



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 27, 2023 – 01:48 PM JST

PDB ID : 7Y0E  
Title : Crystal structure of TMPRSS2 in complex with Camostat  
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Deposited on : 2022-06-04  
Resolution : 2.39 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 2.36  
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) : 2.36

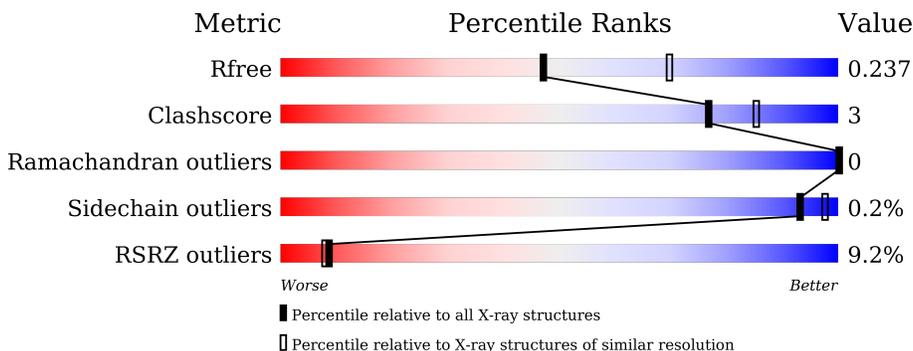
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.39 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	146	 11% 81% 8% 11%
1	B	146	 26% 86% 10%
2	C	249	 2% 91% 5%
2	D	249	 4% 88% 8%

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 5796 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 Transmembrane protease serine 2 catalytic chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	130	966	588	174	192	12	0	0	0
1	B	131	974	597	174	190	13	0	0	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	250	ASP	SER	engineered mutation	UNP O15393
A	251	ASP	SER	engineered mutation	UNP O15393
A	252	ASP	ARG	engineered mutation	UNP O15393
A	253	ASP	GLN	engineered mutation	UNP O15393
A	254	LYS	SER	engineered mutation	UNP O15393
B	250	ASP	SER	engineered mutation	UNP O15393
B	251	ASP	SER	engineered mutation	UNP O15393
B	252	ASP	ARG	engineered mutation	UNP O15393
B	253	ASP	GLN	engineered mutation	UNP O15393
B	254	LYS	SER	engineered mutation	UNP O15393

- Molecule 2 is a protein called Transmembrane protease serine 2 catalytic chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	240	1829	1175	308	330	16	0	1	0
2	D	239	1813	1162	306	329	16	0	1	0

There are 24 discrepancies between the modelled and reference sequences:

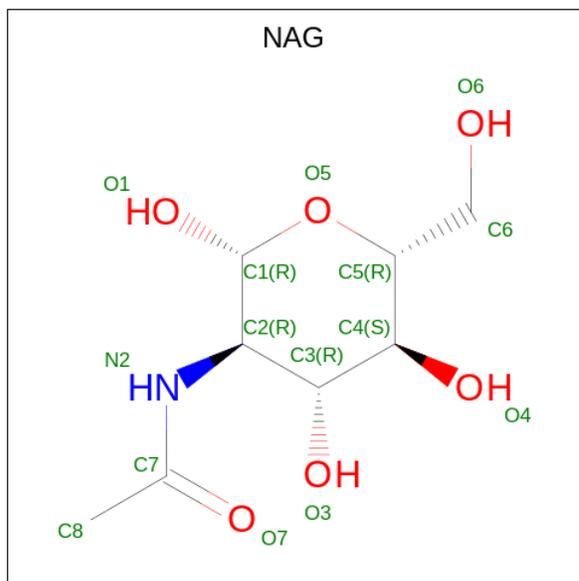
Chain	Residue	Modelled	Actual	Comment	Reference
C	493	GLU	-	expression tag	UNP O15393
C	494	PHE	-	expression tag	UNP O15393

