



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:19 AM UTC

PDB ID : 8J65 / pdb_00008j65
Title : the structure of L,D-transpeptidase LdtMt2 complex with citrate
Authors : Li, D.F.; Qin, Y.L.
Deposited on : 2023-04-24
Resolution : 1.80 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

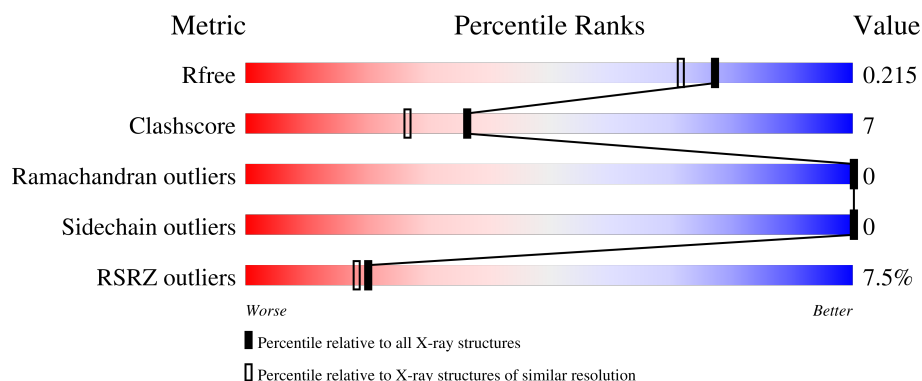
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	293	8% 72% 16% 12%
1	B	293	6% 78% 10% 12%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4345 atoms, of which 10 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 L,D-transpeptidase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	259	Total	C	N	O	S	0	0	0
			1991	1254	345	385	7			
1	B	259	Total	C	N	O	S	0	0	0
			1991	1254	345	385	7			

There are 48 discrepancies between the modelled and reference sequences:

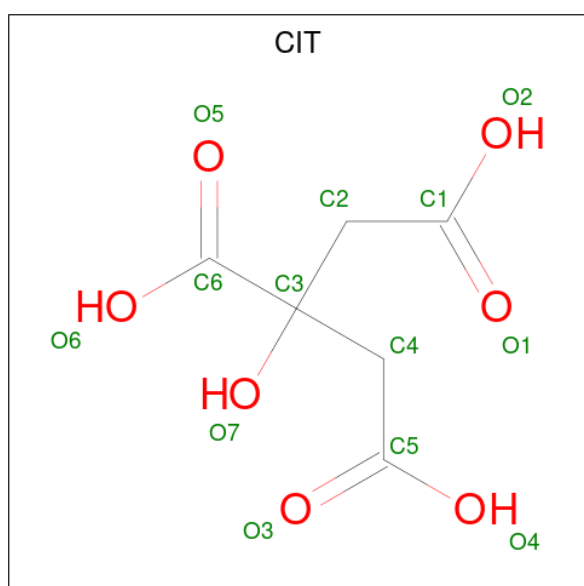
Chain	Residue	Modelled	Actual	Comment	Reference
A	116	MET	-	initiating methionine	UNP O53223
A	117	HIS	-	expression tag	UNP O53223
A	118	HIS	-	expression tag	UNP O53223
A	119	HIS	-	expression tag	UNP O53223
A	120	HIS	-	expression tag	UNP O53223
A	121	HIS	-	expression tag	UNP O53223
A	122	HIS	-	expression tag	UNP O53223
A	123	SER	-	expression tag	UNP O53223
A	124	SER	-	expression tag	UNP O53223
A	125	GLY	-	expression tag	UNP O53223
A	126	VAL	-	expression tag	UNP O53223
A	127	ASP	-	expression tag	UNP O53223
A	128	LEU	-	expression tag	UNP O53223
A	129	GLY	-	expression tag	UNP O53223
A	130	THR	-	expression tag	UNP O53223
A	131	GLU	-	expression tag	UNP O53223
A	132	ASN	-	expression tag	UNP O53223
A	133	LEU	-	expression tag	UNP O53223
A	134	TYR	-	expression tag	UNP O53223
A	135	PHE	-	expression tag	UNP O53223
A	136	GLN	-	expression tag	UNP O53223
A	137	SER	-	expression tag	UNP O53223
A	138	ASN	-	expression tag	UNP O53223
A	139	ALA	-	expression tag	UNP O53223
B	116	MET	-	initiating methionine	UNP O53223

