



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 09:12 AM UTC

PDB ID : 5I6X / pdb_00005i6x
Title : X-ray structure of the ts3 human serotonin transporter complexed with paroxetine at the central site
Authors : Coleman, J.A.; Green, E.M.; Gouaux, E.
Deposited on : 2016-02-16
Resolution : 3.14 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

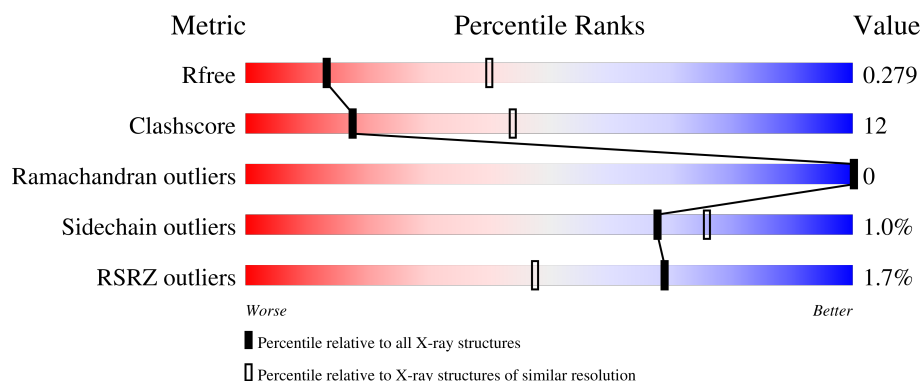
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

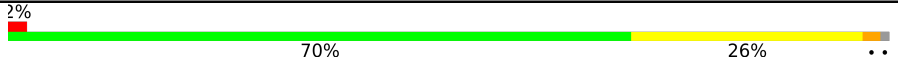
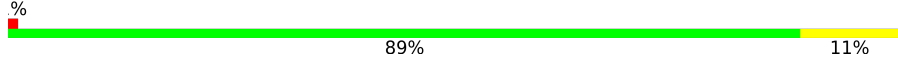
The reported resolution of this entry is 3.14 Å.

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	2351 (3.18-3.10)
Clashscore	190562	2452 (3.18-3.10)
Ramachandran outliers	187476	2324 (3.18-3.10)
Sidechain outliers	187428	2324 (3.18-3.10)
RSRZ outliers	180081	2351 (3.18-3.10)

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	549	
2	B	221	
3	C	214	

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 7631 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 Sodium-dependent serotonin transporter.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	544	Total	C	N	O	S	0	0	0
			4221	2817	656	724	24			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	74	GLY	-	cloning artifact	UNP P31645
A	75	SER	-	cloning artifact	UNP P31645
A	291	ALA	ILE	engineered mutation	UNP P31645
A	439	SER	THR	engineered mutation	UNP P31645
A	554	ALA	CYS	engineered mutation	UNP P31645
A	580	ALA	CYS	engineered mutation	UNP P31645
A	619	LEU	-	cloning artifact	UNP P31645
A	620	VAL	-	cloning artifact	UNP P31645
A	621	PRO	-	cloning artifact	UNP P31645
A	622	ARG	-	cloning artifact	UNP P31645

- Molecule 2 is a protein called 8B6 antibody, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	218	Total	C	N	O	S	0	0	0
			1643	1038	266	331	8			

- Molecule 3 is a protein called 8B6 antibody, light chain.

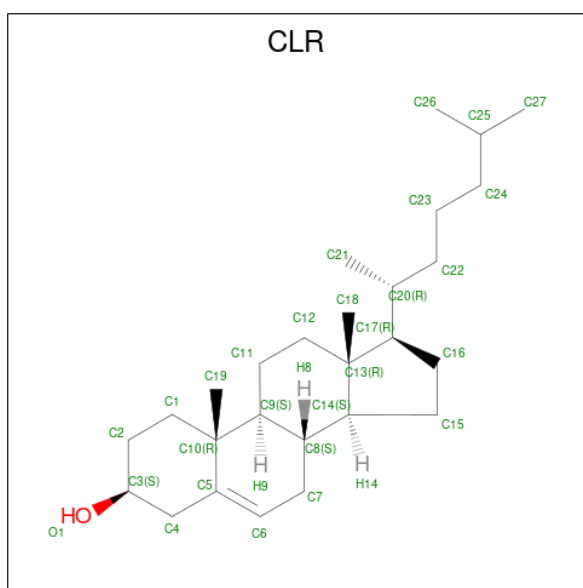
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	214	Total	C	N	O	S	0	0	0
			1662	1037	280	337	8			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 5 is CHOLESTEROL (CCD ID: CLR) (formula: $C_{27}H_{46}O$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			28	27	1		

- Molecule 6 is DODECYL-BETA-D-MALTOSIDE (CCD ID: LMT) (formula: $C_{24}H_{46}O_{11}$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			22	12	10		

- Molecule 7 is Paroxetine (CCD ID: 8PR) (formula: $C_{19}H_{20}FNO_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total 24	C 19	F 1	N 1	O 3	0	0

- Molecule 8 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total	Cl	0	0
			1	1		

- Molecule 9 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	1	Total	Na	0	0
			1	1		