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Chain	Residue	Modelled	Actual	Comment	Reference
C	495	VAL	-	expression tag	UNP O15393
C	496	GLU	-	expression tag	UNP O15393
C	497	HIS	-	expression tag	UNP O15393
C	498	HIS	-	expression tag	UNP O15393
C	499	HIS	-	expression tag	UNP O15393
C	500	HIS	-	expression tag	UNP O15393
C	501	HIS	-	expression tag	UNP O15393
C	502	HIS	-	expression tag	UNP O15393
C	503	HIS	-	expression tag	UNP O15393
C	504	HIS	-	expression tag	UNP O15393
D	493	GLU	-	expression tag	UNP O15393
D	494	PHE	-	expression tag	UNP O15393
D	495	VAL	-	expression tag	UNP O15393
D	496	GLU	-	expression tag	UNP O15393
D	497	HIS	-	expression tag	UNP O15393
D	498	HIS	-	expression tag	UNP O15393
D	499	HIS	-	expression tag	UNP O15393
D	500	HIS	-	expression tag	UNP O15393
D	501	HIS	-	expression tag	UNP O15393
D	502	HIS	-	expression tag	UNP O15393
D	503	HIS	-	expression tag	UNP O15393
D	504	HIS	-	expression tag	UNP O15393

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).

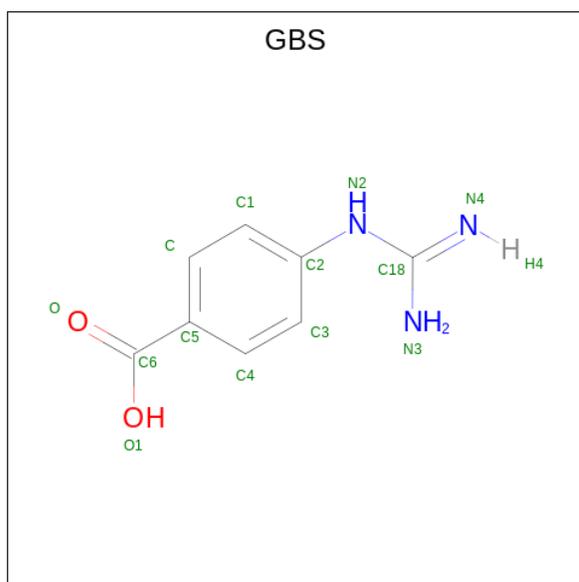


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

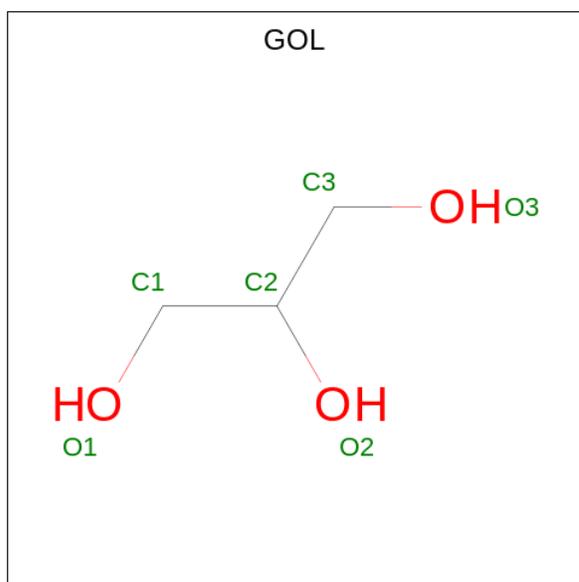
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ca	0	0
			1	1		
4	B	1	Total	Ca	0	0
			1	1		

- Molecule 5 is 4-carbamimidamidobenzoic acid (three-letter code: GBS) (formula: C<sub>8</sub>H<sub>9</sub>N<sub>3</sub>O<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	C	1	Total	C	N	O	0	0
			12	8	3	1		
5	D	1	Total	C	N	O	0	0
			12	8	3	1		

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

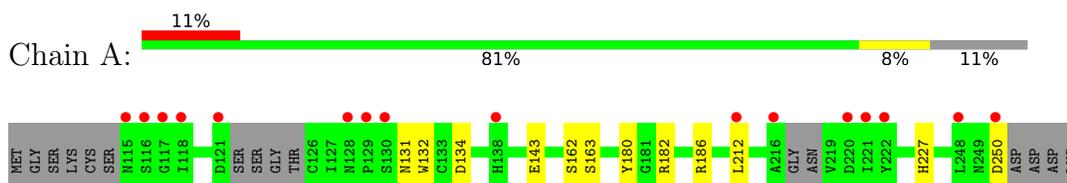
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	27	Total O 27 27	0	0
7	C	72	Total O 72 72	0	0
7	B	7	Total O 7 7	0	0
7	D	48	Total O 48 48	0	0

