



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 15, 2023 – 07:58 AM EDT

PDB ID : 7UVD  
Title : Sco GlgEI-V279S in complex with cyclohexyl carbasugar  
Authors : Jayasinghe, T.D.; Ronning, D.R.  
Deposited on : 2022-04-30  
Resolution : 2.73 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

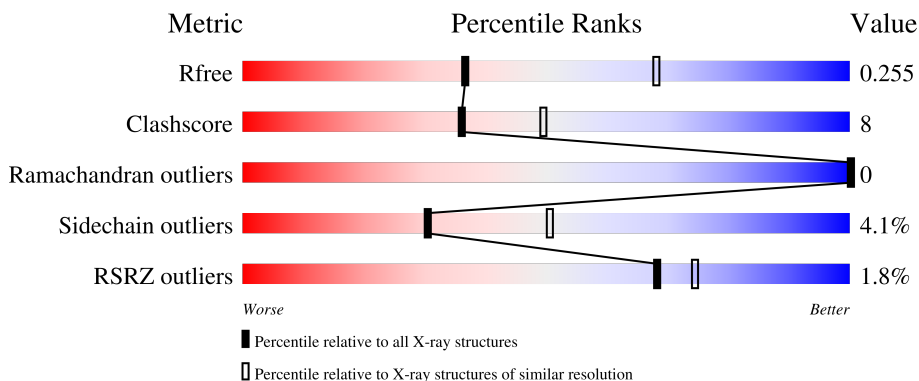
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.73 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	1271 (2.76-2.72)
Clashscore	141614	1322 (2.76-2.72)
Ramachandran outliers	138981	1297 (2.76-2.72)
Sidechain outliers	138945	1298 (2.76-2.72)
RSRZ outliers	127900	1243 (2.76-2.72)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	683	 2% 73% 21% 5%
1	B	683	 % 77% 18% 5%

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10670 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 Alpha-1,4-glucan:maltose-1-phosphate maltosyltransferase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	649	5152	3252	939	949	12	0	2	0
1	A	649	5196	3275	955	956	10	1	5	0

There are 18 discrepancies between the modelled and reference sequences:

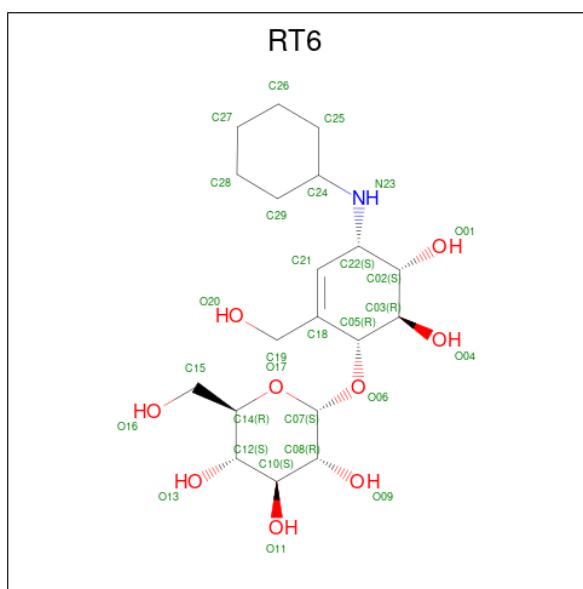
Chain	Residue	Modelled	Actual	Comment	Reference
B	279	SER	VAL	engineered mutation	UNP Q9L1K2
B	676	ALA	-	expression tag	UNP Q9L1K2
B	677	LEU	-	expression tag	UNP Q9L1K2
B	678	HIS	-	expression tag	UNP Q9L1K2
B	679	HIS	-	expression tag	UNP Q9L1K2
B	680	HIS	-	expression tag	UNP Q9L1K2
B	681	HIS	-	expression tag	UNP Q9L1K2
B	682	HIS	-	expression tag	UNP Q9L1K2
B	683	HIS	-	expression tag	UNP Q9L1K2
A	279	SER	VAL	engineered mutation	UNP Q9L1K2
A	676	ALA	-	expression tag	UNP Q9L1K2
A	677	LEU	-	expression tag	UNP Q9L1K2
A	678	HIS	-	expression tag	UNP Q9L1K2
A	679	HIS	-	expression tag	UNP Q9L1K2
A	680	HIS	-	expression tag	UNP Q9L1K2
A	681	HIS	-	expression tag	UNP Q9L1K2
A	682	HIS	-	expression tag	UNP Q9L1K2
A	683	HIS	-	expression tag	UNP Q9L1K2

- Molecule 2 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>).



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

- Molecule 3 is (1R,4S,5S,6R)-4-(cyclohexylamino)-5,6-dihydroxy-2-(hydroxymethyl)cyclohex-2-en-1-yl alpha-D-glucopyranoside (three-letter code: RT6) (formula: C<sub>19</sub>H<sub>33</sub>NO<sub>9</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			29	19	1	9		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	29	19	1	9	0	0

