



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:24 PM UTC

PDB ID : 7LCR / pdb_00007lcr
Title : Improved Feline Drugs as SARS-CoV-2 Mpro Inhibitors: Structure-Activity Studies & Micellar Solubilization for Enhanced Bioavailability
Authors : Khan, M.B.; Arutyunova, E.; Young, H.S.; Lemieux, M.J.
Deposited on : 2021-01-11
Resolution : 1.95 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

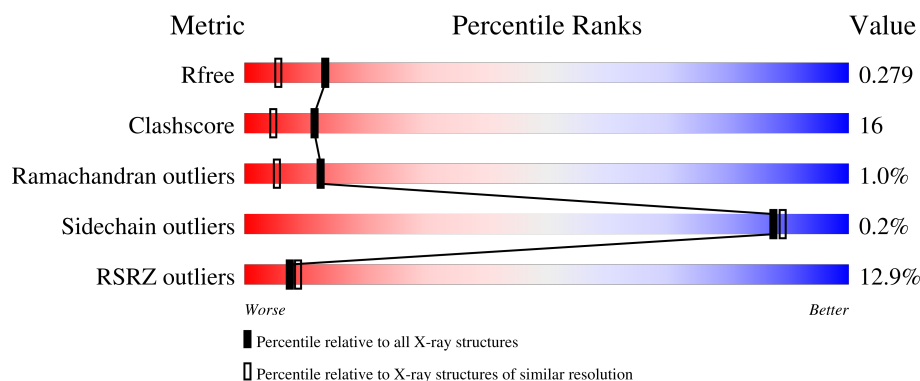
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	306	18% 76% 19% .
1	B	306	8% 76% 23% .

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	XTM	B	401	X	-	-	-

2 Entry composition [i](#)

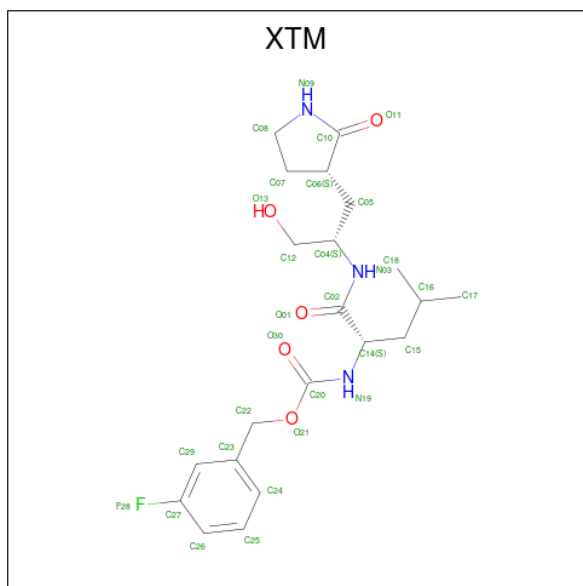
There are 3 unique types of molecules in this entry. The entry contains 9403 atoms, of which 4469 are hydrogens and 0 are deuteriums.

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 3C-like proteinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	306	Total	C	H	N	O	S	0	0	0
			4597	1499	2230	402	444	22			
1	B	306	Total	C	H	N	O	S	0	0	0
			4550	1499	2183	402	444	22			

- Molecule 2 is N 2 -{[(3-fluorophenyl)methoxy]carbonyl}-N-{(2S)-1-hydroxy-3-[(3S)-2-oxopyrrolidin-3-yl]propan-2-yl}-L-leucinamide (CCD ID: XTM) (formula: C₂₁H₃₀FN₃O₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	F	H	N	O	0	0
			58	21	1	28	3	5		
2	B	1	Total	C	F	H	N	O	0	0
			58	21	1	28	3	5		

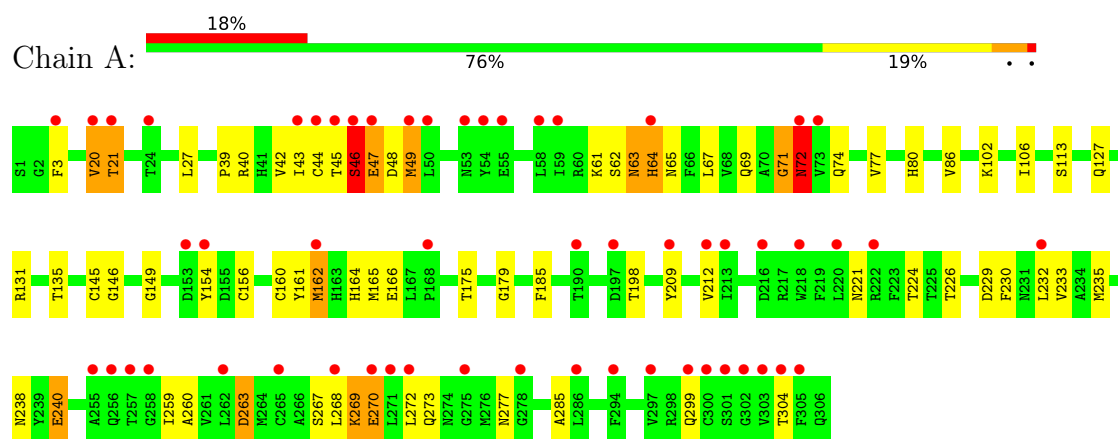
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	49	Total 49	O 49	0	0
3	B	91	Total 91	O 91	0	0

