



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:23 AM UTC

PDB ID : 3LKH / pdb_00003lkh
Title : Inhibitors of Hepatitis C Virus Polymerase: Synthesis and Characterization of Novel 6-Fluoro-N-[2-Hydroxy-1(S)-Benzamides
Authors : Lesburg, C.A.
Deposited on : 2010-01-27
Resolution : 2.05 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

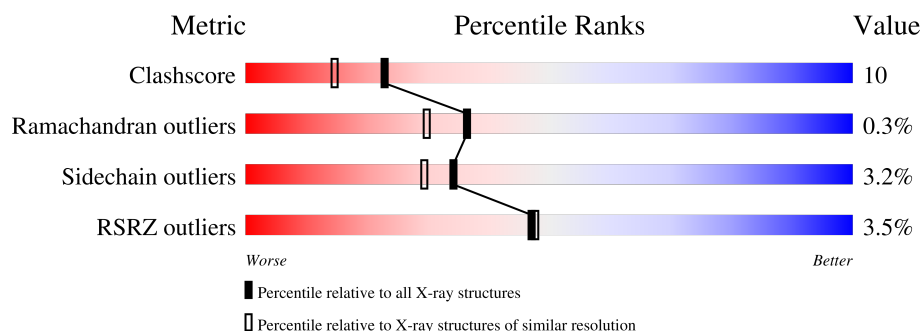
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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))
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	576	2% 77% 17% . .
1	B	576	5% 77% 17% . . .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9891 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.

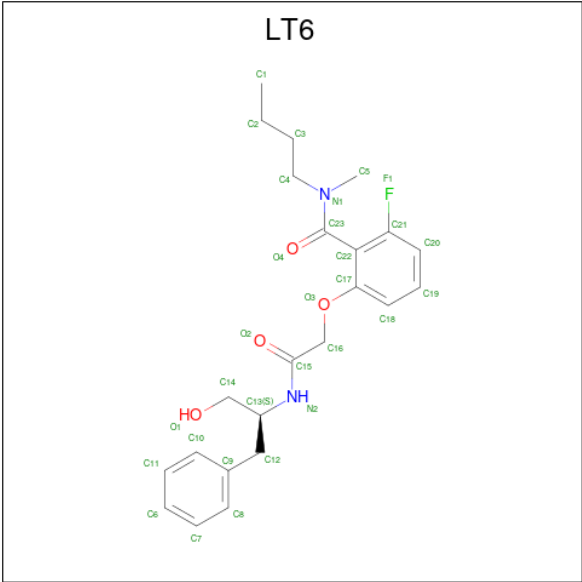
- Molecule 1 is a protein called RNA-directed RNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	554	Total	C	N	O	S	0	3	0
			4340	2734	767	804	35			
1	B	558	Total	C	N	O	S	0	4	0
			4371	2751	773	811	36			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	440	GLY	GLU	conflict	UNP O92972
A	520	ILE	THR	conflict	UNP O92972
A	571	GLU	-	expression tag	UNP O92972
A	572	ASN	-	expression tag	UNP O92972
A	573	LEU	-	expression tag	UNP O92972
A	574	TYR	-	expression tag	UNP O92972
A	575	PHE	-	expression tag	UNP O92972
A	576	GLN	-	expression tag	UNP O92972
B	440	GLY	GLU	conflict	UNP O92972
B	520	ILE	THR	conflict	UNP O92972
B	571	GLU	-	expression tag	UNP O92972
B	572	ASN	-	expression tag	UNP O92972
B	573	LEU	-	expression tag	UNP O92972
B	574	TYR	-	expression tag	UNP O92972
B	575	PHE	-	expression tag	UNP O92972
B	576	GLN	-	expression tag	UNP O92972

- Molecule 2 is 2-(2-{[(1S)-1-benzyl-2-hydroxyethyl]amino}-2-oxoethoxy)-N-butyl-6-fluoro-N-methylbenzamide (CCD ID: LT6) (formula: C₂₃H₂₉FN₂O₄).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			30	23	1	2	4		
2	B	1	Total	C	F	N	O	0	0
			30	23	1	2	4		

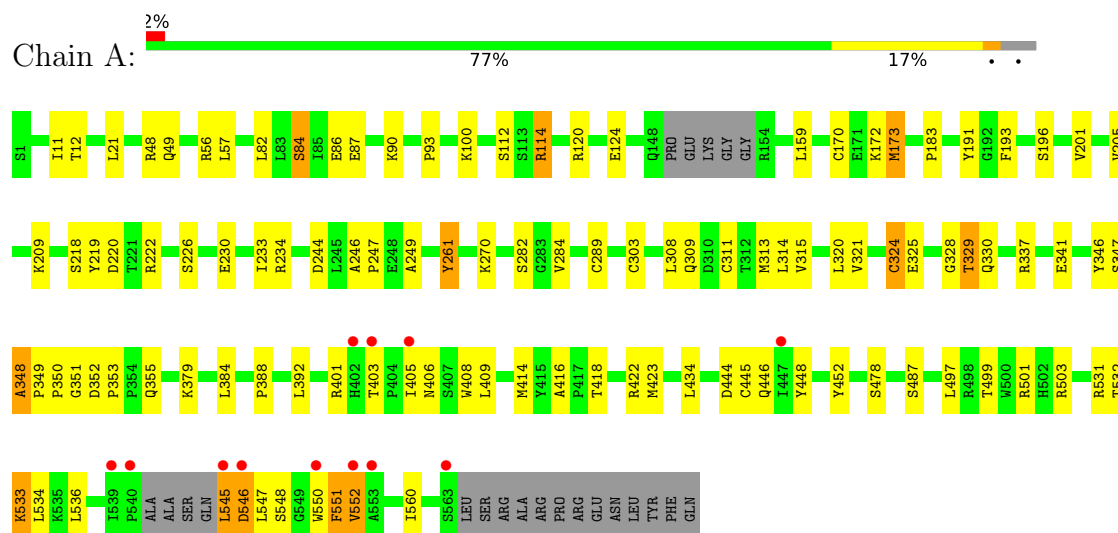
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	613	Total	O	0	0
			613	613		
3	B	507	Total	O	0	0
			507	507		

