



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 17, 2021 – 07:30 PM JST

PDB ID : 7CMD
Title : Crystal structure of the SARS-CoV-2 PLpro with GRL0617
Authors : Gao, X.; Cui, S.
Deposited on : 2020-07-27
Resolution : 2.59 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 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 : 2.18
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) : 2.18

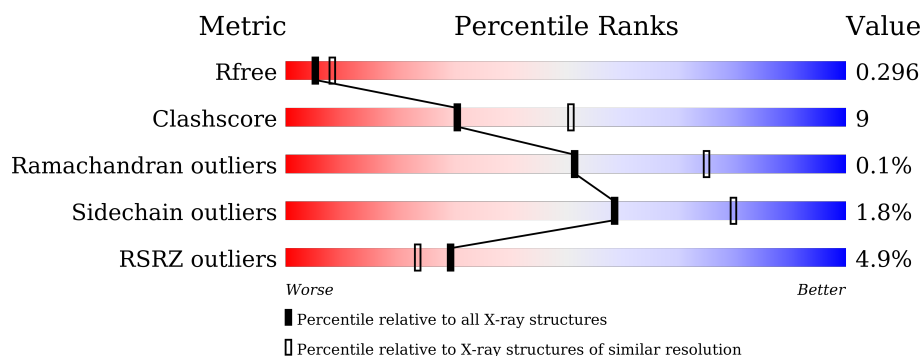
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	319	<div> <div>3%</div> <div>77%</div> <div>18%</div> <div>• •</div> </div>
1	B	319	<div> <div>5%</div> <div>75%</div> <div>22%</div> <div>•</div> </div>
1	C	319	<div> <div>6%</div> <div>76%</div> <div>20%</div> <div>•</div> </div>
1	D	319	<div> <div>4%</div> <div>76%</div> <div>19%</div> <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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ZN	B	503	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 20036 atoms, of which 9830 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.

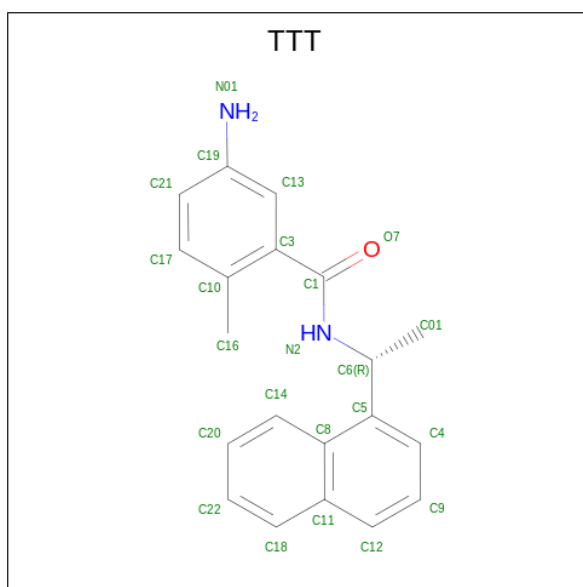
- Molecule 1 is a protein called Non-structural protein 3.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	306	Total	C	H	N	O	S	0	5	0
			4879	1570	2422	402	468	17			
1	B	308	Total	C	H	N	O	S	0	12	0
			5020	1611	2503	412	477	17			
1	C	308	Total	C	H	N	O	S	8	6	0
			4920	1584	2442	405	472	17			
1	D	307	Total	C	H	N	O	S	0	2	0
			4834	1562	2383	403	469	17			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ALA	-	expression tag	UNP P0DTD1
B	0	ALA	-	expression tag	UNP P0DTD1
C	0	ALA	-	expression tag	UNP P0DTD1
D	0	ALA	-	expression tag	UNP P0DTD1

- Molecule 2 is 5-amino-2-methyl-N-[(1R)-1-naphthalen-1-ylethyl]benzamide (three-letter code: TTT) (formula: C₂₀H₂₀N₂O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 20	H 20	N 2	O 1	0	0
2	B	1	Total 43	C 20	H 20	N 2	O 1	0	0
2	C	1	Total 43	C 20	H 20	N 2	O 1	0	0
2	D	1	Total 43	C 20	H 20	N 2	O 1	0	0

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		
3	B	4	Total	Zn	1	0
			4	4		
3	C	1	Total	Zn	0	0
			1	1		
3	D	2	Total	Zn	0	0
			2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	49	Total	O	0	0
			49	49		

Continued on next page...

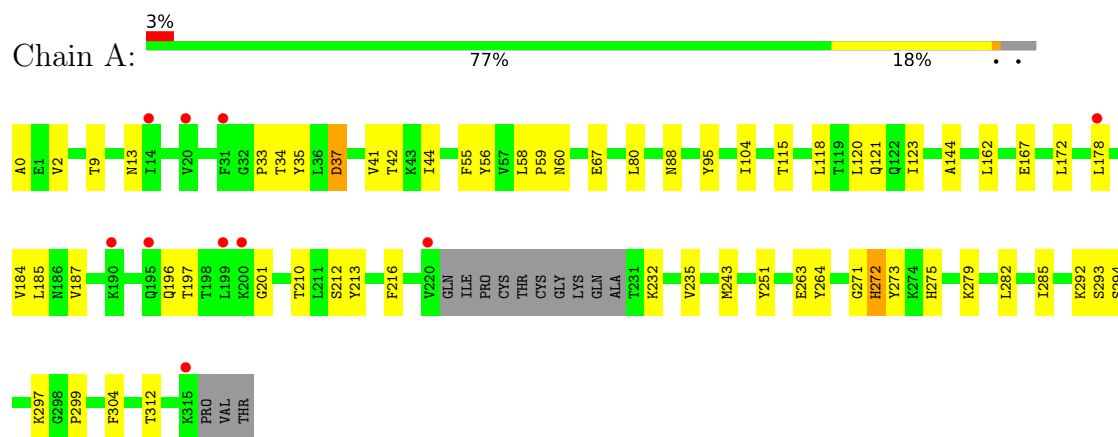
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	51	Total 51	O 51	0	0
4	C	42	Total 42	O 42	0	0
4	D	60	Total 60	O 60	0	0

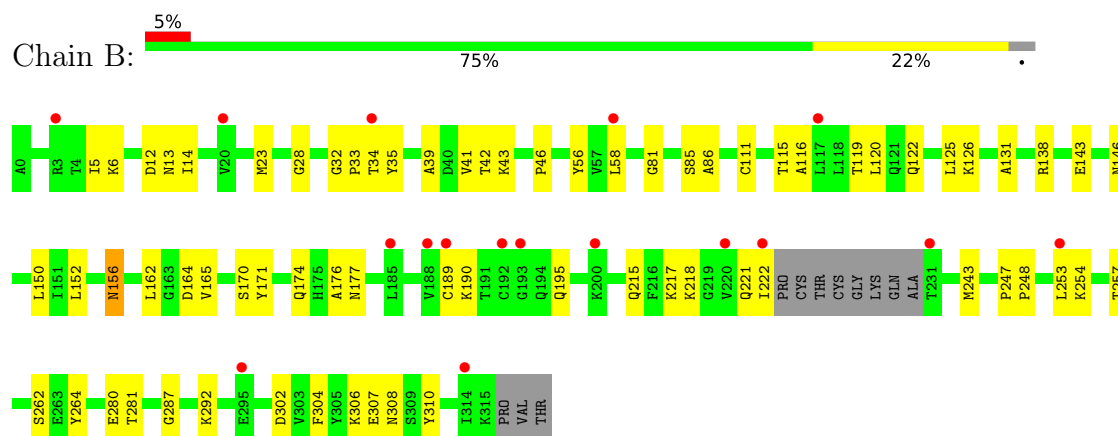
3 Residue-property plots [i](#)

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.

