



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 07:56 PM UTC

PDB ID : 4L8F / pdb_00004l8f
Title : Crystal structure of gamma-glutamyl hydrolase (C108A) complex with MTX
Authors : Chuankhayan, P.; Kao, T.-T.; Chen, C.-J.; Fu, T.-F.
Deposited on : 2013-06-17
Resolution : 1.97 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

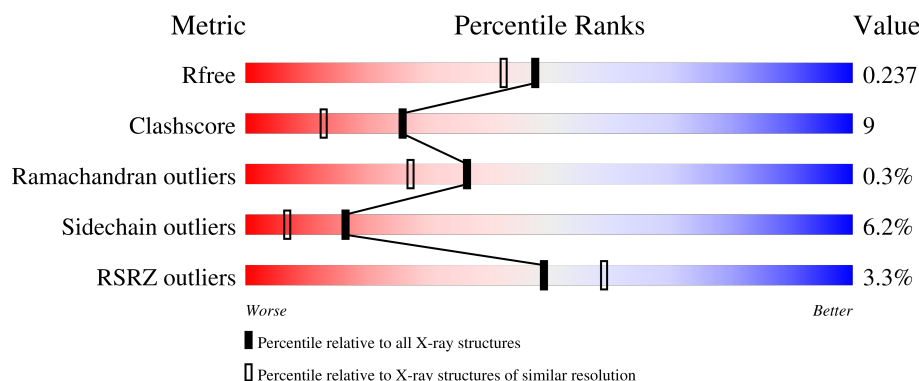
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.97 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	1506 (1.98-1.98)
Clashscore	190562	1534 (1.98-1.98)
Ramachandran outliers	187476	1518 (1.98-1.98)
Sidechain outliers	187428	1518 (1.98-1.98)
RSRZ outliers	180081	1506 (1.98-1.98)

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	312	4% 75% 14% • 8%
1	B	312	3% 72% 15% • • 8%
1	C	312	3% 79% 11% • 8%
1	D	312	2% 76% 11% • • 8%

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	MTX	B	301	-	-	X	-
2	MTX	D	301	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10251 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.

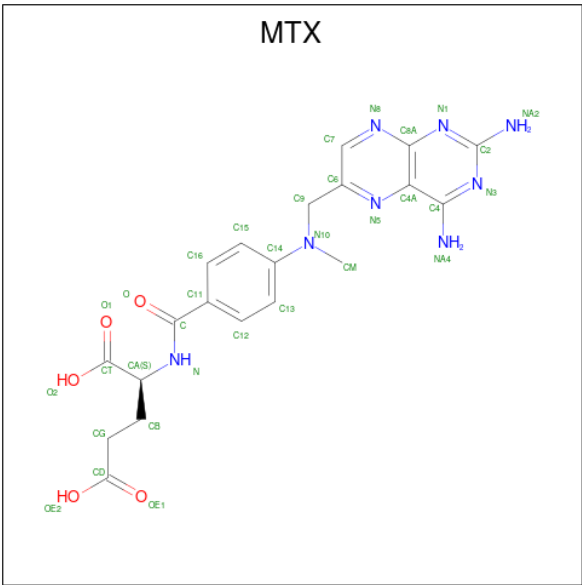
- Molecule 1 is a protein called Gamma-glutamyl hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	287	Total	C	N	O	S	0	0	0
			2308	1492	370	442	4			
1	D	287	Total	C	N	O	S	0	0	0
			2308	1492	370	442	4			
1	B	287	Total	C	N	O	S	0	0	0
			2308	1492	370	442	4			
1	A	287	Total	C	N	O	S	0	0	0
			2308	1492	370	442	4			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	108	ALA	CYS	engineered mutation	UNP Q6NY42
D	108	ALA	CYS	engineered mutation	UNP Q6NY42
B	108	ALA	CYS	engineered mutation	UNP Q6NY42
A	108	ALA	CYS	engineered mutation	UNP Q6NY42

- Molecule 2 is METHOTREXATE (CCD ID: MTX) (formula: C₂₀H₂₂N₈O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	D	1	Total	C	N	O	0	0
			33	20	8	5		
2	B	1	Total	C	N	O	0	0
			33	20	8	5		

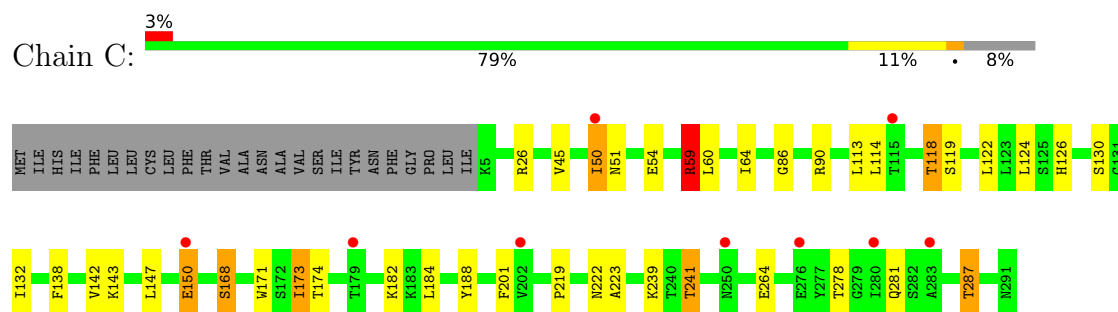
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	265	Total	O	0	0
			265	265		
3	D	266	Total	O	0	0
			266	266		
3	B	271	Total	O	0	0
			271	271		
3	A	151	Total	O	0	0
			151	151		

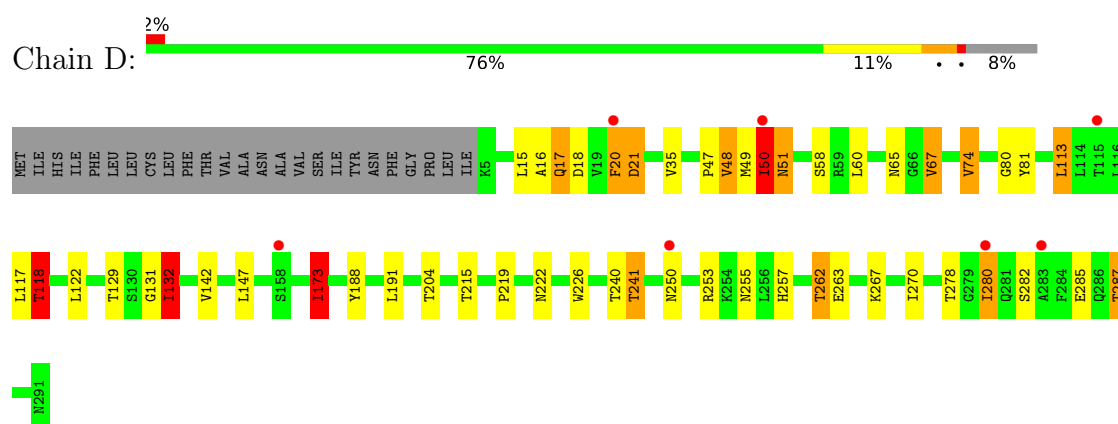
3 Residue-property plots [i](#)

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.