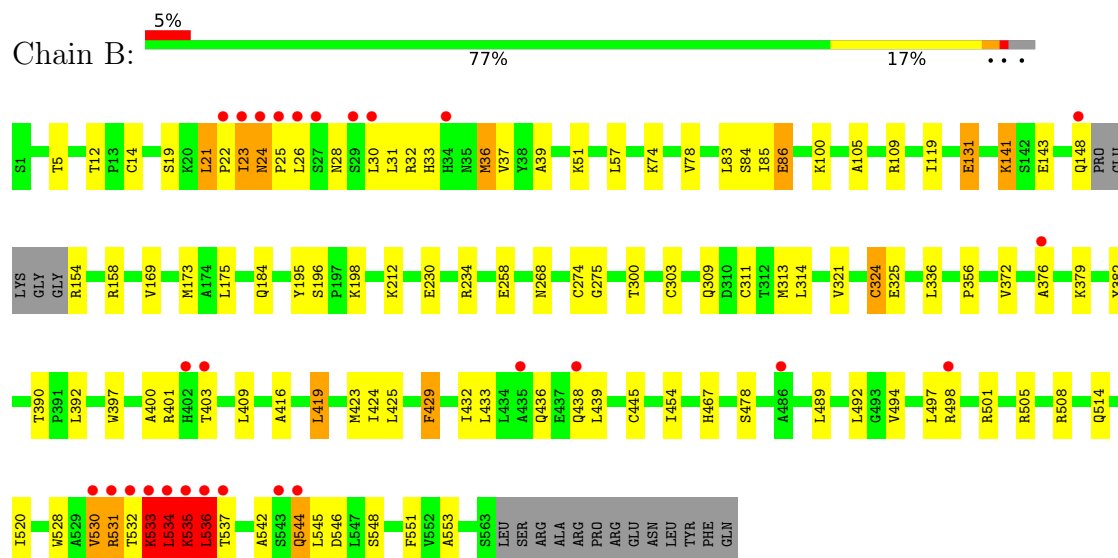
3 Residue-property plots

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: RNA-directed RNA polymerase



• Molecule 1: RNA-directed RNA polymerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.70Å 107.07Å 134.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.05 20.00 – 2.05	Depositor EDS
% Data completeness (in resolution range)	97.8 (20.00-2.05) 94.1 (20.00-2.05)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 2.05Å)	Xtriage
Refinement program	BUSTER-TNT 2.1.1	Depositor
R, R_{free}	0.184 , 0.241 0.186 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	28.6	Xtriage
Anisotropy	0.373	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 57.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9891	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.55% 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: LT6

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	0.91	1/4433 (0.0%)	1.07	28/6013 (0.5%)
1	B	0.83	0/4465	1.12	28/6058 (0.5%)
All	All	0.87	1/8898 (0.0%)	1.09	56/12071 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	173	MET	SD-CE	-8.17	1.59	1.79