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Chain	Residue	Modelled	Actual	Comment	Reference
B	117	HIS	-	expression tag	UNP O53223
B	118	HIS	-	expression tag	UNP O53223
B	119	HIS	-	expression tag	UNP O53223
B	120	HIS	-	expression tag	UNP O53223
B	121	HIS	-	expression tag	UNP O53223
B	122	HIS	-	expression tag	UNP O53223
B	123	SER	-	expression tag	UNP O53223
B	124	SER	-	expression tag	UNP O53223
B	125	GLY	-	expression tag	UNP O53223
B	126	VAL	-	expression tag	UNP O53223
B	127	ASP	-	expression tag	UNP O53223
B	128	LEU	-	expression tag	UNP O53223
B	129	GLY	-	expression tag	UNP O53223
B	130	THR	-	expression tag	UNP O53223
B	131	GLU	-	expression tag	UNP O53223
B	132	ASN	-	expression tag	UNP O53223
B	133	LEU	-	expression tag	UNP O53223
B	134	TYR	-	expression tag	UNP O53223
B	135	PHE	-	expression tag	UNP O53223
B	136	GLN	-	expression tag	UNP O53223
B	137	SER	-	expression tag	UNP O53223
B	138	ASN	-	expression tag	UNP O53223
B	139	ALA	-	expression tag	UNP O53223

- Molecule 2 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			18	6	5	7		
2	B	1	Total	C	H	O	0	0
			18	6	5	7		

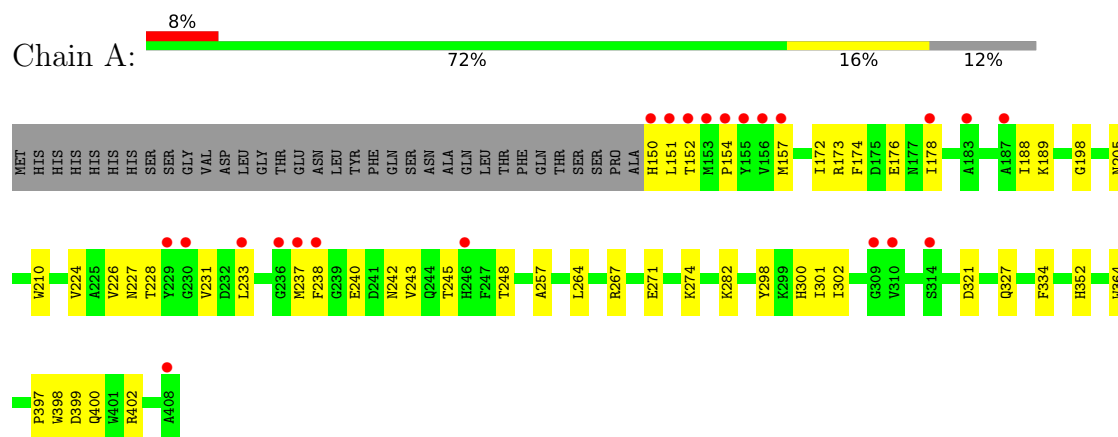
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	166	Total	O	0	0
			166	166		
3	B	161	Total	O	0	0
			161	161		