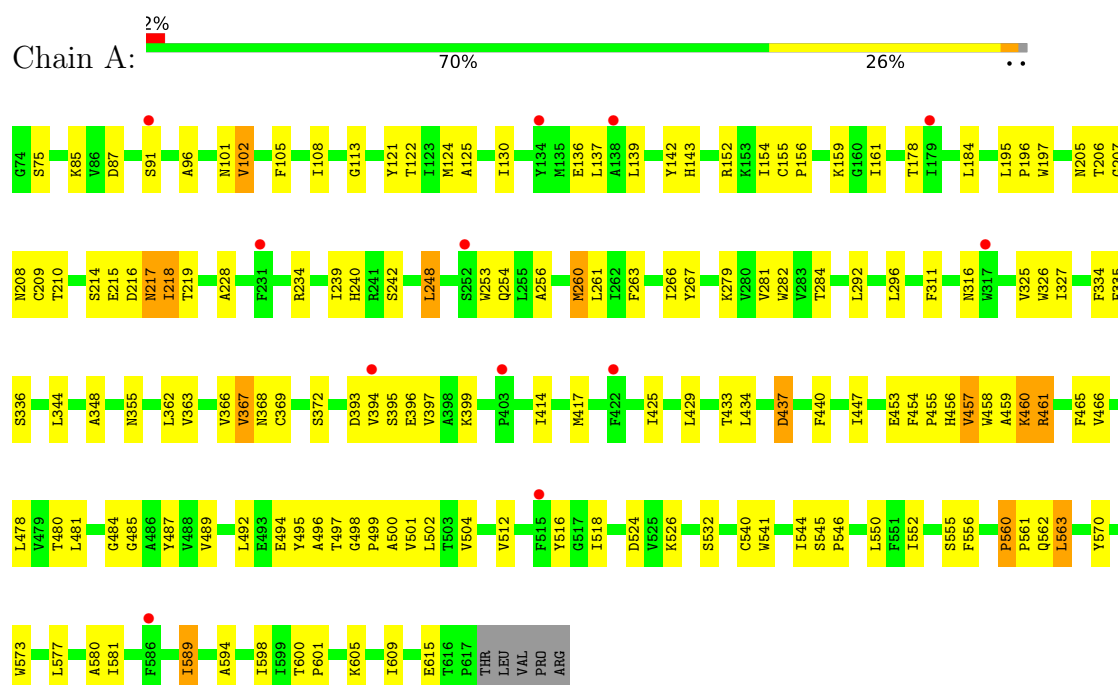
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	1	Total	O	0	0
			1	1		

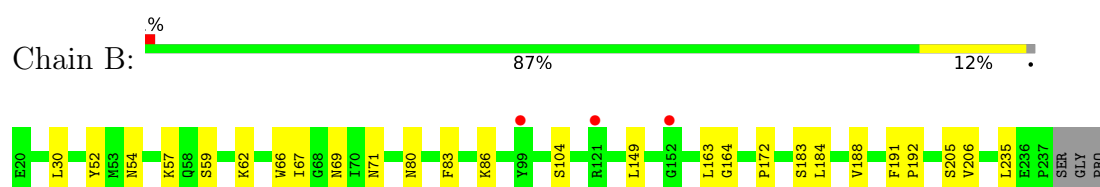
3 Residue-property plots

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.

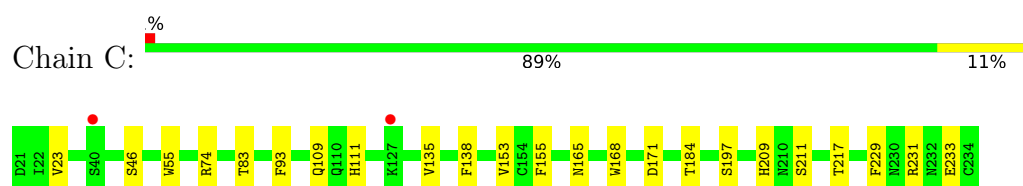
• Molecule 1: Sodium-dependent serotonin transporter



• Molecule 2: 8B6 antibody, heavy chain



• Molecule 3: 8B6 antibody, light chain



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	129.15Å 162.83Å 140.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	53.17 – 3.14 53.17 – 3.14	Depositor EDS
% Data completeness (in resolution range)	99.9 (53.17-3.14) 99.8 (53.17-3.14)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 3.13Å)	Xtriage
Refinement program	PHENIX (DEV_2000: ???)	Depositor
R, R_{free}	0.238 , 0.270 0.241 , 0.279	Depositor DCC
R_{free} test set	1307 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	105.9	Xtriage
Anisotropy	0.134	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 104.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7631	wwPDB-VP
Average B, all atoms (Å ²)	156.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% 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: NA, CLR, CL, 8PR, NAG, LMT

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	0.49	2/4355 (0.0%)	0.82	10/5955 (0.2%)
2	B	0.26	0/1688	0.67	0/2309
3	C	0.25	0/1700	0.64	0/2307
All	All	0.40	2/7743 (0.0%)	0.75	10/10571 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	560	PRO	CA-C	-5.94	1.48	1.51
1	A	560	PRO	C-O	-5.08	1.20	1.25

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	102	VAL	N-CA-C	9.24	120.06	110.72
1	A	369	CYS	N-CA-C	9.09	120.96	111.14
1	A	457	VAL	N-CA-C	-6.50	102.53	111.89
1	A	496	ALA	N-CA-C	6.06	117.76	111.03
1	A	562	GLN	N-CA-C	-5.91	101.10	109.96
1	A	75	SER	CB-CA-C	-5.63	110.06	116.54
1	A	589	ILE	CA-C-N	-5.57	112.88	119.84
1	A	589	ILE	C-N-CA	-5.57	112.88	119.84
1	A	367	VAL	N-CA-C	5.43	115.64	110.53
1	A	218	ILE	O-C-N	5.31	129.21	122.57

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	4221	0	4115	155	0
2	B	1643	0	1589	15	0
3	C	1662	0	1585	12	0
4	A	28	0	26	4	0
5	A	28	0	46	4	0
6	A	22	0	20	5	0
7	A	24	0	20	4	0
8	A	1	0	0	1	0
9	A	1	0	0	0	0
10	A	1	0	0	1	0
All	All	7631	0	7401	186	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 12.