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	351	GLY	N-CA-C	-12.72	98.63	110.21
1	B	535	LYS	CA-C-N	12.70	141.02	122.36
1	B	535	LYS	C-N-CA	12.70	141.02	122.36
1	B	536	LEU	N-CA-C	11.96	128.66	111.56
1	B	534	LEU	N-CA-C	10.74	133.67	110.80
1	B	403	THR	CA-C-N	10.66	130.32	119.24
1	B	403	THR	C-N-CA	10.66	130.32	119.24
1	B	535	LYS	N-CA-C	9.13	123.18	108.76
1	B	537	THR	CA-C-N	-8.20	111.57	119.85
1	B	537	THR	C-N-CA	-8.20	111.57	119.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	348	ALA	CA-C-N	7.62	128.22	120.38
1	A	348	ALA	C-N-CA	7.62	128.22	120.38
1	A	329	THR	N-CA-C	7.16	119.16	111.36
1	B	551	PHE	N-CA-C	6.93	119.86	111.82
1	B	24	ASN	N-CA-C	6.81	115.97	108.14
1	A	308	LEU	N-CA-C	-6.80	101.56	110.53
1	B	429	PHE	N-CA-C	6.80	118.77	111.36
1	B	23	ILE	N-CA-C	6.64	117.31	107.15
1	A	244	ASP	N-CA-C	-6.61	99.25	109.76
1	A	353	PRO	CA-C-N	-6.57	113.18	120.14
1	A	353	PRO	C-N-CA	-6.57	113.18	120.14
1	A	551	PHE	N-CA-C	6.52	120.45	112.23
1	B	545	LEU	N-CA-C	6.49	119.78	110.24
1	B	12	THR	CA-C-N	6.37	126.56	120.31
1	B	12	THR	C-N-CA	6.37	126.56	120.31
1	B	175	LEU	N-CA-C	6.08	122.57	114.12
1	B	195	TYR	N-CA-C	5.93	119.31	109.46
1	B	12	THR	N-CA-C	5.92	118.09	110.39
1	B	533	LYS	CA-C-N	5.92	132.84	121.54
1	B	533	LYS	C-N-CA	5.92	132.84	121.54
1	B	131	GLU	N-CA-C	5.88	123.33	110.80
1	B	478	SER	CA-C-N	5.87	125.08	118.97
1	B	478	SER	C-N-CA	5.87	125.08	118.97
1	A	261	TYR	N-CA-C	5.81	117.28	111.07
1	A	93	PRO	CA-C-N	5.71	126.10	119.47
1	A	93	PRO	C-N-CA	5.71	126.10	119.47
1	A	408	TRP	N-CA-C	5.60	117.38	111.28
1	A	355	GLN	CA-C-N	-5.50	114.31	120.14
1	A	355	GLN	C-N-CA	-5.50	114.31	120.14
1	B	425	LEU	N-CA-C	5.50	117.27	111.28
1	A	84	SER	N-CA-C	-5.45	103.50	110.53
1	A	315	VAL	N-CA-C	5.44	115.73	108.11
1	A	324	CYS	N-CA-C	5.36	115.36	108.34
1	A	328	GLY	N-CA-C	-5.27	106.94	111.95
1	B	196	SER	N-CA-C	-5.26	102.39	110.07
1	B	531	ARG	N-CA-C	5.25	117.79	109.96
1	A	196	SER	N-CA-C	-5.23	102.43	110.07
1	A	218	SER	N-CA-C	-5.14	102.14	110.32
1	A	246	ALA	CA-C-N	5.13	124.58	119.24
1	A	246	ALA	C-N-CA	5.13	124.58	119.24
1	A	478	SER	CA-C-N	5.11	124.56	119.24
1	A	478	SER	C-N-CA	5.11	124.56	119.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	416	ALA	N-CA-C	5.05	120.46	113.45
1	A	21	LEU	CA-C-N	-5.04	114.53	119.92
1	A	21	LEU	C-N-CA	-5.04	114.53	119.92
1	B	78	VAL	N-CA-C	5.02	116.50	108.87

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	535	LYS	Peptide,Mainchain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4340	0	4351	68	1
1	B	4371	0	4379	116	1
2	A	30	0	29	3	0
2	B	30	0	29	0	0
3	A	613	0	0	17	1
3	B	507	0	0	19	1
All	All	9891	0	8788	181	2

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 10.

