



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:17 PM UTC

PDB ID : 6T50 / pdb_00006t50
Title : ROR(γ)t ligand binding domain in complex with 25-hydroxycholesterol and allosteric ligand Glenmark
Authors : de Vries, R.M.J.M.; Meijer, F.A.; Brunsveld, L.
Deposited on : 2019-10-15
Resolution : 1.87 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

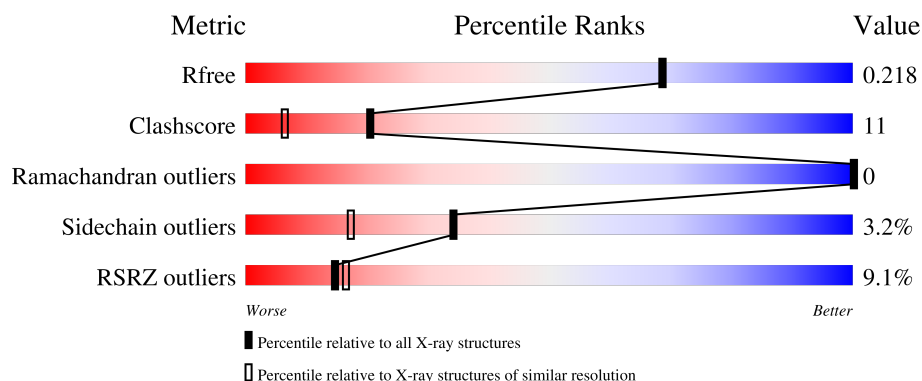
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.87 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	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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	263	8% 81% 8% 8%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2195 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.

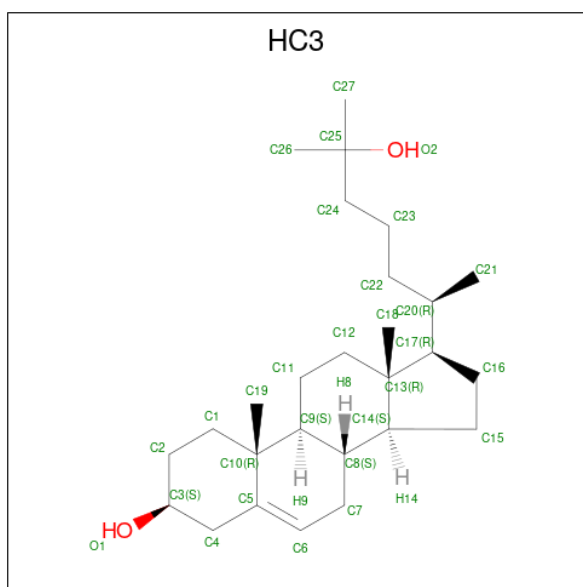
- Molecule 1 is a protein called Nuclear receptor ROR-gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	241	Total	C	N	O	S	0	8	0
			2040	1297	368	360	15			

There are 21 discrepancies between the modelled and reference sequences:

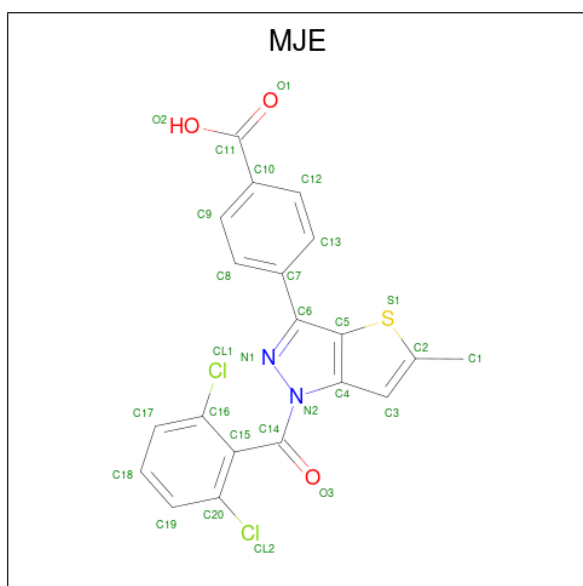
Chain	Residue	Modelled	Actual	Comment	Reference
A	245	GLY	-	expression tag	UNP P51449
A	246	SER	-	expression tag	UNP P51449
A	247	SER	-	expression tag	UNP P51449
A	248	HIS	-	expression tag	UNP P51449
A	249	HIS	-	expression tag	UNP P51449
A	250	HIS	-	expression tag	UNP P51449
A	251	HIS	-	expression tag	UNP P51449
A	252	HIS	-	expression tag	UNP P51449
A	253	HIS	-	expression tag	UNP P51449
A	254	SER	-	expression tag	UNP P51449
A	255	SER	-	expression tag	UNP P51449
A	256	GLY	-	expression tag	UNP P51449
A	257	LEU	-	expression tag	UNP P51449
A	258	VAL	-	expression tag	UNP P51449
A	259	PRO	-	expression tag	UNP P51449
A	260	ARG	-	expression tag	UNP P51449
A	261	GLY	-	expression tag	UNP P51449
A	262	SER	-	expression tag	UNP P51449
A	263	HIS	-	expression tag	UNP P51449
A	264	MET	-	expression tag	UNP P51449
A	455	HIS	CYS	conflict	UNP P51449