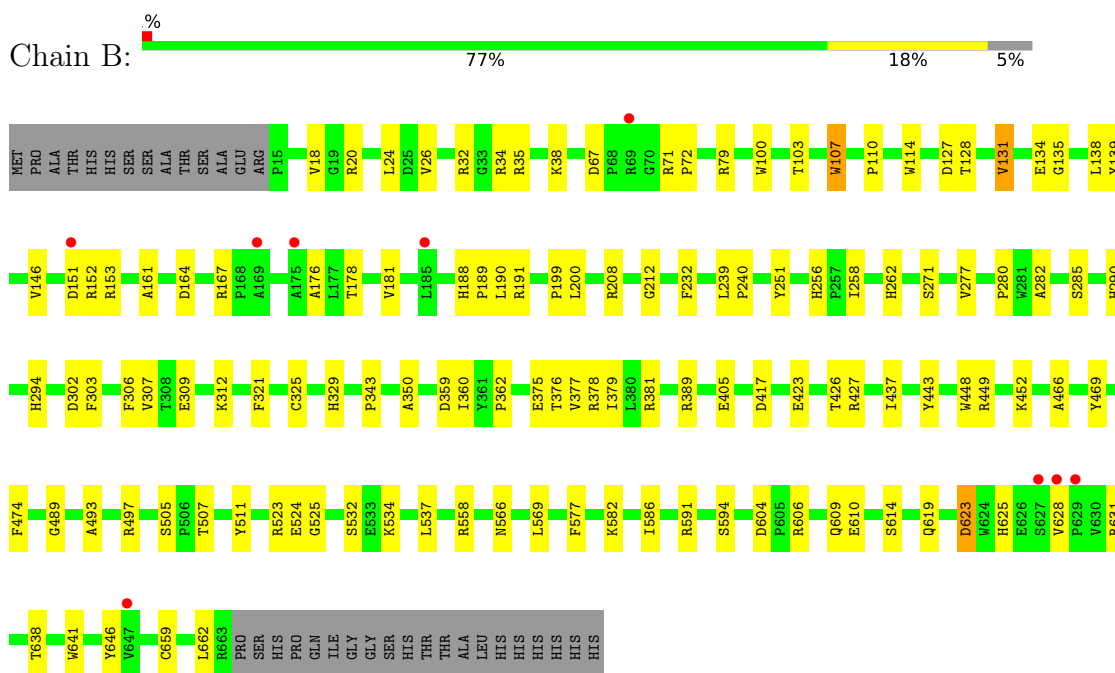
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	128	Total	O	0	0
			128	128		
4	A	122	Total	O	0	0
			122	122		

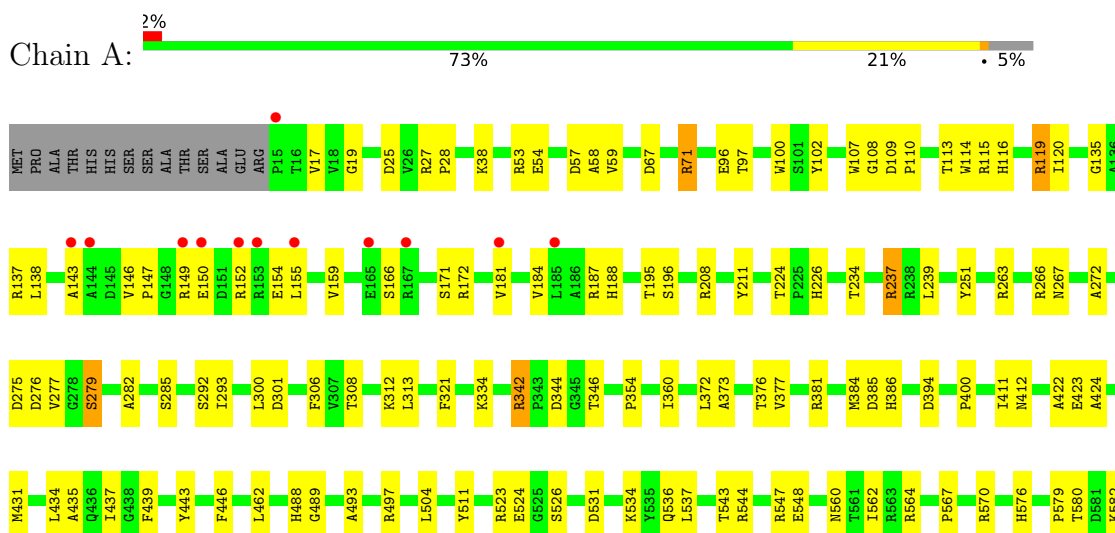
### 3 Residue-property plots [i](#)

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: Alpha-1,4-glucan:maltose-1-phosphate maltosyltransferase 1



- Molecule 1: Alpha-1,4-glucan:maltose-1-phosphate maltosyltransferase 1





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	112.44Å 112.44Å 310.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.40 – 2.73 49.64 – 2.64	Depositor EDS
% Data completeness (in resolution range)	96.3 (49.40-2.73) 86.6 (49.64-2.64)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.76 (at 2.65Å)	Xtrriage
Refinement program	PHENIX 1.19.1_4122	Depositor
R, $R_{free}$	0.190 , 0.259 0.190 , 0.255	Depositor DCC
$R_{free}$ test set	2000 reflections (3.47%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	51.5	Xtrriage
Anisotropy	0.244	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 22.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	10670	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	57.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.60% 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: RT6, PEG

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	0.47	0/5348	0.69	0/7309
1	B	0.46	0/5304	0.68	0/7251
All	All	0.46	0/10652	0.68	0/14560