All (186) 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:461:ARG:HH11	1:A:461:ARG:HG3	1.30	0.97
1:A:563:LEU:HD12	1:A:563:LEU:H	1.32	0.94
1:A:573:TRP:CE3	5:A:702:CLR:H71	2.03	0.93
1:A:393:ASP:OD2	1:A:396:GLU:OE2	1.91	0.88
1:A:453:GLU:HG3	1:A:454:PHE:CE1	2.10	0.86
1:A:560:PRO:HA	6:A:703:LMT:H6'2	1.61	0.81
1:A:215:GLU:O	1:A:218:ILE:HG23	1.80	0.81
1:A:453:GLU:HG3	1:A:454:PHE:CD1	2.14	0.81
1:A:219:THR:OG1	4:A:704:NAG:H61	1.81	0.80
1:A:253:TRP:CZ2	1:A:254:GLN:NE2	2.52	0.78
1:A:327:ILE:HD13	1:A:555:SER:CB	2.14	0.78
1:A:393:ASP:OD2	1:A:396:GLU:HG2	1.83	0.77
1:A:327:ILE:HD13	1:A:555:SER:HB2	1.67	0.76
2:B:54:ASN:HD22	2:B:66:TRP:HE1	1.31	0.76
1:A:461:ARG:HG3	1:A:461:ARG:NH1	1.97	0.73
1:A:563:LEU:HD12	1:A:563:LEU:N	2.03	0.73
1:A:393:ASP:O	1:A:395:SER:N	2.22	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:458:TRP:CZ3	1:A:465:PHE:HB2	2.26	0.71
1:A:334:PHE:CE2	1:A:502:LEU:HD23	2.26	0.70
7:A:705:8PR:CAF	7:A:705:8PR:HAM	2.20	0.70
1:A:216:ASP:CG	1:A:217:ASN:H	1.99	0.70
1:A:142:TYR:HD2	1:A:143:HIS:CE1	2.10	0.70
1:A:208:ASN:O	1:A:210:THR:HG23	1.92	0.69
1:A:205:ASN:ND2	1:A:209:CYS:SG	2.66	0.69
1:A:101:ASN:ND2	1:A:372:SER:OG	2.26	0.69
1:A:142:TYR:HD2	1:A:143:HIS:ND1	1.92	0.68
1:A:334:PHE:CZ	1:A:502:LEU:HD23	2.29	0.68
6:A:703:LMT:O3'	6:A:703:LMT:H6D	1.93	0.67
1:A:498:GLY:N	1:A:499:PRO:CD	2.57	0.67
1:A:453:GLU:C	1:A:454:PHE:HD1	2.04	0.66
1:A:499:PRO:CD	1:A:500:ALA:H	2.09	0.66
1:A:335:PHE:HD1	7:A:705:8PR:HAB	1.61	0.65
1:A:453:GLU:HG3	1:A:454:PHE:HE1	1.62	0.64
1:A:437:ASP:OD1	1:A:437:ASP:N	2.29	0.64
1:A:497:THR:C	1:A:499:PRO:HD2	2.23	0.64
1:A:437:ASP:OD2	10:A:1001:HOH:O	2.14	0.64
1:A:215:GLU:O	1:A:218:ILE:CG2	2.47	0.63
2:B:192:PRO:HD3	3:C:184:THR:HG22	1.81	0.63
1:A:393:ASP:OD2	1:A:396:GLU:CG	2.47	0.63
1:A:142:TYR:CD2	1:A:143:HIS:CE1	2.87	0.62
1:A:363:VAL:O	1:A:367:VAL:HG23	1.98	0.62
1:A:573:TRP:CZ3	5:A:702:CLR:H71	2.33	0.62
1:A:393:ASP:OD2	1:A:396:GLU:CD	2.43	0.62
3:C:171:ASP:OD2	3:C:209:HIS:ND1	2.30	0.61
1:A:499:PRO:HD2	1:A:500:ALA:H	1.65	0.61
1:A:334:PHE:CE2	1:A:502:LEU:CD2	2.83	0.61
1:A:396:GLU:OE1	1:A:399:LYS:HE3	2.01	0.61
1:A:447:ILE:HD13	1:A:466:VAL:HG22	1.83	0.60
1:A:113:GLY:HA2	1:A:316:ASN:HB3	1.82	0.60
1:A:552:ILE:HG22	1:A:556:PHE:CE1	2.36	0.60
1:A:454:PHE:O	1:A:458:TRP:HB2	2.01	0.60
1:A:455:PRO:O	1:A:457:VAL:N	2.27	0.60
1:A:91:SER:HA	1:A:281:VAL:HG11	1.83	0.59
1:A:429:LEU:O	1:A:433:THR:OG1	2.16	0.59
1:A:526:LYS:HD3	1:A:532:SER:HB2	1.85	0.58
1:A:85:LYS:NZ	1:A:355:ASN:OD1	2.35	0.57
3:C:74:ARG:HH21	3:C:83:THR:HG22	1.70	0.57
1:A:499:PRO:CG	1:A:500:ALA:N	2.67	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:LEU:HD21	1:A:261:LEU:HD23	1.86	0.56
1:A:546:PRO:O	1:A:550:LEU:HB2	2.05	0.56
1:A:292:LEU:O	1:A:296:LEU:HB2	2.06	0.56
1:A:394:VAL:O	1:A:397:VAL:HG12	2.07	0.55
1:A:216:ASP:OD1	1:A:217:ASN:N	2.37	0.55
2:B:86:LYS:NZ	2:B:104:SER:O	2.39	0.55
2:B:30:LEU:HD22	2:B:172:PRO:HD3	1.89	0.55
1:A:334:PHE:CD2	1:A:502:LEU:HD21	2.43	0.54
1:A:495:TYR:HE1	1:A:561:PRO:HD2	1.71	0.54
1:A:556:PHE:CD2	6:A:703:LMT:H4B	2.41	0.54
2:B:188:VAL:HG22	2:B:206:VAL:HG23	1.89	0.54
1:A:502:LEU:HD13	1:A:552:ILE:HB	1.91	0.53
1:A:125:ALA:O	1:A:130:ILE:HG12	2.08	0.53
1:A:501:VAL:HA	1:A:504:VAL:HG12	1.90	0.53
1:A:498:GLY:N	1:A:499:PRO:HD3	2.24	0.53
1:A:85:LYS:HE2	1:A:355:ASN:HD21	1.74	0.53
1:A:499:PRO:CG	1:A:500:ALA:H	2.22	0.53
1:A:453:GLU:HG3	1:A:454:PHE:HD1	1.70	0.52
1:A:492:LEU:O	1:A:497:THR:HG23	2.10	0.52
1:A:195:LEU:HD12	1:A:195:LEU:H	1.73	0.52
6:A:703:LMT:O2B	6:A:703:LMT:H4'	2.09	0.52
7:A:705:8PR:HAH	7:A:705:8PR:HAF	1.90	0.52
3:C:55:TRP:CD2	3:C:93:PHE:HB2	2.45	0.52
1:A:216:ASP:CG	1:A:217:ASN:N	2.67	0.52
1:A:499:PRO:HG2	1:A:500:ALA:N	2.25	0.52
1:A:396:GLU:OE1	1:A:396:GLU:HA	2.10	0.52
1:A:362:LEU:O	1:A:366:VAL:HG23	2.10	0.51
1:A:453:GLU:C	1:A:454:PHE:CD1	2.86	0.51
1:A:178:THR:HG21	1:A:480:THR:HB	1.92	0.51
1:A:393:ASP:O	1:A:394:VAL:C	2.51	0.51
1:A:136:GLU:HG2	1:A:344:LEU:HB2	1.92	0.51
3:C:111:HIS:O	3:C:111:HIS:ND1	2.43	0.51
1:A:414:ILE:HA	1:A:417:MET:HE2	1.93	0.51
1:A:516:TYR:OH	1:A:524:ASP:OD2	2.24	0.51
1:A:499:PRO:CD	1:A:500:ALA:N	2.73	0.50
1:A:253:TRP:CE2	1:A:254:GLN:NE2	2.79	0.50
1:A:458:TRP:O	1:A:459:ALA:HB3	2.11	0.50
3:C:135:VAL:HA	3:C:155:PHE:O	2.10	0.50
3:C:211:SER:HA	3:C:229:PHE:O	2.12	0.50
1:A:494:GLU:O	6:A:703:LMT:O3B	2.26	0.49