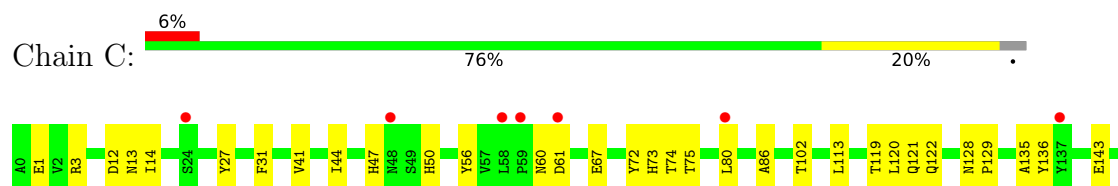
• Molecule 1: Non-structural protein 3

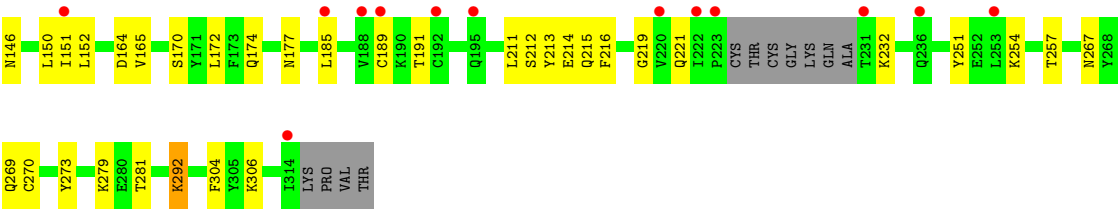


• Molecule 1: Non-structural protein 3

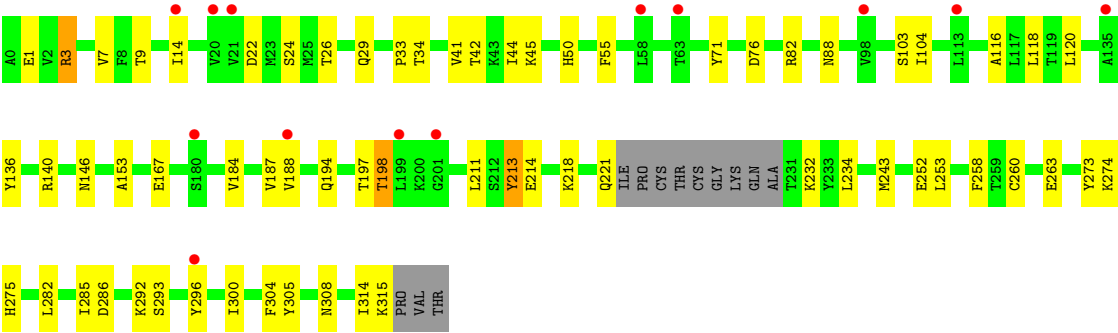
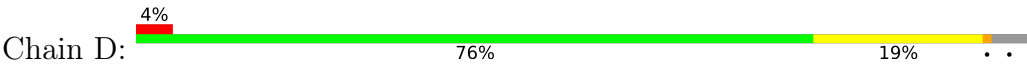


• Molecule 1: Non-structural protein 3





● Molecule 1: Non-structural protein 3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.47Å 123.51Å 146.81Å 90.00° 90.06° 90.00°	Depositor
Resolution (Å)	94.51 – 2.59 94.51 – 2.59	Depositor EDS
% Data completeness (in resolution range)	99.0 (94.51-2.59) 99.4 (94.51-2.59)	Depositor EDS
R_{merge}	0.32	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.58Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.244 , 0.298 0.240 , 0.296	Depositor DCC
R_{free} test set	3383 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	36.3	Xtriage
Anisotropy	0.950	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 39.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.439 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	20036	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.57 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.2967e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.44	0/2529	0.61	0/3429
1	B	0.44	0/2610	0.58	0/3535
1	C	0.43	0/2554	0.58	0/3465
1	D	0.46	0/2514	0.59	0/3409
All	All	0.44	0/10207	0.59	0/13838

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	2457	2422	2422	43	0
1	B	2517	2503	2505	51	0
1	C	2478	2442	2445	49	0
1	D	2451	2383	2400	41	0
2	A	23	20	20	1	0
2	B	23	20	20	4	0
2	C	23	20	20	0	0
2	D	23	20	20	0	0
3	A	2	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	4	0	0	0	0
3	C	1	0	0	0	0
3	D	2	0	0	0	0
4	A	49	0	0	10	0
4	B	51	0	0	17	0
4	C	42	0	0	17	0
4	D	60	0	0	6	0
All	All	10206	9830	9852	182	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 9.

All (182) 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:B:111:CYS:SG	4:B:642:HOH:O	2.09	1.11
1:A:292:LYS:NZ	4:A:702:HOH:O	1.93	0.98
1:C:189:CYS:SG	4:C:707:HOH:O	2.21	0.98
1:D:82:ARG:NH2	1:D:153:ALA:O	2.00	0.95
1:D:167:GLU:OE1	4:D:701:HOH:O	1.88	0.92
1:C:143:GLU:O	4:C:701:HOH:O	1.87	0.92
1:B:189:CYS:SG	4:B:608:HOH:O	2.28	0.89
1:A:294:SER:O	4:A:701:HOH:O	1.91	0.88
1:B:23:MET:O	4:B:601:HOH:O	1.93	0.85
1:C:279:LYS:NZ	4:C:703:HOH:O	1.97	0.84
1:B:162:LEU:CD1	2:B:502:TTT:H16	2.08	0.84
1:B:253:LEU:HA	4:B:603:HOH:O	1.76	0.83
1:D:308:ASN:OD1	4:D:702:HOH:O	1.95	0.82
1:C:164:ASP:OD2	4:C:702:HOH:O	1.97	0.81
1:D:184:VAL:HG22	1:D:198:THR:HG22	1.61	0.81
1:C:177:ASN:O	4:C:704:HOH:O	2.00	0.78
1:A:123:ILE:O	4:A:704:HOH:O	2.03	0.76
1:D:76:ASP:OD2	1:D:82:ARG:NH1	2.19	0.76
1:B:164:ASP:OD2	4:B:604:HOH:O	2.03	0.75
1:A:13:ASN:O	1:A:13:ASN:ND2	2.20	0.75
1:A:33:PRO:O	4:A:703:HOH:O	2.03	0.75
1:B:190:LYS:N	4:B:608:HOH:O	2.20	0.74
1:A:56:TYR:N	4:A:706:HOH:O	2.19	0.73
1:B:177:ASN:O	4:B:606:HOH:O	2.07	0.73
1:C:213:TYR:N	4:C:709:HOH:O	2.22	0.73
1:B:126[A]:LYS:NZ	4:B:605:HOH:O	2.07	0.72