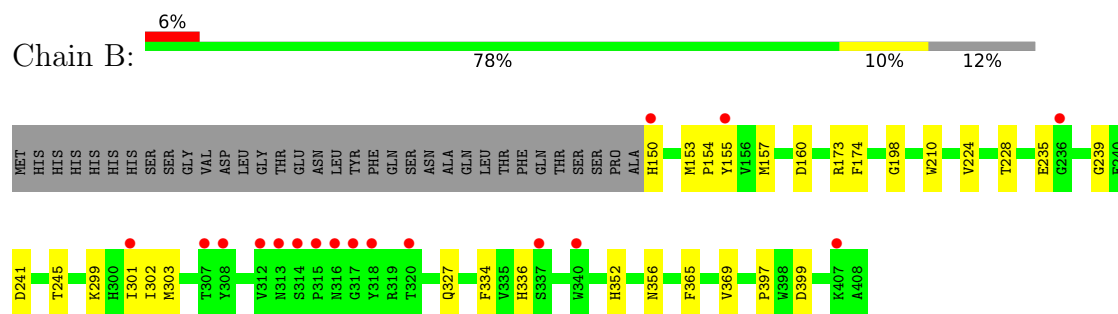
3 Residue-property plots [i](#)

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: L,D-transpeptidase 2



• Molecule 1: L,D-transpeptidase 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	51.18Å 75.89Å 67.07Å 90.00° 111.92° 90.00°	Depositor
Resolution (Å)	48.12 – 1.80 48.12 – 1.80	Depositor EDS
% Data completeness (in resolution range)	95.3 (48.12-1.80) 89.3 (48.12-1.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 1.79Å)	Xtriage
Refinement program	PHENIX (1.14_3260: ???)	Depositor
R, R_{free}	0.212 , 0.236 (Not available) , 0.215	Depositor DCC
R_{free} test set	2008 reflections (4.55%)	wwPDB-VP
Wilson B-factor (Å ²)	21.8	Xtriage
Anisotropy	0.592	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 46.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.046 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4345	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 25.80 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.9617e-03.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.18	0/2047	0.36	0/2799
1	B	0.19	0/2047	0.43	1/2799 (0.0%)
All	All	0.18	0/4094	0.40	1/5598 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	299	LYS	N-CA-CB	-5.54	101.97	110.16

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1991	0	1885	35	0
1	B	1991	0	1885	21	0
2	A	13	5	5	1	0
2	B	13	5	5	3	0
3	A	166	0	0	8	1
3	B	161	0	0	4	0
All	All	4335	10	3780	56	1

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 7.

