1:A:247:SER:OG	2:A:601:G3L:O1	2.16	0.56
1:A:304:LEU:HD23	1:A:305:SER:N	2.22	0.55
1:B:230:ASP:CG	1:B:231:SER:H	2.10	0.55
1:A:407:HIS:CE1	2:A:601:G3L:C16	2.92	0.52
1:B:270:GLU:HB3	1:B:448:LEU:HD23	1.91	0.52
1:A:213:LEU:HD11	1:A:304:LEU:HG	1.92	0.51
1:B:281:PHE:HZ	1:B:323:MET:HE2	1.75	0.50
1:A:233:GLY:O	1:A:236:ILE:HG12	2.13	0.49
1:A:381:ARG:O	1:A:383:GLN:N	2.45	0.49
1:B:214:GLN:HB3	1:B:305:SER:HB2	1.96	0.48
1:B:217:GLY:C	1:B:219:ASP:H	2.17	0.48
1:B:408:THR:O	1:B:412:LEU:HG	2.13	0.48
1:B:254:ILE:HD12	1:B:283:LEU:HB3	1.95	0.48
1:A:352:ASP:OD1	1:A:352:ASP:N	2.39	0.48
1:B:213:LEU:HD21	1:B:215:LEU:HD11	1.96	0.47
1:A:323:MET:O	1:A:323:MET:HE2	2.15	0.46
1:A:213:LEU:HD11	1:A:304:LEU:CG	2.46	0.46
1:B:406:GLN:NE2	3:B:502:HOH:O	2.35	0.45
1:A:304:LEU:HD23	1:A:304:LEU:C	2.37	0.45
1:A:281:PHE:HD1	2:A:601:G3L:O6	2.00	0.45
1:A:205:ASP:HB3	1:A:410:ARG:HG3	1.98	0.45
1:A:319:LEU:O	1:A:319:LEU:HG	2.17	0.45
1:A:213:LEU:HD13	1:A:306:TYR:CE1	2.52	0.45
1:B:209:LEU:HD11	1:B:410:ARG:HD3	2.00	0.44
1:A:193:ARG:NH1	1:A:193:ARG:CB	2.63	0.43
1:A:209:LEU:O	2:A:601:G3L:C28	2.66	0.43
1:B:205:ASP:HB3	1:B:410:ARG:HG3	2.00	0.43
1:A:410:ARG:HE	1:A:410:ARG:HB2	1.56	0.43
1:B:209:LEU:HD21	1:B:407[B]:HIS:HE1	1.83	0.43
1:A:209:LEU:CD2	2:A:601:G3L:C28	2.94	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:LYS:HD2	1:A:459:LEU:HA	2.01	0.43
1:A:323:MET:HE2	1:A:327:HIS:CD2	2.54	0.42
1:B:349:PHE:O	1:B:366:GLN:HB2	2.20	0.42
1:A:304:LEU:HD21	1:A:306:TYR:CE1	2.54	0.41
1:A:349:PHE:O	1:A:366:GLN:HB2	2.20	0.41
1:B:216:ARG:NH2	1:B:302:GLY:O	2.44	0.41
1:A:318:LEU:N	1:A:319:LEU:HD23	2.36	0.41
1:A:209:LEU:CB	2:A:601:G3L:C28	2.98	0.41
1:A:323:MET:HE2	1:A:327:HIS:HD2	1.86	0.41
1:A:208:SER:HB2	1:A:321:GLU:OE1	2.21	0.41
1:A:384:PRO:O	1:A:385:ALA:HB3	2.21	0.41
1:A:318:LEU:O	1:A:318:LEU:HG	2.20	0.40

There are no symmetry-related clashes.

## 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	276/337 (82%)	264 (96%)	9 (3%)	3 (1%)	14	5
1	B	276/337 (82%)	268 (97%)	5 (2%)	3 (1%)	14	5
All	All	552/674 (82%)	532 (96%)	14 (2%)	6 (1%)	14	5

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	231	SER
1	A	382	PRO
1	A	383	GLN
1	B	231	SER
1	B	218	GLU
1	B	232	GLY

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	245/299 (82%)	239 (98%)	6 (2%)	49	36
1	B	247/299 (83%)	246 (100%)	1 (0%)	91	91
All	All	492/598 (82%)	485 (99%)	7 (1%)	67	58

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	162	PHE
1	A	199	TRP
1	A	319	LEU
1	A	352	ASP
1	A	360	ARG
1	A	447	SER
1	B	402	SER

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

Mol	Chain	Res	Type
1	A	406	GLN
1	A	407	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

1 ligand is 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	$\# Z  > 2$
2	G3L	A	601	-	36,42,42	1.17	4 (11%)	57,72,72	1.38	8 (14%)

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	G3L	A	601	-	-	4/4/103/103	0/6/6/6

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	G3L	C6-C4	2.66	1.61	1.56
2	A	601	G3L	C3-C11	2.43	1.58	1.55
2	A	601	G3L	C1-C2	2.27	1.57	1.54
2	A	601	G3L	C1-C5	2.17	1.52	1.48

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	G3L	C1-O1-C5	4.13	63.66	61.43
2	A	601	G3L	C22-C17-C21	-2.95	105.42	109.92
2	A	601	G3L	C15-C3-C7	2.88	112.75	107.84
2	A	601	G3L	C17-O4-C23	2.74	130.72	126.77
2	A	601	G3L	O2-C16-O6	2.64	122.32	118.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	G3L	C8-C7-C3	2.49	117.13	112.84
2	A	601	G3L	O1-C1-C5	-2.37	57.30	58.80
2	A	601	G3L	O1-C1-C2	-2.13	113.04	114.90

There are no chirality outliers.