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

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	5196	0	5037	89	0
1	B	5152	0	4992	77	0
2	A	7	0	10	0	0
2	B	7	0	10	0	0
3	A	29	0	0	1	0
3	B	29	0	0	1	0
4	A	122	0	0	3	0
4	B	128	0	0	6	0
All	All	10670	0	10049	163	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 (163) 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)
3:A:701:RT6:C03	3:A:701:RT6:C05	1.74	1.62
1:A:580:THR:HG22	1:A:613:VAL:HG23	1.67	0.77
1:B:309:GLU:OE2	1:B:312:LYS:NZ	2.17	0.76
1:A:146:VAL:HG23	1:A:152:ARG:HG2	1.67	0.75
1:A:154:GLU:HG2	1:A:184:VAL:HG21	1.70	0.73
1:B:146:VAL:HG13	1:B:152:ARG:HG2	1.71	0.72
1:A:25[B]:ASP:OD2	1:A:27[B]:ARG:NH1	2.22	0.72
1:A:58:ALA:H	1:A:113:THR:HG22	1.56	0.70
1:B:161:ALA:HB1	1:B:167:ARG:HH12	1.58	0.69
1:B:79:ARG:NH2	4:B:805:HOH:O	2.26	0.68
1:A:623:ASP:N	1:A:626:GLU:OE1	2.24	0.68
1:A:544:ARG:NH1	1:A:548:GLU:OE2	2.27	0.68
1:B:423:GLU:OE2	3:B:702:RT6:N23	2.28	0.67
1:B:290:HIS:HB2	1:B:379:ILE:HD13	1.77	0.66
1:A:146:VAL:HG12	1:A:188:HIS:ND1	2.11	0.66
1:B:107:TRP:CD2	1:B:191:ARG:HG2	2.30	0.66
1:B:452:LYS:NZ	4:B:804:HOH:O	2.26	0.65
1:B:375:GLU:OE2	1:B:378:ARG:NH2	2.30	0.64
1:B:110:PRO:HG2	1:B:190:LEU:HD23	1.80	0.64
1:A:109:ASP:O	1:A:113:THR:HG23	1.98	0.63
1:A:275:ASP:N	1:A:275:ASP:OD1	2.33	0.61
1:B:146:VAL:HG23	1:B:188:HIS:CD2	2.36	0.61
1:B:628:VAL:HG23	1:B:641:TRP:HB2	1.81	0.61
1:B:405:GLU:HG3	1:B:437:ILE:HB	1.81	0.60
1:A:372:LEU:O	1:A:376:THR:OG1	2.17	0.60
1:A:38:LYS:NZ	1:A:435:ALA:O	2.32	0.59
1:B:110:PRO:HB2	1:B:138:LEU:HD11	1.84	0.59
1:A:143:ALA:HA	1:A:146:VAL:HG22	1.85	0.58
1:A:489:GLY:HA3	1:A:493:ALA:HB2	1.83	0.58
1:A:543:THR:HG22	1:A:547:ARG:HE	1.69	0.58
1:B:359:ASP:OD1	1:B:359:ASP:N	2.34	0.58
1:B:134:GLU:OE2	1:A:342:ARG:HG2	2.04	0.57
1:A:560:ASN:O	1:A:564:ARG:HG3	2.05	0.56
1:A:28:PRO:O	1:A:102:TYR:OH	2.24	0.56
1:B:610:GLU:OE1	1:B:646:TYR:OH	2.23	0.56
1:A:147:PRO:HD3	1:A:188:HIS:HE1	1.70	0.56
1:A:58:ALA:H	1:A:113:THR:CG2	2.19	0.56
1:B:282:ALA:HA	1:B:360:ILE:HG22	1.88	0.55
1:B:631:ARG:HB2	1:B:662:LEU:HD21	1.88	0.55
1:A:277:VAL:HG11	1:A:537:LEU:HD21	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:452:LYS:HG3	1:B:586:ILE:HD12	1.88	0.55
1:A:116:HIS:O	1:A:120:ILE:HD12	2.07	0.55
1:B:604:ASP:OD2	1:B:609:GLN:NE2	2.40	0.54
1:A:266:ARG:NH2	1:A:536:GLN:OE1	2.40	0.54
1:B:423:GLU:HA	1:B:443:TYR:CD1	2.43	0.54
1:B:164:ASP:OD2	1:B:167:ARG:HD2	2.08	0.54
1:B:489:GLY:HA3	1:B:493:ALA:HB2	1.89	0.54
1:B:631:ARG:NH1	4:B:815:HOH:O	2.41	0.54
1:B:623:ASP:HB3	1:B:625:HIS:CE1	2.43	0.54
1:B:325[B]:CYS:HB2	1:B:329:HIS:HB2	1.89	0.53
1:A:266:ARG:NH1	4:A:808:HOH:O	2.37	0.53
1:A:321:PHE:CZ	1:A:376:THR:HG23	2.43	0.53
1:A:488:HIS:O	1:A:606:ARG:NH2	2.41	0.53
1:A:300:LEU:HD21	1:A:386:HIS:NE2	2.24	0.53
1:B:591:ARG:HD3	4:B:910:HOH:O	2.08	0.52
1:A:239:LEU:HD21	1:A:306:PHE:CZ	2.45	0.52
1:B:474:PHE:CD2	1:B:505:SER:HB2	2.45	0.52
1:A:187:ARG:O	1:A:188:HIS:HD2	1.91	0.52
1:A:321:PHE:HZ	1:A:376:THR:HG23	1.75	0.52
1:B:32:ARG:HD2	4:B:823:HOH:O	2.09	0.51
1:B:146:VAL:HG21	1:B:151:ASP:HB3	1.91	0.51
1:A:377:VAL:O	1:A:381:ARG:HG3	2.11	0.51
1:B:466:ALA:HA	1:B:469:TYR:CE1	2.46	0.51
1:B:38:LYS:NZ	1:B:469:TYR:O	2.41	0.50
1:A:613:VAL:HG12	1:A:645:ASN:HB2	1.94	0.50
1:B:178:THR:OG1	1:B:181:VAL:HG23	2.11	0.50
1:B:303:PHE:O	1:B:307:VAL:HG23	2.12	0.50
1:A:446:PHE:O	1:A:497:ARG:NH2	2.45	0.50
1:A:344:ASP:OD1	1:A:346:THR:OG1	2.23	0.49
1:A:579:PRO:HD2	1:A:614:SER:O	2.12	0.49
1:B:26:VAL:HG12	1:B:200:LEU:HD12	1.95	0.49
1:A:263:ARG:HG2	1:A:272:ALA:HB2	1.94	0.49
1:A:628:VAL:HG13	1:A:641:TRP:C	2.34	0.49
1:B:139:TYR:HE2	1:B:181:VAL:HG11	1.78	0.48
1:B:523:ARG:HG3	1:B:524:GLU:O	2.12	0.48
1:B:577:PHE:HE2	1:B:586:ILE:HD11	1.78	0.48
1:B:18:VAL:HG22	1:B:24:LEU:HD21	1.96	0.48
1:A:276:ASP:OD2	4:A:801:HOH:O	2.19	0.48
1:B:212:GLY:O	1:B:507:THR:HA	2.14	0.48
1:B:524:GLU:HG2	1:B:525:GLY:N	2.29	0.48