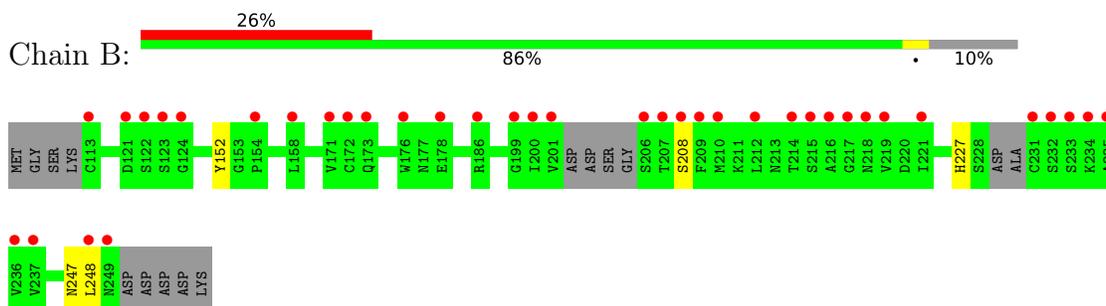
### 3 Residue-property plots [i](#)

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.

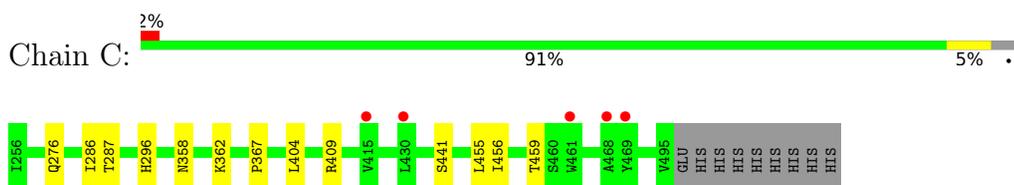
- Molecule 1: Transmembrane protease serine 2 catalytic chain



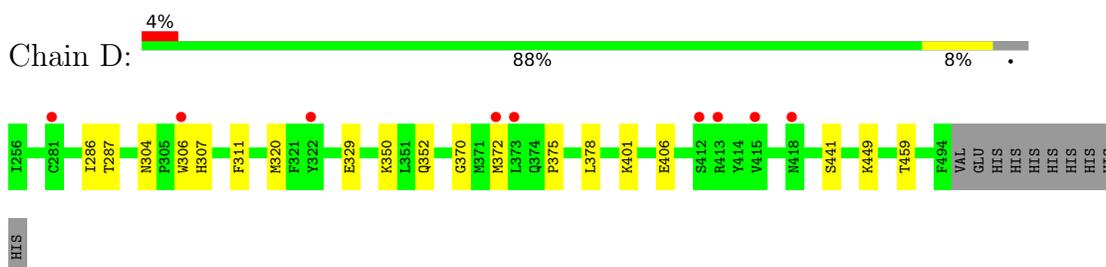
- Molecule 1: Transmembrane protease serine 2 catalytic chain



- Molecule 2: Transmembrane protease serine 2 catalytic chain



- Molecule 2: Transmembrane protease serine 2 catalytic chain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	49.16Å 93.54Å 92.35Å 90.00° 100.08° 90.00°	Depositor
Resolution (Å)	45.46 – 2.39 45.46 – 2.39	Depositor EDS
% Data completeness (in resolution range)	99.5 (45.46-2.39) 99.5 (45.46-2.39)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.03 (at 2.39Å)	Xtrriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, $R_{free}$	0.195 , 0.239 0.198 , 0.237	Depositor DCC
$R_{free}$ test set	1998 reflections (6.13%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	42.0	Xtrriage
Anisotropy	0.189	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 45.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5796	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	48.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.93% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, GBS, NAG, 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	0.32	0/985	0.52	0/1335
1	B	0.27	0/994	0.49	0/1346
2	C	0.27	0/1886	0.47	0/2579
2	D	0.32	0/1869	0.52	0/2558
All	All	0.30	0/5734	0.50	0/7818

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	966	0	834	8	0
1	B	974	0	858	3	0
2	C	1829	0	1742	9	0
2	D	1813	0	1720	13	0
3	A	14	0	13	0	0
3	B	14	0	13	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	C	12	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	12	0	0	0	0
6	D	6	0	8	0	0
7	A	27	0	0	0	0
7	B	7	0	0	0	0
7	C	72	0	0	1	0
7	D	48	0	0	0	0
All	All	5796	0	5188	32	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 3.