All (56) 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:151:LEU:HB3	1:A:240:GLU:HG3	1.52	0.91
1:A:157:MET:HE1	1:A:173:ARG:HD2	1.56	0.86
1:A:152:THR:HB	3:A:708:HOH:O	1.87	0.74
1:A:237:MET:SD	3:A:708:HOH:O	2.46	0.72
1:B:327:GLN:OE1	3:B:601:HOH:O	2.09	0.69
1:A:176:GLU:HG3	3:A:708:HOH:O	1.95	0.67
1:A:302:ILE:HG12	1:A:321:ASP:OD2	1.97	0.64
1:B:356:ASN:ND2	3:B:604:HOH:O	2.30	0.63
1:B:155:TYR:OH	1:B:157:MET:SD	2.56	0.62
1:B:365:PHE:O	1:B:369:VAL:HG22	1.99	0.62
1:B:235:GLU:HA	1:B:235:GLU:OE1	2.01	0.58
1:B:224:VAL:HB	1:B:245:THR:HG22	1.87	0.56
1:B:365:PHE:CE1	1:B:369:VAL:HG21	2.42	0.55
1:A:150:HIS:HA	3:A:738:HOH:O	2.08	0.54
1:B:301:ILE:HD12	1:B:302:ILE:H	1.73	0.54
1:B:160:ASP:N	3:B:605:HOH:O	2.33	0.53
1:B:198:GLY:HA3	1:B:210:TRP:CZ2	2.45	0.52
1:A:151:LEU:HB3	1:A:240:GLU:CG	2.35	0.52
1:A:151:LEU:HD13	1:A:240:GLU:CG	2.40	0.52
1:B:327:GLN:HB2	1:B:334:PHE:CE2	2.46	0.51
1:B:173:ARG:HD3	1:B:174:PHE:N	2.25	0.50
1:A:224:VAL:HB	1:A:245:THR:HG22	1.92	0.50
1:B:154:PRO:HD2	1:B:241:ASP:OD1	2.12	0.49
1:A:298:TYR:HB3	1:A:301:ILE:HG23	1.93	0.49
1:B:336:HIS:NE2	2:B:501:CIT:O6	2.43	0.49
1:A:157:MET:CE	1:A:173:ARG:HD2	2.37	0.48
1:A:198:GLY:HA3	1:A:210:TRP:CZ2	2.49	0.48
1:B:352:HIS:HB2	2:B:501:CIT:C5	2.44	0.47
1:A:205:ASN:HA	3:A:622:HOH:O	2.15	0.46
1:A:300:HIS:O	1:A:301:ILE:HG23	2.16	0.46
1:B:303:MET:HE1	2:B:501:CIT:O6	2.15	0.45
1:A:274:LYS:HE3	1:A:364:TRP:CD1	2.51	0.45
1:A:172:ILE:HD13	1:A:188:ILE:HD13	1.98	0.45
1:B:397:PRO:HB2	1:B:399:ASP:OD1	2.16	0.45
1:A:174:PHE:CG	1:A:178:ILE:HD11	2.52	0.44
1:A:189:LYS:NZ	3:A:604:HOH:O	2.31	0.44
1:A:227:ASN:HA	1:A:242:ASN:CG	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:VAL:O	1:A:238:PHE:HA	2.18	0.43
1:A:157:MET:HE3	1:A:157:MET:HB2	1.71	0.43
1:A:257:ALA:HB2	1:A:264:LEU:HD13	1.99	0.43
1:A:352:HIS:HB2	2:A:501:CIT:C5	2.49	0.43
1:B:173:ARG:HD3	1:B:174:PHE:H	1.83	0.43
1:B:153:MET:HE2	1:B:153:MET:HB2	1.69	0.43
1:A:154:PRO:HG3	1:A:228:THR:OG1	2.19	0.42
1:A:152:THR:HG21	1:A:233:LEU:CD1	2.49	0.42
1:B:150:HIS:N	3:B:612:HOH:O	2.52	0.42
1:A:398:TRP:HE1	1:A:402:ARG:NH2	2.17	0.42
1:A:399:ASP:OD1	1:A:400:GLN:N	2.53	0.42
1:A:248:THR:HG23	3:A:687:HOH:O	2.18	0.42
1:B:228:THR:O	1:B:239:GLY:HA3	2.20	0.42
1:A:267:ARG:HA	1:A:271:GLU:O	2.20	0.42
1:A:298:TYR:CD2	1:A:301:ILE:HG21	2.54	0.42
1:A:327:GLN:HB2	1:A:334:PHE:CE2	2.54	0.41
1:A:397:PRO:HB2	1:A:399:ASP:OD1	2.20	0.41
1:A:226:VAL:HB	1:A:243:VAL:HG22	2.03	0.41
1:A:282:LYS:NZ	3:A:613:HOH:O	2.50	0.41

All (1) 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)
3:A:621:HOH:O	3:A:723:HOH:O[1_655]	1.58	0.62

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	257/293 (88%)	250 (97%)	7 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	257/293 (88%)	251 (98%)	6 (2%)	0	100	100
All	All	514/586 (88%)	501 (98%)	13 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	212/242 (88%)	212 (100%)	0	100	100
1	B	212/242 (88%)	212 (100%)	0	100	100
All	All	424/484 (88%)	424 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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	316	ASN
1	B	316	ASN
1	B	347	HIS
1	B	356	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

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	CIT	B	501	-	12,12,12	1.04	0	17,17,17	1.43	1 (5%)
2	CIT	A	501	-	12,12,12	1.10	0	17,17,17	1.90	6 (35%)

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	CIT	B	501	-	-	6/16/16/16	-
2	CIT	A	501	-	-	8/16/16/16	-

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	CIT	O6-C6-C3	3.97	120.76	113.14
2	B	501	CIT	O6-C6-C3	3.70	120.24	113.14
2	A	501	CIT	C2-C3-C6	-2.76	103.92	110.03
2	A	501	CIT	C4-C3-C2	2.43	115.55	109.31
2	A	501	CIT	O2-C1-C2	2.41	121.98	114.35
2	A	501	CIT	O4-C5-C4	2.15	121.17	114.35
2	A	501	CIT	O4-C5-O3	-2.12	117.89	123.33

There are no chirality outliers.