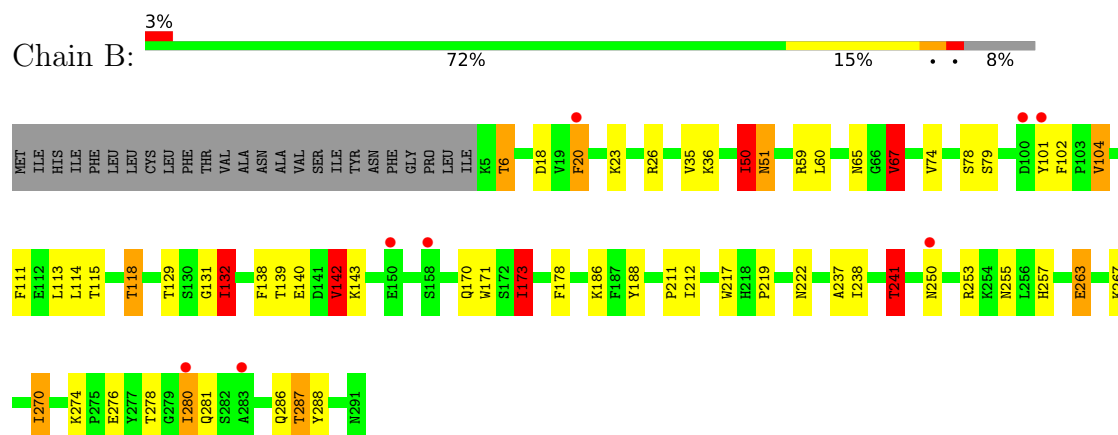
• Molecule 1: Gamma-glutamyl hydrolase



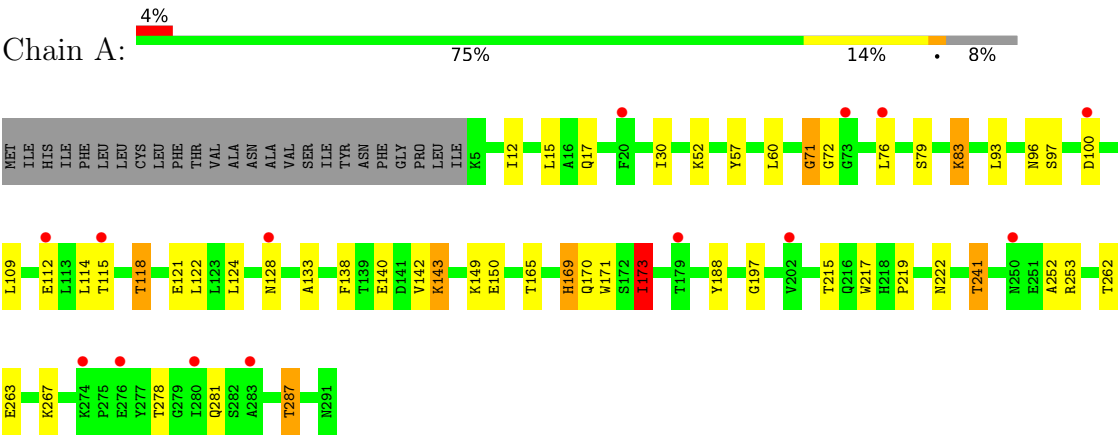
• Molecule 1: Gamma-glutamyl hydrolase



• Molecule 1: Gamma-glutamyl hydrolase



● Molecule 1: Gamma-glutamyl hydrolase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	53.40Å 198.10Å 65.11Å 90.00° 113.82° 90.00°	Depositor
Resolution (Å)	30.00 – 1.97 30.00 – 1.97	Depositor EDS
% Data completeness (in resolution range)	99.2 (30.00-1.97) 99.2 (30.00-1.97)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.01 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.175 , 0.239 0.177 , 0.237	Depositor DCC
R_{free} test set	4330 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	20.0	Xtriage
Anisotropy	0.130	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 38.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.048 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10251	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% 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: MTX

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	1.31	8/2373 (0.3%)	1.16	4/3220 (0.1%)
1	B	1.30	3/2373 (0.1%)	1.27	16/3220 (0.5%)
1	C	1.24	3/2373 (0.1%)	1.16	5/3220 (0.2%)
1	D	1.29	5/2373 (0.2%)	1.26	18/3220 (0.6%)
All	All	1.29	19/9492 (0.2%)	1.21	43/12880 (0.3%)

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	51	ASN	N-CA	7.53	1.55	1.46
1	B	51	ASN	N-CA	7.48	1.55	1.46
1	A	71	GLY	N-CA	6.91	1.53	1.45
1	D	17	GLN	CB-CG	-6.34	1.33	1.52
1	C	45	VAL	C-O	6.14	1.30	1.24
1	A	133	ALA	CA-CB	-5.82	1.45	1.53
1	C	51	ASN	N-CA	5.68	1.53	1.46
1	A	215	THR	C-O	5.56	1.30	1.24
1	A	12	ILE	C-O	5.43	1.29	1.24
1	A	97	SER	N-CA	5.42	1.52	1.46
1	B	104	VAL	C-O	5.30	1.29	1.24
1	D	262	THR	CA-CB	5.25	1.61	1.53
1	A	93	LEU	N-CA	5.22	1.52	1.46
1	C	223	ALA	CA-CB	-5.17	1.45	1.53
1	A	252	ALA	CA-CB	5.17	1.61	1.53
1	D	122	LEU	CA-C	-5.13	1.46	1.52
1	A	30	ILE	C-O	-5.12	1.18	1.24
1	D	48	VAL	CA-CB	-5.12	1.47	1.53
1	B	237	ALA	CA-CB	5.04	1.61	1.53