All (181) 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:423:MET:HE2	1:B:497:LEU:HD22	1.25	1.07
1:B:376:ALA:HB3	1:B:535:LYS:NZ	1.71	1.06
1:B:530:VAL:HG11	1:B:533:LYS:HE3	1.52	0.91
1:B:119:ILE:CG2	1:B:173:MET:HE1	2.03	0.89
1:B:22:PRO:HD3	1:B:401:ARG:NE	1.88	0.88
1:B:501:ARG:NH2	1:B:531:ARG:HD3	1.88	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:376:ALA:HB3	1:B:535:LYS:HZ2	1.32	0.86
1:B:119:ILE:HG21	1:B:173:MET:HE1	1.57	0.85
1:B:85:ILE:HA	1:B:173:MET:HE2	1.58	0.84
1:B:409:LEU:HD23	1:B:445:CYS:SG	2.16	0.84
1:B:169:VAL:HG12	1:B:173:MET:HE3	1.58	0.83
1:B:530:VAL:CG1	1:B:533:LYS:HE3	2.09	0.83
1:A:124:GLU:HG3	3:A:1074:HOH:O	1.78	0.82
1:B:535:LYS:HE3	1:B:536:LEU:HG	1.62	0.81
1:B:303[B]:CYS:SG	1:B:313:MET:HE1	2.21	0.81
1:B:423:MET:HE2	1:B:497:LEU:CD2	2.08	0.81
1:B:234:ARG:NH1	1:B:258:GLU:OE2	2.15	0.80
1:B:501:ARG:NH2	1:B:531:ARG:HH11	1.80	0.80
1:B:501:ARG:NE	1:B:505:ARG:NH2	2.31	0.80
1:B:21:LEU:CD1	1:B:37:VAL:HB	2.12	0.79
1:B:86:GLU:H	1:B:86:GLU:CD	1.91	0.79
1:B:423:MET:CE	1:B:497:LEU:HB3	2.13	0.79
1:B:230:GLU:HG2	3:B:909:HOH:O	1.80	0.78
1:B:22:PRO:HG3	1:B:401:ARG:HG2	1.65	0.78
1:A:414:MET:HE3	2:A:577:LT6:C9	2.15	0.77
1:A:534:LEU:HD21	3:A:1011:HOH:O	1.83	0.77
1:B:542:ALA:HB1	1:B:544:GLN:OE1	1.84	0.77
1:A:230:GLU:HG2	3:A:612:HOH:O	1.85	0.76
1:B:553:ALA:HB3	3:B:755:HOH:O	1.85	0.76
1:A:170:CYS:HA	1:A:173:MET:HE3	1.67	0.75
1:B:501:ARG:HH22	1:B:531:ARG:HH11	1.32	0.75
1:B:24:ASN:HB3	1:B:25:PRO:HD2	1.69	0.75
1:B:376:ALA:HB3	1:B:535:LYS:HZ3	1.50	0.75
1:B:501:ARG:HH21	1:B:531:ARG:HD3	1.50	0.74
1:A:531:ARG:O	1:A:533:LYS:HG2	1.87	0.74
1:A:409:LEU:HD23	1:A:445:CYS:SG	2.28	0.72
1:B:534:LEU:C	1:B:535:LYS:HG2	2.14	0.71
1:A:100:LYS:HE3	3:A:703:HOH:O	1.89	0.71
1:B:154:ARG:HD2	3:B:1065:HOH:O	1.90	0.70
1:B:21:LEU:HD12	1:B:37:VAL:HB	1.73	0.70
1:A:311[A]:CYS:HA	1:A:324:CYS:HB3	1.71	0.70
1:B:498:ARG:NH1	3:B:954:HOH:O	2.25	0.69
1:B:33:HIS:HB3	1:B:36:MET:HG3	1.75	0.69
1:B:423:MET:HE1	1:B:497:LEU:HB3	1.76	0.68
1:B:131:GLU:HG2	3:B:983:HOH:O	1.94	0.67
1:A:172:LYS:HE3	1:A:560:ILE:HD13	1.75	0.67
1:A:247:PRO:HG3	1:B:234:ARG:HD3	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:501:ARG:HD3	3:A:775:HOH:O	1.95	0.66
1:B:21:LEU:HD11	1:B:37:VAL:HB	1.79	0.65
1:A:309:GLN:HG3	3:A:762:HOH:O	1.97	0.64
1:B:22:PRO:HD3	1:B:401:ARG:CZ	2.28	0.64
1:B:31:LEU:HD12	1:B:492:LEU:HB3	1.80	0.64
1:B:24:ASN:HB3	1:B:25:PRO:CD	2.28	0.64
1:A:303[B]:CYS:SG	1:A:313:MET:HE1	2.38	0.63
1:B:530:VAL:CG1	1:B:533:LYS:HB3	2.28	0.63
1:B:429:PHE:O	1:B:433:LEU:HG	1.99	0.63
1:B:501:ARG:HH22	1:B:531:ARG:NH1	1.96	0.62
1:A:550:TRP:CD1	2:A:577:LT6:H1A	2.34	0.62
1:B:22:PRO:HG2	1:B:400:ALA:HB1	1.82	0.62
1:B:74:LYS:NZ	3:B:953:HOH:O	2.27	0.62
1:B:379:LYS:HE2	3:B:606:HOH:O	1.99	0.61
1:A:325:GLU:HG3	3:A:638:HOH:O	2.00	0.61
1:A:547:LEU:HD23	1:A:547:LEU:N	2.14	0.61
1:B:433:LEU:HB3	1:B:439:LEU:HD23	1.83	0.60
1:A:531:ARG:O	1:A:533:LYS:N	2.35	0.60
1:B:501:ARG:NE	1:B:505:ARG:HH22	1.99	0.59
1:B:51:LYS:NZ	3:B:1010:HOH:O	2.29	0.59
1:B:419:LEU:HD12	1:B:423:MET:SD	2.42	0.59
1:A:547:LEU:O	1:A:551:PHE:HD1	1.86	0.58
1:A:309:GLN:O	1:A:324:CYS:HB2	2.03	0.58
1:B:534:LEU:O	1:B:535:LYS:HG2	2.03	0.58
1:B:433:LEU:CB	1:B:439:LEU:HD23	2.34	0.58
1:A:84:SER:OG	1:A:87:GLU:HG3	2.03	0.57
1:B:303[B]:CYS:HG	1:B:311[B]:CYS:HG	0.59	0.56
1:B:530:VAL:HG13	1:B:533:LYS:HB2	1.87	0.56
1:B:184:GLN:HB2	3:B:587:HOH:O	2.05	0.56
1:A:330:GLN:NE2	3:A:1110:HOH:O	2.29	0.55
1:A:337:ARG:O	1:A:341:GLU:HG3	2.06	0.55
1:A:56:ARG:HD2	1:A:226:SER:O	2.06	0.55
1:A:405:ILE:HG22	1:A:406:ASN:N	2.21	0.55
1:B:492:LEU:HB2	1:B:494:VAL:HG23	1.88	0.55
1:B:530:VAL:HG13	1:B:533:LYS:CB	2.37	0.55
1:A:170:CYS:HA	1:A:173:MET:CE	2.36	0.54
1:B:309:GLN:O	1:B:324:CYS:HB3	2.07	0.54
1:A:434:LEU:HD21	3:A:935:HOH:O	2.07	0.54
1:B:33:HIS:CB	1:B:36:MET:HG3	2.37	0.54
1:B:432:ILE:O	1:B:436:GLN:HG3	2.08	0.54
1:B:336:LEU:HD23	1:B:356:PRO:HD3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:501:ARG:NH2	1:B:531:ARG:NH1	2.53	0.53
1:B:534:LEU:HD12	3:B:820:HOH:O	2.08	0.53