All (14) torsion outliers are listed below:

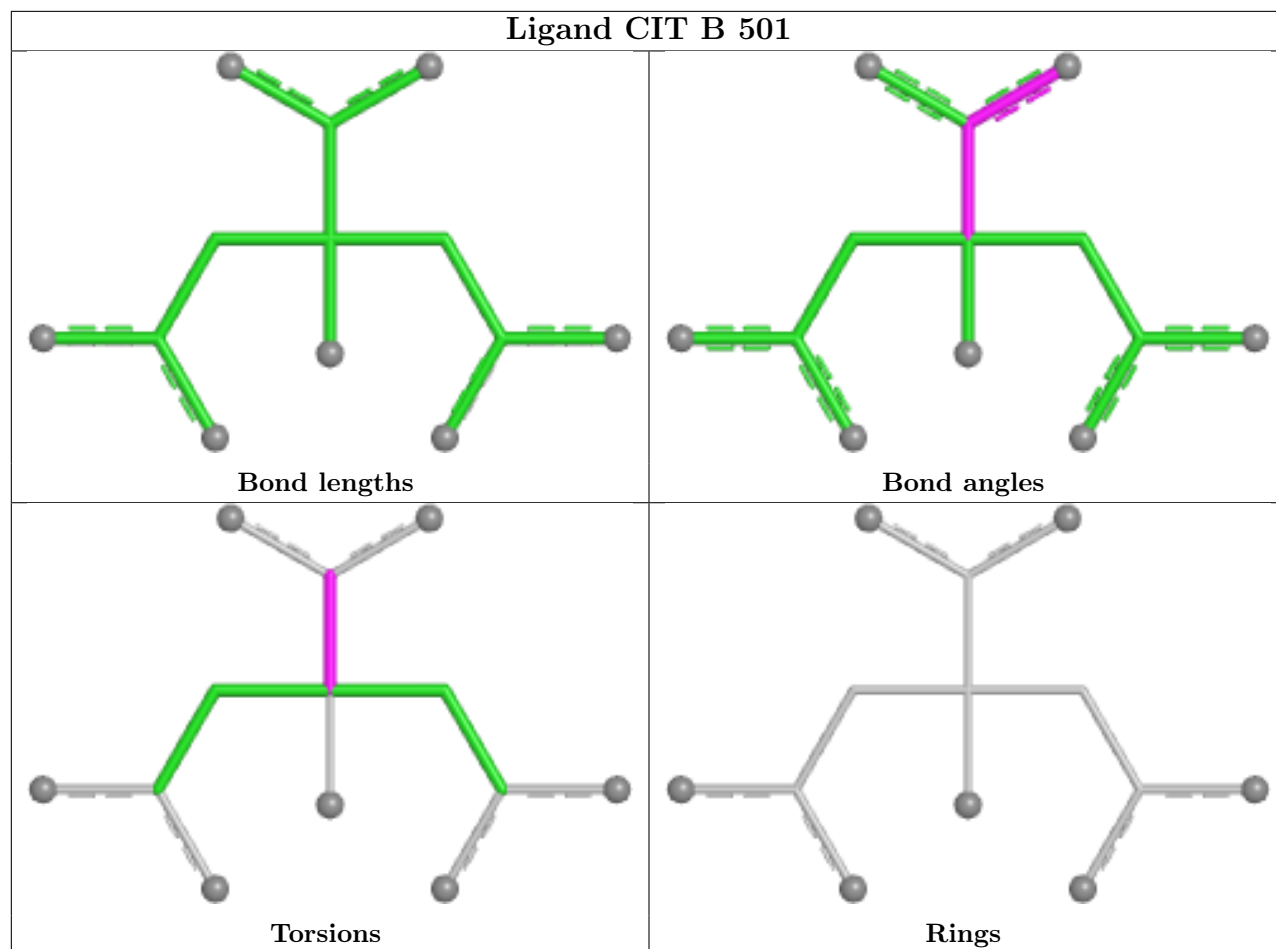
Mol	Chain	Res	Type	Atoms
2	A	501	CIT	O7-C3-C6-O5
2	A	501	CIT	O7-C3-C6-O6
2	A	501	CIT	C4-C3-C6-O5
2	A	501	CIT	C4-C3-C6-O6
2	B	501	CIT	O7-C3-C6-O6
2	B	501	CIT	C4-C3-C6-O6
2	B	501	CIT	C4-C3-C6-O5
2	A	501	CIT	C3-C4-C5-O4
2	A	501	CIT	C3-C4-C5-O3
2	B	501	CIT	C2-C3-C6-O5
2	B	501	CIT	C2-C3-C6-O6
2	B	501	CIT	O7-C3-C6-O5
2	A	501	CIT	O2-C1-C2-C3
2	A	501	CIT	O1-C1-C2-C3