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	50	ILE	N-CA-C	13.08	123.93	110.72
1	D	50	ILE	CA-C-N	-12.36	97.94	121.54
1	D	50	ILE	C-N-CA	-12.36	97.94	121.54
1	B	50	ILE	CA-C-N	-11.08	100.38	121.54
1	B	50	ILE	C-N-CA	-11.08	100.38	121.54
1	D	67	VAL	CB-CA-C	-10.17	92.87	111.18
1	B	132	ILE	CB-CA-C	-9.86	98.09	111.80
1	C	50	ILE	CA-C-N	-9.55	103.30	121.54
1	C	50	ILE	C-N-CA	-9.55	103.30	121.54
1	B	173	ILE	CB-CA-C	-9.49	96.68	110.83
1	B	67	VAL	CB-CA-C	-8.79	95.36	111.18
1	A	173	ILE	CB-CA-C	-8.32	98.43	110.83
1	B	142	VAL	CB-CA-C	-8.23	99.40	112.16
1	A	71	GLY	N-CA-C	-8.05	99.62	112.66
1	D	173	ILE	CB-CA-C	-7.07	100.30	110.83
1	D	132	ILE	CB-CA-C	-6.80	98.93	111.18
1	B	50	ILE	N-CA-C	6.79	120.54	111.44
1	D	74	VAL	N-CA-CB	-6.51	100.48	111.23
1	D	280	ILE	CB-CA-C	-6.36	103.55	112.14
1	B	50	ILE	O-C-N	-6.29	113.20	121.82
1	B	20	PHE	N-CA-C	6.24	118.17	111.36
1	D	20	PHE	N-CA-C	6.12	118.74	111.71
1	C	50	ILE	O-C-N	-5.98	115.67	121.83
1	B	241	THR	CA-CB-OG1	5.69	118.14	109.60
1	D	50	ILE	O-C-N	-5.57	114.19	121.82
1	B	274	LYS	CA-C-N	-5.46	114.95	120.85
1	B	274	LYS	C-N-CA	-5.46	114.95	120.85
1	D	67	VAL	CG1-CB-CG2	5.45	122.80	110.80
1	D	21	ASP	CA-C-N	5.43	125.34	119.85
1	D	21	ASP	C-N-CA	5.43	125.34	119.85
1	D	118	THR	N-CA-C	5.27	118.18	111.69
1	D	240	THR	CA-C-N	-5.24	113.26	120.28
1	D	240	THR	C-N-CA	-5.24	113.26	120.28
1	B	67	VAL	CG1-CB-CG2	5.23	122.31	110.80
1	C	59	ARG	CB-CG-CD	5.22	123.30	111.30
1	B	280	ILE	CB-CA-C	-5.21	105.10	112.14
1	A	169	HIS	CA-C-N	-5.20	114.36	122.37
1	A	169	HIS	C-N-CA	-5.20	114.36	122.37
1	D	67	VAL	CA-CB-CG2	5.20	119.24	110.40
1	B	67	VAL	CA-CB-CG2	5.20	119.24	110.40
1	D	173	ILE	N-CA-CB	5.20	118.32	111.25
1	B	241	THR	OG1-CB-CG2	5.14	119.58	109.30
1	D	50	ILE	N-CA-C	5.01	118.15	111.44