1:A:206:THR:HG23	1:A:234:ARG:NH2	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:THR:O	1:A:125:ALA:HB3	2.13	0.49
2:B:59:SER:HB2	2:B:62:LYS:HB2	1.95	0.49
1:A:219:THR:OG1	4:A:704:NAG:C6	2.56	0.49
1:A:263:PHE:O	1:A:267:TYR:HB2	2.13	0.49
1:A:478:LEU:HD23	1:A:481:LEU:HD12	1.95	0.49
1:A:425:ILE:O	1:A:429:LEU:HB2	2.13	0.48
1:A:197:TRP:HB2	1:A:228:ALA:HA	1.95	0.48
1:A:137:LEU:HD12	1:A:348:ALA:HB2	1.96	0.48
1:A:605:LYS:O	1:A:609:ILE:CB	2.62	0.47
1:A:152:ARG:O	1:A:615:GLU:HG2	2.14	0.47
1:A:495:TYR:CE1	1:A:561:PRO:HD2	2.48	0.47
3:C:138:PHE:HB2	3:C:153:VAL:HB	1.95	0.47
1:A:573:TRP:CZ3	5:A:702:CLR:C7	2.98	0.47
1:A:455:PRO:O	1:A:456:HIS:HB3	2.15	0.47
1:A:455:PRO:C	1:A:457:VAL:N	2.72	0.46
1:A:256:ALA:O	1:A:260:MET:HB2	2.15	0.46
1:A:336:SER:O	7:A:705:8PR:HAIA	2.15	0.46
1:A:485:GLY:O	1:A:489:VAL:HG23	2.16	0.46
1:A:393:ASP:C	1:A:395:SER:N	2.71	0.46
1:A:87:ASP:OD1	1:A:282:TRP:NE1	2.39	0.46
1:A:455:PRO:C	1:A:457:VAL:H	2.19	0.46
1:A:594:ALA:O	1:A:598:ILE:CB	2.64	0.46
1:A:552:ILE:CG2	1:A:556:PHE:CE1	2.98	0.46
3:C:168:TRP:NE1	3:C:197:SER:OG	2.41	0.46
1:A:396:GLU:HA	1:A:399:LYS:HG3	1.98	0.45
1:A:447:ILE:HA	1:A:465:PHE:HE2	1.81	0.45
1:A:121:TYR:OH	8:A:706:CL:CL	2.56	0.45
1:A:459:ALA:O	1:A:460:LYS:C	2.59	0.45
1:A:484:GLY:O	1:A:487:TYR:HB2	2.16	0.45
1:A:497:THR:C	1:A:499:PRO:CD	2.89	0.45
2:B:80:ASN:HB3	2:B:83:PHE:HD2	1.81	0.45
3:C:23:VAL:H	3:C:46:SER:HB3	1.81	0.45
3:C:231:ARG:O	3:C:233:GLU:N	2.47	0.45
1:A:497:THR:O	1:A:499:PRO:HD2	2.16	0.45
1:A:207:GLY:HA3	4:A:701:NAG:H82	1.98	0.45
1:A:577:LEU:O	1:A:581:ILE:HG13	2.17	0.45
1:A:456:HIS:C	1:A:458:TRP:N	2.69	0.45
2:B:149:LEU:HB2	2:B:164:GLY:C	2.42	0.44
1:A:266:ILE:HG22	1:A:440:PHE:HE1	1.82	0.44
1:A:499:PRO:HG2	1:A:500:ALA:H	1.81	0.44
2:B:163:LEU:O	2:B:205:SER:HA	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:CYS:HA	1:A:156:PRO:HD2	1.88	0.44
1:A:266:ILE:HG22	1:A:440:PHE:CE1	2.53	0.43
1:A:396:GLU:OE1	1:A:399:LYS:CE	2.66	0.43
1:A:540:CYS:HA	1:A:544:ILE:HB	1.99	0.43
1:A:219:THR:CG2	4:A:704:NAG:H61	2.48	0.43
1:A:279:LYS:O	1:A:282:TRP:HB2	2.19	0.43
1:A:545:SER:HB2	1:A:546:PRO:HD3	2.00	0.43
1:A:216:ASP:C	1:A:218:ILE:H	2.26	0.43
2:B:57:LYS:HB2	2:B:67:ILE:HD11	2.01	0.43
1:A:124:MET:HE2	1:A:326:TRP:HE3	1.83	0.43
5:A:702:CLR:H162	5:A:702:CLR:H231	2.00	0.43
1:A:454:PHE:CD1	1:A:454:PHE:N	2.87	0.43
1:A:460:LYS:HG2	1:A:461:ARG:H	1.83	0.43
1:A:214:SER:OG	1:A:215:GLU:N	2.51	0.42
1:A:139:LEU:HD11	1:A:512:VAL:HG11	2.01	0.42
1:A:195:LEU:HA	1:A:196:PRO:HD3	1.84	0.42
1:A:253:TRP:CH2	1:A:254:GLN:NE2	2.88	0.42
2:B:52:TYR:HE1	2:B:71:ASN:HB2	1.85	0.42
1:A:263:PHE:HD1	1:A:266:ILE:HD11	1.85	0.42
1:A:325:VAL:HG23	1:A:326:TRP:N	2.35	0.42
3:C:165:ASN:HB3	3:C:217:THR:HB	2.01	0.42
1:A:156:PRO:O	1:A:159:LYS:HB2	2.20	0.41
1:A:281:VAL:HA	1:A:284:THR:OG1	2.20	0.41
1:A:425:ILE:O	1:A:429:LEU:CB	2.67	0.41
2:B:191:PHE:HA	2:B:192:PRO:HD3	1.87	0.41
1:A:102:VAL:HG21	1:A:434:LEU:CD1	2.50	0.41
2:B:183:SER:HA	2:B:184:LEU:HA	1.51	0.41
1:A:96:ALA:O	1:A:368:ASN:OD1	2.39	0.41
1:A:242:SER:HB2	1:A:248:LEU:HA	2.03	0.41
1:A:234:ARG:HE	1:A:240:HIS:CG	2.38	0.41
1:A:461:ARG:HH11	1:A:461:ARG:CG	2.12	0.41
1:A:518:ILE:HD11	1:A:541:TRP:CE3	2.55	0.41
1:A:105:PHE:HA	1:A:108:ILE:HG22	2.02	0.41
1:A:577:LEU:O	1:A:580:ALA:HB3	2.21	0.41
1:A:600:THR:HA	1:A:601:PRO:HD3	1.85	0.41
2:B:54:ASN:ND2	2:B:69:ASN:HB3	2.35	0.41
1:A:143:HIS:NE2	1:A:154:ILE:HD11	2.36	0.41
1:A:239:ILE:HD11	1:A:487:TYR:CE1	2.56	0.41
1:A:263:PHE:HA	1:A:266:ILE:HG12	2.03	0.40
2:B:163:LEU:HB3	2:B:235:LEU:HD22	2.03	0.40
1:A:161:ILE:HG13	1:A:589:ILE:HD12	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:454:PHE:HD1	1:A:454:PHE:N	2.19	0.40
1:A:101:ASN:HD21	1:A:368:ASN:ND2	2.18	0.40
1:A:458:TRP:O	1:A:459:ALA:CB	2.70	0.40
1:A:311:PHE:CZ	1:A:395:SER:HB2	2.56	0.40
1:A:563:LEU:HD13	1:A:570:TYR:CG	2.57	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	542/549 (99%)	521 (96%)	21 (4%)	0	100	100
2	B	216/221 (98%)	207 (96%)	9 (4%)	0	100	100
3	C	212/214 (99%)	200 (94%)	12 (6%)	0	100	100
All	All	970/984 (99%)	928 (96%)	42 (4%)	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	430/462 (93%)	423 (98%)	7 (2%)	55	71
2	B	190/193 (98%)	190 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	C	189/190 (100%)	188 (100%)	1 (0%)	81	82
All	All	809/845 (96%)	801 (99%)	8 (1%)	68	76