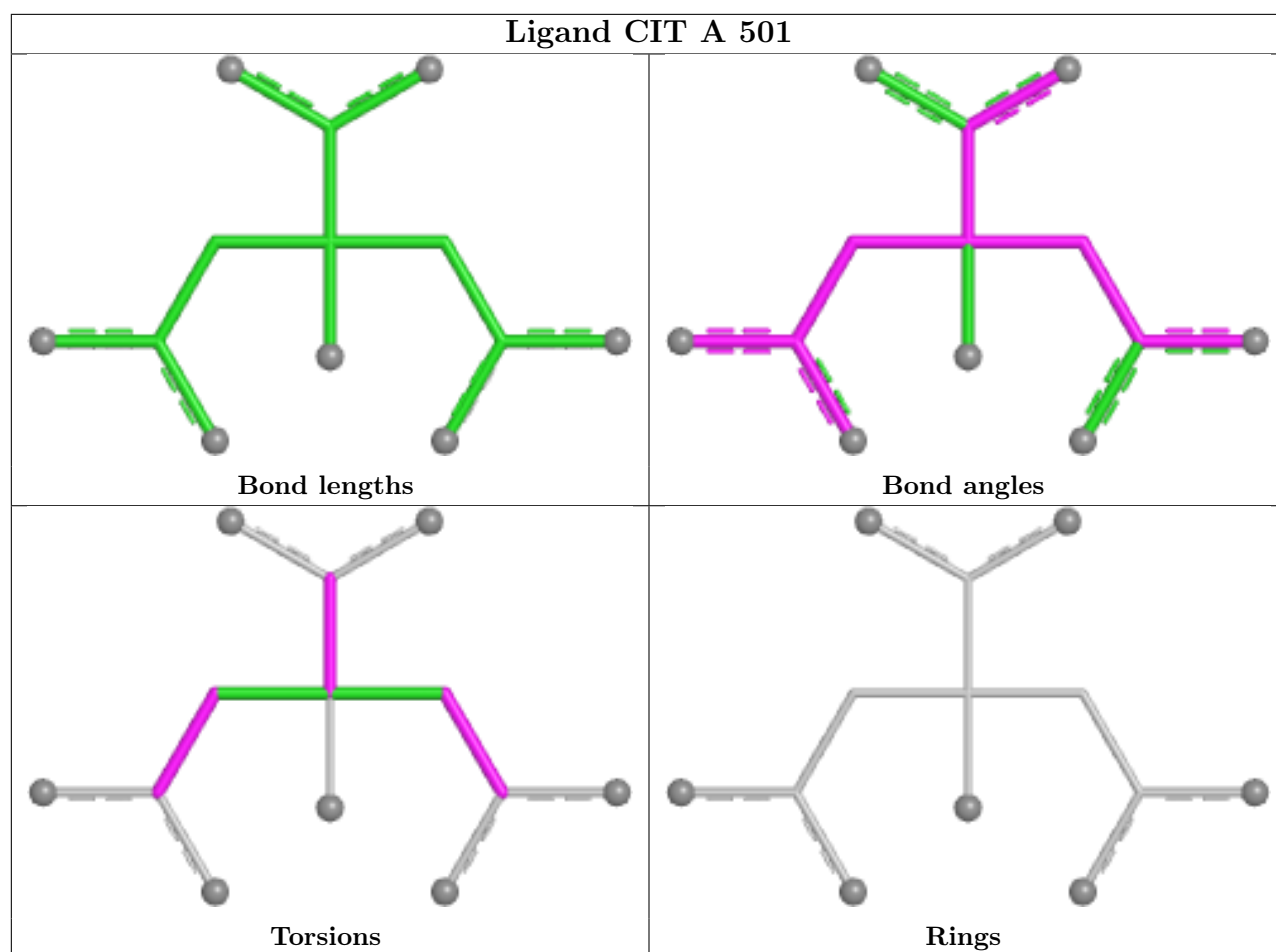
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	CIT	3	0
2	A	501	CIT	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	259/293 (88%)	0.69	22 (8%) 16 14	17, 26, 50, 63	0
1	B	259/293 (88%)	0.66	17 (6%) 24 23	16, 26, 43, 59	0
All	All	518/586 (88%)	0.68	39 (7%) 20 18	16, 26, 46, 63	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	155	TYR	6.0
1	B	150	HIS	4.9
1	B	155	TYR	4.8
1	B	314	SER	4.6
1	B	316	ASN	4.0
1	B	315	PRO	3.8
1	A	150	HIS	3.7
1	A	238	PHE	3.5
1	A	229	TYR	3.4
1	A	233	LEU	3.4
1	A	154	PRO	3.3
1	B	308	TYR	3.2
1	B	313	ASN	3.1
1	B	317	GLY	3.1
1	A	156	VAL	2.9
1	A	237	MET	2.9
1	A	408	ALA	2.8
1	A	153	MET	2.8
1	B	307	THR	2.7
1	A	246	HIS	2.5
1	A	151	LEU	2.4
1	A	310	VAL	2.4
1	A	236	GLY	2.4
1	A	230	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	320	THR	2.3
1	B	337	SER	2.3
1	B	340	TRP	2.3
1	A	152	THR	2.3
1	A	183	ALA	2.2
1	B	318	TYR	2.2
1	B	407	LYS	2.2
1	B	301	ILE	2.1
1	A	187	ALA	2.1
1	A	157	MET	2.1
1	B	312	VAL	2.1
1	A	178	ILE	2.0
1	A	309	GLY	2.0
1	A	314	SER	2.0
1	B	236	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](#)