1:A:452:TYR:CE2	1:A:550:TRP:NE1	2.75	0.52
1:B:520:ILE:HD13	3:B:1054:HOH:O	2.08	0.52
1:B:86:GLU:OE1	1:B:86:GLU:N	2.32	0.52
1:B:30:LEU:O	1:B:494:VAL:HG13	2.10	0.52
1:A:548:SER:HA	3:A:1115:HOH:O	2.09	0.52
1:B:5:THR:HG22	3:B:737:HOH:O	2.08	0.52
1:A:445:CYS:HB3	3:A:877:HOH:O	2.09	0.51
1:A:311[B]:CYS:HA	1:A:324:CYS:HB3	1.91	0.51
1:B:501:ARG:CZ	1:B:505:ARG:HH22	2.23	0.51
1:A:403:THR:O	1:A:405:ILE:O	2.29	0.51
1:B:86:GLU:CD	1:B:86:GLU:N	2.67	0.50
1:B:535:LYS:CE	1:B:536:LEU:HG	2.36	0.50
1:A:120:ARG:HD2	3:A:1109:HOH:O	2.12	0.50
1:B:445:CYS:HB2	1:B:454:ILE:HD12	1.94	0.50
1:A:444:ASP:HB3	3:A:931:HOH:O	2.12	0.49
1:B:31:LEU:CD1	1:B:492:LEU:HB3	2.43	0.49
1:B:530:VAL:CG1	1:B:533:LYS:CB	2.90	0.49
1:B:501:ARG:NH1	3:B:930:HOH:O	2.45	0.49
1:A:219:TYR:HB3	1:A:320:LEU:HD23	1.95	0.49
1:A:548:SER:O	1:A:552:VAL:HG13	2.13	0.49
1:B:268:ASN:HB3	1:B:274:CYS:SG	2.53	0.49
1:A:205:VAL:O	1:A:209:LYS:HG3	2.13	0.49
1:A:499:THR:O	1:A:503:ARG:HG3	2.13	0.49
1:B:424:ILE:CD1	1:B:489:LEU:HD21	2.44	0.48
1:B:546:ASP:OD1	1:B:548:SER:HB2	2.13	0.48
1:A:247:PRO:CG	1:B:234:ARG:NH1	2.77	0.48
1:B:5:THR:O	1:B:275:GLY:HA3	2.13	0.48
1:A:233:ILE:HD13	1:A:261:TYR:O	2.14	0.48
1:A:546:ASP:C	1:A:547:LEU:HD23	2.39	0.48
1:B:141:LYS:CD	1:B:158:ARG:NH2	2.77	0.48
1:B:303[B]:CYS:SG	1:B:311[B]:CYS:CB	3.02	0.48
1:B:498:ARG:NH1	3:B:684:HOH:O	2.46	0.47
1:B:21:LEU:HD11	1:B:37:VAL:CB	2.44	0.47
1:B:141:LYS:CE	1:B:158:ARG:NH2	2.78	0.47
1:B:33:HIS:O	1:B:36:MET:HG3	2.15	0.47
1:B:22:PRO:HG3	1:B:401:ARG:CG	2.41	0.47
1:A:82:LEU:HD13	1:A:249:ALA:HB2	1.97	0.46
1:A:314:LEU:HB3	1:A:321:VAL:CG1	2.46	0.46
1:B:19:SER:HB3	3:B:1004:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:372:VAL:HG22	1:B:382:TYR:CD1	2.50	0.46
1:B:530:VAL:HG13	1:B:533:LYS:HE3	1.94	0.46
1:A:282:SER:HB3	3:A:669:HOH:O	2.14	0.46
1:A:388:PRO:O	1:A:392:LEU:HG	2.15	0.46
1:A:401:ARG:HD3	1:A:401:ARG:HA	1.78	0.46
1:A:446:GLN:NE2	1:A:448:TYR:O	2.48	0.46
1:B:300:THR:HA	1:B:313:MET:HE3	1.97	0.46
1:B:100:LYS:HD2	3:B:830:HOH:O	2.15	0.45
1:B:436:GLN:OE1	1:B:438:GLN:OE1	2.33	0.45
1:B:148:GLN:O	1:B:148:GLN:HG3	2.17	0.45
1:B:84:SER:HB2	1:B:86:GLU:OE1	2.16	0.45
1:A:86:GLU:O	1:A:90:LYS:HG3	2.16	0.45
1:B:119:ILE:HG21	1:B:173:MET:CE	2.39	0.45
1:B:501:ARG:CZ	1:B:505:ARG:NH2	2.79	0.44
1:B:508:ARG:HH22	1:B:533:LYS:NZ	2.16	0.44
1:B:423:MET:HG2	1:B:528:TRP:CZ3	2.53	0.44
1:B:24:ASN:HB2	3:B:1085:HOH:O	2.16	0.44
1:A:405:ILE:CG2	1:A:406:ASN:N	2.80	0.44
1:B:31:LEU:HG	1:B:492:LEU:O	2.18	0.43
1:B:372:VAL:HG22	1:B:382:TYR:CE1	2.53	0.43
1:A:183:PRO:HG3	1:A:289:CYS:SG	2.59	0.43
1:B:314:LEU:HB3	1:B:321:VAL:HG13	2.01	0.43
1:B:544:GLN:HB2	3:B:741:HOH:O	2.19	0.43
1:B:212:LYS:HG3	1:B:325:GLU:OE1	2.18	0.42
1:B:36:MET:HE2	1:B:36:MET:HB3	1.80	0.42
1:B:392:LEU:HD21	1:B:424:ILE:HG21	2.02	0.42
1:B:531:ARG:O	1:B:532:THR:HB	2.19	0.42
1:A:201:VAL:CG2	1:A:384:LEU:HG	2.49	0.42
1:A:403:THR:C	1:A:405:ILE:N	2.76	0.42
1:A:414:MET:HE3	2:A:577:LT6:C12	2.49	0.42
1:A:423:MET:HE1	1:A:497:LEU:HB3	2.01	0.42
1:B:141:LYS:HE2	1:B:158:ARG:NH2	2.34	0.42
1:A:545:LEU:HB3	1:A:546:ASP:H	1.54	0.41
1:B:39:ALA:HA	1:B:143:GLU:O	2.20	0.41
1:B:198:LYS:NZ	3:B:1079:HOH:O	2.52	0.41
1:A:12:THR:HG21	1:A:270:LYS:HD2	2.03	0.41
1:A:329:THR:HG22	3:A:1110:HOH:O	2.21	0.41
1:A:220:ASP:O	1:A:350:PRO:HA	2.21	0.41
1:B:105:ALA:O	1:B:109:ARG:HG3	2.21	0.41
1:A:48:ARG:HG2	1:A:159:LEU:HG	2.03	0.41
1:B:83:LEU:HB2	1:B:173:MET:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:397:TRP:CE2	1:B:401:ARG:HD2	2.56	0.41
1:B:416:ALA:HB3	1:B:467:HIS:CD2	2.56	0.41
1:A:418:THR:O	1:A:422:ARG:HG3	2.20	0.40
1:A:222:ARG:NH2	3:A:948:HOH:O	2.54	0.40
1:A:403:THR:HB	1:A:405:ILE:O	2.21	0.40
1:A:247:PRO:HG3	1:B:234:ARG:HH11	1.86	0.40
1:A:284:VAL:CG2	3:A:1054:HOH:O	2.69	0.40
1:A:346:TYR:O	1:A:347:SER:HB3	2.22	0.40
1:A:11:ILE:HD12	1:A:49:GLN:CD	2.46	0.40
1:A:191:TYR:CZ	1:A:193:PHE:HB2	2.57	0.40
1:A:348:ALA:N	1:A:349:PRO:CD	2.85	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:ARG:NH2	1:B:514:GLN:O[4_455]	2.16	0.04
3:A:1190:HOH:O	3:B:599:HOH:O[4_455]	2.18	0.02