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

Mol	Chain	Res	Type
1	A	217	ASN
1	A	248	LEU
1	A	260	MET
1	A	437	ASP
1	A	460	LYS
1	A	461	ARG
1	A	563	LEU
3	C	109	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	141	GLN
1	A	351	ASN
2	B	58	GLN
3	C	28	HIS
3	C	58	GLN
3	C	157	ASN
3	C	158	ASN
3	C	230	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

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
4	NAG	A	701	1	14,14,15	0.27	0	17,19,21	0.63	0
5	CLR	A	702	-	31,31,31	0.86	1 (3%)	48,48,48	1.36	6 (12%)
4	NAG	A	704	1	14,14,15	0.25	0	17,19,21	0.39	0
6	LMT	A	703	-	23,23,36	0.42	0	33,33,47	1.34	4 (12%)
7	8PR	A	705	-	27,27,27	1.94	6 (22%)	33,37,37	1.71	7 (21%)

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
4	NAG	A	701	1	-	2/6/23/26	0/1/1/1
5	CLR	A	702	-	-	6/10/68/68	0/4/4/4
4	NAG	A	704	1	-	1/6/23/26	0/1/1/1
6	LMT	A	703	-	-	7/8/45/61	0/2/2/2
7	8PR	A	705	-	-	4/9/26/26	0/4/4/4

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	705	8PR	CAT-CAX	4.66	1.58	1.51
7	A	705	8PR	CAJ-CAX	-3.92	1.48	1.53
7	A	705	8PR	OAO-CAS	3.87	1.46	1.37
7	A	705	8PR	OAP-CAU	3.37	1.43	1.38
7	A	705	8PR	CAX-CAW	-3.02	1.47	1.53
7	A	705	8PR	OAQ-CAV	2.95	1.43	1.38
5	A	702	CLR	C13-C14	-2.06	1.51	1.55

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	705	8PR	OAQ-CAS-CAH	-3.82	101.34	119.74
7	A	705	8PR	OAQ-CAS-CAD	3.58	137.66	120.01
6	A	703	LMT	C1B-O5B-C5B	3.48	120.51	113.72
7	A	705	8PR	OAQ-CAV-CAH	3.41	132.39	127.86
5	A	702	CLR	C8-C7-C6	-3.21	108.31	112.76
6	A	703	LMT	C1B-C2B-C3B	3.18	116.70	110.01
5	A	702	CLR	C9-C10-C5	3.14	114.25	109.65
6	A	703	LMT	O5B-C1B-C2B	2.77	116.06	110.37
5	A	702	CLR	C13-C14-C8	-2.75	110.50	114.41
7	A	705	8PR	CAC-CAR-CAB	-2.62	119.36	122.80
7	A	705	8PR	CAE-CAB-CAR	2.58	121.03	118.38
5	A	702	CLR	C4-C5-C10	2.42	119.52	116.42
7	A	705	8PR	OAP-CAU-CAG	2.22	132.00	127.72
7	A	705	8PR	OAQ-CAV-CAU	-2.19	107.14	109.79
6	A	703	LMT	C3B-C4B-C5B	-2.17	106.30	110.23
5	A	702	CLR	C13-C17-C20	-2.02	116.38	119.50
5	A	702	CLR	C16-C17-C20	-2.01	109.14	112.18

There are no chirality outliers.