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:221:GLN:N	1:C:221:GLN:OE1	2.23	0.71
1:B:156[A]:ASN:OD1	4:B:607:HOH:O	2.09	0.70
1:D:263:GLU:OE2	1:D:296:TYR:OH	2.04	0.70
1:C:279:LYS:HD2	4:C:703:HOH:O	1.92	0.69
1:A:167:GLU:OE1	4:A:705:HOH:O	2.10	0.69
1:D:221:GLN:O	4:D:704:HOH:O	2.11	0.69
1:A:35:TYR:O	4:A:706:HOH:O	2.11	0.67
1:A:104:ILE:O	4:A:707:HOH:O	2.13	0.67
1:D:234:LEU:O	4:D:705:HOH:O	2.14	0.66
1:B:287:GLY:N	4:B:609:HOH:O	2.30	0.65
1:B:138:ARG:CG	1:B:143:GLU:HG3	2.27	0.64
1:A:282:LEU:HB2	1:A:293:SER:O	1.98	0.64
1:C:67:GLU:OE1	4:C:706:HOH:O	2.14	0.64
1:B:264:TYR:CE2	2:B:502:TTT:H14	2.32	0.64
1:C:221:GLN:N	4:C:708:HOH:O	2.31	0.63
1:A:185:LEU:HD21	1:A:216:PHE:HZ	1.62	0.63
1:B:254:LYS:N	4:B:603:HOH:O	2.02	0.63
1:B:171:TYR:OH	4:B:607:HOH:O	2.10	0.63
1:B:138:ARG:HG2	1:B:143:GLU:HG3	1.83	0.60
1:D:188:VAL:HG13	1:D:194:GLN:HG2	1.83	0.59
1:D:211:LEU:HD13	1:D:300:ILE:HG22	1.85	0.58
1:C:135:ALA:N	4:C:711:HOH:O	2.36	0.58
1:B:81:GLY:O	1:B:85[A]:SER:OG	2.18	0.57
1:C:27:TYR:OH	1:C:50:HIS:ND1	2.19	0.57
1:B:42:THR:HG22	1:B:43:LYS:HD2	1.85	0.57
1:D:34:THR:HG22	1:D:41:VAL:CG2	2.35	0.57
1:C:113:LEU:HD11	1:C:152:LEU:HD21	1.87	0.56
1:C:122:GLN:O	1:C:306:LYS:NZ	2.40	0.55
1:B:33:PRO:HG2	1:B:58:LEU:HD12	1.89	0.55
1:A:210:THR:HG22	1:A:212:SER:H	1.72	0.54
1:B:5:ILE:HD11	1:B:46:PRO:HB3	1.90	0.54
1:B:162:LEU:HD13	2:B:502:TTT:H16	1.90	0.53
1:C:41:VAL:O	1:C:44:ILE:HG22	2.09	0.53
1:C:279:LYS:CE	4:C:703:HOH:O	2.54	0.53
1:C:191:THR:N	4:C:707:HOH:O	2.17	0.53
1:A:41:VAL:O	1:A:44:ILE:HG22	2.08	0.53
1:D:282:LEU:HB2	1:D:293:SER:O	2.08	0.52
1:A:59:PRO:HD3	1:A:80[A]:LEU:HD12	1.90	0.52
1:C:221:GLN:O	4:C:708:HOH:O	2.19	0.52
1:C:120:LEU:HD11	1:C:172:LEU:HD13	1.91	0.52
1:B:35:TYR:HA	1:B:39:ALA:O	2.10	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:264:TYR:OH	1:A:271:GLY:N	2.40	0.52
1:B:195:GLN:NE2	4:B:615:HOH:O	2.43	0.52
1:B:254:LYS:HB2	1:B:257:THR:OG1	2.10	0.52
1:D:116:ALA:O	1:D:120:LEU:HD23	2.11	0.51
1:C:136:TYR:N	4:C:711:HOH:O	2.35	0.51
1:A:185:LEU:HD21	1:A:216:PHE:CZ	2.44	0.51
1:C:146:ASN:O	1:C:150:LEU:HD12	2.11	0.51
1:A:121:GLN:O	1:C:267:ASN:ND2	2.44	0.50
1:A:37:ASP:N	1:A:37:ASP:OD1	2.43	0.50
1:C:151:ILE:HG12	1:C:172:LEU:HD21	1.93	0.50
1:B:177:ASN:HB3	4:B:602:HOH:O	2.11	0.50
1:A:187:VAL:HG22	1:A:232:LYS:HG3	1.94	0.50
1:B:162:LEU:HD11	2:B:502:TTT:H16	1.89	0.49
1:C:13:ASN:HB2	1:C:56:TYR:OH	2.12	0.49
1:D:1:GLU:HG2	1:D:3:ARG:HD3	1.94	0.49
1:C:279:LYS:CD	4:C:703:HOH:O	2.57	0.49
1:A:178:LEU:O	1:A:201:GLY:HA2	2.13	0.49
1:B:116:ALA:O	1:B:120:LEU:HD23	2.12	0.49
1:D:103:SER:OG	1:D:104:ILE:N	2.46	0.49
1:A:115[B]:THR:HG23	1:A:275:HIS:HB2	1.95	0.49
1:B:34:THR:HG22	1:B:41:VAL:HG23	1.94	0.48
1:B:138:ARG:HG2	1:B:143:GLU:O	2.13	0.48
1:C:12:ASP:OD1	1:C:14:ILE:N	2.37	0.48
1:D:9:THR:HA	1:D:55:PHE:O	2.14	0.48
1:D:88:ASN:ND2	4:D:713:HOH:O	2.44	0.48
1:C:254:LYS:HB2	1:C:257:THR:OG1	2.13	0.47
1:B:243:MET:HG3	1:B:304:PHE:CE2	2.49	0.47
1:C:185:LEU:HD21	1:C:216:PHE:HZ	1.79	0.47
1:D:188:VAL:HG22	1:D:194:GLN:HG2	1.96	0.47
1:B:262:SER:OG	1:B:302:ASP:HB2	2.14	0.47
1:D:7:VAL:HG12	1:D:50:HIS:O	2.15	0.47
1:D:260:CYS:HB2	1:D:275:HIS:CE1	2.50	0.47
1:D:214:GLU:OE1	1:D:218:LYS:HE2	2.15	0.46
1:B:217:LYS:HE2	1:B:310:TYR:CE1	2.50	0.46
1:C:170:SER:O	1:C:174:GLN:HG2	2.16	0.46
1:C:281:THR:HG21	1:C:292:LYS:HD3	1.97	0.46
1:B:221:GLN:OE1	1:B:221:GLN:N	2.49	0.46
1:C:185:LEU:HD21	1:C:216:PHE:CZ	2.51	0.46
1:A:33:PRO:HA	1:A:42:THR:OG1	2.16	0.46
1:B:28:GLY:HA3	1:B:42:THR:HG23	1.98	0.46