1:A:373:ALA:O	1:A:377:VAL:HG22	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:309:GLU:HA	1:B:312:LYS:HE2	1.96	0.47
1:A:293:ILE:HD11	1:A:300:LEU:HB2	1.96	0.47
1:A:544:ARG:NH1	4:A:817:HOH:O	2.47	0.47
1:A:57:ASP:HB3	1:A:113:THR:HG21	1.96	0.47
1:B:377:VAL:O	1:B:381:ARG:HG3	2.15	0.47
1:A:267:ASN:ND2	1:A:531:ASP:OD1	2.40	0.47
1:A:544:ARG:O	1:A:548:GLU:HG3	2.15	0.46
1:B:146:VAL:HG23	1:B:188:HIS:HD2	1.78	0.46
1:A:71:ARG:HE	1:A:71:ARG:HB2	1.63	0.46
1:A:279:SER:HA	1:A:534:LYS:O	2.15	0.46
1:A:19:GLY:HA3	1:A:195:THR:HG21	1.96	0.46
1:A:272:ALA:HB1	1:A:276:ASP:HB2	1.97	0.46
1:A:67:ASP:HA	1:A:100:TRP:CD2	2.50	0.46
1:B:532:SER:OG	1:B:534:LYS:HG3	2.16	0.46
1:B:566:ASN:HB2	1:B:569:LEU:HD12	1.97	0.46
1:B:239:LEU:HD21	1:B:306:PHE:CZ	2.51	0.46
1:A:96:GLU:HG3	1:A:97:THR:O	2.16	0.46
1:A:422:ALA:HB2	1:A:439:PHE:CD1	2.51	0.46
1:A:234:THR:O	1:A:237:ARG:HB2	2.16	0.46
1:A:293:ILE:H	1:A:293:ILE:HD12	1.81	0.46
1:A:308:THR:O	1:A:312:LYS:HG2	2.16	0.45
1:A:434:LEU:O	1:A:437:ILE:HG13	2.16	0.45
1:B:282:ALA:HA	1:B:360:ILE:CG2	2.46	0.45
1:B:321:PHE:CZ	1:B:376:THR:HG23	2.52	0.45
1:A:334:LYS:HA	1:A:334:LYS:HD3	1.70	0.45
1:A:423:GLU:HA	1:A:443:TYR:CD1	2.51	0.45
1:B:103:THR:HG22	1:B:199:PRO:HA	1.98	0.45
1:A:567:PRO:HG3	1:A:570:ARG:NH2	2.32	0.45
1:B:448:TRP:O	1:B:449:ARG:NH1	2.47	0.45
1:B:67:ASP:HA	1:B:100:TRP:CD2	2.52	0.45
1:A:208:ARG:HA	1:A:211:TYR:O	2.17	0.45
1:A:576:HIS:NE2	1:A:619:GLN:HG3	2.32	0.45
1:A:282:ALA:HA	1:A:360:ILE:HG22	1.98	0.45
1:A:54:GLU:OE1	1:A:54:GLU:N	2.41	0.44
1:A:621:GLY:O	1:A:622:LEU:HD23	2.18	0.44
1:A:53:ARG:HG2	1:A:59:VAL:HG13	1.98	0.44
1:B:239:LEU:HB2	1:B:240:PRO:HD3	2.00	0.44
1:B:619:GLN:HA	1:B:619:GLN:OE1	2.17	0.44
1:A:114:TRP:CZ2	1:A:135:GLY:HA3	2.52	0.44
1:A:282:ALA:HA	1:A:360:ILE:CG2	2.47	0.43
1:A:635:THR:OG1	1:A:637:GLU:HG3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:ARG:HE	1:A:53:ARG:HB2	1.70	0.43
1:A:562:ILE:HG12	1:A:633:GLU:HB3	1.99	0.43
1:B:389:ARG:HD2	1:B:417:ASP:OD2	2.18	0.43
1:A:110:PRO:HB2	1:A:138:LEU:HD11	2.00	0.43
1:A:596:THR:HB	1:A:661:VAL:CG2	2.48	0.43
1:B:258:ILE:HG12	1:B:280:PRO:HA	1.99	0.43
1:B:232:PHE:HB2	1:B:302:ASP:HB3	2.00	0.43
1:B:631:ARG:O	1:B:659:CYS:HA	2.19	0.43
1:B:146:VAL:CG2	1:B:151:ASP:HB3	2.48	0.42
1:B:35:ARG:HD2	1:A:17:VAL:HG21	2.01	0.42
1:A:576:HIS:CD2	1:A:619:GLN:HG3	2.54	0.42
1:A:582:LYS:HA	1:A:582:LYS:HD2	1.87	0.42
1:B:350:ALA:HB2	1:B:362:PRO:HD3	2.02	0.42
1:B:189:PRO:O	4:B:801:HOH:O	2.21	0.42
1:A:155:LEU:O	1:A:159:VAL:HG23	2.20	0.42
1:B:71:ARG:HA	1:B:72:PRO:HD3	1.95	0.42
1:B:497:ARG:HD2	1:B:497:ARG:HA	1.88	0.41
1:B:114:TRP:CZ2	1:B:135:GLY:HA3	2.56	0.41
1:B:277:VAL:HG11	1:B:537:LEU:HD21	2.02	0.41
1:A:615:LEU:HD22	1:A:620:LEU:HD11	2.02	0.41
1:B:350:ALA:HB3	1:B:360:ILE:HD11	2.03	0.41
1:B:426:THR:OG1	1:B:427:ARG:N	2.53	0.41
1:B:577:PHE:CE2	1:B:586:ILE:HD11	2.56	0.41
1:B:139:TYR:CZ	1:B:176:ALA:HB1	2.56	0.41
1:B:256:HIS:O	1:B:294:HIS:HB2	2.21	0.41
1:B:262:HIS:O	1:B:359:ASP:HB3	2.21	0.41
1:B:343:PRO:HB3	1:A:137:ARG:NH2	2.36	0.41
1:A:313:LEU:HD23	1:A:313:LEU:HA	1.83	0.41
1:A:462:LEU:HD23	1:A:462:LEU:HA	1.79	0.41
1:A:504:LEU:HA	1:A:504:LEU:HD23	1.82	0.41
1:B:127:ASP:O	1:B:131:VAL:HG13	2.21	0.41
1:A:108:GLY:O	1:A:110:PRO:HD3	2.21	0.41
1:A:628:VAL:O	1:A:628:VAL:HG22	2.21	0.41
1:A:424:ALA:O	1:A:431:MET:HG2	2.22	0.40
1:A:115:ARG:O	1:A:119:ARG:HG2	2.20	0.40
1:A:523:ARG:HG2	1:A:526:SER:HB2	2.02	0.40
1:B:20:ARG:NH1	1:A:400:PRO:HB3	2.37	0.40
1:A:384:MET:HE1	1:A:411:ILE:HD12	2.04	0.40
1:A:412:ASN:HD22	1:A:412:ASN:HA	1.69	0.40
1:A:643:ARG:HE	1:A:643:ARG:HB2	1.64	0.40
1:B:152:ARG:HH21	1:B:153:ARG:HE	1.69	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	653/683 (96%)	630 (96%)	23 (4%)	0	100	100
1	B	649/683 (95%)	636 (98%)	13 (2%)	0	100	100
All	All	1302/1366 (95%)	1266 (97%)	36 (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	543/565 (96%)	514 (95%)	29 (5%)	22	39
1	B	539/565 (95%)	523 (97%)	16 (3%)	41	61
All	All	1082/1130 (96%)	1037 (96%)	45 (4%)	30	49