All (20) torsion outliers are listed below:

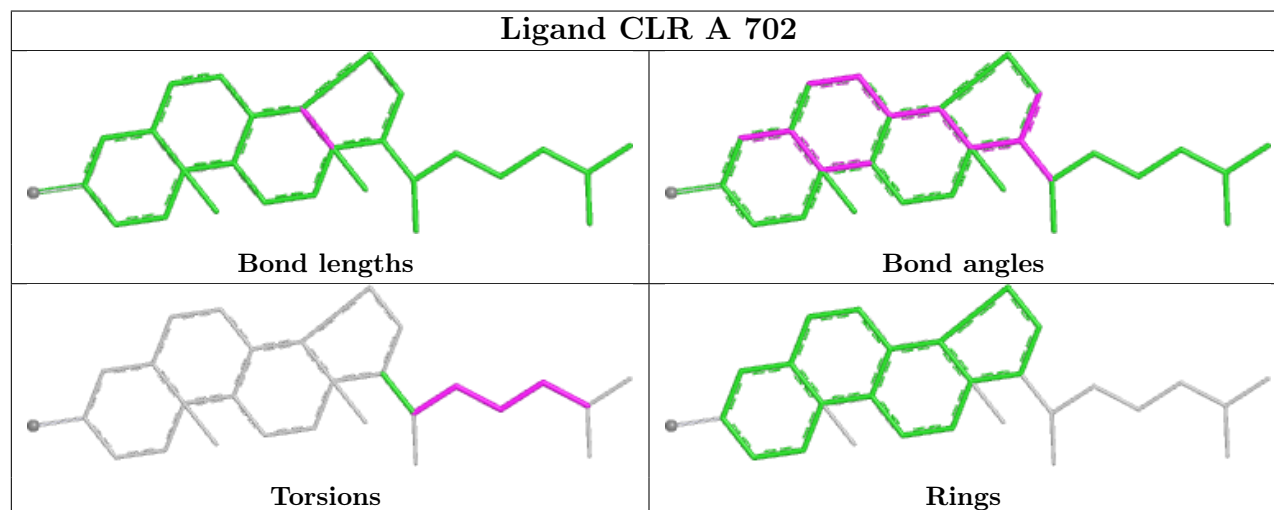
Mol	Chain	Res	Type	Atoms
7	A	705	8PR	OAQ-CAM-CAW-CAK
6	A	703	LMT	O5B-C1B-O1B-C4'
6	A	703	LMT	C2B-C1B-O1B-C4'
4	A	701	NAG	O5-C5-C6-O6
6	A	703	LMT	O5'-C5'-C6'-O6'
4	A	701	NAG	C4-C5-C6-O6
6	A	703	LMT	C3'-C4'-O1B-C1B
7	A	705	8PR	CAD-CAS-OAQ-CAM
7	A	705	8PR	CAH-CAS-OAQ-CAM
6	A	703	LMT	C4'-C5'-C6'-O6'
5	A	702	CLR	C17-C20-C22-C23
6	A	703	LMT	C5'-C4'-O1B-C1B
5	A	702	CLR	C21-C20-C22-C23
6	A	703	LMT	O5B-C5B-C6B-O6B
4	A	704	NAG	O5-C5-C6-O6
5	A	702	CLR	C20-C22-C23-C24
5	A	702	CLR	C23-C24-C25-C26
5	A	702	CLR	C23-C24-C25-C27
7	A	705	8PR	OAQ-CAM-CAW-CAX
5	A	702	CLR	C22-C23-C24-C25