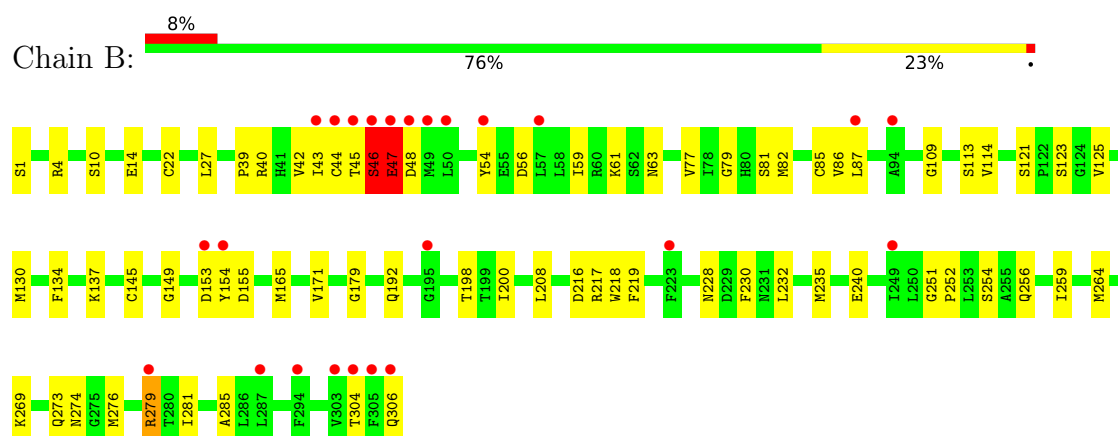
3 Residue-property plots [i](#)

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: 3C-like proteinase



• Molecule 1: 3C-like proteinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.74Å 101.51Å 102.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.00 – 1.95 38.00 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.8 (38.00-1.95) 99.9 (38.00-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.95 (at 1.89Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.238 , 0.277 0.243 , 0.279	Depositor DCC
R_{free} test set	1999 reflections (3.51%)	wwPDB-VP
Wilson B-factor (Å ²)	35.8	Xtriage
Anisotropy	0.483	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 33.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9403	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.33% 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: XTM

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	0.80	9/2420 (0.4%)	1.27	34/3289 (1.0%)
1	B	0.48	4/2420 (0.2%)	0.78	11/3289 (0.3%)
All	All	0.66	13/4840 (0.3%)	1.06	45/6578 (0.7%)

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	A	0	6
1	B	1	1
All	All	1	7

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	162	MET	CB-CG	21.83	2.17	1.52
1	B	47	GLU	CB-CG	13.81	1.93	1.52
1	A	21	THR	CB-CG2	12.84	1.95	1.52
1	A	162	MET	SD-CE	12.10	2.09	1.79
1	A	47	GLU	CD-OE2	7.73	1.40	1.25
1	A	161	TYR	C-N	-7.47	1.22	1.33
1	B	47	GLU	CG-CD	7.40	1.70	1.52
1	A	47	GLU	CB-CG	7.35	1.74	1.52
1	A	240	GLU	CD-OE2	6.20	1.37	1.25
1	A	21	THR	CA-CB	6.17	1.63	1.53
1	A	161	TYR	C-O	-6.06	1.17	1.23
1	B	46	SER	CA-C	5.39	1.59	1.52
1	B	46	SER	C-N	5.19	1.41	1.33

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	47	GLU	N-CA-CB	20.48	145.10	110.49
1	A	47	GLU	CG-CD-OE2	-19.28	74.05	118.40
1	A	47	GLU	CG-CD-OE1	18.09	160.02	118.40
1	A	240	GLU	CG-CD-OE1	16.19	155.64	118.40
1	A	47	GLU	OE1-CD-OE2	-15.16	86.53	122.90
1	A	72	ASN	N-CA-CB	14.38	131.32	110.47
1	A	240	GLU	CG-CD-OE2	-14.15	85.86	118.40
1	A	71	GLY	CA-C-N	-13.79	98.84	122.56
1	A	71	GLY	C-N-CA	-13.79	98.84	122.56
1	B	47	GLU	CB-CG-CD	12.40	133.68	112.60
1	A	240	GLU	OE1-CD-OE2	-11.66	94.90	122.90
1	A	21	THR	OG1-CB-CG2	11.34	131.98	109.30
1	B	47	GLU	CA-C-O	-10.85	105.00	120.51
1	A	162	MET	CB-CG-SD	-10.74	80.47	112.70
1	A	47	GLU	N-CA-C	-10.41	101.08	113.88
1	A	72	ASN	N-CA-C	-10.36	100.58	113.02
1	A	263	ASP	CB-CA-C	9.19	128.13	109.67
1	A	72	ASN	CB-CA-C	-9.15	94.07	110.37
1	A	161	TYR	O-C-N	-8.72	112.71	123.44
1	A	263	ASP	N-CA-CB	-8.46	97.63	110.33
1	B	279	ARG	CB-CG-CD	7.95	129.58	111.30
1	A	47	GLU	CB-CA-C	-7.81	96.37	109.03
1	A	64	HIS	N-CA-CB	-7.81	96.67	110.42
1	A	47	GLU	N-CA-CB	7.59	120.69	110.59
1	A	161	TYR	CA-C-N	-7.31	109.43	122.74
1	A	161	TYR	C-N-CA	-7.31	109.43	122.74
1	B	279	ARG	CG-CD-NE	7.05	127.52	112.00
1	A	71	GLY	CA-C-O	-6.88	113.02	121.68
1	A	64	HIS	CB-CG-CD2	-6.53	122.72	131.20
1	A	162	MET	CB-CA-C	-6.50	99.07	109.80
1	B	46	SER	CA-C-N	-6.37	109.37	121.54
1	B	46	SER	C-N-CA	-6.37	109.37	121.54
1	B	47	GLU	O-C-N	-6.20	114.34	122.59
1	B	46	SER	N-CA-C	6.17	120.89	112.68
1	A	21	THR	CB-CA-C	6.14	122.65	110.42
1	A	47	GLU	CB-CG-CD	6.10	122.97	112.60
1	B	47	GLU	CG-CD-OE1	5.83	131.80	118.40
1	A	20	VAL	CA-C-N	5.80	132.61	121.54
1	A	20	VAL	C-N-CA	5.80	132.61	121.54
1	B	279	ARG	CA-CB-CG	-5.27	103.56	114.10
1	A	263	ASP	N-CA-C	-5.17	106.13	112.90
1	A	63	ASN	CA-C-N	5.17	133.61	121.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	63	ASN	C-N-CA	5.17	133.61	121.52
1	A	64	HIS	ND1-CG-CD2	-5.16	100.94	106.10
1	A	71	GLY	O-C-N	-5.02	116.94	122.61