- Molecule 2 is 25-HYDROXYCHOLESTEROL (CCD ID: HC3) (formula: C₂₇H₄₆O₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			29	27	2		

- Molecule 3 is 4-[1-[2,6-bis(chloranyl)phenyl]carbonyl-5-methyl-thieno[3,2-c]pyrazol-3-yl]benzoic acid (CCD ID: MJE) (formula: $C_{20}H_{12}Cl_2N_2O_3S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	S	0	0
			28	20	2	2	3	1		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 5 is water.

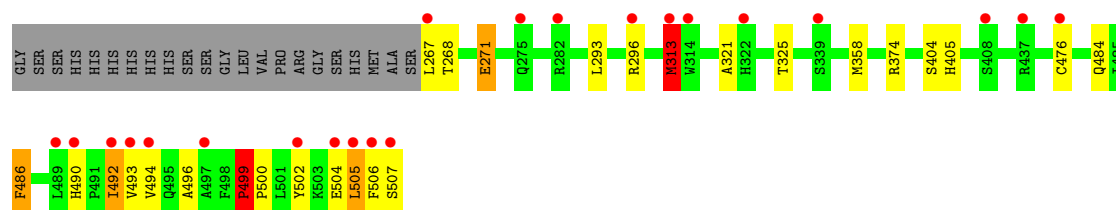
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	92	Total	O	0	0
			92	92		

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: Nuclear receptor ROR-gamma

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	108.33Å 108.33Å 108.51Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	93.81 – 1.87 93.81 – 1.87	Depositor EDS
% Data completeness (in resolution range)	100.0 (93.81-1.87) 93.5 (93.81-1.87)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.95 (at 1.87Å)	Xtriage
Refinement program	PHENIX 1.16_3549, REFMAC 5	Depositor
R, R_{free}	0.186 , 0.210 0.192 , 0.218	Depositor DCC
R_{free} test set	1578 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	41.7	Xtriage
Anisotropy	0.367	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 35.0	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	2195	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.32% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HC3, GOL, MJE

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.13	5/2088 (0.2%)	1.33	5/2810 (0.2%)

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

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	490	HIS	C-N	13.29	1.48	1.33
1	A	499	PRO	C-O	-6.59	1.17	1.24
1	A	321	ALA	C-O	-5.36	1.18	1.24
1	A	486	PHE	C-O	-5.29	1.18	1.24
1	A	405	HIS	CE1-NE2	5.02	1.37	1.32