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	2308	0	2229	35	0
1	B	2308	0	2229	63	0
1	C	2308	0	2229	34	0
1	D	2308	0	2229	41	0
2	B	33	0	20	15	0
2	D	33	0	20	14	0
3	A	151	0	0	6	0
3	B	271	0	0	5	0
3	C	265	0	0	5	0
3	D	266	0	0	1	0
All	All	10251	0	8956	173	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (173) 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)
2:D:301:MTX:H13	2:D:301:MTX:H7	1.28	1.09
1:B:6:THR:HG23	3:B:654:HOH:O	1.49	1.09
1:B:59:ARG:HD3	3:B:633:HOH:O	1.54	1.06
1:C:59:ARG:HD2	3:C:317:HOH:O	1.57	1.04
2:D:301:MTX:H7	2:D:301:MTX:C13	1.96	0.96
2:B:301:MTX:C13	2:B:301:MTX:H7	1.98	0.91
1:B:65:ASN:HD22	1:B:255:ASN:HD22	1.18	0.90
2:B:301:MTX:H7	2:B:301:MTX:H13	1.55	0.89
1:A:287:THR:HG21	3:A:423:HOH:O	1.76	0.85
1:B:270:ILE:HD11	1:B:288:TYR:CZ	2.12	0.84
1:D:65:ASN:HD22	1:D:255:ASN:HD22	1.24	0.83
1:C:150:GLU:HA	1:C:150:GLU:OE1	1.79	0.83
1:D:222:ASN:HB2	1:D:241:THR:CG2	2.09	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:222:ASN:HB2	1:B:241:THR:CG2	2.08	0.82
1:C:287:THR:HG21	3:C:304:HOH:O	1.80	0.81
1:A:76:LEU:HD11	3:A:435:HOH:O	1.81	0.80
1:B:129:THR:HA	1:B:132:ILE:HD11	1.65	0.77
1:A:222:ASN:HB2	1:A:241:THR:CG2	2.15	0.77
1:C:222:ASN:HB2	1:C:241:THR:CG2	2.16	0.76
1:B:101:TYR:CZ	1:B:211:PRO:HB3	2.20	0.76
1:B:287:THR:HG21	3:B:606:HOH:O	1.85	0.76
1:C:124:LEU:HD23	1:C:173:ILE:CD1	2.17	0.74
1:C:54:GLU:HG3	3:C:525:HOH:O	1.87	0.74
1:D:20:PHE:CD1	2:D:301:MTX:C2	2.71	0.73
2:D:301:MTX:H13	2:D:301:MTX:C7	2.14	0.73
1:B:278:THR:HG21	1:B:287:THR:HG23	1.71	0.71
1:B:104:VAL:HB	1:B:212:ILE:HD13	1.72	0.71
1:A:149:LYS:HB3	1:A:150:GLU:OE2	1.90	0.71
1:D:20:PHE:CD1	2:D:301:MTX:N1	2.58	0.71
1:D:20:PHE:HB3	2:D:301:MTX:N3	2.07	0.69
1:A:112:GLU:OE1	3:A:435:HOH:O	2.09	0.69
1:B:101:TYR:OH	1:B:211:PRO:HB3	1.92	0.69
1:D:35:VAL:HG21	1:D:270:ILE:CD1	2.24	0.68
1:B:74:VAL:HG22	1:B:79:SER:HB3	1.78	0.65
1:B:20:PHE:CZ	2:B:301:MTX:C8A	2.80	0.65
1:B:20:PHE:CD2	2:B:301:MTX:C2	2.80	0.65
1:D:20:PHE:CZ	2:D:301:MTX:N8	2.66	0.64
1:C:168:SER:HB3	1:A:128:ASN:ND2	2.13	0.64
1:B:6:THR:CG2	3:B:654:HOH:O	2.24	0.63
1:C:278:THR:HG21	1:C:287:THR:HG23	1.81	0.62
1:B:6:THR:HG21	3:B:539:HOH:O	1.99	0.62
1:D:129:THR:HG22	1:D:132:ILE:HD11	1.80	0.62
1:D:20:PHE:CE1	2:D:301:MTX:N1	2.68	0.62
1:B:222:ASN:HB2	1:B:241:THR:HG21	1.81	0.62
1:C:168:SER:HB3	1:A:128:ASN:HD21	1.64	0.61
1:A:278:THR:HG21	1:A:287:THR:HG23	1.82	0.61
1:B:20:PHE:CD1	2:B:301:MTX:C4A	2.84	0.61
1:D:20:PHE:CZ	2:D:301:MTX:C8A	2.84	0.61
1:B:65:ASN:HD22	1:B:255:ASN:ND2	1.96	0.61
1:A:96:ASN:HA	1:A:100:ASP:O	2.00	0.61
1:C:138:PHE:CD2	1:C:142:VAL:HG11	2.35	0.61
1:B:238:ILE:HD13	1:A:253:ARG:NH2	2.16	0.60
1:D:50:ILE:HD11	1:D:81:TYR:HA	1.84	0.60
1:C:239:LYS:NZ	3:C:537:HOH:O	2.30	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:VAL:HB	1:B:212:ILE:CD1	2.31	0.60
1:A:52:LYS:HB2	1:A:57:TYR:CZ	2.37	0.60
1:C:219:PRO:O	1:C:241:THR:HB	2.02	0.59
1:D:35:VAL:HG11	1:D:270:ILE:HD13	1.85	0.58
1:B:255:ASN:HD21	1:B:257:HIS:CG	2.21	0.58
1:B:140:GLU:O	1:B:143:LYS:HE3	2.04	0.58
1:A:115:THR:HG21	1:A:173:ILE:HD11	1.85	0.58
1:A:219:PRO:O	1:A:241:THR:HB	2.03	0.57
1:C:138:PHE:HB3	1:C:142:VAL:HG11	1.86	0.57
1:D:173:ILE:HD12	1:D:173:ILE:O	2.04	0.57
1:D:222:ASN:HB2	1:D:241:THR:HG21	1.84	0.57
1:B:278:THR:CG2	1:B:287:THR:HG23	2.34	0.57
1:A:138:PHE:HB3	1:A:142:VAL:HG11	1.86	0.57
1:B:20:PHE:CE2	2:B:301:MTX:N1	2.72	0.56
1:D:16:ALA:O	1:D:50:ILE:HD13	2.05	0.56
1:B:250:ASN:ND2	1:B:253:ARG:HH11	2.04	0.56
1:D:20:PHE:CG	2:D:301:MTX:C2	2.88	0.56
1:B:255:ASN:ND2	1:B:257:HIS:H	2.04	0.56
1:C:59:ARG:HH22	1:C:264:GLU:HB3	1.71	0.56
1:A:169:HIS:HD2	3:A:302:HOH:O	1.89	0.56
1:B:170:GLN:HG3	1:B:171:TRP:CD1	2.41	0.55
2:B:301:MTX:H13	2:B:301:MTX:C7	2.33	0.55
1:B:20:PHE:CG	2:B:301:MTX:C4	2.90	0.54
1:A:222:ASN:HB2	1:A:241:THR:HG21	1.87	0.54
1:D:118:THR:CG2	1:D:188:TYR:OH	2.55	0.54
1:D:219:PRO:O	1:D:241:THR:HB	2.07	0.53
1:D:20:PHE:CE1	2:D:301:MTX:C8A	2.92	0.53
1:D:50:ILE:HG23	1:D:51:ASN:ND2	2.23	0.53
1:B:129:THR:HG22	1:B:132:ILE:HD11	1.92	0.52
1:D:255:ASN:HD21	1:D:257:HIS:CG	2.27	0.52
1:D:250:ASN:ND2	1:D:253:ARG:HH11	2.08	0.52
1:B:270:ILE:HD11	1:B:288:TYR:CE2	2.44	0.52