All (32) 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)
2:C:409:ARG:HH11	2:C:409:ARG:HG3	1.36	0.89
2:C:409:ARG:HG3	2:C:409:ARG:NH1	2.04	0.68
2:C:276:GLN:NE2	7:C:701:HOH:O	2.30	0.64
2:D:378:LEU:HD11	2:D:401:LYS:HD2	1.85	0.58
1:A:134:ASP:OD1	1:A:134:ASP:N	2.41	0.53
1:A:143:GLU:OE2	1:A:143:GLU:N	2.35	0.52
1:B:247:ASN:OD1	1:B:248:LEU:N	2.40	0.51
2:C:286:ILE:HG13	2:C:287:THR:HG23	1.92	0.51
1:B:208:SER:HB3	1:B:227:HIS:CE1	2.46	0.51
2:D:286:ILE:HG13	2:D:287:THR:HG23	1.95	0.48
1:A:182:ARG:O	1:A:186:ARG:HG3	2.15	0.47
1:B:152:TYR:CZ	2:D:370:GLY:HA3	2.50	0.47
2:D:375:PRO:HB3	2:D:406:GLU:OE1	2.15	0.47
2:C:367:PRO:HD3	2:C:455:LEU:O	2.15	0.46
2:D:329:GLU:N	2:D:350:LYS:O	2.42	0.46
2:D:329:GLU:HB2	2:D:352:GLN:HG2	1.98	0.45
2:C:404:LEU:HD21	2:C:456:ILE:HD11	1.99	0.45
2:D:304:ASN:HB2	2:D:307:HIS:CD2	2.52	0.45
2:C:358:ASN:O	2:C:362:LYS:HG3	2.16	0.45
2:C:441:SER:HA	2:C:459:THR:HB	1.99	0.45
2:C:296:HIS:CE1	2:C:441:SER:HB3	2.52	0.45
1:A:162:SER:HB2	1:A:227:HIS:HE1	1.83	0.44
1:A:131:ASN:HA	1:A:134:ASP:OD2	2.18	0.43
2:D:441:SER:HA	2:D:459:THR:HB	2.00	0.43
2:D:306:TRP:NE1	2:D:307:HIS:CE1	2.87	0.43
2:D:372:MET:O	2:D:449:LYS:NZ	2.50	0.43
2:D:304:ASN:HB2	2:D:307:HIS:NE2	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:304:ASN:O	2:D:307:HIS:HD2	2.01	0.42
1:A:180:TYR:CE2	1:A:212:LEU:HD21	2.53	0.42
2:D:311:PHE:CG	2:D:320:MET:HG2	2.54	0.42
1:A:132:TRP:O	1:A:163:SER:HB3	2.19	0.42
1:A:162:SER:CB	1:A:227:HIS:HE1	2.34	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	124/146 (85%)	122 (98%)	2 (2%)	0	100	100
1	B	125/146 (86%)	120 (96%)	5 (4%)	0	100	100
2	C	239/249 (96%)	231 (97%)	8 (3%)	0	100	100
2	D	238/249 (96%)	231 (97%)	7 (3%)	0	100	100
All	All	726/790 (92%)	704 (97%)	22 (3%)	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	101/127 (80%)	100 (99%)	1 (1%)	76	88

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	104/127 (82%)	104 (100%)	0	100	100
2	C	191/210 (91%)	191 (100%)	0	100	100
2	D	189/210 (90%)	189 (100%)	0	100	100
All	All	585/674 (87%)	584 (100%)	1 (0%)	93	97

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	250	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	131	ASN
1	A	249	ASN
1	B	173	GLN
2	D	307	HIS
2	D	398	ASN
2	D	487	GLN

### 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](#)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	GBS	C	601	2	12,12,13	0.93	1 (8%)	15,15,17	0.56	0
6	GOL	D	602	-	5,5,5	0.08	0	5,5,5	0.35	0
3	NAG	B	301	1	14,14,15	0.40	0	17,19,21	0.36	0
5	GBS	D	601	2	12,12,13	0.81	0	15,15,17	0.61	0
3	NAG	A	301	1	14,14,15	0.39	0	17,19,21	0.95	1 (5%)