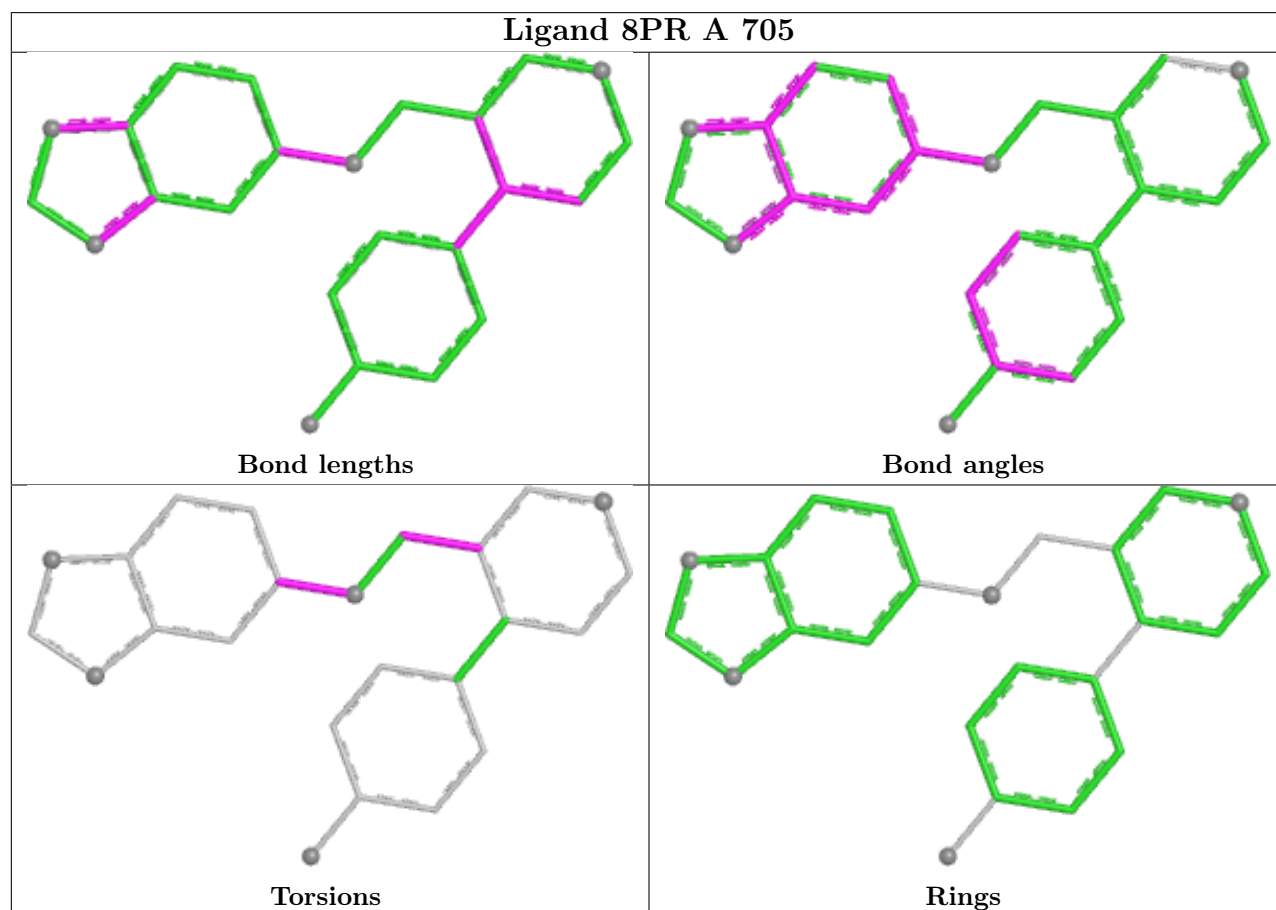
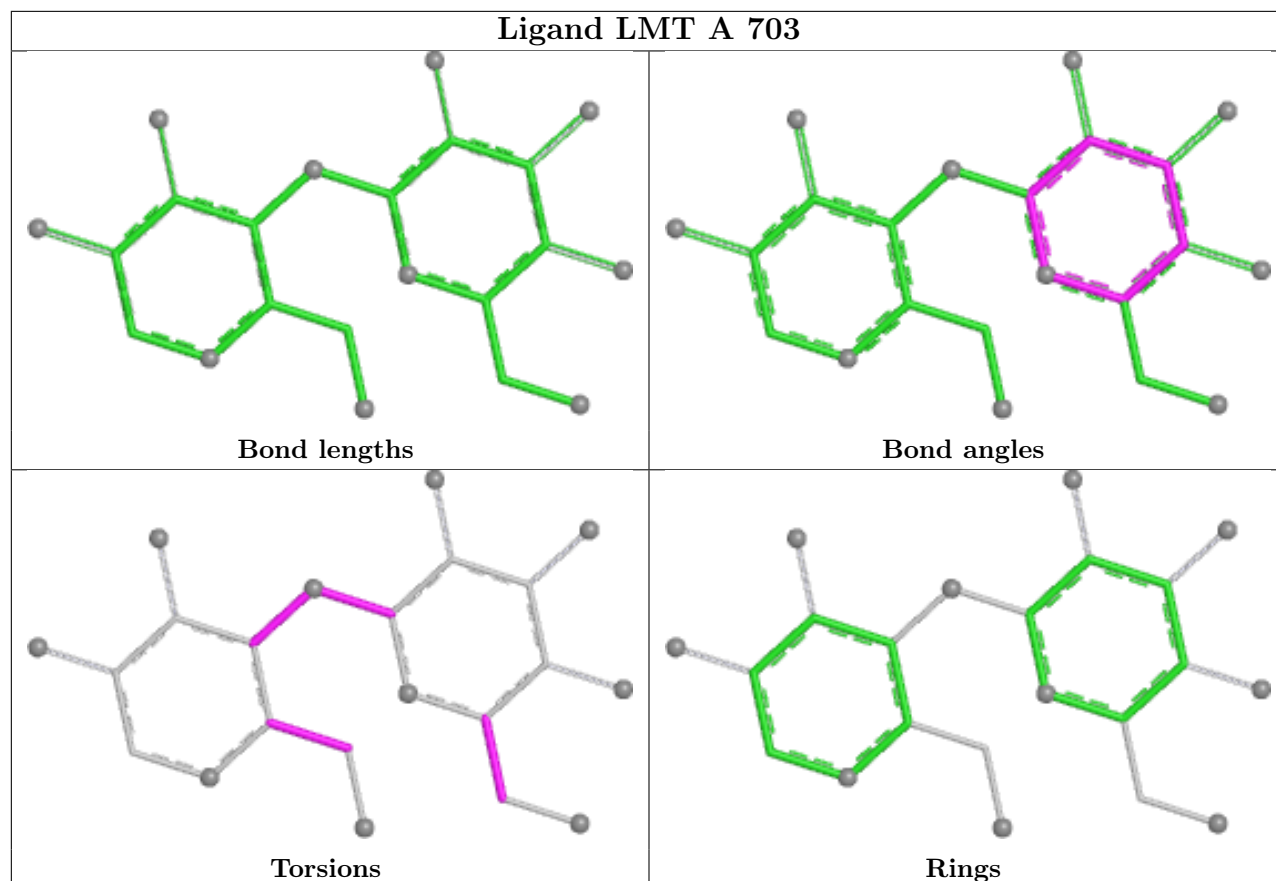
There are no ring outliers.