1:A:115:THR:HG21	1:A:173:ILE:CD1	2.40	0.52
1:D:20:PHE:CD2	2:D:301:MTX:C4A	2.93	0.52
1:B:238:ILE:CD1	1:A:253:ARG:NH2	2.72	0.51
1:D:118:THR:HG21	1:D:188:TYR:OH	2.11	0.51
1:B:20:PHE:CZ	2:B:301:MTX:N8	2.79	0.51
1:A:263:GLU:O	1:A:267:LYS:HG3	2.11	0.50
1:D:142:VAL:CG1	1:D:191:LEU:O	2.60	0.50
1:C:168:SER:CB	1:A:128:ASN:HD21	2.24	0.50
1:C:278:THR:CG2	1:C:287:THR:HG23	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:118:THR:HG21	1:C:188:TYR:OH	2.12	0.50
1:C:124:LEU:HD23	1:C:173:ILE:HD11	1.93	0.49
1:B:20:PHE:CE1	2:B:301:MTX:C4A	2.95	0.49
1:C:119:SER:HB3	1:C:184:LEU:HG	1.93	0.49
1:C:124:LEU:CD2	1:C:173:ILE:HD11	2.43	0.49
1:B:67:VAL:HG22	1:B:102:PHE:HE2	1.77	0.49
1:B:20:PHE:CE1	2:B:301:MTX:C8A	2.96	0.48
1:C:124:LEU:HD23	1:C:173:ILE:HD13	1.92	0.48
1:B:238:ILE:HD13	1:A:253:ARG:HH22	1.79	0.48
1:D:278:THR:HG21	1:D:287:THR:HG23	1.95	0.48
1:B:270:ILE:CD1	1:B:288:TYR:CZ	2.93	0.48
1:A:71:GLY:HA2	1:A:109:LEU:H	1.78	0.48
1:D:222:ASN:HB2	1:D:241:THR:HG22	1.92	0.47
1:C:130:SER:OG	1:A:170:GLN:NE2	2.46	0.47
1:B:118:THR:HG21	1:B:188:TYR:OH	2.14	0.47
1:A:165:THR:HB	1:A:217:TRP:CE3	2.49	0.47
1:D:255:ASN:ND2	1:D:257:HIS:H	2.13	0.47
1:A:79:SER:O	1:A:83:LYS:HD3	2.14	0.47
1:A:118:THR:HG21	1:A:188:TYR:OH	2.14	0.47
1:C:126:HIS:CE1	1:C:171:TRP:CH2	3.03	0.47
1:B:114:LEU:O	1:B:118:THR:HG23	2.14	0.46
1:B:263:GLU:H	1:B:263:GLU:CD	2.23	0.46
1:B:138:PHE:CD2	1:B:142:VAL:HG11	2.51	0.46
1:D:113:LEU:HD22	1:D:117:LEU:HG	1.98	0.46
1:B:74:VAL:HG22	1:B:79:SER:CB	2.45	0.46
1:B:270:ILE:HD13	3:A:347:HOH:O	2.16	0.46
1:A:169:HIS:CD2	3:A:302:HOH:O	2.65	0.46
1:C:118:THR:CG2	1:C:188:TYR:OH	2.63	0.46
1:A:217:TRP:H	1:A:217:TRP:CD1	2.32	0.46
1:D:15:LEU:HD23	1:D:48:VAL:HB	1.98	0.45
1:D:20:PHE:CE2	2:D:301:MTX:C8A	2.99	0.45
1:D:21:ASP:OD1	3:D:488:HOH:O	2.21	0.45
1:D:35:VAL:CG2	1:D:270:ILE:CD1	2.94	0.45
1:A:17:GLN:OE1	1:A:72:GLY:HA3	2.17	0.45
1:D:50:ILE:HG13	1:D:80:GLY:C	2.42	0.45
1:A:124:LEU:CD2	1:A:173:ILE:HD13	2.47	0.45
1:B:20:PHE:CE2	2:B:301:MTX:C8A	3.01	0.44
1:B:139:THR:O	1:B:142:VAL:HG22	2.18	0.44
1:D:204:THR:HG23	1:D:215:THR:HG22	2.00	0.44
1:A:114:LEU:O	1:A:118:THR:HG23	2.18	0.44
1:B:36:LYS:NZ	1:B:286:GLN:NE2	2.65	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:255:ASN:HD21	1:B:257:HIS:CB	2.31	0.44
1:C:122:LEU:HD13	1:C:124:LEU:H	1.83	0.43
1:B:101:TYR:CE2	1:B:211:PRO:HB3	2.51	0.43
1:D:117:LEU:HD23	1:D:117:LEU:HA	1.87	0.43
1:B:118:THR:CG2	1:B:188:TYR:OH	2.67	0.43
1:D:35:VAL:HG21	1:D:270:ILE:HD11	1.96	0.43
2:B:301:MTX:C13	2:B:301:MTX:C7	2.86	0.43
1:B:111:PHE:CD2	1:B:111:PHE:C	2.96	0.43
1:B:219:PRO:O	1:B:241:THR:HB	2.18	0.43
1:A:278:THR:CG2	1:A:287:THR:HG23	2.48	0.43
1:A:140:GLU:O	1:A:143:LYS:HB2	2.19	0.43
1:C:114:LEU:O	1:C:118:THR:HG23	2.20	0.42
1:C:174:THR:HA	1:C:201:PHE:HA	2.01	0.42
1:A:124:LEU:HB3	1:A:171:TRP:HB3	2.01	0.42
1:D:20:PHE:CB	2:D:301:MTX:N3	2.81	0.42
1:C:222:ASN:HB2	1:C:241:THR:HG21	1.96	0.42
1:B:20:PHE:CD1	2:B:301:MTX:C4	3.03	0.42
1:B:36:LYS:HZ1	1:B:286:GLN:NE2	2.18	0.42
2:B:301:MTX:HM1	2:B:301:MTX:H15	1.67	0.42
1:C:281:GLN:O	1:A:197:GLY:HA2	2.20	0.41
1:D:47:PRO:HB2	1:D:49:MET:CE	2.50	0.41
1:B:26:ARG:NH2	1:B:276:GLU:OE2	2.53	0.41
1:B:115:THR:HG21	1:B:173:ILE:HD13	2.02	0.41
1:B:178:PHE:HZ	1:B:188:TYR:HB2	1.85	0.41
1:C:278:THR:OG1	1:C:287:THR:CG2	2.68	0.41
1:B:35:VAL:HG11	1:B:270:ILE:HD12	2.02	0.41
1:B:217:TRP:CD1	1:B:217:TRP:H	2.39	0.41
1:B:278:THR:OG1	1:B:287:THR:CG2	2.69	0.41
1:B:50:ILE:HG23	1:B:51:ASN:ND2	2.36	0.41
1:B:115:THR:HG21	1:B:173:ILE:CD1	2.51	0.41
1:C:138:PHE:CB	1:C:142:VAL:HG11	2.50	0.41
1:C:86:GLY:O	1:C:90:ARG:HG3	2.22	0.40
1:D:35:VAL:CG1	1:D:270:ILE:HD13	2.52	0.40
1:D:226:TRP:HB2	1:D:285:GLU:OE1	2.21	0.40
1:C:60:LEU:CD1	1:C:64:ILE:HD13	2.51	0.40
1:C:182:LYS:HG3	3:C:322:HOH:O	2.21	0.40
1:B:263:GLU:O	1:B:267:LYS:HG3	2.21	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	285/312 (91%)	276 (97%)	8 (3%)	1 (0%)	30	20
1	B	285/312 (91%)	276 (97%)	8 (3%)	1 (0%)	30	20
1	C	285/312 (91%)	278 (98%)	7 (2%)	0	100	100
1	D	285/312 (91%)	279 (98%)	5 (2%)	1 (0%)	30	20
All	All	1140/1248 (91%)	1109 (97%)	28 (2%)	3 (0%)	36	27