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	GBS	C	601	2	-	2/6/6/8	0/1/1/1
6	GOL	D	602	-	-	2/4/4/4	-
3	NAG	B	301	1	-	2/6/23/26	0/1/1/1
5	GBS	D	601	2	-	3/6/6/8	0/1/1/1
3	NAG	A	301	1	-	3/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	601	GBS	C18-N2	-2.07	1.35	1.37

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	301	NAG	O5-C1-C2	-2.96	106.61	111.29

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	301	NAG	C3-C2-N2-C7
6	D	602	GOL	O1-C1-C2-C3

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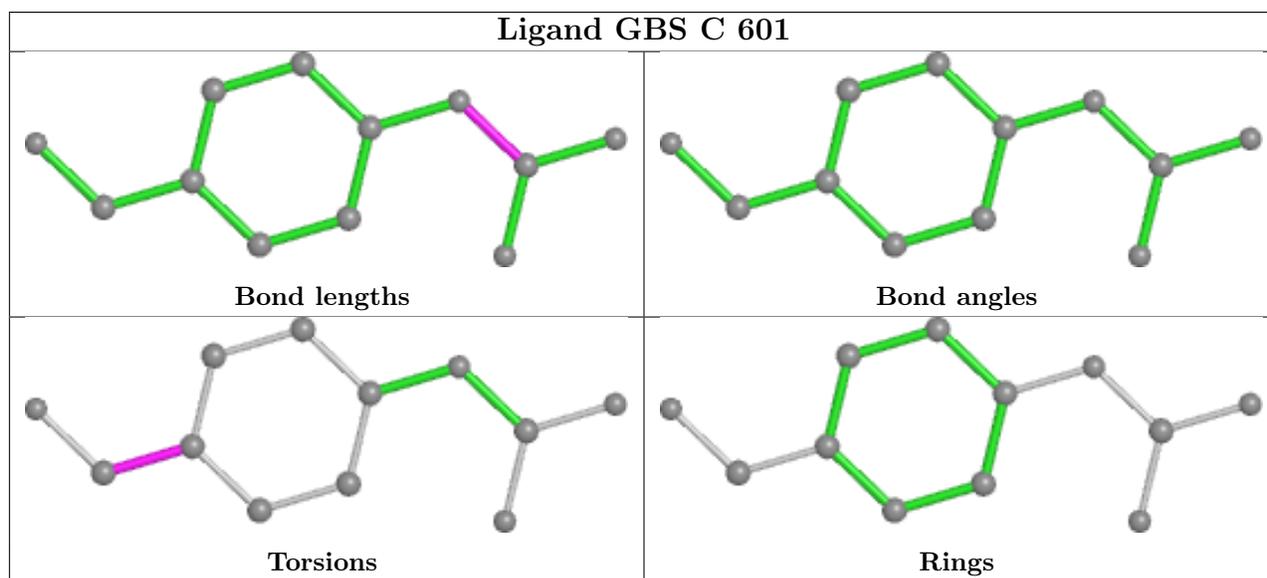
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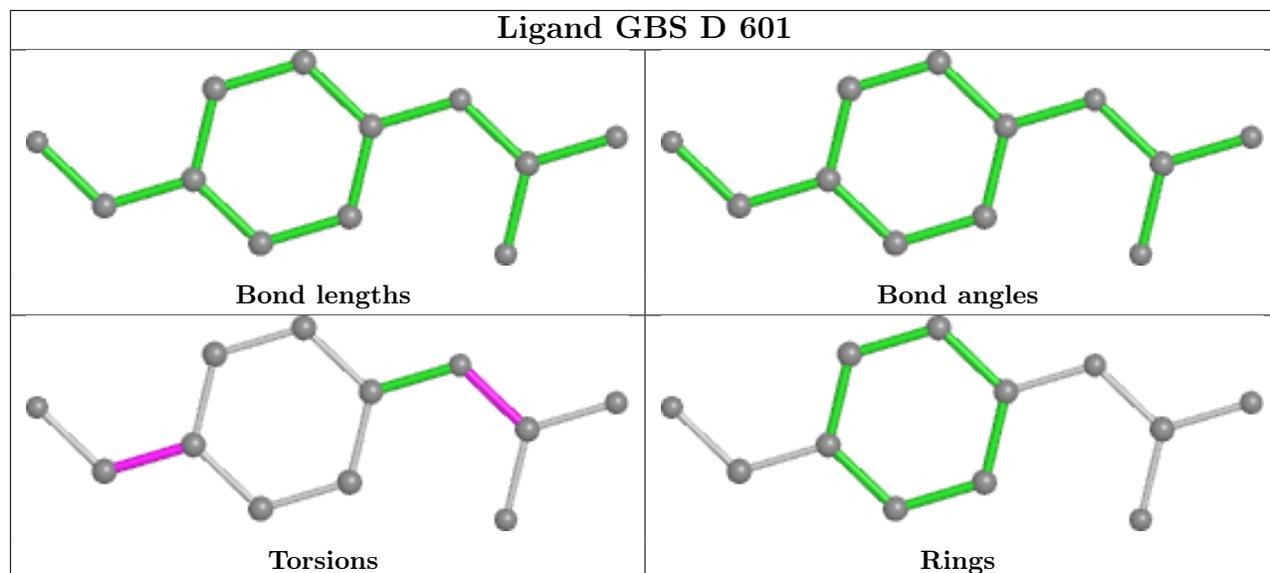
Mol	Chain	Res	Type	Atoms
3	A	301	NAG	C8-C7-N2-C2
3	A	301	NAG	O7-C7-N2-C2
3	B	301	NAG	C8-C7-N2-C2
3	B	301	NAG	O7-C7-N2-C2
5	D	601	GBS	C-C5-C6-O
6	D	602	GOL	O1-C1-C2-O2
5	D	601	GBS	C4-C5-C6-O
5	C	601	GBS	C-C5-C6-O
5	C	601	GBS	C4-C5-C6-O
5	D	601	GBS	N3-C18-N2-C2