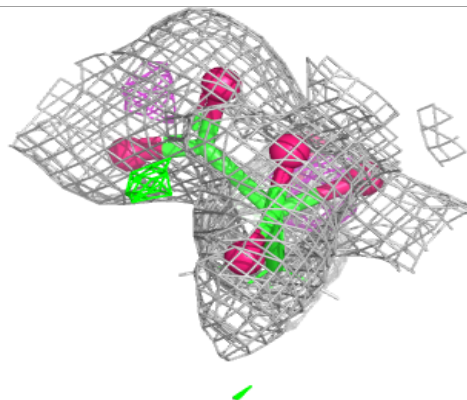
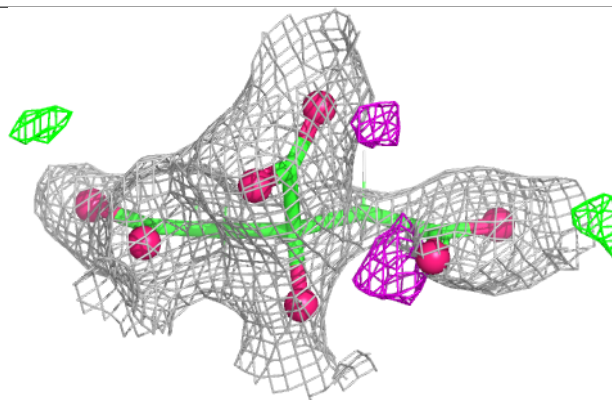
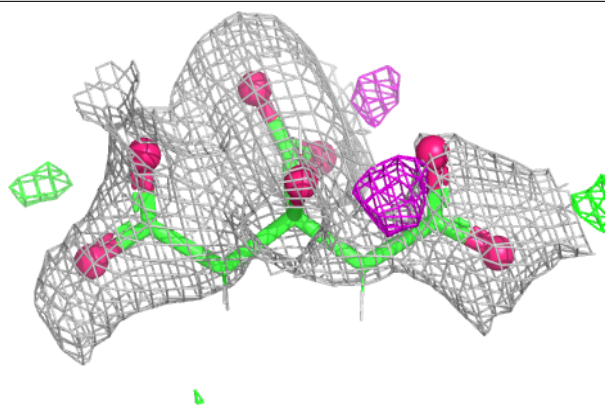
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	CIT	B	501	13/13	0.82	0.13	32,40,54,54	0
2	CIT	A	501	13/13	0.92	0.08	19,25,33,35	0