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	490	HIS	CA-C-N	8.76	130.14	119.98
1	A	490	HIS	C-N-CA	8.76	130.14	119.98
1	A	490	HIS	O-C-N	-7.63	114.42	121.36
1	A	499	PRO	N-CA-C	6.45	118.57	110.70
1	A	271	GLU	CB-CG-CD	5.56	122.05	112.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	313[A]	MET	Mainchain
1	A	313[B]	MET	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	2040	0	2034	47	0
2	A	29	0	46	0	0
3	A	28	0	0	7	0
4	A	6	0	8	0	0
5	A	92	0	0	4	0
All	All	2195	0	2088	47	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (47) 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:502:TYR:CE1	1:A:506:PHE:CG	2.21	1.28
1:A:502:TYR:CE1	1:A:506:PHE:CD1	2.28	1.20
1:A:502:TYR:CE1	1:A:506:PHE:CB	2.25	1.20
1:A:502:TYR:HE1	1:A:506:PHE:CB	1.57	1.18
1:A:502:TYR:CE1	1:A:506:PHE:HB2	1.81	1.15
1:A:494:VAL:HB	1:A:502:TYR:CE2	1.93	1.03
1:A:313[B]:MET:HE3	1:A:486:PHE:CE1	1.97	1.00
1:A:374:ARG:NH2	5:A:702:HOH:O	1.95	0.99
1:A:494:VAL:HB	1:A:502:TYR:HE2	1.29	0.95
1:A:313[B]:MET:CE	1:A:486:PHE:CE1	2.49	0.94
1:A:313[B]:MET:CE	1:A:486:PHE:HE1	1.81	0.93
1:A:502:TYR:CD1	1:A:506:PHE:HB2	2.04	0.92
1:A:271:GLU:OE1	5:A:701:HOH:O	1.89	0.90
1:A:502:TYR:CZ	1:A:506:PHE:CD1	2.61	0.89
1:A:502:TYR:CD1	1:A:506:PHE:CD1	2.60	0.88
1:A:476:CYS:SG	5:A:780:HOH:O	2.29	0.88
1:A:499:PRO:HB2	1:A:500:PRO:HD3	1.56	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:492:ILE:O	1:A:492:ILE:HD13	1.79	0.82
1:A:502:TYR:CD1	1:A:506:PHE:HD1	1.98	0.79
1:A:496:ALA:HB2	1:A:502:TYR:CD2	2.18	0.78
1:A:502:TYR:CZ	1:A:506:PHE:CG	2.75	0.75
1:A:313[B]:MET:HE3	1:A:486:PHE:CZ	2.23	0.74
1:A:494:VAL:CB	1:A:502:TYR:HE2	2.01	0.72
1:A:313[B]:MET:HE2	1:A:486:PHE:HE1	1.53	0.71
1:A:505:LEU:HD11	3:A:602:MJE:C4	2.22	0.70
1:A:313[B]:MET:HE2	1:A:486:PHE:CE1	2.24	0.69
1:A:494:VAL:HB	1:A:502:TYR:CZ	2.27	0.69
1:A:313[B]:MET:HE3	1:A:486:PHE:HE1	1.43	0.63
1:A:494:VAL:HB	1:A:502:TYR:OH	1.99	0.63
1:A:502:TYR:HE1	1:A:506:PHE:HB3	1.60	0.62
1:A:492:ILE:HD13	1:A:492:ILE:C	2.29	0.57
1:A:267:LEU:HD22	1:A:267:LEU:N	2.21	0.56
1:A:505:LEU:CD1	3:A:602:MJE:C4	2.83	0.56
1:A:268:THR:HG23	5:A:701:HOH:O	2.05	0.55
1:A:494:VAL:CG1	1:A:502:TYR:HE2	2.19	0.55
1:A:484[A]:GLN:HA	1:A:484[A]:GLN:OE1	2.06	0.55
1:A:502:TYR:O	1:A:505:LEU:N	2.37	0.54
1:A:505:LEU:HD12	3:A:602:MJE:C6	2.40	0.52
1:A:496:ALA:CB	1:A:502:TYR:CD2	2.90	0.50
1:A:507:SER:OXT	1:A:507:SER:OG	2.15	0.49
1:A:293:LEU:O	1:A:296:ARG:HG2	2.13	0.48
1:A:505:LEU:CD1	3:A:602:MJE:N2	2.76	0.48
1:A:358:MET:HB3	1:A:476:CYS:SG	2.54	0.48
1:A:499:PRO:HB2	1:A:500:PRO:CD	2.35	0.47
1:A:505:LEU:HD12	3:A:602:MJE:N1	2.31	0.45
1:A:325[B]:THR:HG23	3:A:602:MJE:C17	2.47	0.45
1:A:505:LEU:HD11	3:A:602:MJE:N2	2.33	0.43

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	247/263 (94%)	242 (98%)	5 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	224/234 (96%)	216 (96%)	8 (4%)	31	14