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	551/576 (96%)	537 (98%)	12 (2%)	2 (0%)	30	22
1	B	558/576 (97%)	541 (97%)	16 (3%)	1 (0%)	43	37
All	All	1109/1152 (96%)	1078 (97%)	28 (2%)	3 (0%)	36	30

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	534	LEU

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Mol	Chain	Res	Type
1	A	532	THR
1	A	546	ASP

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	476/490 (97%)	465 (98%)	11 (2%)	44	42
1	B	479/490 (98%)	460 (96%)	19 (4%)	28	22
All	All	955/980 (97%)	925 (97%)	30 (3%)	34	30

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	LEU
1	A	112	SER
1	A	114	ARG
1	A	234	ARG
1	A	352	ASP
1	A	379	LYS
1	A	487	SER
1	A	533	LYS
1	A	536	LEU
1	A	545	LEU
1	A	552	VAL
1	B	14	CYS
1	B	21	LEU
1	B	23	ILE
1	B	26	LEU
1	B	28	ASN
1	B	32	ARG
1	B	36	MET
1	B	57	LEU
1	B	86	GLU
1	B	141	LYS

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Mol	Chain	Res	Type
1	B	324	CYS
1	B	390	THR
1	B	419	LEU
1	B	530	VAL
1	B	533	LYS
1	B	534	LEU
1	B	535	LYS
1	B	536	LEU
1	B	544	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	514	GLN
1	A	562	HIS
1	B	438	GLN
1	B	527	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	LT6	A	577	-	31,31,31	1.17	2 (6%)	37,40,40	0.96	3 (8%)
2	LT6	B	577	-	31,31,31	1.05	2 (6%)	37,40,40	0.94	2 (5%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	LT6	A	577	-	-	2/27/27/27	0/2/2/2
2	LT6	B	577	-	-	0/27/27/27	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	577	LT6	C23-N1	3.91	1.41	1.34
2	B	577	LT6	C23-N1	2.75	1.39	1.34
2	A	577	LT6	C15-N2	2.66	1.39	1.34
2	B	577	LT6	C15-N2	2.42	1.39	1.34

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	577	LT6	C16-C15-N2	-2.59	111.94	116.41
2	A	577	LT6	C16-O3-C17	2.36	122.54	117.60
2	B	577	LT6	C16-C15-N2	-2.13	112.73	116.41
2	B	577	LT6	F1-C21-C22	2.05	121.75	118.15
2	A	577	LT6	O4-C23-C22	-2.02	115.64	119.49

There are no chirality outliers.