1:D:136:TYR:CZ	1:D:140:ARG:HD3	2.51	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:22:ASP:OD1	1:D:24:SER:N	2.43	0.45
1:A:120:LEU:HD11	1:A:172:LEU:CD1	2.46	0.45
1:B:131:ALA:HB1	1:B:150:LEU:HD21	1.98	0.45
1:D:243:MET:HG3	1:D:304:PHE:CZ	2.52	0.45
1:C:41:VAL:HB	1:C:44:ILE:CG2	2.47	0.45
1:C:215:GLN:NE2	1:C:219:GLY:O	2.49	0.45
1:D:14:ILE:HD12	1:D:71:TYR:CE1	2.51	0.45
1:B:307:GLU:HG3	1:B:308:ASN:N	2.32	0.45
1:B:115[A]:THR:CG2	1:B:165:VAL:HG21	2.47	0.45
1:C:86:ALA:HB1	1:C:152:LEU:HD12	2.00	0.44
1:A:118:LEU:HD11	1:A:285:ILE:HG21	1.99	0.44
1:A:118:LEU:HD11	1:A:285:ILE:CG2	2.47	0.44
1:A:162:LEU:HD23	2:A:601:TTT:H16	2.00	0.44
1:C:165:VAL:HG23	1:C:273:TYR:CE2	2.53	0.44
1:B:126[A]:LYS:HG3	4:B:602:HOH:O	2.16	0.44
1:C:214:GLU:N	4:C:709:HOH:O	2.20	0.44
1:D:263:GLU:O	1:D:273:TYR:HA	2.18	0.44
1:A:9:THR:HA	1:A:55:PHE:O	2.17	0.44
1:B:28:GLY:HA2	1:B:32:GLY:O	2.18	0.44
1:C:211:LEU:HD12	1:C:251:TYR:CD1	2.53	0.44
1:D:211:LEU:CD1	1:D:300:ILE:HG22	2.47	0.44
1:B:170:SER:O	1:B:174:GLN:HG2	2.18	0.43
1:B:280:GLU:HG3	1:B:281:THR:HG22	2.00	0.43
1:D:253:LEU:HB3	1:D:258:PHE:CE1	2.54	0.43
1:A:185:LEU:HB2	1:A:197:THR:OG1	2.18	0.43
1:A:95:TYR:CD2	1:A:144:ALA:HB3	2.53	0.43
1:B:86:ALA:HB1	1:B:152:LEU:HD12	2.00	0.43
1:C:41:VAL:HB	1:C:44:ILE:HG22	2.00	0.43
1:D:314:ILE:O	1:D:315:LYS:HE2	2.19	0.43
1:A:88:ASN:ND2	4:A:710:HOH:O	2.37	0.43
1:D:33:PRO:HA	1:D:42:THR:OG1	2.18	0.43
1:B:122:GLN:O	1:B:306:LYS:NZ	2.52	0.43
1:D:146:ASN:ND2	4:D:703:HOH:O	2.09	0.43
1:C:102:THR:O	1:C:121:GLN:NE2	2.50	0.43
1:C:219:GLY:HA2	1:C:232:LYS:O	2.19	0.43
1:D:118:LEU:HD11	1:D:285:ILE:CG2	2.49	0.43
1:B:217:LYS:HE2	1:B:310:TYR:CD1	2.54	0.42
1:A:58:LEU:O	1:A:60:ASN:N	2.52	0.42
1:A:185:LEU:O	1:A:196:GLN:HA	2.19	0.42
1:A:0:ALA:N	1:D:252:GLU:O	2.53	0.42
1:C:72:TYR:CE2	1:C:80[B]:LEU:HD21	2.55	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:12:ASP:OD1	1:B:14:ILE:N	2.37	0.42
1:B:146:ASN:O	1:B:150:LEU:CD1	2.67	0.42
1:B:125:LEU:HD11	1:B:176:ALA:CB	2.49	0.42
1:D:41:VAL:O	1:D:44:ILE:HG22	2.20	0.42
1:A:34:THR:HG22	1:A:41:VAL:CG2	2.49	0.42
1:A:279:LYS:HD3	1:C:269:GLN:OE1	2.19	0.42
1:B:6:LYS:HD2	4:B:619:HOH:O	2.20	0.42
1:D:213:TYR:HB2	1:D:305:TYR:CE2	2.54	0.42
1:D:184:VAL:HA	1:D:197:THR:O	2.19	0.42
1:D:274:LYS:HD3	1:D:286:ASP:HB2	2.01	0.42
1:A:2:VAL:HG13	1:A:2:VAL:O	2.20	0.42
1:A:67:GLU:OE2	4:A:708:HOH:O	2.22	0.42
1:B:119:THR:HG21	1:B:304:PHE:CZ	2.55	0.41
1:C:191:THR:OG1	4:C:707:HOH:O	2.19	0.41
1:A:263:GLU:HG3	1:A:299:PRO:O	2.20	0.41
1:A:251:TYR:O	1:A:297:LYS:HA	2.21	0.41
1:A:184:VAL:HB	1:A:235:VAL:HG13	2.03	0.41
1:B:247:PRO:O	1:B:248:PRO:C	2.59	0.41
1:A:243:MET:HG3	1:A:304:PHE:CZ	2.56	0.41
1:B:13:ASN:HB2	1:B:56:TYR:OH	2.21	0.41
1:C:119:THR:HG21	1:C:304:PHE:CZ	2.55	0.41
1:D:26:THR:HA	1:D:45:LYS:HA	2.02	0.41
1:D:187:VAL:HG22	1:D:232:LYS:HG3	2.02	0.41
1:D:213:TYR:HB2	1:D:305:TYR:CD2	2.56	0.41
1:C:31:PHE:O	1:C:60:ASN:ND2	2.54	0.40
1:C:74:THR:OG1	1:C:75:THR:N	2.54	0.40
1:A:210:THR:HG22	1:A:212:SER:N	2.35	0.40
1:A:272:HIS:ND1	1:A:273:TYR:N	2.69	0.40
1:D:118:LEU:HD11	1:D:285:ILE:HG21	2.03	0.40
1:C:1:GLU:OE2	1:C:3:ARG:HD2	2.21	0.40
1:C:146:ASN:O	1:C:150:LEU:CD1	2.70	0.40
1:C:128:ASN:HB2	1:C:129:PRO:HD3	2.03	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	307/319 (96%)	290 (94%)	17 (6%)	0	100	100
1	B	316/319 (99%)	293 (93%)	22 (7%)	1 (0%)	41	64
1	C	310/319 (97%)	289 (93%)	21 (7%)	0	100	100
1	D	305/319 (96%)	285 (93%)	20 (7%)	0	100	100
All	All	1238/1276 (97%)	1157 (94%)	80 (6%)	1 (0%)	51	75