All (4) torsion outliers are listed below:

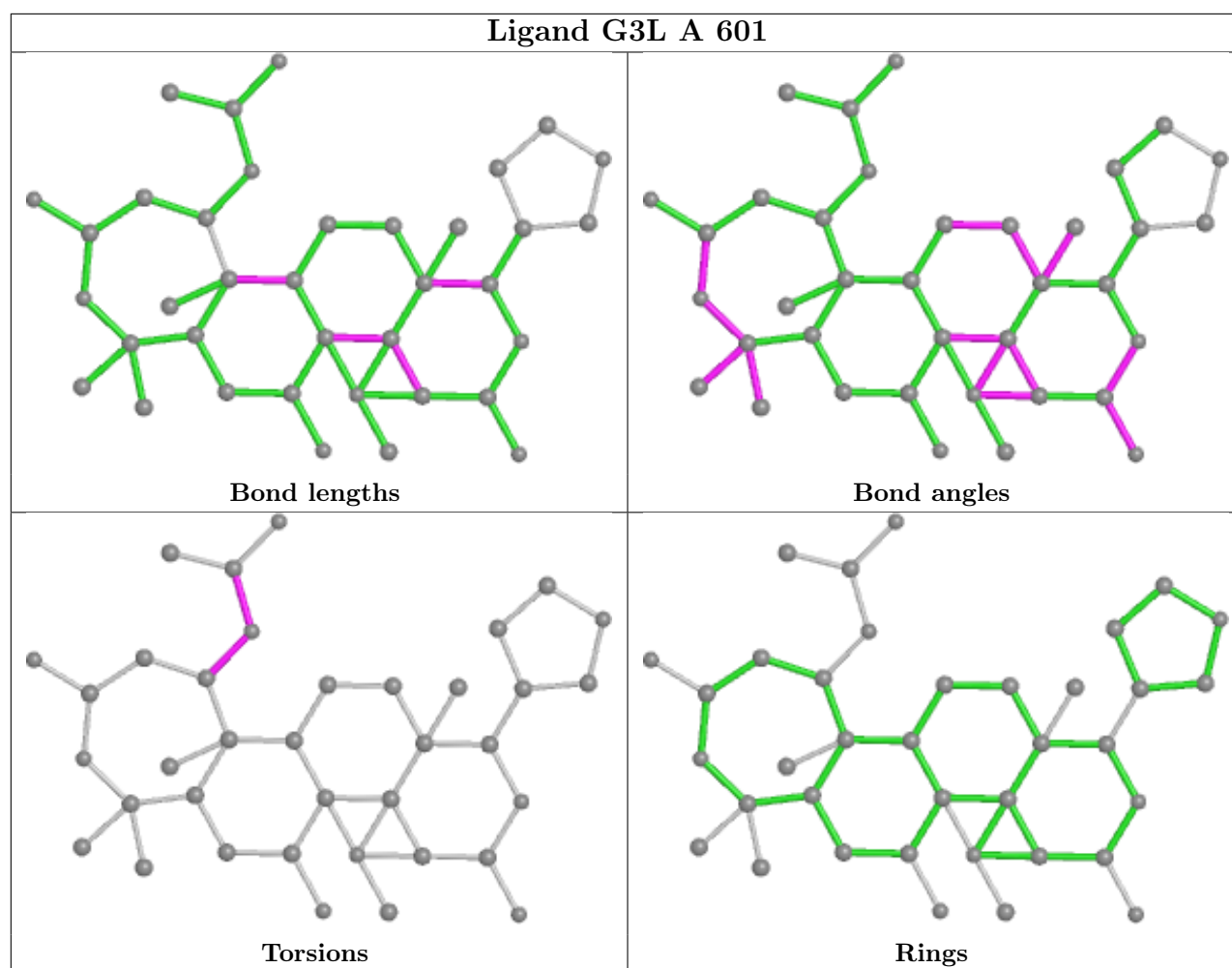
Mol	Chain	Res	Type	Atoms
2	A	601	G3L	C6-C14-O5-C26
2	A	601	G3L	C20-C14-O5-C26
2	A	601	G3L	C28-C26-O5-C14
2	A	601	G3L	O9-C26-O5-C14

There are no ring outliers.