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	143	LYS
1	B	131	GLY
1	D	131	GLY

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	249/271 (92%)	238 (96%)	11 (4%)	25	14
1	B	249/271 (92%)	230 (92%)	19 (8%)	12	4
1	C	249/271 (92%)	236 (95%)	13 (5%)	21	9
1	D	249/271 (92%)	230 (92%)	19 (8%)	12	4
All	All	996/1084 (92%)	934 (94%)	62 (6%)	16	7

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	26	ARG
1	C	50	ILE
1	C	59	ARG
1	C	113	LEU
1	C	118	THR
1	C	132	ILE
1	C	143	LYS
1	C	147	LEU
1	C	150	GLU
1	C	168	SER
1	C	173	ILE
1	C	241	THR
1	C	287	THR
1	D	17	GLN
1	D	18	ASP
1	D	50	ILE
1	D	58	SER
1	D	60	LEU
1	D	67	VAL
1	D	74	VAL
1	D	113	LEU
1	D	118	THR
1	D	132	ILE
1	D	147	LEU
1	D	173	ILE
1	D	241	THR
1	D	262	THR
1	D	263	GLU
1	D	267	LYS
1	D	280	ILE
1	D	282	SER
1	D	287	THR
1	B	6	THR
1	B	18	ASP
1	B	23	LYS
1	B	50	ILE
1	B	60	LEU
1	B	67	VAL
1	B	78	SER
1	B	113	LEU
1	B	118	THR
1	B	132	ILE
1	B	142	VAL

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Mol	Chain	Res	Type
1	B	173	ILE
1	B	186	LYS
1	B	241	THR
1	B	263	GLU
1	B	270	ILE
1	B	280	ILE
1	B	281	GLN
1	B	287	THR
1	A	15	LEU
1	A	60	LEU
1	A	83	LYS
1	A	118	THR
1	A	121	GLU
1	A	122	LEU
1	A	173	ILE
1	A	241	THR
1	A	262	THR
1	A	281	GLN
1	A	287	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	128	ASN
1	C	250	ASN
1	C	272	ASN
1	D	51	ASN
1	D	216	GLN
1	D	250	ASN
1	D	255	ASN
1	D	286	GLN
1	B	51	ASN
1	B	98	ASN
1	B	167	ASN
1	B	170	GLN
1	B	218	HIS
1	B	250	ASN
1	B	255	ASN
1	B	286	GLN
1	A	169	HIS
1	A	170	GLN
1	A	246	ASN

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Mol	Chain	Res	Type
1	A	250	ASN
1	A	281	GLN

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 ⓘ

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	MTX	D	301	-	35,35,35	2.11	7 (20%)	47,49,49	2.20	12 (25%)
2	MTX	B	301	-	35,35,35	2.22	11 (31%)	47,49,49	2.13	15 (31%)

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	MTX	D	301	-	-	1/25/25/25	0/3/3/3
2	MTX	B	301	-	-	1/25/25/25	0/3/3/3

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	301	MTX	C9-C6	-6.32	1.40	1.51
2	B	301	MTX	C7-N8	5.94	1.41	1.31
2	B	301	MTX	C9-C6	-5.81	1.41	1.51
2	B	301	MTX	O-C	5.37	1.35	1.23
2	D	301	MTX	O-C	5.26	1.35	1.23
2	D	301	MTX	C7-N8	4.76	1.39	1.31
2	D	301	MTX	C11-C	-4.22	1.41	1.50
2	B	301	MTX	C11-C	-3.47	1.42	1.50
2	B	301	MTX	O2-CT	-2.65	1.22	1.30
2	D	301	MTX	C13-C14	2.41	1.43	1.39
2	B	301	MTX	C13-C12	2.40	1.42	1.38
2	B	301	MTX	C6-N5	2.35	1.36	1.32
2	D	301	MTX	C16-C11	2.29	1.42	1.39
2	D	301	MTX	O2-CT	-2.24	1.23	1.30
2	B	301	MTX	C13-C14	2.16	1.43	1.39
2	B	301	MTX	C12-C11	2.11	1.42	1.39
2	B	301	MTX	CA-CT	2.10	1.58	1.52
2	B	301	MTX	C16-C11	2.07	1.42	1.39

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	MTX	C2-N1-C8A	6.24	122.21	115.48
2	B	301	MTX	N1-C2-N3	-5.90	119.70	127.21
2	B	301	MTX	C2-N1-C8A	5.88	121.82	115.48
2	D	301	MTX	N1-C2-N3	-5.80	119.84	127.21
2	B	301	MTX	C4-C4A-N5	4.08	123.47	120.33
2	D	301	MTX	C4-C4A-N5	3.77	123.23	120.33
2	B	301	MTX	C7-C6-N5	-3.75	118.44	120.87
2	D	301	MTX	N8-C8A-N1	3.72	119.82	115.77
2	D	301	MTX	C13-C12-C11	-3.70	116.84	120.80
2	B	301	MTX	N8-C8A-N1	3.70	119.80	115.77
2	B	301	MTX	C12-C13-C14	3.55	124.80	120.30
2	D	301	MTX	C12-C13-C14	3.46	124.68	120.30
2	D	301	MTX	C6-C7-N8	-3.24	120.03	123.14
2	B	301	MTX	C13-C12-C11	-3.23	117.35	120.80
2	D	301	MTX	C6-C9-N10	-3.19	107.49	112.95
2	D	301	MTX	CB-CA-N	-3.10	104.78	110.91
2	D	301	MTX	C15-C16-C11	2.94	123.94	120.80
2	B	301	MTX	C15-C16-C11	2.80	123.79	120.80
2	D	301	MTX	C8A-C4A-N5	-2.71	119.32	122.35
2	B	301	MTX	C9-C6-N5	2.71	121.49	117.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	MTX	C9-N10-C14	2.56	123.27	120.17
2	B	301	MTX	NA2-C2-N3	2.41	120.84	117.22
2	B	301	MTX	C6-N5-C4A	2.33	121.46	118.17
2	D	301	MTX	C9-C6-N5	2.30	120.83	117.09
2	B	301	MTX	C8A-C4A-N5	-2.17	119.92	122.35
2	B	301	MTX	OE1-CD-CG	-2.14	116.30	123.09
2	B	301	MTX	C6-C7-N8	-2.06	121.16	123.14

There are no chirality outliers.