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	218	LYS

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	272/278 (98%)	268 (98%)	4 (2%)	65	83
1	B	281/278 (101%)	276 (98%)	5 (2%)	59	80
1	C	275/278 (99%)	269 (98%)	6 (2%)	52	76
1	D	270/278 (97%)	265 (98%)	5 (2%)	57	79
All	All	1098/1112 (99%)	1078 (98%)	20 (2%)	59	80

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ASP
1	A	213	TYR
1	A	272	HIS
1	A	312	THR
1	B	156[A]	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	156[B]	ASN
1	B	215	GLN
1	B	222	ILE
1	B	292	LYS
1	C	47	HIS
1	C	61	ASP
1	C	73	HIS
1	C	212	SER
1	C	270	CYS
1	C	292	LYS
1	D	3	ARG
1	D	29	GLN
1	D	198	THR
1	D	213	TYR
1	D	292	LYS

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	13	ASN
1	D	308	ASN

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 13 ligands modelled in this entry, 9 are monoatomic - leaving 4 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	$\# Z > 2$
2	TTT	D	601	-	25,25,25	3.51	12 (48%)	35,35,35	1.34	2 (5%)
2	TTT	B	502	-	25,25,25	3.53	12 (48%)	35,35,35	1.39	6 (17%)
2	TTT	C	601	-	25,25,25	3.48	11 (44%)	35,35,35	1.26	4 (11%)
2	TTT	A	601	-	25,25,25	3.58	12 (48%)	35,35,35	1.44	5 (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	TTT	D	601	-	-	6/12/12/12	0/3/3/3
2	TTT	B	502	-	-	6/12/12/12	0/3/3/3
2	TTT	C	601	-	-	4/12/12/12	0/3/3/3
2	TTT	A	601	-	-	4/12/12/12	0/3/3/3

All (47) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	TTT	C22-C20	8.35	1.60	1.38
2	D	601	TTT	C22-C20	8.24	1.59	1.38
2	A	601	TTT	C4-C5	8.18	1.51	1.37
2	C	601	TTT	C22-C20	8.17	1.59	1.38
2	C	601	TTT	C4-C5	8.16	1.51	1.37
2	B	502	TTT	C22-C20	8.12	1.59	1.38
2	D	601	TTT	C4-C5	8.11	1.51	1.37
2	B	502	TTT	C4-C5	8.10	1.51	1.37
2	C	601	TTT	C9-C12	6.92	1.52	1.36
2	B	502	TTT	C9-C12	6.65	1.51	1.36
2	A	601	TTT	C9-C12	6.64	1.51	1.36
2	D	601	TTT	C9-C12	6.48	1.51	1.36
2	A	601	TTT	C5-C8	-5.94	1.33	1.43
2	D	601	TTT	C5-C8	-5.65	1.33	1.43

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	502	TTT	C5-C8	-5.31	1.34	1.43
2	A	601	TTT	C1-N2	5.31	1.45	1.34
2	C	601	TTT	C1-N2	5.17	1.45	1.34
2	B	502	TTT	C1-N2	5.07	1.45	1.34
2	D	601	TTT	C1-N2	5.01	1.45	1.34
2	D	601	TTT	C9-C4	-4.70	1.29	1.38
2	A	601	TTT	C9-C4	-4.60	1.29	1.38
2	C	601	TTT	C5-C8	-4.53	1.35	1.43
2	B	502	TTT	C9-C4	-4.45	1.29	1.38
2	C	601	TTT	C9-C4	-4.41	1.29	1.38
2	B	502	TTT	C12-C11	-3.84	1.32	1.41
2	A	601	TTT	C12-C11	-3.82	1.32	1.41
2	B	502	TTT	C14-C8	3.76	1.50	1.42
2	D	601	TTT	C12-C11	-3.71	1.33	1.41
2	C	601	TTT	C12-C11	-3.42	1.33	1.41
2	C	601	TTT	C14-C8	3.40	1.49	1.42
2	A	601	TTT	C14-C8	3.11	1.48	1.42
2	D	601	TTT	C14-C8	3.11	1.48	1.42
2	C	601	TTT	C18-C11	2.84	1.48	1.41
2	D	601	TTT	O7-C1	-2.64	1.17	1.23
2	B	502	TTT	C18-C11	2.60	1.48	1.41
2	D	601	TTT	C18-C11	2.58	1.48	1.41
2	A	601	TTT	C18-C11	2.54	1.48	1.41
2	A	601	TTT	C8-C11	2.46	1.47	1.43
2	C	601	TTT	C8-C11	2.43	1.47	1.43
2	B	502	TTT	C8-C11	2.36	1.47	1.43
2	A	601	TTT	O7-C1	-2.30	1.18	1.23
2	D	601	TTT	C8-C11	2.24	1.47	1.43
2	A	601	TTT	C19-N01	2.13	1.45	1.38
2	C	601	TTT	C19-N01	2.08	1.45	1.38
2	D	601	TTT	C19-N01	2.07	1.45	1.38
2	B	502	TTT	C19-N01	2.02	1.45	1.38
2	B	502	TTT	O7-C1	-2.02	1.19	1.23

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	601	TTT	C5-C6-N2	-4.46	101.19	110.89
2	A	601	TTT	C5-C6-N2	-4.24	101.67	110.89
2	B	502	TTT	C3-C1-N2	2.83	122.05	116.80
2	D	601	TTT	C4-C5-C6	-2.71	115.48	120.64
2	B	502	TTT	O7-C1-C3	-2.68	116.12	121.01

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	TTT	C16-C10-C3	-2.66	118.66	122.79
2	B	502	TTT	C5-C6-N2	-2.64	105.15	110.89
2	C	601	TTT	C16-C10-C3	-2.58	118.79	122.79
2	B	502	TTT	C10-C3-C1	-2.57	118.64	120.56
2	C	601	TTT	C3-C1-N2	2.53	121.49	116.80
2	B	502	TTT	C16-C10-C3	-2.51	118.90	122.79
2	A	601	TTT	C14-C8-C5	-2.40	120.61	123.40
2	C	601	TTT	O7-C1-C3	-2.38	116.67	121.01
2	C	601	TTT	C4-C5-C6	-2.34	116.19	120.64
2	A	601	TTT	C01-C6-N2	2.23	112.98	109.05
2	B	502	TTT	C13-C3-C10	2.22	121.53	119.57
2	A	601	TTT	C4-C5-C6	-2.09	116.67	120.64

There are no chirality outliers.