1 monomer is involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	G3L	16	0

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



## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	282/337 (83%)	0.75	40 (14%) 2 4	30, 50, 81, 97	0
1	B	283/337 (83%)	0.84	40 (14%) 2 4	31, 52, 83, 102	0
All	All	565/674 (83%)	0.79	80 (14%) 2 4	30, 51, 82, 102	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	196	ALA	10.4
1	B	178	LEU	8.5
1	B	432	THR	7.4
1	B	179	SER	7.2
1	B	320	LEU	5.7
1	B	318	LEU	5.4
1	A	319	LEU	5.3
1	B	197	ALA	5.2
1	B	319	LEU	4.9
1	A	233	GLY	4.8
1	A	229	ALA	4.8
1	B	142	GLY	4.7
1	B	384	PRO	4.4
1	A	142	GLY	4.4
1	A	232	GLY	4.4
1	B	229	ALA	4.3
1	A	448	LEU	4.2
1	B	235	GLU	4.2
1	B	171	ASN	4.1
1	A	283	LEU	4.1
1	A	281	PHE	4.1
1	A	199	TRP	3.8
1	B	209	LEU	3.8
1	A	447	SER	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	359	HIS	3.7
1	B	195	GLU	3.7
1	A	360	ARG	3.6
1	B	283	LEU	3.6
1	A	230	ASP	3.5
1	A	209	LEU	3.5
1	B	232	GLY	3.4
1	B	199	TRP	3.4
1	B	451	ARG	3.4
1	B	233	GLY	3.3
1	A	194	GLU	3.3
1	B	317	GLN	3.3
1	B	220	GLY	3.3
1	A	254	ILE	3.2
1	B	348	LEU	3.2
1	B	207	CYS	3.1
1	A	346	ILE	3.0
1	A	284[A]	CYS	3.0
1	A	286	LEU	3.0
1	A	280	ALA	2.9
1	B	286	LEU	2.9
1	A	279	ALA	2.9
1	B	431	ILE	2.9
1	B	448	LEU	2.9
1	A	348	LEU	2.7
1	A	345	ALA	2.7
1	A	449	THR	2.7
1	A	343	MET	2.7
1	A	318	LEU	2.7
1	A	341	VAL	2.6
1	A	383	GLN	2.6
1	A	177	VAL	2.6
1	B	295	GLU	2.6
1	B	357	LEU	2.5
1	B	219	ASP	2.5
1	A	145	GLU	2.5
1	A	382	PRO	2.5
1	A	195	GLU	2.5
1	B	279	ALA	2.4
1	A	171	ASN	2.4
1	B	450	GLU	2.4
1	B	177	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	193	ARG	2.4
1	A	231	SER	2.4
1	B	385	ALA	2.4
1	B	234	LYS	2.3
1	A	235	GLU	2.3
1	A	349	PHE	2.3
1	A	432	THR	2.3
1	B	360	ARG	2.2
1	A	143	LEU	2.1
1	A	288	PHE	2.1
1	B	309	GLU	2.1
1	A	342	LEU	2.1
1	B	168	HIS	2.1
1	B	407[A]	HIS	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

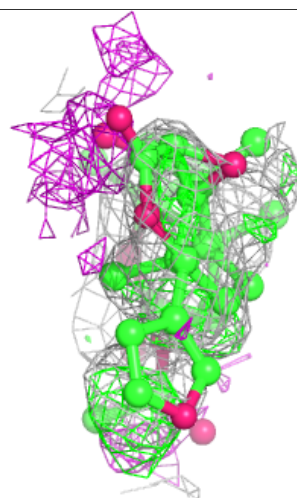
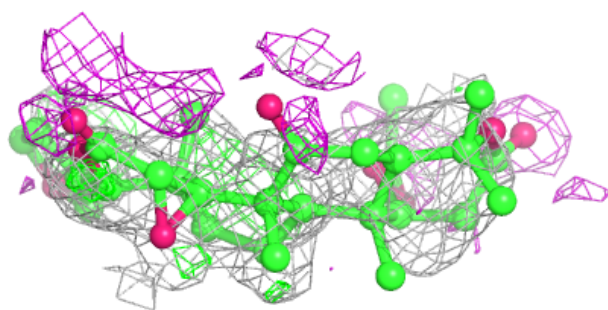
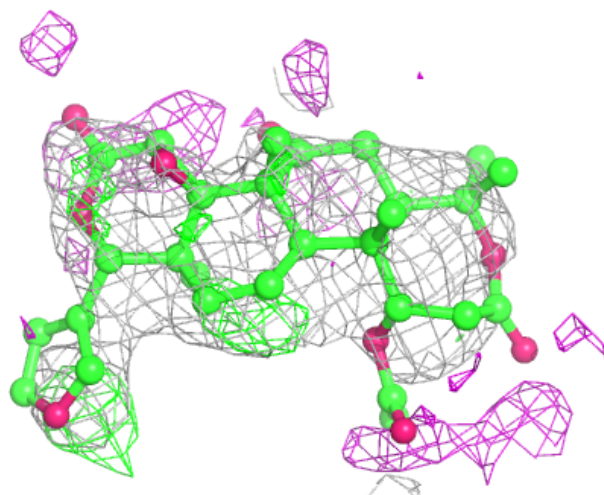
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	G3L	A	601	37/37	0.62	0.56	160,160,160,160	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 G3L A 601:**

$2mF_o-DF_c$  (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.