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	34	ARG
1	B	107	TRP
1	B	128	THR
1	B	131	VAL
1	B	208	ARG

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Mol	Chain	Res	Type
1	B	251	TYR
1	B	271	SER
1	B	285	SER
1	B	511	TYR
1	B	558	ARG
1	B	582	LYS
1	B	594	SER
1	B	606	ARG
1	B	614	SER
1	B	623	ASP
1	B	638	THR
1	A	71	ARG
1	A	107	TRP
1	A	119	ARG
1	A	149	ARG
1	A	150	GLU
1	A	166	SER
1	A	171	SER
1	A	172	ARG
1	A	181	VAL
1	A	196	SER
1	A	224	THR
1	A	226	HIS
1	A	237	ARG
1	A	251	TYR
1	A	279	SER
1	A	285	SER
1	A	292	SER
1	A	301	ASP
1	A	342	ARG
1	A	354	PRO
1	A	385	ASP
1	A	394	ASP
1	A	511	TYR
1	A	524	GLU
1	A	628	VAL
1	A	648	ARG
1	A	653[A]	ARG
1	A	653[B]	ARG
1	A	663	ARG

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

Mol	Chain	Res	Type
1	B	188	HIS
1	A	188	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](#)

4 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
2	PEG	A	702	-	6,6,6	1.37	0	5,5,5	0.90	0
3	RT6	B	702	-	31,31,31	3.91	15 (48%)	32,44,44	2.03	9 (28%)
2	PEG	B	701	-	6,6,6	1.65	2 (33%)	5,5,5	0.55	0
3	RT6	A	701	-	31,31,31	3.91	15 (48%)	32,44,44	2.20	11 (34%)

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	PEG	A	702	-	-	1/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	RT6	B	702	-	-	6/12/60/60	0/3/3/3
2	PEG	B	701	-	-	1/4/4/4	-
3	RT6	A	701	-	-	6/12/60/60	0/3/3/3