There are no ring outliers.

No monomer is involved in short contacts.

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





## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

## 6 Fit of model and data [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, 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	130/146 (89%)	0.94	16 (12%) <b>4</b> <b>3</b>	33, 48, 92, 102	0
1	B	131/146 (89%)	1.43	38 (29%) <b>0</b> <b>0</b>	39, 60, 93, 103	0
2	C	240/249 (96%)	0.46	5 (2%) 63 61	30, 40, 55, 79	0
2	D	239/249 (95%)	0.56	9 (3%) 40 39	33, 44, 64, 76	0
All	All	740/790 (93%)	0.75	68 (9%) <b>9</b> <b>8</b>	30, 45, 80, 103	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	129	PRO	7.1
1	B	207	THR	6.9
1	A	116	SER	6.5
1	B	206	SER	6.0
1	B	201	VAL	5.8
1	A	222	TYR	5.1
1	A	221	ILE	4.9
1	B	217	GLY	4.8
1	A	130	SER	4.7
1	B	232	SER	4.4
1	A	138	HIS	4.3
1	B	234	LYS	4.3
1	A	115	ASN	4.3
1	A	118	ILE	3.8
2	C	430	LEU	3.8
2	D	306	TRP	3.8
1	B	209	PHE	3.7
1	B	208	SER	3.7
1	B	123	SER	3.7
2	D	413	ARG	3.6
1	B	221	ILE	3.6

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	B	231	CYS	3.5
1	B	233	SER	3.5
1	B	121	ASP	3.5
1	B	176	TRP	3.3
1	A	216	ALA	3.3
2	C	415	VAL	3.3
1	B	215	SER	3.3
1	B	248	LEU	3.3
2	D	372	MET	3.3
1	B	124	GLY	3.2
1	B	171	VAL	3.2
1	A	220	ASP	3.1
1	B	200	ILE	3.1
2	D	415	VAL	3.0
1	B	235	ALA	3.0
1	B	210	MET	2.8
1	B	178	GLU	2.8
1	A	117	GLY	2.7
2	D	418	ASN	2.7
2	C	469	TYR	2.7
1	B	172	CYS	2.6
1	B	212	LEU	2.6
1	B	173	GLN	2.6
1	B	154	PRO	2.6
1	B	218	ASN	2.6
1	B	113	CYS	2.5
1	B	158	LEU	2.5
1	B	219	VAL	2.5
2	D	412	SER	2.5
1	B	249	ASN	2.5
1	A	212	LEU	2.4
1	A	121	ASP	2.4
2	D	281[A]	CYS	2.3
1	B	122	SER	2.3
1	A	128	ASN	2.3
1	B	214	THR	2.2
2	C	468	ALA	2.2
2	C	461	TRP	2.2
1	A	248	LEU	2.2
2	D	373	LEU	2.2
1	B	216	ALA	2.1
1	A	250	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	199	GLY	2.1
2	D	322	TYR	2.1
1	B	186	ARG	2.1
1	B	237	VAL	2.1
1	B	236	VAL	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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