All (2) torsion outliers are listed below:

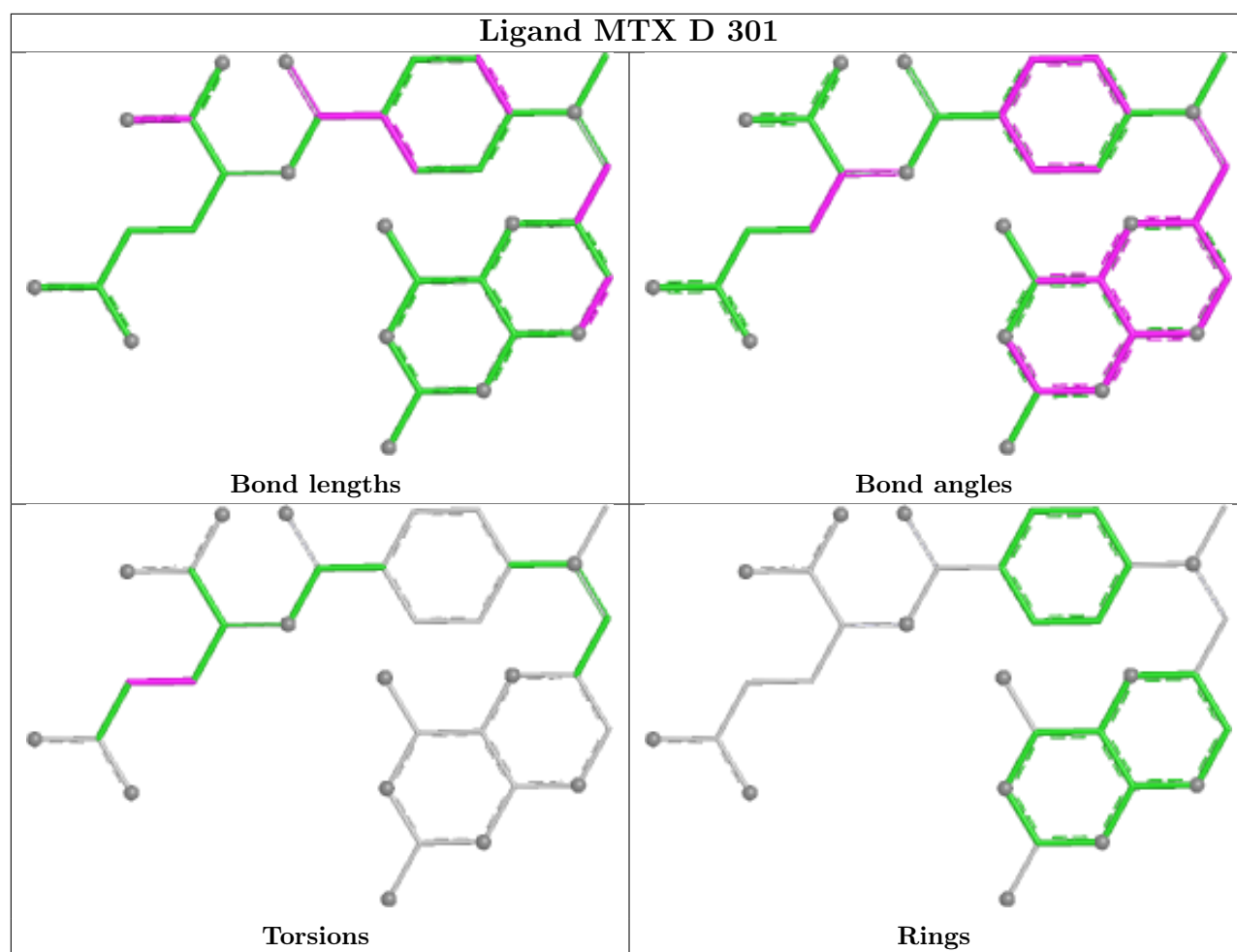
Mol	Chain	Res	Type	Atoms
2	B	301	MTX	CA-CB-CG-CD
2	D	301	MTX	CA-CB-CG-CD

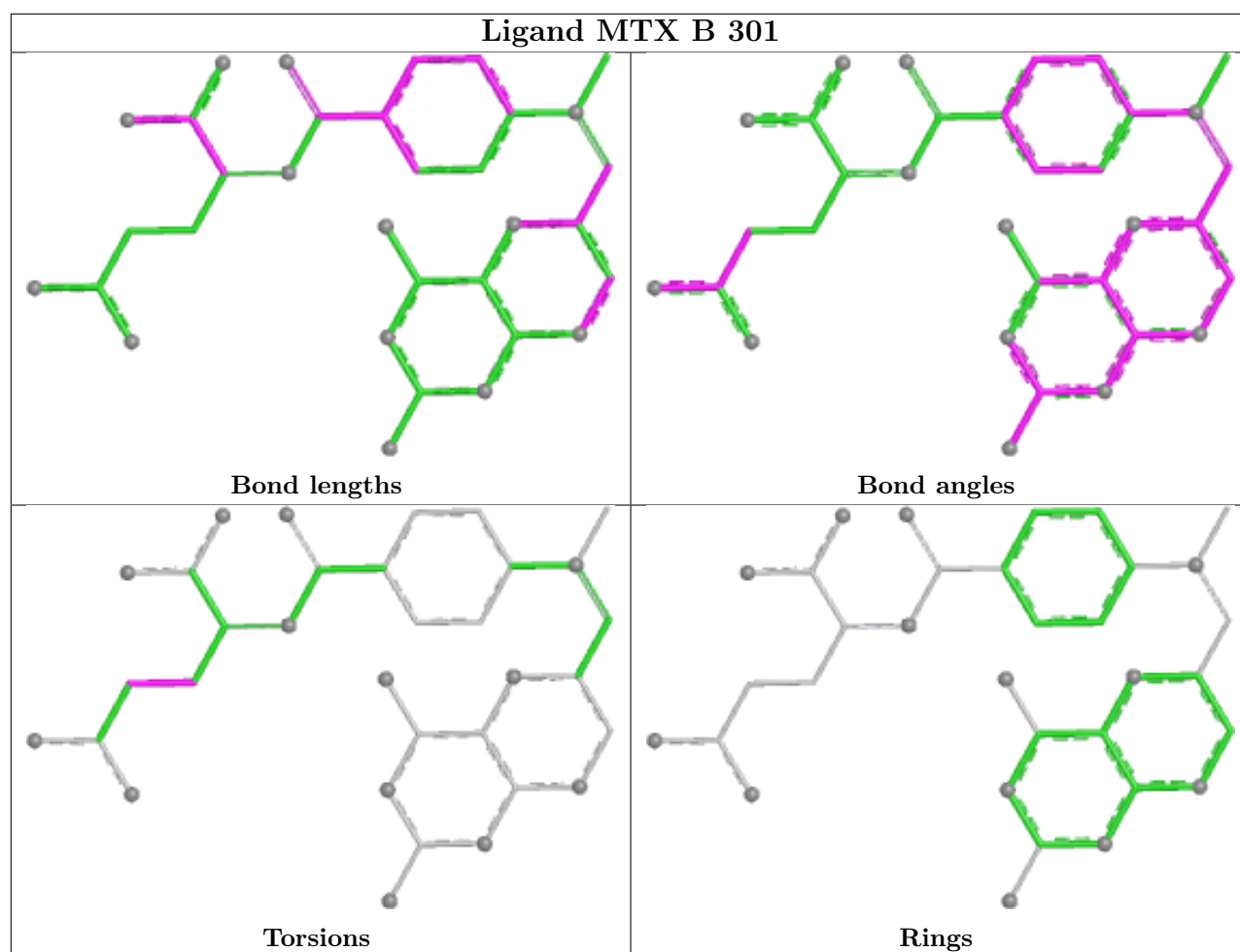
There are no ring outliers.