All (20) torsion outliers are listed below:

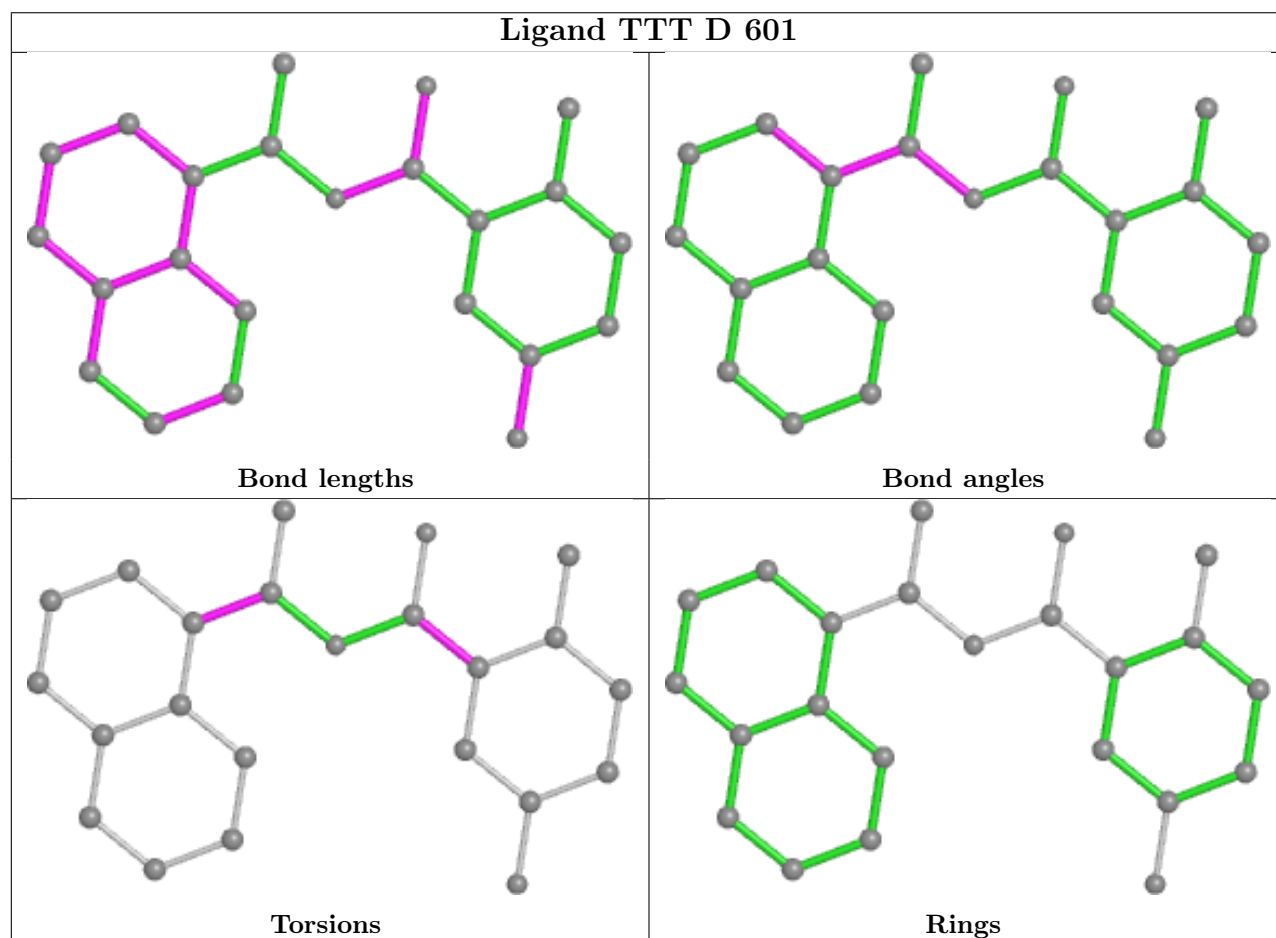
Mol	Chain	Res	Type	Atoms
2	A	601	TTT	C4-C5-C6-N2
2	B	502	TTT	C4-C5-C6-N2
2	B	502	TTT	O7-C1-C3-C13
2	B	502	TTT	N2-C1-C3-C13
2	C	601	TTT	C8-C5-C6-N2
2	D	601	TTT	C4-C5-C6-N2
2	A	601	TTT	C8-C5-C6-C01
2	B	502	TTT	C8-C5-C6-C01
2	B	502	TTT	C4-C5-C6-C01
2	C	601	TTT	C8-C5-C6-C01
2	C	601	TTT	C4-C5-C6-C01
2	D	601	TTT	C8-C5-C6-C01
2	D	601	TTT	C4-C5-C6-C01
2	A	601	TTT	C8-C5-C6-N2
2	B	502	TTT	C8-C5-C6-N2
2	D	601	TTT	C8-C5-C6-N2
2	D	601	TTT	N2-C1-C3-C13
2	A	601	TTT	C4-C5-C6-C01
2	C	601	TTT	C4-C5-C6-N2
2	D	601	TTT	O7-C1-C3-C13