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	313[A]	MET
1	A	313[B]	MET
1	A	404	SER
1	A	492	ILE
1	A	493	VAL
1	A	499	PRO
1	A	504	GLU
1	A	505	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	275	GLN
1	A	286	GLN
1	A	441	GLN
1	A	452	HIS
1	A	453	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	MJE	A	602	-	31,31,31	0.73	0	40,46,46	1.25	3 (7%)
2	HC3	A	601	-	32,32,32	0.38	0	51,51,51	0.48	0
4	GOL	A	603	-	5,5,5	0.14	0	5,5,5	0.19	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
3	MJE	A	602	-	-	0/16/16/16	0/4/4/4
2	HC3	A	601	-	-	0/11/69/69	0/4/4/4
4	GOL	A	603	-	-	2/4/4/4	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	MJE	C4-N2-N1	-5.28	109.50	112.92
3	A	602	MJE	C6-N1-N2	2.90	107.06	103.97
3	A	602	MJE	C14-N2-N1	2.42	124.95	119.78

There are no chirality outliers.

All (2) torsion outliers are listed below:

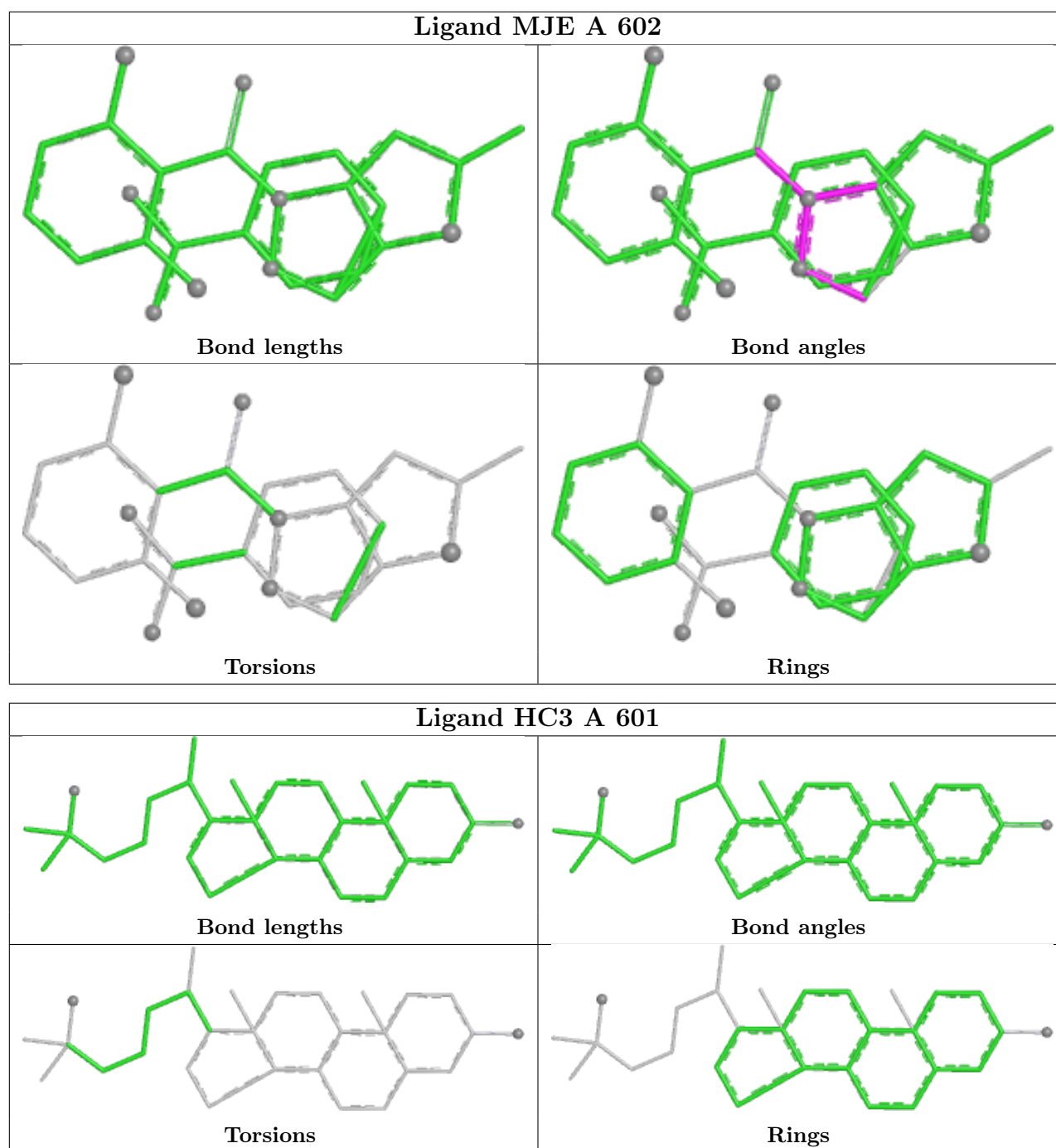
Mol	Chain	Res	Type	Atoms
4	A	603	GOL	C1-C2-C3-O3
4	A	603	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	MJE	7	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, 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	241/263 (91%)	0.58	22 (9%) 15 16	21, 50, 82, 142	8 (3%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	267	LEU	8.8
1	A	506	PHE	6.9
1	A	314	TRP	4.8
1	A	505	LEU	4.6
1	A	490	HIS	4.3
1	A	502	TYR	4.2
1	A	507	SER	3.8
1	A	322[A]	HIS	3.6
1	A	296	ARG	3.3
1	A	489	LEU	3.2
1	A	492	ILE	3.1
1	A	275	GLN	3.1
1	A	313[A]	MET	2.8
1	A	282	ARG	2.8
1	A	494	VAL	2.8
1	A	493	VAL	2.6
1	A	408[A]	SER	2.4
1	A	437	ARG	2.2
1	A	504	GLU	2.1
1	A	476	CYS	2.1
1	A	497	ALA	2.1
1	A	339	SER	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 [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