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	701	RT6	C03-C05	12.67	1.74	1.53
3	B	702	RT6	C03-C05	11.45	1.72	1.53
3	B	702	RT6	C02-C22	11.08	1.67	1.52
3	A	701	RT6	C02-C22	9.70	1.65	1.52
3	B	702	RT6	C22-C21	7.13	1.60	1.50
3	A	701	RT6	C22-C21	6.19	1.59	1.50
3	A	701	RT6	C19-C18	5.77	1.63	1.50
3	A	701	RT6	C05-C18	5.58	1.56	1.51
3	B	702	RT6	C19-C18	5.37	1.62	1.50
3	B	702	RT6	C05-C18	5.23	1.56	1.51
3	B	702	RT6	C12-C14	4.51	1.62	1.53
3	B	702	RT6	C22-N23	4.30	1.55	1.47
3	A	701	RT6	C12-C14	4.10	1.61	1.53
3	A	701	RT6	C07-C08	3.98	1.63	1.52
3	A	701	RT6	C22-N23	3.79	1.54	1.47
3	B	702	RT6	C24-N23	3.26	1.53	1.48
3	B	702	RT6	C25-C24	3.20	1.59	1.52
3	A	701	RT6	C15-C14	3.06	1.62	1.51
3	B	702	RT6	C07-C08	3.01	1.61	1.52
3	A	701	RT6	C29-C24	2.99	1.59	1.52
3	A	701	RT6	C24-N23	2.95	1.52	1.48
3	A	701	RT6	C25-C24	2.93	1.58	1.52
3	B	702	RT6	C15-C14	2.80	1.61	1.51
3	B	702	RT6	C12-C10	2.73	1.59	1.52
3	A	701	RT6	C12-C10	2.53	1.58	1.52
3	B	702	RT6	C29-C24	2.39	1.57	1.52
3	B	702	RT6	C26-C25	2.23	1.59	1.53
3	A	701	RT6	C10-C08	2.22	1.58	1.52
2	B	701	PEG	C2-C1	2.20	1.61	1.49
3	B	702	RT6	C10-C08	2.19	1.57	1.52
3	A	701	RT6	O17-C07	2.09	1.47	1.41
2	B	701	PEG	O2-C2	2.02	1.50	1.42

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	701	RT6	C07-O17-C14	5.84	125.16	113.69
3	A	701	RT6	C02-C03-C05	5.62	118.86	109.25
3	B	702	RT6	C26-C25-C24	5.20	120.89	111.11
3	B	702	RT6	O01-C02-C22	4.91	118.97	109.12
3	A	701	RT6	O04-C03-C02	-4.32	100.35	110.35
3	B	702	RT6	C27-C26-C25	3.83	119.22	111.42
3	A	701	RT6	O01-C02-C22	3.61	116.35	109.12
3	A	701	RT6	O17-C14-C12	3.12	115.36	109.69
3	B	702	RT6	O20-C19-C18	-3.11	105.05	112.50
3	B	702	RT6	O04-C03-C02	-2.84	103.77	110.35
3	B	702	RT6	C07-O17-C14	2.78	119.14	113.69
3	B	702	RT6	O17-C14-C15	2.75	113.27	106.44
3	A	701	RT6	C15-C14-C12	-2.74	106.59	113.00
3	B	702	RT6	O01-C02-C03	-2.67	104.18	110.35
3	A	701	RT6	C28-C29-C24	2.57	115.95	111.11
3	A	701	RT6	O01-C02-C03	-2.39	104.82	110.35
3	A	701	RT6	O17-C14-C15	2.37	112.33	106.44
3	A	701	RT6	O06-C07-C08	2.25	113.92	108.10
3	B	702	RT6	O04-C03-C05	2.24	114.76	109.74
3	A	701	RT6	O20-C19-C18	-2.13	107.41	112.50

There are no chirality outliers.

All (14) torsion outliers are listed below:

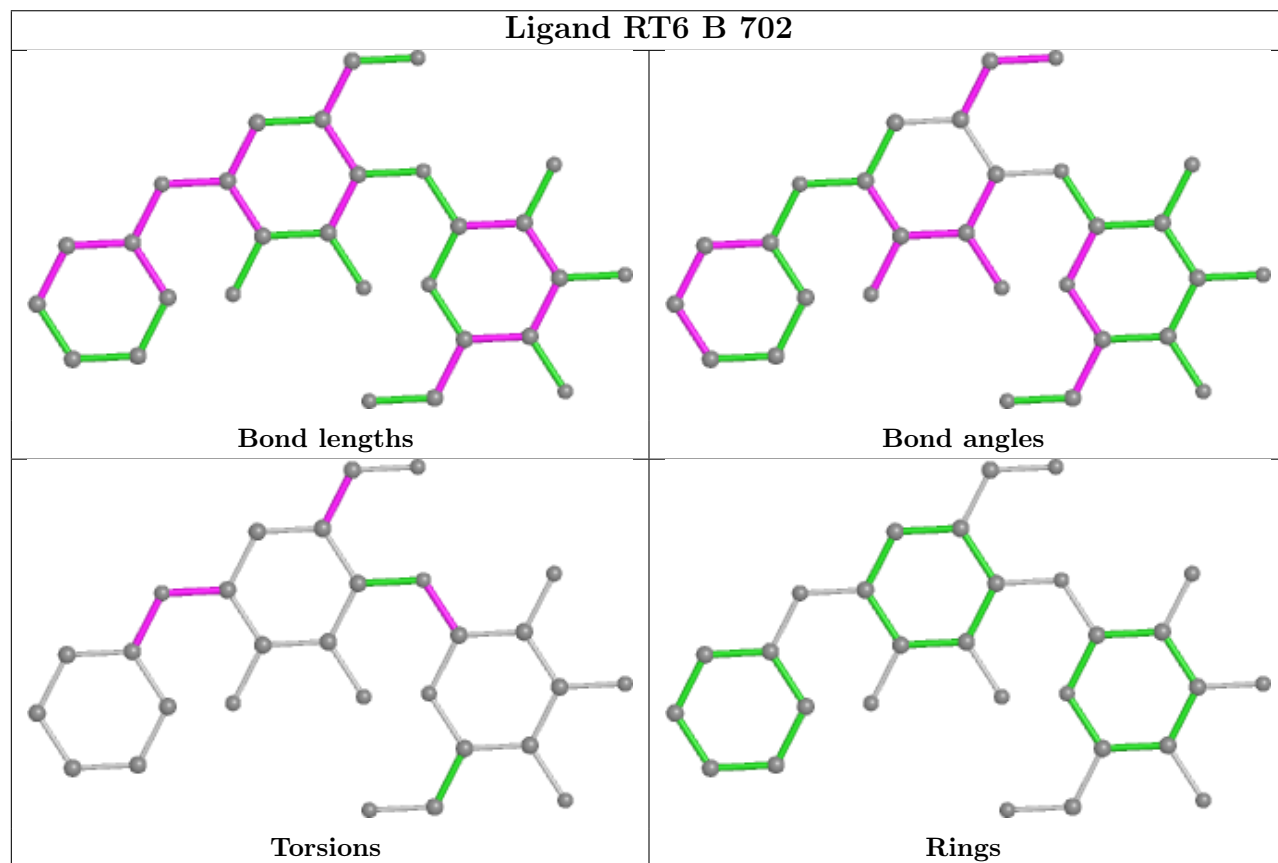
Mol	Chain	Res	Type	Atoms
3	B	702	RT6	C21-C22-N23-C24
3	B	702	RT6	C25-C24-N23-C22
3	B	702	RT6	C29-C24-N23-C22
3	A	701	RT6	C05-C18-C19-O20
3	A	701	RT6	C21-C22-N23-C24
2	B	701	PEG	O2-C3-C4-O4
3	A	701	RT6	O17-C07-O06-C05
3	B	702	RT6	O17-C07-O06-C05
3	A	701	RT6	C25-C24-N23-C22
3	A	701	RT6	C29-C24-N23-C22
2	A	702	PEG	C4-C3-O2-C2
3	A	701	RT6	C08-C07-O06-C05
3	B	702	RT6	C08-C07-O06-C05
3	B	702	RT6	C21-C18-C19-O20

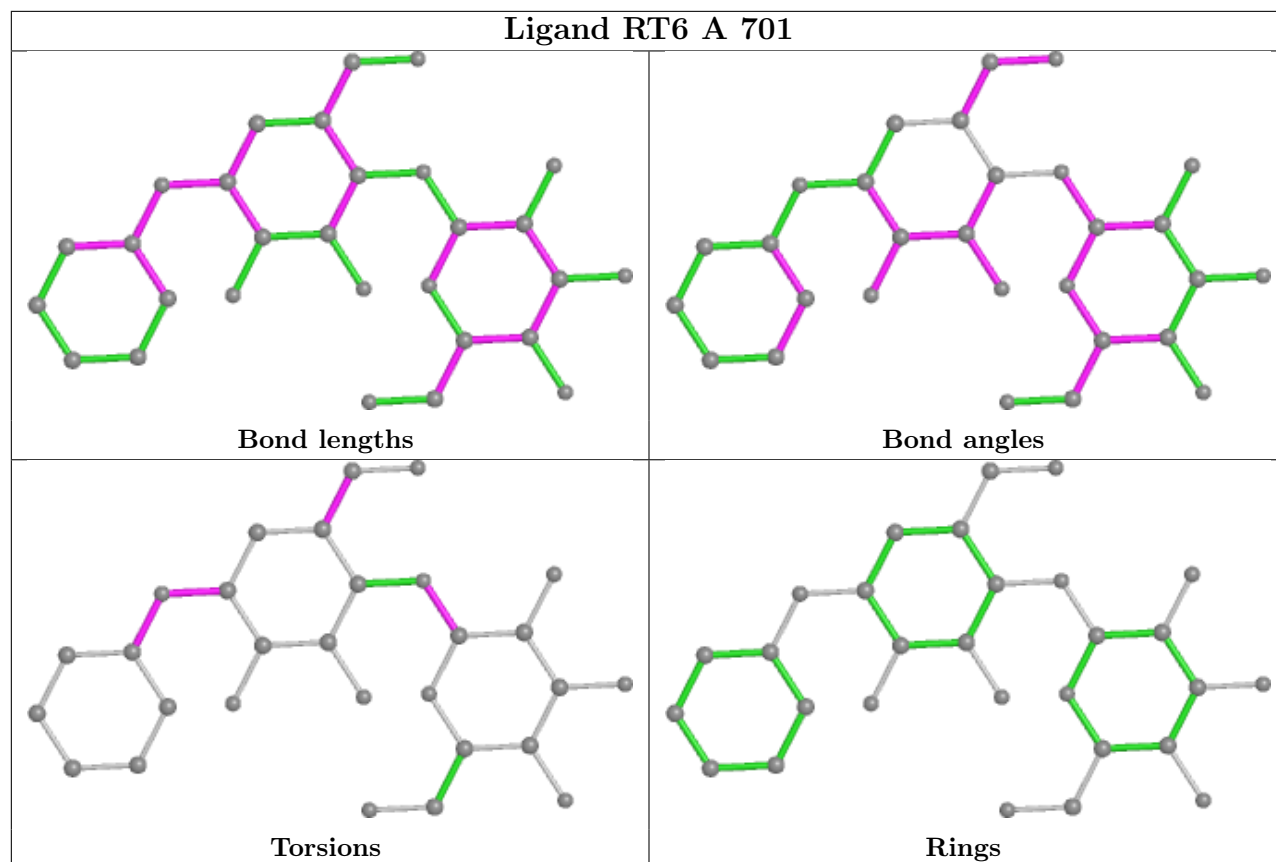
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	702	RT6	1	0
3	A	701	RT6	1	0