All (2) torsion outliers are listed below:

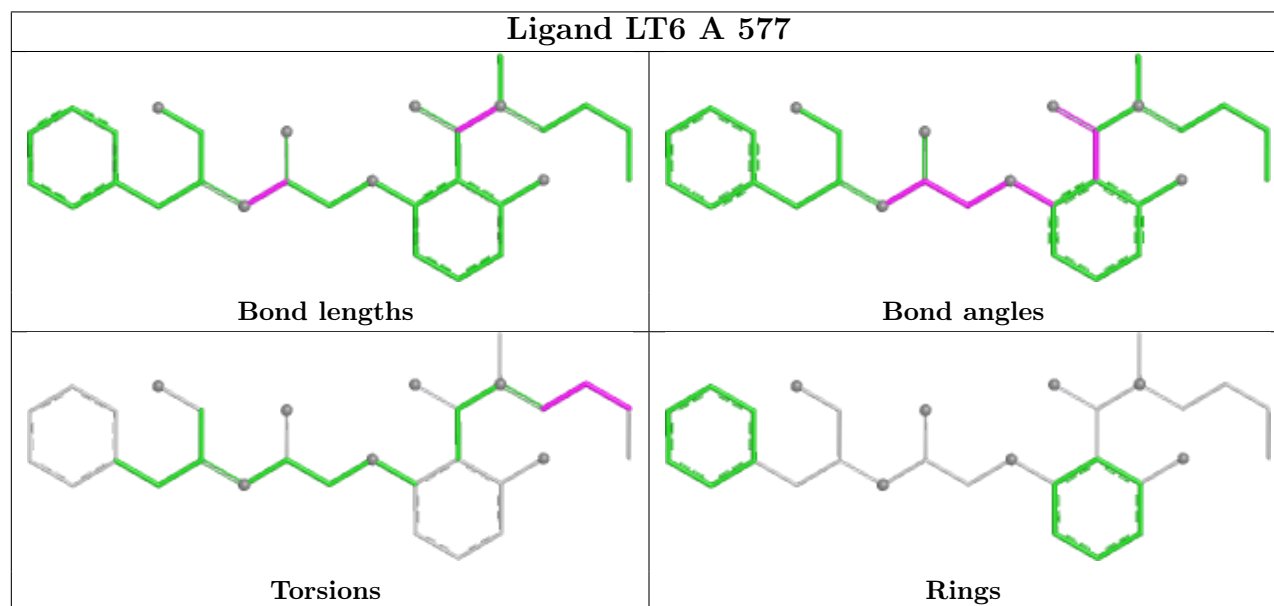
Mol	Chain	Res	Type	Atoms
2	A	577	LT6	C1-C2-C3-C4
2	A	577	LT6	C2-C3-C4-N1

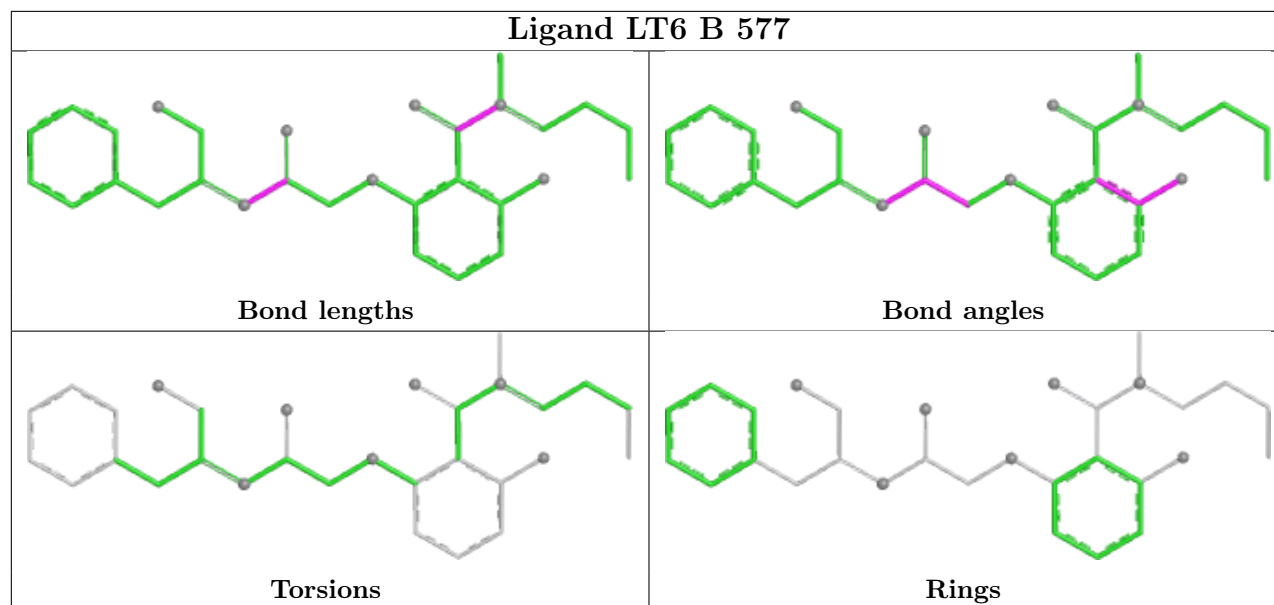
There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	577	LT6	3	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	554/576 (96%)	-0.17	12 (2%) 62 64	15, 31, 56, 83	3 (0%)
1	B	558/576 (96%)	0.11	27 (4%) 35 36	16, 36, 66, 107	4 (0%)
All	All	1112/1152 (96%)	-0.03	39 (3%) 47 48	15, 33, 60, 107	7 (0%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	26	LEU	7.4
1	B	532	THR	6.0
1	A	546	ASP	4.3
1	B	534	LEU	4.1
1	B	23	ILE	4.0
1	B	25	PRO	3.9
1	A	563	SER	3.6
1	B	536	LEU	3.4
1	B	543	SER	3.2
1	A	539	ILE	3.1
1	B	34	HIS	3.1
1	A	553	ALA	3.1
1	A	540	PRO	3.0
1	B	24	ASN	2.9
1	A	552	VAL	2.9
1	A	403	THR	2.8
1	B	486	ALA	2.7
1	B	376	ALA	2.6
1	B	27	SER	2.6
1	B	530	VAL	2.6
1	A	550	TRP	2.5
1	B	533	LYS	2.5
1	B	402	HIS	2.5
1	B	22	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	545	LEU	2.4
1	A	402	HIS	2.4
1	B	498	ARG	2.4
1	B	544	GLN	2.4
1	A	447	ILE	2.3
1	B	30	LEU	2.3
1	B	29	SER	2.2
1	B	148	GLN	2.1
1	B	435	ALA	2.1
1	B	403	THR	2.1
1	B	537	THR	2.1
1	A	405	ILE	2.1
1	B	531	ARG	2.0
1	B	535	LYS	2.0
1	B	438	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