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	47	GLU	CA

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	20	VAL	Peptide
1	A	240	GLU	Sidechain
1	A	46	SER	Peptide
1	A	64	HIS	Sidechain
1	A	72	ASN	Mainchain,Sidechain
1	B	46	SER	Peptide

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	2367	2230	2313	89	0
1	B	2367	2183	2314	65	0
2	A	30	28	0	3	0
2	B	30	28	0	2	0
3	A	49	0	0	2	0
3	B	91	0	0	9	0
All	All	4934	4469	4627	149	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 16.

All (149) 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:47:GLU:CB	1:A:47:GLU:CG	1.74	1.62
1:B:47:GLU:CB	1:B:47:GLU:CG	1.93	1.43
1:A:21:THR:CB	1:A:21:THR:CG2	1.94	1.42
1:A:162:MET:SD	1:A:162:MET:CE	2.09	1.40
1:A:162:MET:CB	1:A:162:MET:CG	2.17	1.22
1:A:45:THR:HB	1:A:49:MET:HE2	1.30	1.12
1:B:47:GLU:OE1	1:B:48:ASP:N	1.83	1.12
1:B:47:GLU:OE1	1:B:48:ASP:HB3	1.50	1.11
1:A:47:GLU:CB	1:A:47:GLU:OE2	1.99	1.07
1:A:233:VAL:HG21	1:A:269:LYS:HD3	1.36	1.07
1:A:162:MET:HE2	1:A:162:MET:HB2	1.34	1.06
1:B:47:GLU:OE1	1:B:48:ASP:CB	2.05	1.04
1:A:47:GLU:OE1	1:A:48:ASP:HB2	1.63	0.98
1:A:162:MET:SD	1:A:162:MET:CB	2.52	0.97
1:A:21:THR:HG23	1:A:67:LEU:O	1.63	0.96
1:A:226:THR:OG1	1:A:229:ASP:OD1	1.85	0.94
1:B:85:CYS:SG	3:B:503:HOH:O	2.26	0.93
1:B:79:GLY:N	3:B:501:HOH:O	2.02	0.93
1:B:47:GLU:OE1	1:B:48:ASP:CA	2.18	0.92
1:B:145:CYS:SG	2:B:401:XTM:C04	2.58	0.91
1:B:77:VAL:HG12	3:B:501:HOH:O	1.71	0.91
1:A:162:MET:CE	1:A:162:MET:HB2	2.02	0.88
1:A:45:THR:HG22	1:A:49:MET:HB3	1.57	0.87
1:B:54:TYR:HB3	1:B:82:MET:HE1	1.55	0.86
1:A:233:VAL:HG11	1:A:269:LYS:HG2	1.56	0.85
1:A:102:LYS:NZ	1:A:156:CYS:SG	2.50	0.85
1:A:62:SER:N	1:A:65:ASN:OD1	2.12	0.83
1:A:45:THR:HG22	1:A:49:MET:CB	2.12	0.80
1:A:221:ASN:OD1	3:A:501:HOH:O	1.98	0.79
1:A:47:GLU:CG	1:A:47:GLU:C	2.55	0.79
1:B:208:LEU:HB3	1:B:264:MET:HE2	1.66	0.78
1:A:39:PRO:HG3	1:A:162:MET:HG2	1.65	0.78
1:A:21:THR:CG2	1:A:21:THR:HB	2.11	0.78
1:A:46:SER:HB2	1:A:49:MET:H	1.50	0.76
1:A:47:GLU:C	1:A:47:GLU:HG2	2.12	0.75
1:B:198:THR:OG1	1:B:240:GLU:OE2	2.03	0.75
1:A:47:GLU:OE2	1:A:47:GLU:HB3	1.88	0.73
1:B:27:LEU:HD21	1:B:42:VAL:HB	1.73	0.70
1:B:153:ASP:O	1:B:154:TYR:HB2	1.91	0.70
1:A:45:THR:CB	1:A:49:MET:HE2	2.18	0.70
1:A:86:VAL:HG23	1:A:179:GLY:HA2	1.74	0.69
1:A:162:MET:CB	1:A:162:MET:HE2	2.18	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:VAL:HG21	1:A:269:LYS:CD	2.19	0.69
1:A:47:GLU:HG2	1:A:47:GLU:O	1.92	0.69
1:A:162:MET:CE	1:A:162:MET:CB	2.70	0.68
1:A:277:ASN:OD1	3:A:502:HOH:O	2.10	0.68
1:A:269:LYS:O	1:A:272:LEU:N	2.27	0.67
1:A:27:LEU:HD21	1:A:42:VAL:HB	1.77	0.66
1:A:145:CYS:SG	2:A:401:XTM:C04	2.82	0.65
1:A:45:THR:HG21	1:A:49:MET:HG3	1.79	0.65
1:A:198:THR:HG22	1:A:238:ASN:OD1	1.97	0.65
1:B:47:GLU:CD	1:B:48:ASP:N	2.54	0.65
1:A:49:MET:O	1:A:49:MET:HG2	1.95	0.65
1:A:165:MET:HE1	1:A:185:PHE:HB3	1.79	0.65
1:A:260:ALA:HB3	1:A:263:ASP:CB	2.28	0.64
1:B:273:GLN:NE2	1:B:274:ASN:OD1	2.30	0.63
1:B:306:GLN:NE2	3:B:504:HOH:O	2.30	0.63
1:B:45:THR:OG1	1:B:47:GLU:OE2	2.16	0.62
1:B:86:VAL:HG23	1:B:179:GLY:HA2	1.82	0.61
1:A:47:GLU:OE2	1:A:48:ASP:N	2.34	0.61
1:B:130:MET:HE2	1:B:134:PHE:HA	1.82	0.61
1:A:69:GLN:OE1	1:A:74:GLN:OE1	2.19	0.59
1:B:40:ARG:NH2	3:B:503:HOH:O	2.28	0.59
1:B:217:ARG:NH2	3:B:505:HOH:O	2.37	0.58
1:B:145:CYS:SG	2:B:401:XTM:N03	2.77	0.57
1:A:45:THR:CG2	1:A:49:MET:CB	2.81	0.57
1:A:162:MET:CE	1:A:162:MET:CG	2.81	0.56
1:A:260:ALA:HB3	1:A:263:ASP:HB3	1.90	0.54
1:A:285:ALA:HB3	1:B:285:ALA:HB3	1.89	0.54
1:B:219:PHE:HE1	1:B:264:MET:HE1	1.73	0.54
1:B:228:ASN:O	1:B:232:LEU:HD13	2.07	0.54