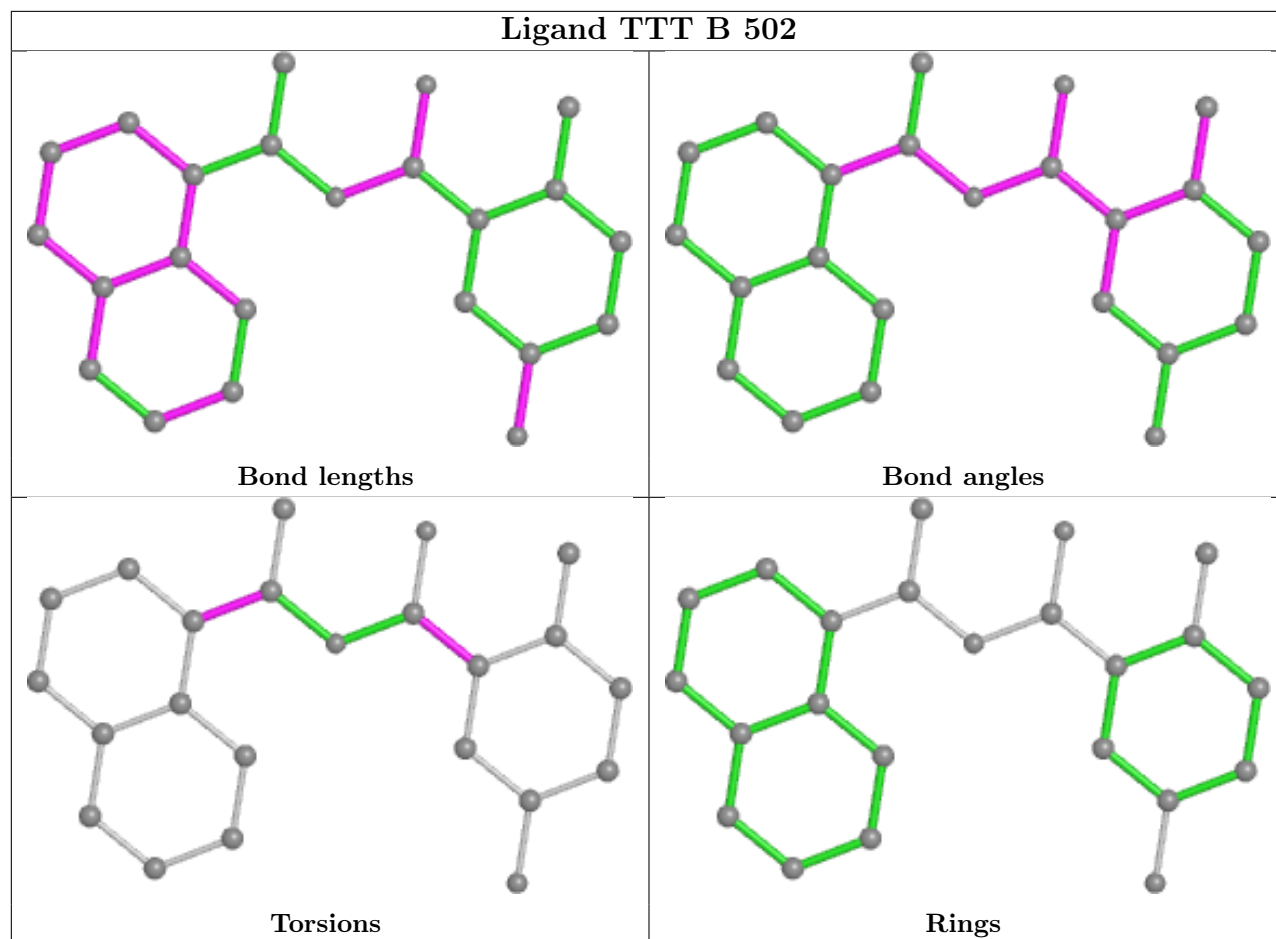
There are no ring outliers.

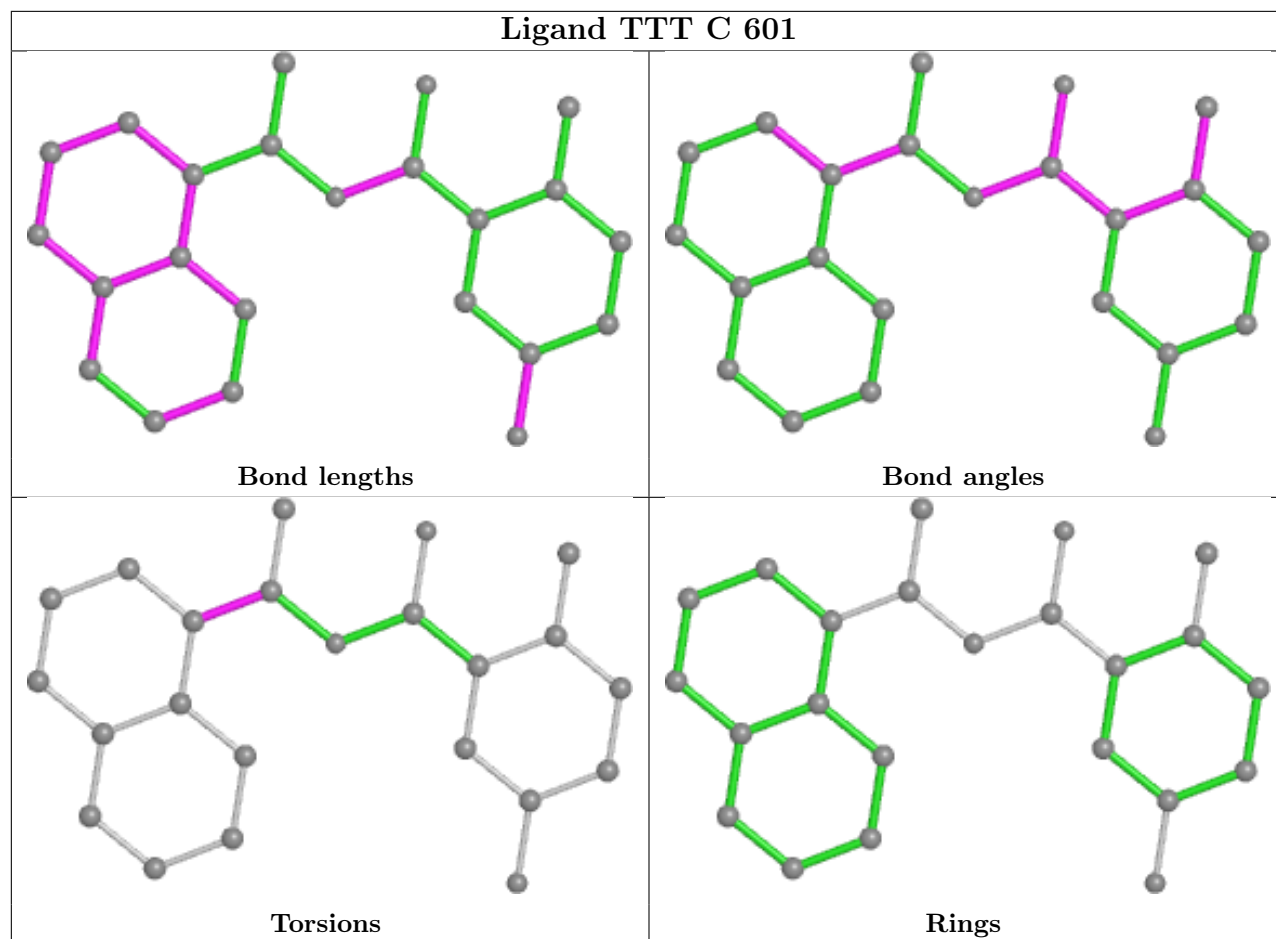
2 monomers are involved in 5 short contacts:

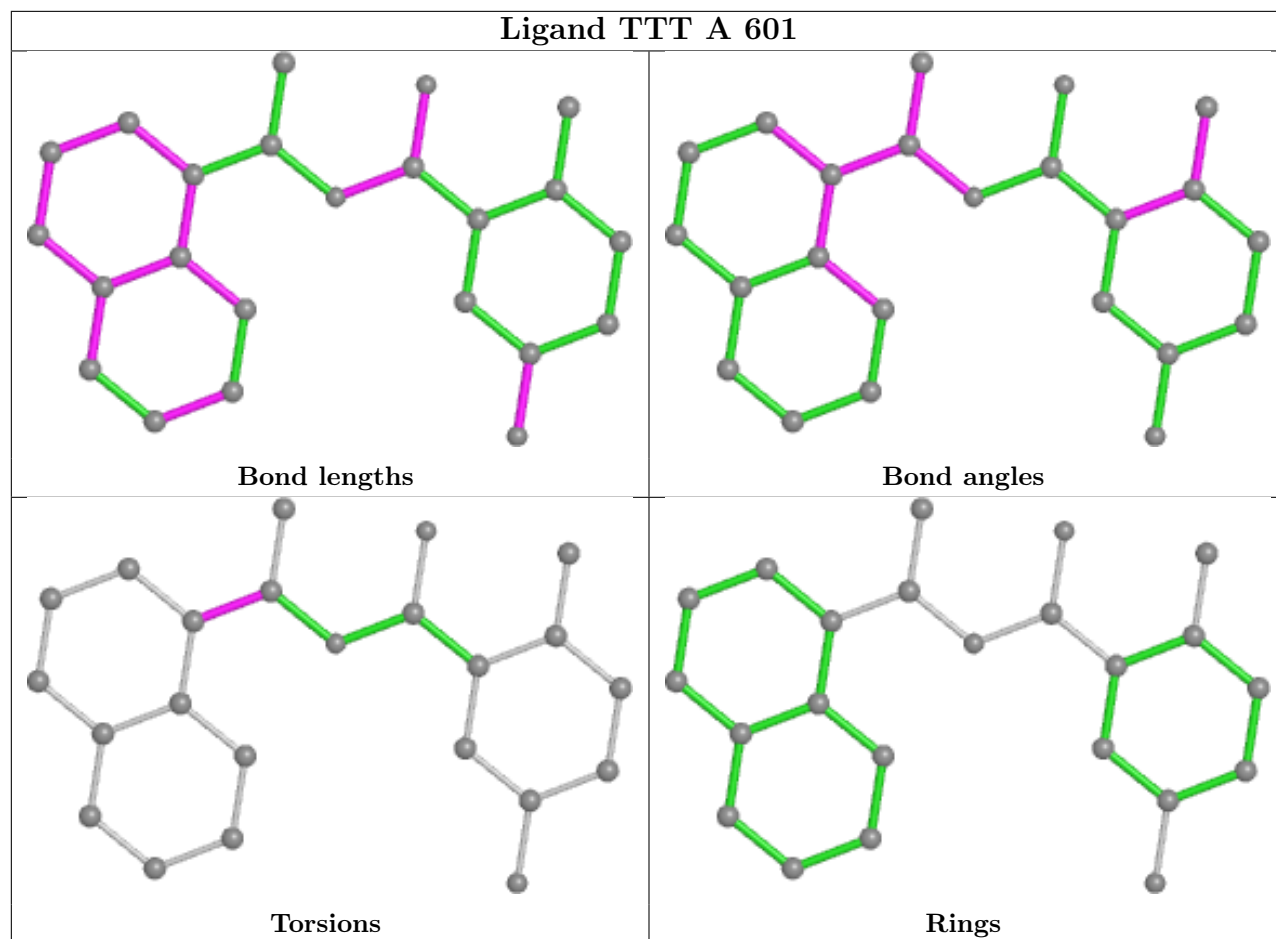
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	502	TTT	4	0
2	A	601	TTT	1	0

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, 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	306/319 (95%)	0.61	10 (3%) 46 39	23, 45, 67, 97	0
1	B	308/319 (96%)	0.71	17 (5%) 25 19	24, 49, 84, 115	0
1	C	308/319 (96%)	0.70	20 (6%) 18 14	26, 48, 82, 115	0
1	D	307/319 (96%)	0.61	13 (4%) 36 29	22, 46, 71, 107	0
All	All	1229/1276 (96%)	0.66	60 (4%) 29 23	22, 47, 76, 115	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	192	CYS	6.5
1	B	192	CYS	5.8
1	B	222	ILE	5.5
1	B	220	VAL	5.5
1	C	314	ILE	5.1
1	B	314	ILE	4.8
1	C	220	VAL	4.6
1	C	189	CYS	4.4
1	B	193	GLY	4.3
1	D	188	VAL	4.2
1	B	188	VAL	4.2
1	B	189	CYS	4.0
1	A	31	PHE	3.7
1	C	222	ILE	3.7
1	C	223	PRO	3.5
1	D	199	LEU	3.1
1	D	296	TYR	2.9
1	A	20	VAL	2.8
1	D	58	LEU	2.8
1	B	20	VAL	2.7
1	B	200[A]	LYS	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	80[A]	LEU	2.6
1	D	63	THR	2.5
1	C	137	TYR	2.5
1	A	199	LEU	2.5
1	B	58	LEU	2.5
1	C	61	ASP	2.5
1	B	185	LEU	2.5
1	D	21	VAL	2.5
1	A	220	VAL	2.4
1	D	180[A]	SER	2.4
1	B	3[A]	ARG	2.4
1	B	117	LEU	2.4
1	B	253	LEU	2.4
1	C	185	LEU	2.4
1	C	48	ASN	2.4
1	C	231	THR	2.4
1	A	178	LEU	2.4
1	C	188	VAL	2.3
1	C	59	PRO	2.2
1	B	295[A]	GLU	2.2
1	A	14	ILE	2.2
1	B	34	THR	2.2
1	C	24	SER	2.2
1	A	190	LYS	2.2
1	D	201	GLY	2.2
1	A	200[A]	LYS	2.1
1	A	315	LYS	2.1
1	B	231	THR	2.1
1	D	14	ILE	2.1
1	C	236[A]	GLN	2.1
1	D	20	VAL	2.1
1	D	98	VAL	2.1
1	C	253	LEU	2.1
1	C	151	ILE	2.1
1	C	58	LEU	2.0
1	A	195	GLN	2.0
1	D	135	ALA	2.0
1	D	113	LEU	2.0
1	C	195	GLN	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](#)

LIGAND-RSR INFOmissingINFO

6.5 Other polymers [i](#)

There are no such residues in this entry.