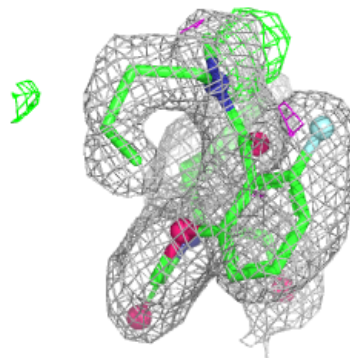
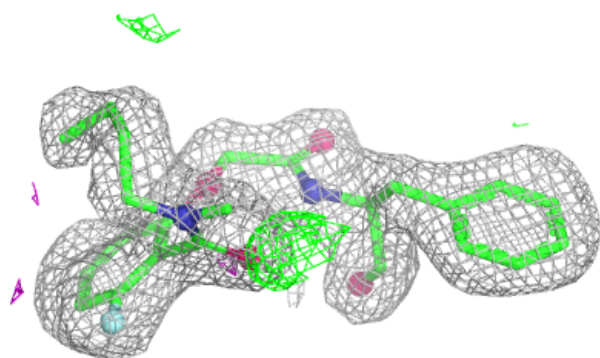
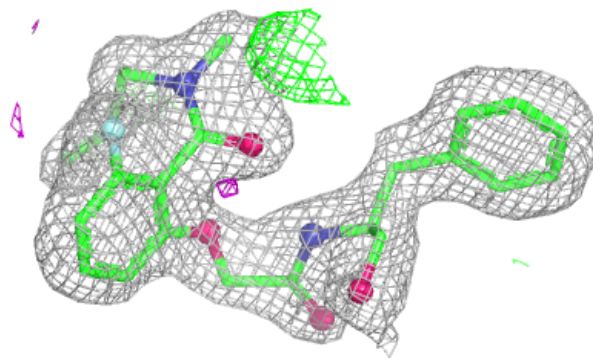
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	LT6	A	577	30/30	0.90	0.09	31,35,40,44	0
2	LT6	B	577	30/30	0.93	0.07	24,29,32,34	0

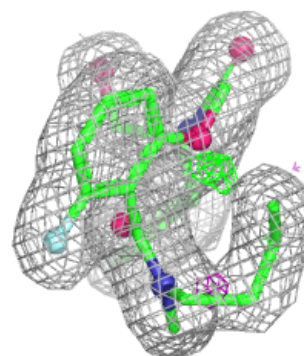
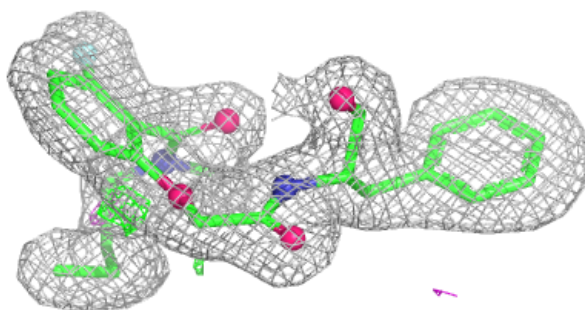
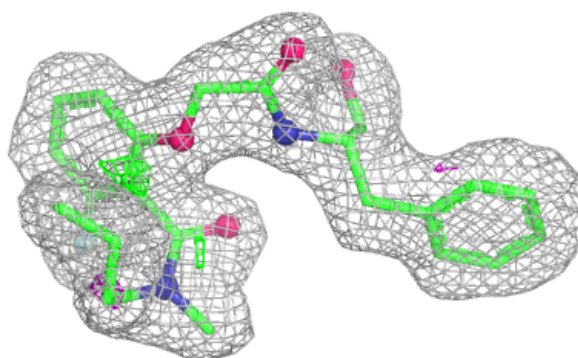
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around LT6 A 577:

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

**Electron density around LT6 B 577:**

$2mF_o-DF_c$ (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.