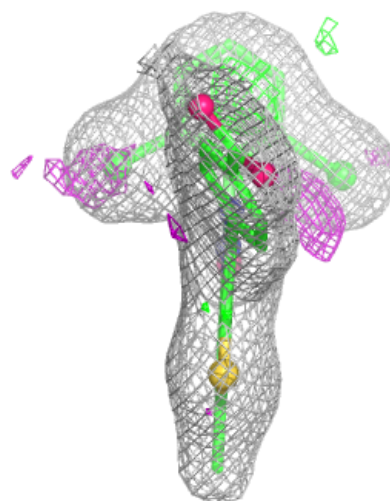
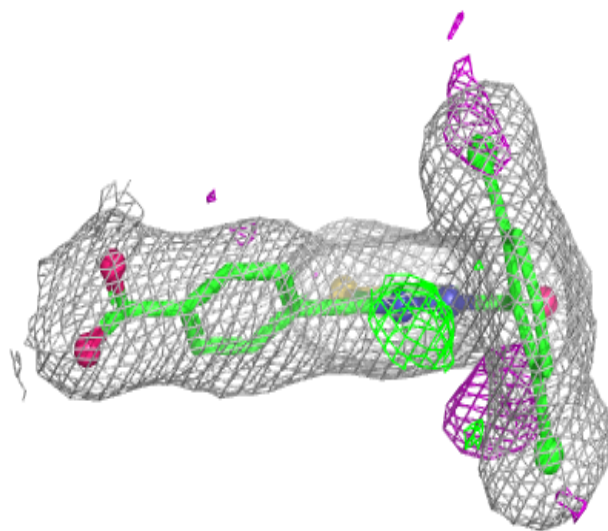
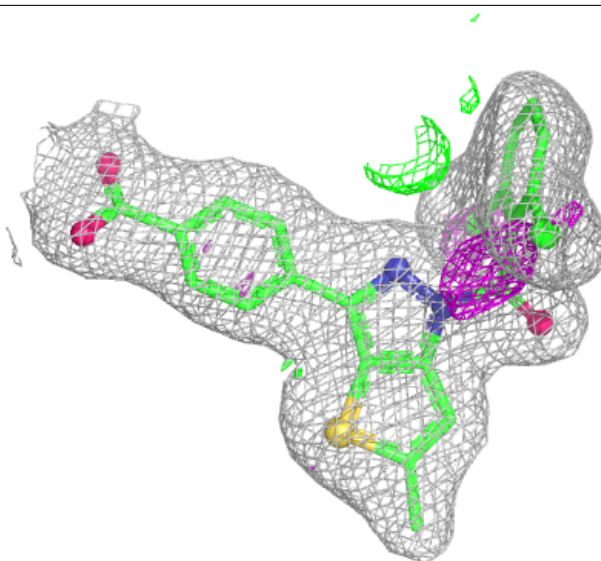
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