5 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	701	NAG	1	0
5	A	702	CLR	4	0
4	A	704	NAG	3	0
6	A	703	LMT	5	0
7	A	705	8PR	4	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 [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, 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	544/549 (99%)	-0.14	12 (2%) 62 41	72, 131, 218, 322	0
2	B	218/221 (98%)	-0.25	3 (1%) 73 53	73, 161, 330, 391	0
3	C	214/214 (100%)	-0.14	2 (0%) 81 63	85, 176, 282, 408	0
All	All	976/984 (99%)	-0.17	17 (1%) 69 48	72, 145, 269, 408	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	134	TYR	4.5
1	A	138	ALA	4.0
2	B	99	TYR	3.9
3	C	40	SER	3.6
1	A	515	PHE	2.8
1	A	422	PHE	2.7
1	A	91	SER	2.5
3	C	127	LYS	2.5
1	A	586	PHE	2.4
1	A	403	PRO	2.4
1	A	231	PHE	2.4
1	A	252	SER	2.3
1	A	394	VAL	2.3
2	B	152	GLY	2.2
2	B	121	ARG	2.1
1	A	179	ILE	2.1
1	A	317	TRP	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

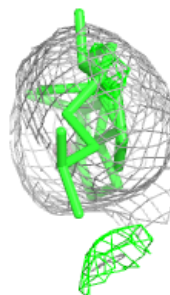
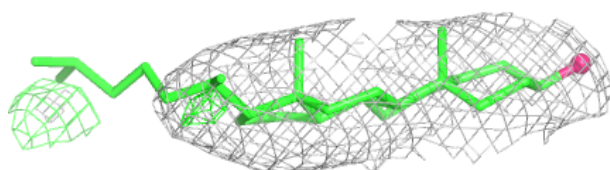
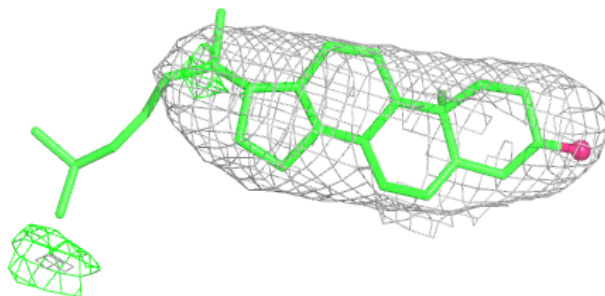
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAG	A	704	14/15	0.85	0.15	274,279,284,284	0
5	CLR	A	702	28/28	0.88	0.17	184,199,204,209	0
6	LMT	A	703	22/35	0.88	0.08	113,132,162,167	0
4	NAG	A	701	14/15	0.90	0.10	64,83,132,137	0
8	CL	A	706	1/1	0.93	0.10	162,162,162,162	0
7	8PR	A	705	24/24	0.94	0.08	85,98,114,119	0
9	NA	A	707	1/1	0.95	0.07	51,51,51,51	1

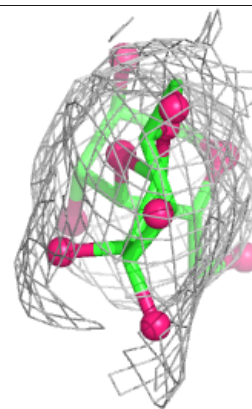
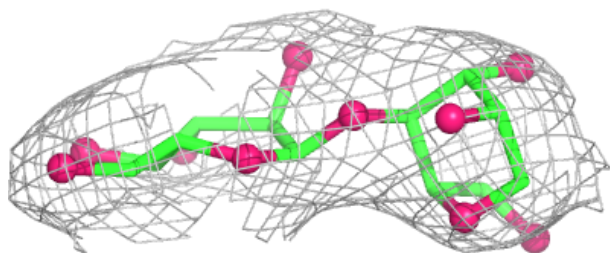
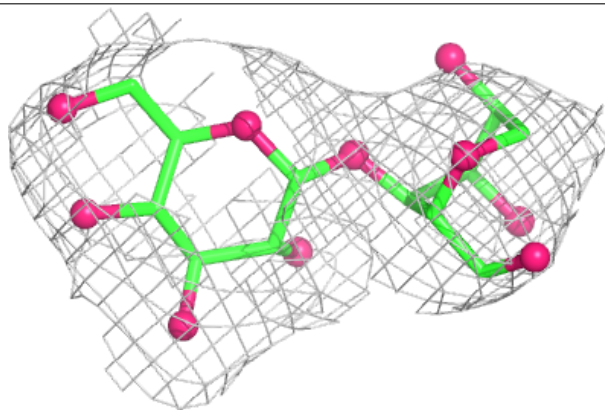
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around CLR A 702:

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

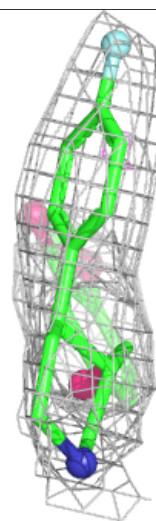
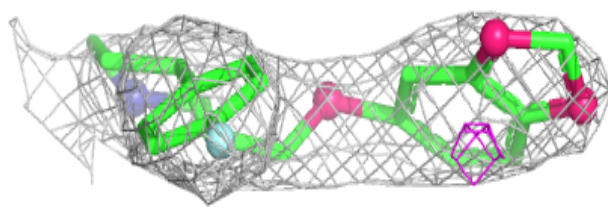
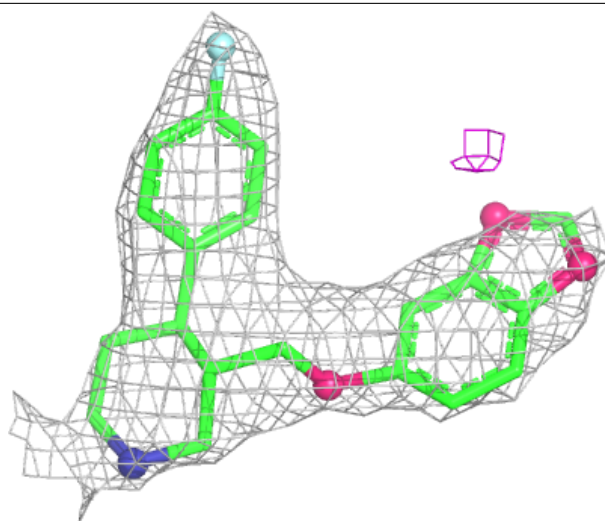
**Electron density around LMT A 703:**

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



Electron density around 8PR A 705:

$2mF_o-DF_c$ (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.