1:A:21:THR:CG2	1:A:67:LEU:O	2.48	0.54
1:A:46:SER:HB2	1:A:49:MET:N	2.23	0.53
1:A:48:ASP:O	1:A:49:MET:C	2.52	0.53
1:B:40:ARG:O	1:B:43:ILE:HG12	2.08	0.53
1:B:276:MET:HE1	1:B:281:ILE:HG13	1.90	0.52
1:A:145:CYS:CB	2:A:401:XTM:C12	2.87	0.52
1:B:44:CYS:HG	1:B:54:TYR:HE1	1.57	0.52
1:A:162:MET:N	1:A:175:THR:O	2.37	0.51
1:A:260:ALA:HB3	1:A:263:ASP:HB2	1.92	0.51
1:A:45:THR:CG2	1:A:49:MET:HG3	2.40	0.51
1:A:106:ILE:HG23	1:A:160:CYS:HB2	1.93	0.51
1:A:165:MET:CE	1:A:185:PHE:HB3	2.39	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:GLY:HA2	1:A:162:MET:CG	2.41	0.50
1:B:165:MET:HE1	1:B:192:GLN:NE2	2.26	0.50
1:A:233:VAL:CG2	1:A:269:LYS:HD3	2.26	0.50
1:A:47:GLU:O	1:A:49:MET:SD	2.70	0.50
1:B:46:SER:O	1:B:47:GLU:C	2.53	0.50
1:B:10:SER:O	1:B:14:GLU:HG3	2.12	0.49
1:A:230:PHE:CZ	1:A:268:LEU:HD23	2.48	0.49
1:A:40:ARG:O	1:A:43:ILE:HD12	2.12	0.49
1:B:56:ASP:O	1:B:59:ILE:HG22	2.11	0.49
1:A:209:TYR:O	1:A:212:VAL:HG22	2.12	0.49
1:B:27:LEU:HD13	1:B:39:PRO:HD2	1.95	0.48
1:A:47:GLU:OE1	1:A:48:ASP:CB	2.51	0.48
1:A:131:ARG:HG2	1:A:135:THR:O	2.13	0.48
1:A:304:THR:HG22	1:B:123:SER:HB2	1.95	0.48
1:A:224:THR:HG22	1:A:263:ASP:CB	2.44	0.48
1:A:267:SER:O	1:A:270:GLU:HB2	2.14	0.48
1:B:154:TYR:O	1:B:155:ASP:OD1	2.31	0.47
1:B:43:ILE:HD11	1:B:54:TYR:HD1	1.78	0.47
1:A:166:GLU:OE2	1:B:1:SER:N	2.48	0.47
1:B:254:SER:OG	1:B:259:ILE:O	2.27	0.47
1:A:269:LYS:HD2	1:A:273:GLN:OE1	2.15	0.46
1:A:259:ILE:HG22	1:A:260:ALA:O	2.16	0.46
1:B:130:MET:HE2	1:B:134:PHE:C	2.40	0.46
1:B:63:ASN:CG	3:B:501:HOH:O	2.59	0.46
1:B:208:LEU:CB	1:B:264:MET:HE2	2.41	0.46
1:A:127:GLN:O	1:B:4:ARG:NH2	2.39	0.45
1:A:260:ALA:O	1:A:263:ASP:HB3	2.17	0.45
1:B:22:CYS:SG	1:B:61:LYS:NZ	2.88	0.45
1:B:130:MET:HE2	1:B:134:PHE:CA	2.46	0.45
1:A:164:HIS:CD2	1:A:175:THR:HG23	2.52	0.45
1:B:155:ASP:OD1	1:B:155:ASP:C	2.60	0.45
1:A:45:THR:O	1:A:46:SER:O	2.35	0.45
1:B:230:PHE:HA	1:B:269:LYS:HD2	1.98	0.44
1:A:45:THR:C	1:A:49:MET:HE3	2.43	0.44
1:A:145:CYS:SG	2:A:401:XTM:N03	2.91	0.44
1:A:232:LEU:HD13	1:A:235:MET:HE3	2.00	0.44
1:B:276:MET:O	1:B:279:ARG:HB2	2.18	0.44
1:A:113:SER:O	1:A:149:GLY:HA2	2.18	0.44
1:A:3:PHE:CD2	1:A:299:GLN:HG2	2.53	0.44
1:B:256:GLN:HG2	1:B:304:THR:HG23	2.00	0.43
1:B:232:LEU:HA	1:B:235:MET:HE3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:HIS:O	1:A:80:HIS:ND1	2.50	0.43
1:A:44:CYS:HB2	1:A:61:LYS:NZ	2.33	0.43
1:A:63:ASN:O	1:A:77:VAL:HG11	2.18	0.43
1:B:81:SER:O	1:B:87:LEU:HD12	2.19	0.43
1:B:137:LYS:HG2	1:B:171:VAL:HG12	2.01	0.43
1:B:44:CYS:SG	1:B:54:TYR:CE1	3.12	0.43
1:A:71:GLY:O	1:A:72:ASN:C	2.57	0.42
1:B:109:GLY:HA2	1:B:200:ILE:HD13	2.02	0.42
1:A:304:THR:CB	1:B:121:SER:OG	2.68	0.42
1:A:45:THR:C	1:A:49:MET:CE	2.93	0.42
1:B:45:THR:OG1	1:B:47:GLU:CD	2.63	0.41
1:A:162:MET:SD	1:A:162:MET:HB2	2.52	0.41
1:B:81:SER:N	3:B:512:HOH:O	2.54	0.41
1:B:153:ASP:C	1:B:154:TYR:HD2	2.29	0.41
1:B:114:VAL:O	1:B:125:VAL:HA	2.21	0.41
1:B:54:TYR:CB	1:B:82:MET:HE1	2.39	0.41
1:A:45:THR:CG2	1:A:49:MET:CG	2.99	0.41
1:A:232:LEU:HA	1:A:235:MET:HE2	2.02	0.41
1:B:216:ASP:OD1	1:B:218:TRP:HD1	2.03	0.41
1:B:251:GLY:N	1:B:252:PRO:HD2	2.36	0.41
1:A:47:GLU:OE2	1:A:47:GLU:C	2.64	0.40
1:B:130:MET:CE	1:B:134:PHE:C	2.94	0.40
1:B:113:SER:O	1:B:149:GLY:HA2	2.21	0.40
1:B:87:LEU:CD1	3:B:512:HOH:O	2.68	0.40
1:A:269:LYS:O	1:A:270:GLU:C	2.65	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	304/306 (99%)	291 (96%)	8 (3%)	5 (2%)	7 2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	304/306 (99%)	295 (97%)	8 (3%)	1 (0%)	36	28
All	All	608/612 (99%)	586 (96%)	16 (3%)	6 (1%)	12	5