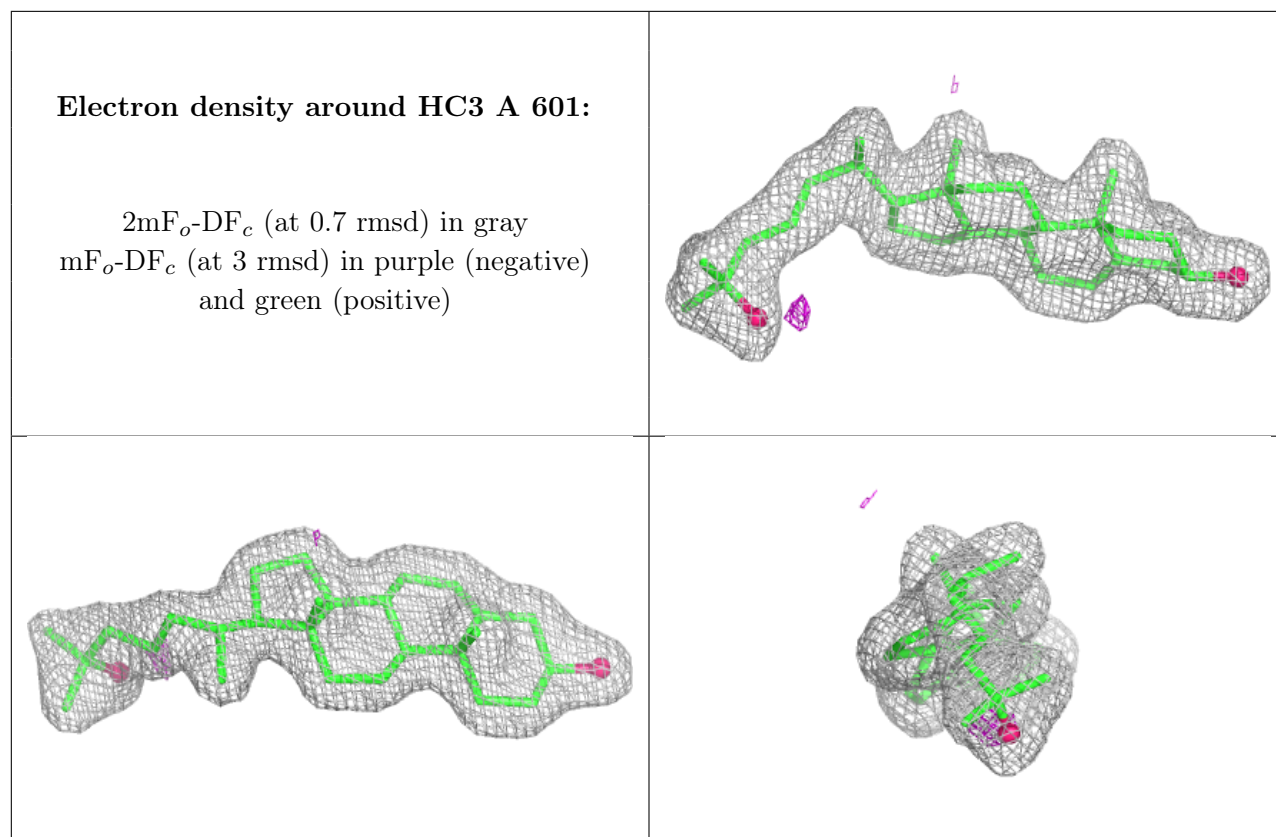
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	GOL	A	603	6/6	0.89	0.12	53,57,59,61	0
3	MJE	A	602	28/28	0.96	0.08	37,44,53,55	0
2	HC3	A	601	29/29	0.97	0.07	30,35,42,49	0

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 MJE 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)





6.5 Other polymers [i](#)

There are no such residues in this entry.