2 monomers are involved in 29 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	301	MTX	14	0
2	B	301	MTX	15	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	287/312 (91%)	-0.07	14 (4%)	35 44	2, 18, 34, 47	9 (3%)
1	B	287/312 (91%)	-0.23	8 (2%)	55 65	3, 17, 31, 41	5 (1%)
1	C	287/312 (91%)	-0.18	9 (3%)	51 61	3, 18, 32, 41	8 (2%)
1	D	287/312 (91%)	-0.21	7 (2%)	59 69	3, 17, 31, 41	5 (1%)
All	All	1148/1248 (91%)	-0.17	38 (3%)	49 59	2, 17, 32, 47	27 (2%)

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	179	THR	6.7
1	A	283	ALA	6.3
1	C	179	THR	5.9
1	C	283	ALA	4.9
1	D	283	ALA	4.6
1	B	101	TYR	4.6
1	B	250	ASN	4.5
1	D	158	SER	4.5
1	A	250	ASN	4.4
1	C	250	ASN	4.4
1	B	158	SER	4.1
1	B	283	ALA	4.0
1	D	250	ASN	3.7
1	A	20	PHE	3.6
1	A	128	ASN	3.2
1	B	150	GLU	3.2
1	A	280	ILE	3.1
1	A	76	LEU	3.1
1	C	202	VAL	3.0
1	A	276	GLU	3.0
1	C	150	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	280	ILE	2.9
1	A	115	THR	2.9
1	D	20	PHE	2.8
1	D	50	ILE	2.7
1	A	100	ASP	2.6
1	C	280	ILE	2.6
1	C	50	ILE	2.5
1	D	280	ILE	2.5
1	C	115	THR	2.5
1	A	202	VAL	2.5
1	B	100	ASP	2.4
1	B	20	PHE	2.3
1	C	276	GLU	2.3
1	D	115	THR	2.2
1	A	274	LYS	2.1
1	A	112	GLU	2.0
1	A	73	GLY	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](#)

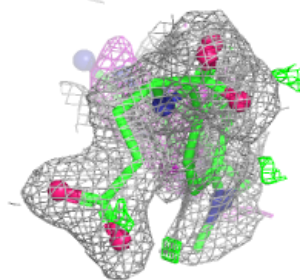
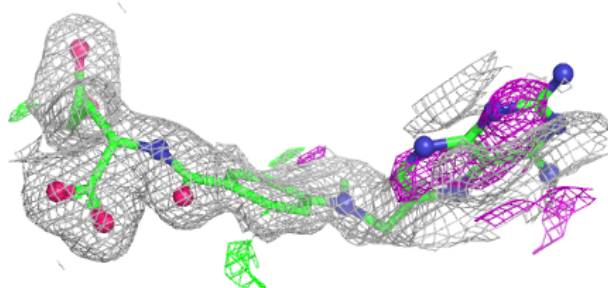
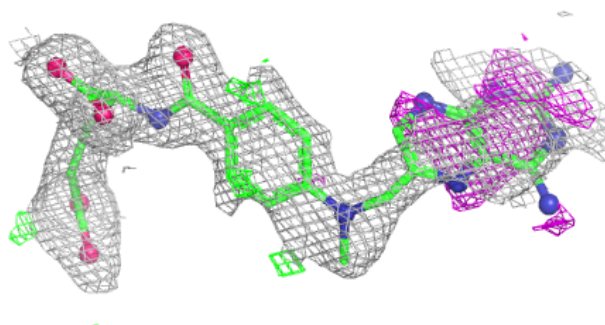
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	MTX	B	301	33/33	0.86	0.15	17,42,86,88	0
2	MTX	D	301	33/33	0.87	0.14	16,39,80,80	0

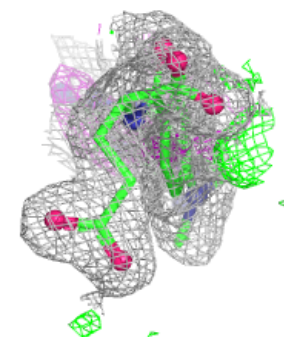
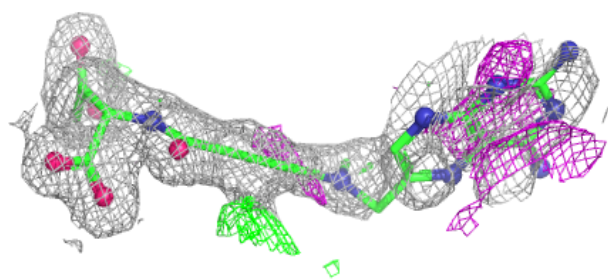
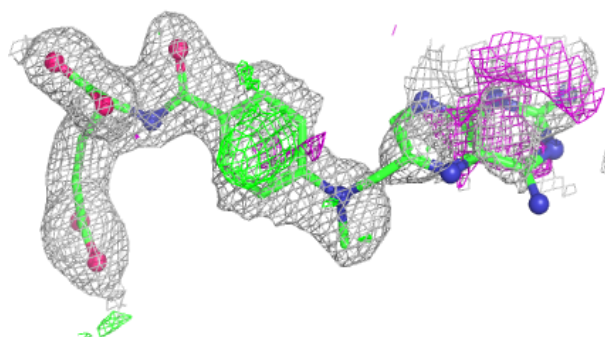
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around MTX B 301:

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

**Electron density around MTX D 301:**

$2mF_o-DF_c$ (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.