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	46	SER
1	A	270	GLU
1	B	47	GLU
1	A	49	MET
1	A	154	TYR
1	A	269	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	263/263 (100%)	263 (100%)	0	100	100
1	B	263/263 (100%)	262 (100%)	1 (0%)	84	85
All	All	526/526 (100%)	525 (100%)	1 (0%)	87	89

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

Mol	Chain	Res	Type
1	B	47	GLU

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	84	ASN
1	A	119	ASN
1	A	274	ASN
1	B	51	ASN

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Mol	Chain	Res	Type
1	B	119	ASN
1	B	142	ASN
1	B	180	ASN
1	B	228	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	XTM	B	401	1	31,31,31	4.32	15 (48%)	37,41,41	1.65	7 (18%)
2	XTM	A	401	1	31,31,31	4.33	15 (48%)	37,41,41	1.51	6 (16%)

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	XTM	B	401	1	1/1/10/10	5/27/37/37	0/2/2/2
2	XTM	A	401	1	-	0/27/37/37	0/2/2/2

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	XTM	C10-N09	14.87	1.49	1.33
2	B	401	XTM	C10-N09	14.58	1.49	1.33
2	A	401	XTM	C25-C24	8.41	1.53	1.38
2	B	401	XTM	C25-C24	8.35	1.53	1.38
2	A	401	XTM	C29-C23	8.30	1.53	1.39
2	B	401	XTM	C29-C23	8.28	1.53	1.39
2	B	401	XTM	C26-C27	8.24	1.53	1.37
2	A	401	XTM	C26-C27	8.24	1.53	1.37
2	B	401	XTM	C02-N03	5.72	1.46	1.34
2	A	401	XTM	C02-N03	5.58	1.46	1.34
2	B	401	XTM	O13-C12	-5.51	1.19	1.42
2	A	401	XTM	O13-C12	-5.49	1.19	1.42
2	B	401	XTM	C20-N19	4.28	1.45	1.34
2	A	401	XTM	C20-N19	4.08	1.44	1.34
2	B	401	XTM	C25-C26	-4.00	1.32	1.38
2	A	401	XTM	C25-C26	-3.93	1.32	1.38
2	B	401	XTM	C24-C23	-3.43	1.32	1.38
2	A	401	XTM	C24-C23	-3.41	1.32	1.38
2	B	401	XTM	C29-C27	-3.31	1.32	1.37
2	A	401	XTM	C29-C27	-3.26	1.32	1.37
2	B	401	XTM	O21-C20	2.67	1.40	1.35
2	B	401	XTM	O11-C10	-2.60	1.18	1.23
2	A	401	XTM	O21-C20	2.57	1.40	1.35
2	B	401	XTM	O01-C02	-2.54	1.18	1.23
2	A	401	XTM	O11-C10	-2.51	1.18	1.23
2	A	401	XTM	O21-C22	-2.49	1.40	1.45
2	B	401	XTM	O21-C22	-2.46	1.40	1.45
2	A	401	XTM	O01-C02	-2.30	1.19	1.23
2	A	401	XTM	C08-N09	2.16	1.50	1.46
2	B	401	XTM	C08-N09	2.14	1.50	1.46