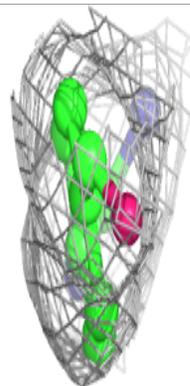
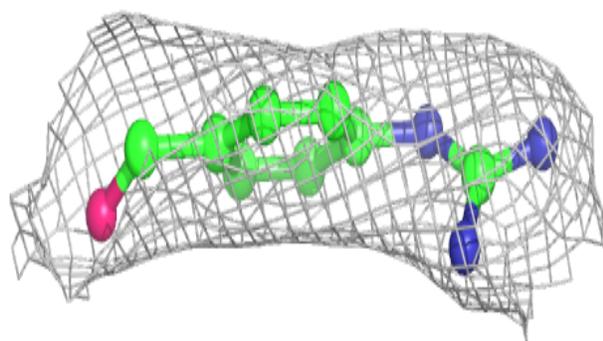
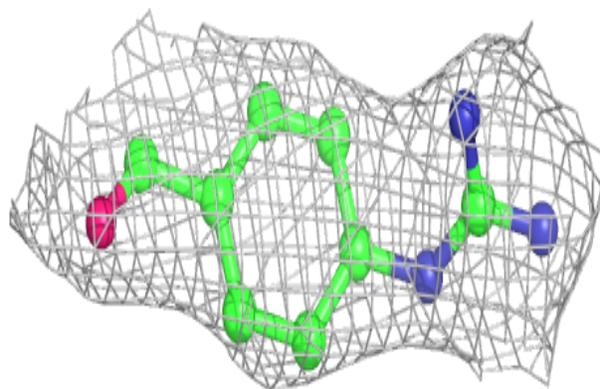
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
3	NAG	A	301	14/15	0.58	0.36	72,92,98,100	0
3	NAG	B	301	14/15	0.77	0.23	83,90,93,94	0
4	CA	A	302	1/1	0.81	0.07	66,66,66,66	0
6	GOL	D	602	6/6	0.84	0.23	41,47,52,54	0
5	GBS	D	601	12/13	0.95	0.18	41,45,48,50	0
5	GBS	C	601	12/13	0.96	0.20	35,42,49,49	0
4	CA	B	302	1/1	0.98	0.12	46,46,46,46	0

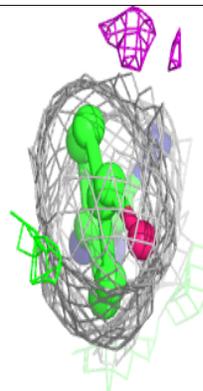
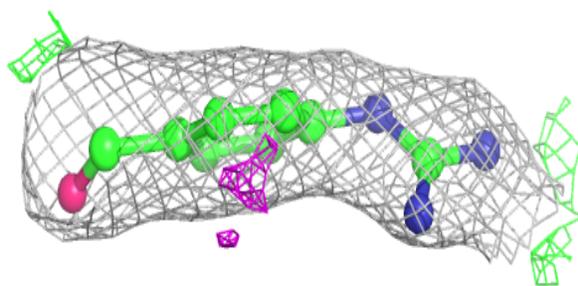
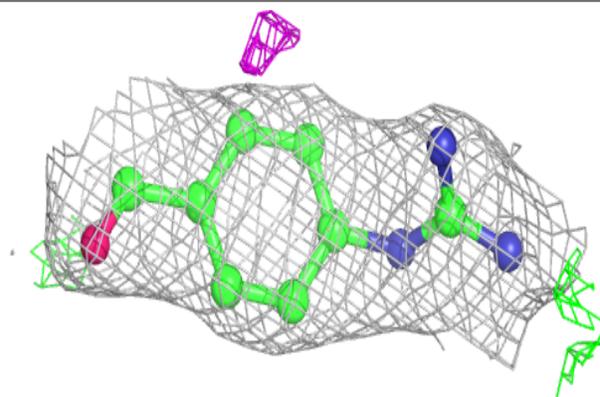
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around GBS D 601:**

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

**Electron density around GBS C 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.