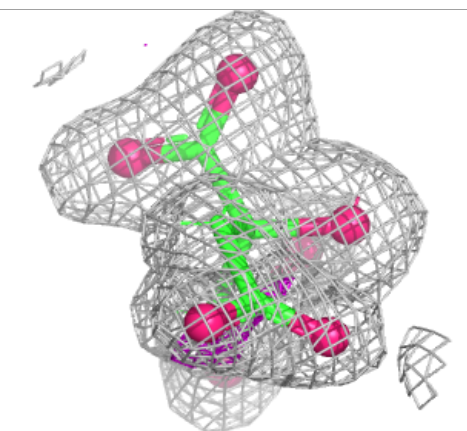
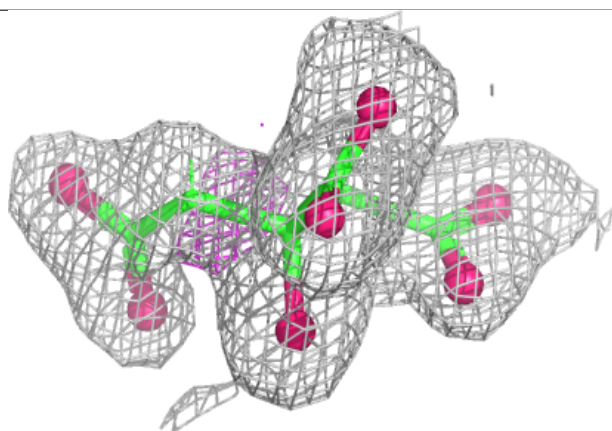
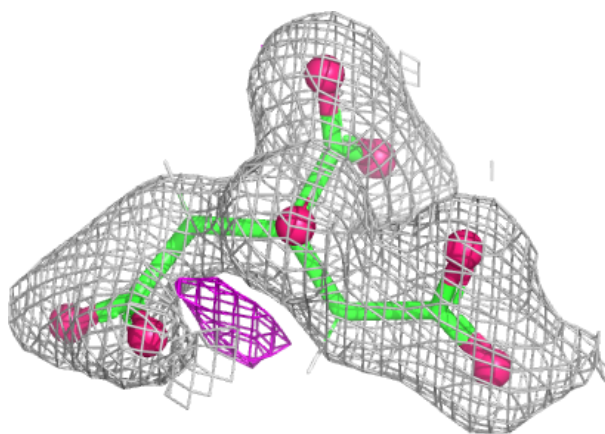
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around CIT B 501:

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

**Electron density around CIT A 501:**

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



6.5 Other polymers ⓘ

There are no such residues in this entry.