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	XTM	O21-C20-N19	4.19	119.40	110.45
2	A	401	XTM	O21-C20-N19	3.79	118.56	110.45
2	B	401	XTM	O13-C12-C04	3.77	121.22	111.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	XTM	O13-C12-C04	3.50	120.54	111.73
2	A	401	XTM	C07-C06-C10	3.47	107.63	102.87
2	B	401	XTM	C07-C06-C10	3.40	107.54	102.87
2	B	401	XTM	C08-N09-C10	-3.31	106.77	113.79
2	B	401	XTM	O30-C20-N19	-3.28	119.48	124.86
2	A	401	XTM	O30-C20-N19	-3.26	119.52	124.86
2	A	401	XTM	C08-N09-C10	-2.85	107.74	113.79
2	B	401	XTM	C16-C15-C14	-2.68	108.19	115.40
2	B	401	XTM	C26-C27-C29	-2.39	120.08	123.23
2	A	401	XTM	C26-C27-C29	-2.37	120.11	123.23

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	401	XTM	C04

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	401	XTM	C04-C05-C06-C07
2	B	401	XTM	C05-C04-N03-C02
2	B	401	XTM	C14-C15-C16-C17
2	B	401	XTM	C04-C05-C06-C10
2	B	401	XTM	C14-C15-C16-C18

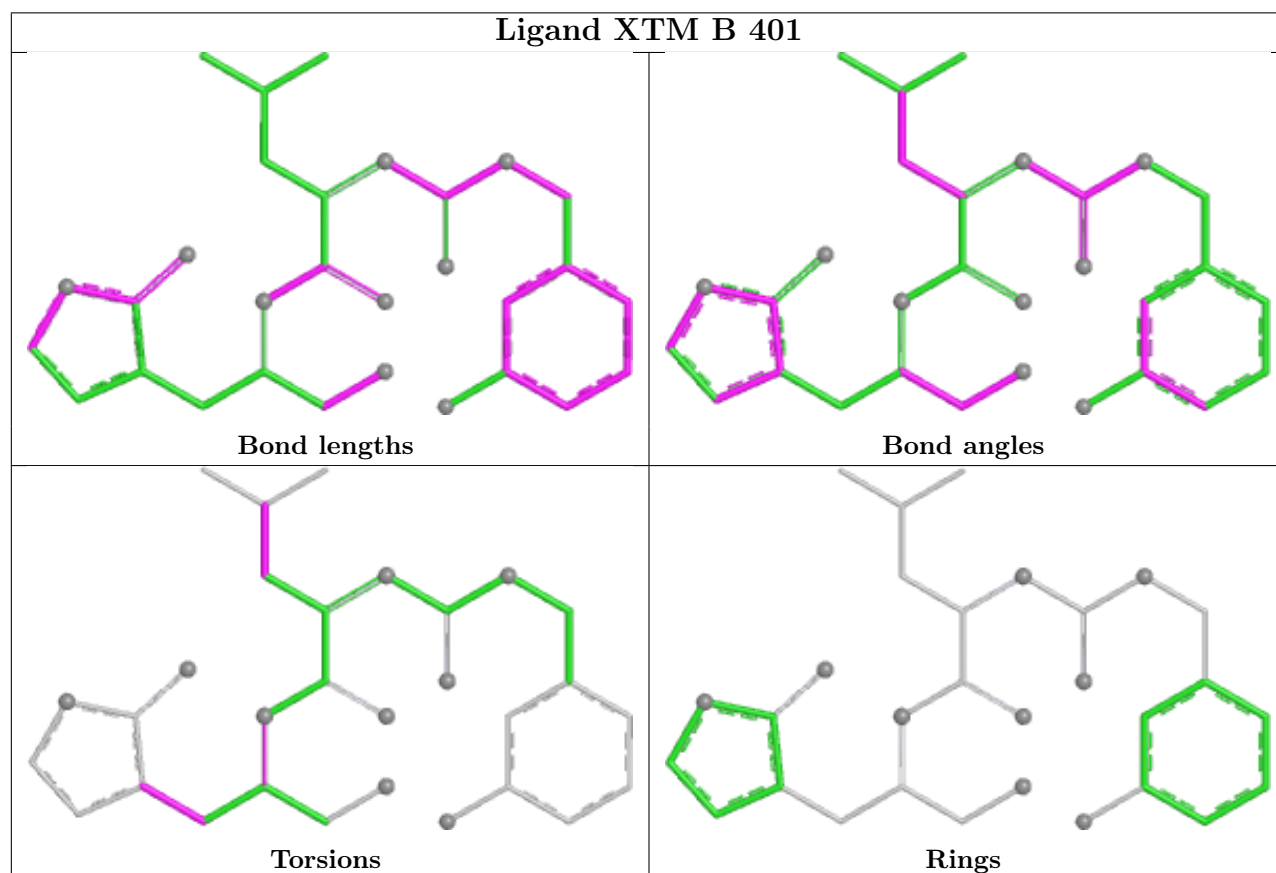
There are no ring outliers.