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





## 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	649/683 (95%)	-0.19	15 (2%) 60 67	41, 53, 87, 123	1 (0%)
1	B	649/683 (95%)	-0.14	9 (1%) 75 80	40, 54, 80, 108	1 (0%)
All	All	1298/1366 (95%)	-0.16	24 (1%) 68 74	40, 54, 83, 123	2 (0%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	627	SER	3.9
1	B	627	SER	3.5
1	A	153	ARG	3.3
1	B	169	ALA	3.2
1	A	144	ALA	3.2
1	A	150	GLU	3.0
1	A	15	PRO	2.8
1	A	143	ALA	2.6
1	A	149	ARG	2.6
1	A	185	LEU	2.6
1	B	69	ARG	2.6
1	A	181	VAL	2.5
1	B	175	ALA	2.5
1	B	628	VAL	2.4
1	A	152	ARG	2.4
1	A	165	GLU	2.3
1	B	151	ASP	2.3
1	A	662	LEU	2.3
1	A	167	ARG	2.3
1	B	647	VAL	2.3
1	B	185	LEU	2.2
1	A	155	LEU	2.2
1	B	629	PRO	2.1
1	A	620	LEU	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](#)

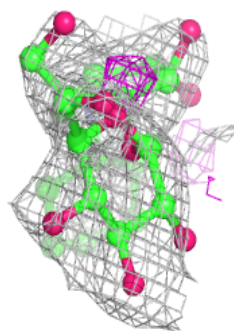
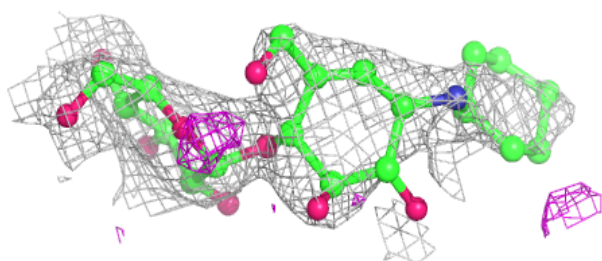
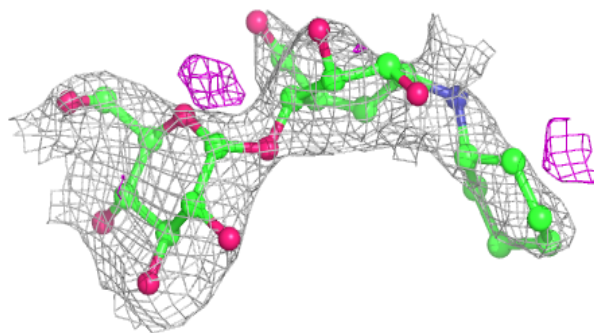
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
2	PEG	B	701	7/7	0.87	0.13	46,52,62,68	0
2	PEG	A	702	7/7	0.87	0.16	63,68,69,73	0
3	RT6	A	701	29/29	0.90	0.30	56,71,76,81	3
3	RT6	B	702	29/29	0.93	0.21	59,72,87,88	3

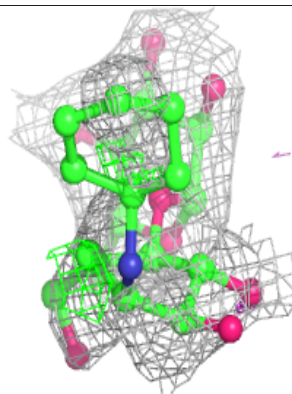
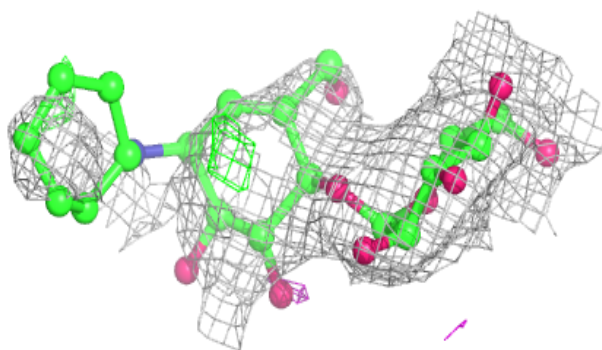
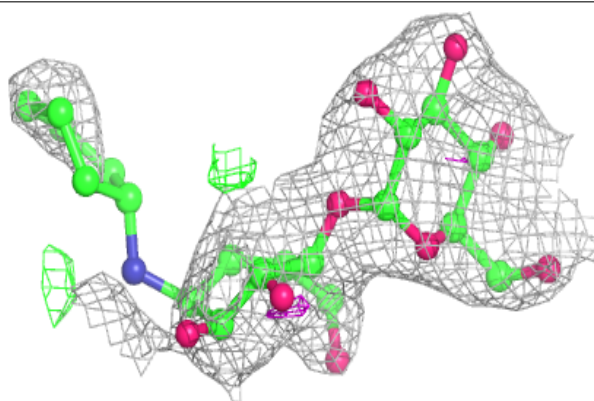
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around RT6 A 701:**

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

**Electron density around RT6 B 702:**

$2mF_o-DF_c$  (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.