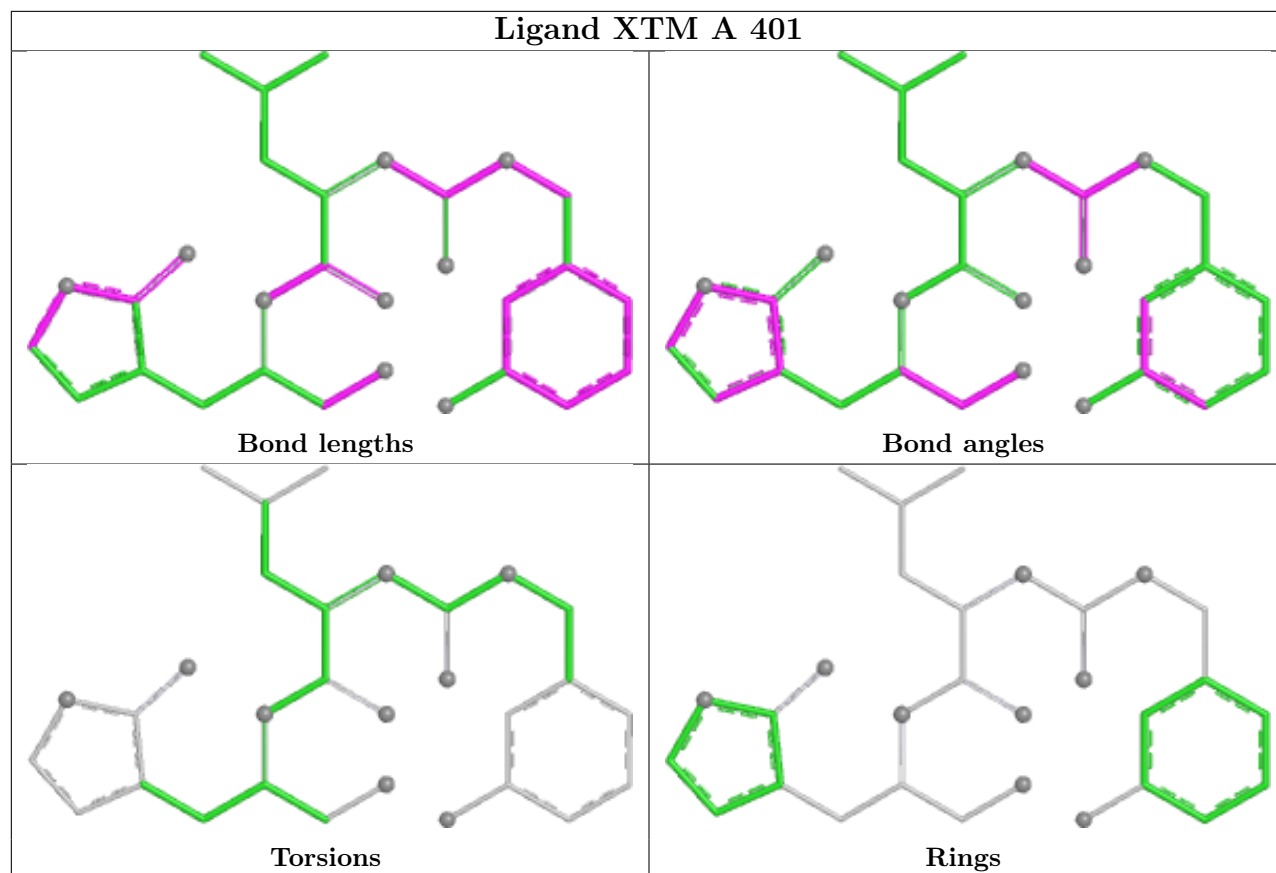
2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	XTM	2	0
2	A	401	XTM	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	306/306 (100%)	1.11	55 (17%) 3 3	45, 71, 135, 211	0
1	B	306/306 (100%)	0.72	24 (7%) 19 22	45, 60, 97, 211	0
All	All	612/612 (100%)	0.92	79 (12%) 7 9	45, 64, 120, 211	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	162	MET	7.0
1	A	305	PHE	6.6
1	A	303	VAL	4.8
1	A	300	CYS	4.0
1	A	294	PHE	3.9
1	B	47	GLU	3.8
1	A	299	GLN	3.8
1	B	303	VAL	3.8
1	A	302	GLY	3.6
1	B	294	PHE	3.6
1	A	50	LEU	3.5
1	A	222	ARG	3.4
1	A	21	THR	3.4
1	A	301	SER	3.3
1	B	154	TYR	3.2
1	A	278	GLY	3.1
1	A	45	THR	3.1
1	A	275	GLY	3.1
1	A	286	LEU	3.1
1	A	257	THR	3.1
1	A	297	VAL	3.0
1	B	44	CYS	3.0
1	A	46	SER	3.0
1	A	258	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	20	VAL	2.9
1	B	45	THR	2.9
1	A	64	HIS	2.8
1	A	24	THR	2.8
1	A	209	TYR	2.8
1	A	59	ILE	2.8
1	A	304	THR	2.8
1	A	268	LEU	2.7
1	A	255	ALA	2.7
1	B	249	ILE	2.7
1	A	270	GLU	2.6
1	B	304	THR	2.6
1	A	44	CYS	2.6
1	A	43	ILE	2.6
1	A	232	LEU	2.6
1	A	73	VAL	2.6
1	B	50	LEU	2.5
1	B	87	LEU	2.5
1	A	271	LEU	2.5
1	B	287	LEU	2.5
1	B	43	ILE	2.5
1	A	72	ASN	2.4
1	A	154	TYR	2.4
1	A	218	TRP	2.4
1	B	153	ASP	2.4
1	A	3	PHE	2.3
1	A	220	LEU	2.3
1	B	57	LEU	2.3
1	B	54	TYR	2.3
1	A	213	ILE	2.3
1	A	212	VAL	2.3
1	A	190	THR	2.3
1	B	305	PHE	2.2
1	A	153	ASP	2.2
1	A	55	GLU	2.2
1	A	216	ASP	2.2
1	B	46	SER	2.2
1	A	168	PRO	2.2
1	A	58	LEU	2.2
1	A	54	TYR	2.1
1	A	272	LEU	2.1
1	A	53	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	47	GLU	2.1
1	B	94	ALA	2.1
1	A	262	LEU	2.1
1	A	265	CYS	2.1
1	A	197	ASP	2.1
1	B	279	ARG	2.1
1	B	223	PHE	2.0
1	A	49	MET	2.0
1	B	49	MET	2.0
1	B	195	GLY	2.0
1	A	256	GLN	2.0
1	B	48	ASP	2.0
1	B	306	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](#)

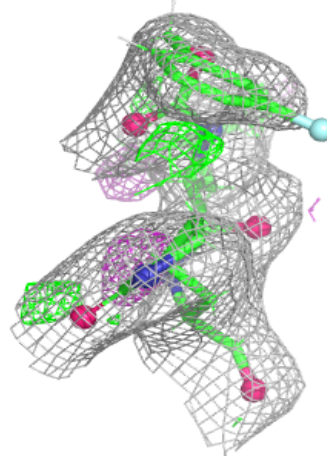
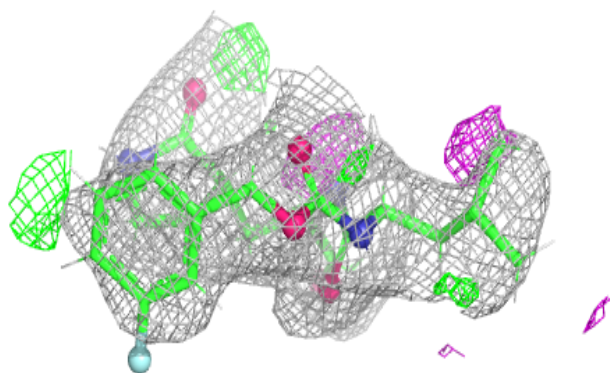
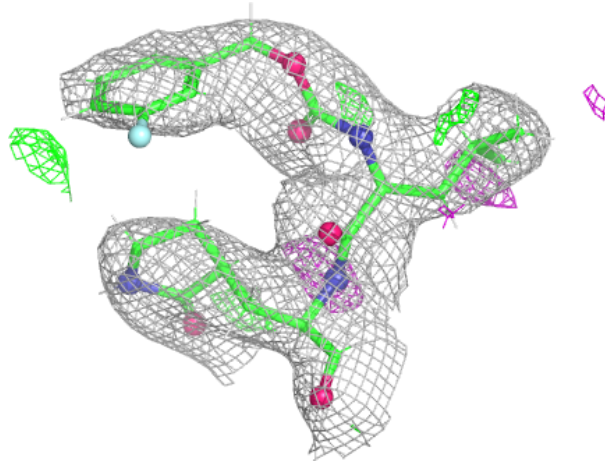
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	XTM	B	401	30/30	0.88	0.15	47,80,112,118	0
2	XTM	A	401	30/30	0.94	0.10	43,68,94,113	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.

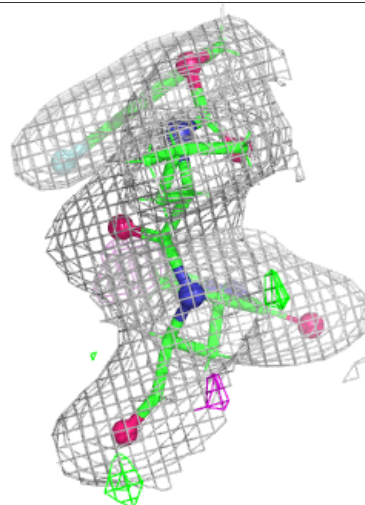
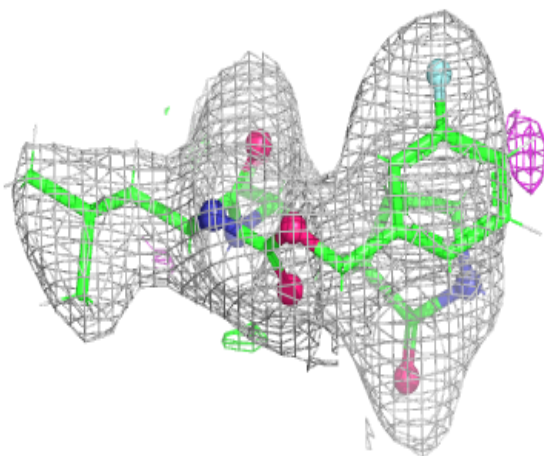
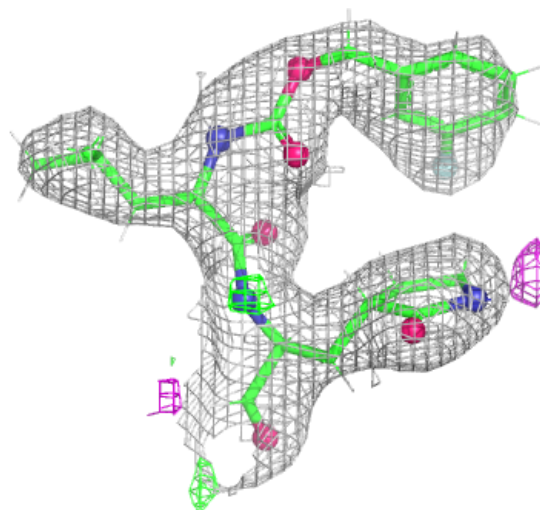
Electron density around XTM B 401:

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



Electron density around XTM A 401:

$2mF_o-DF_c$ (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.