WORLDWIDE PDB PROTEIN DATA BANK

wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 5, 2024 – 01:01 am GMT

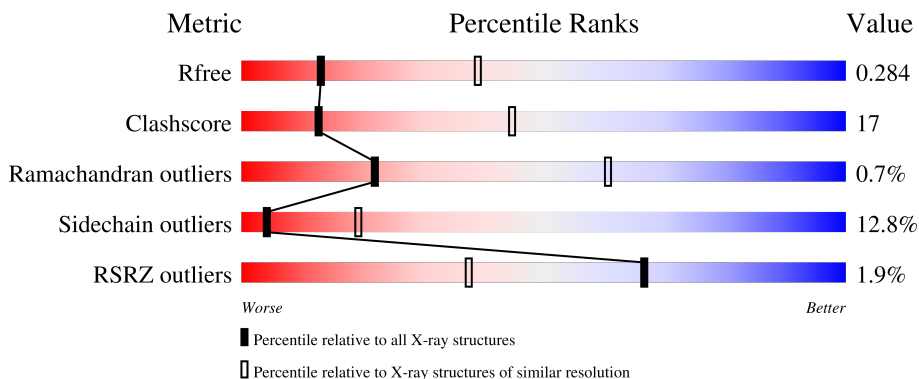|		
PDB ID	:	5A0Z
Title	:	STRUCTURE OF CUTC CHOLINE LYASE CHOLINE FREE FORM FROM KLEBSIELLA PNEUMONIAE
Authors	:	Kalnins, G.; Tars, K.
Deposited on	:	2015-04-27
Resolution	:	3.00 Å(reported)

1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	Percentile Ranks	Value
Rfree		0.284
Clashscore		17
Ramachandran outliers		0.7%
Sidechain outliers		12.8%
RSRZ outliers		1.9%

Worse   *Better*

■ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R$_{free}$	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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 <=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	792	% 64% 25% 5% • 5%
1	B	792	65% 25% • 6%
1	C	792	3% 66% 24% • 6%
1	D	792	4% 62% 22% 5% 10%

2   Entry composition ⓘ

There is only 1 type of molecule in this entry. The entry contains 23177 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 CHOLINE TRIMETHYLAMINE LYASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	750	Total	C	N	O	S	0	0	0
			5919	3747	1020	1114	38			
1	B	746	Total	C	N	O	S	0	0	0
			5868	3715	1011	1103	39			
1	C	746	Total	C	N	O	S	0	0	0
			5856	3707	1005	1107	37			
1	D	710	Total	C	N	O	S	0	0	0
			5534	3507	947	1046	34			

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.

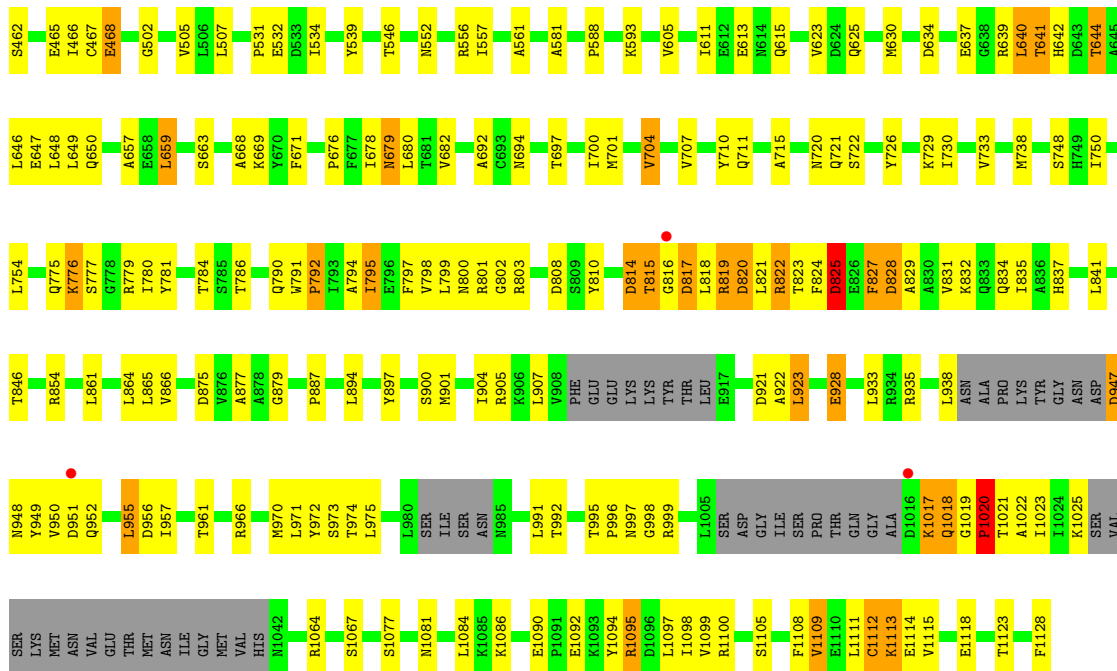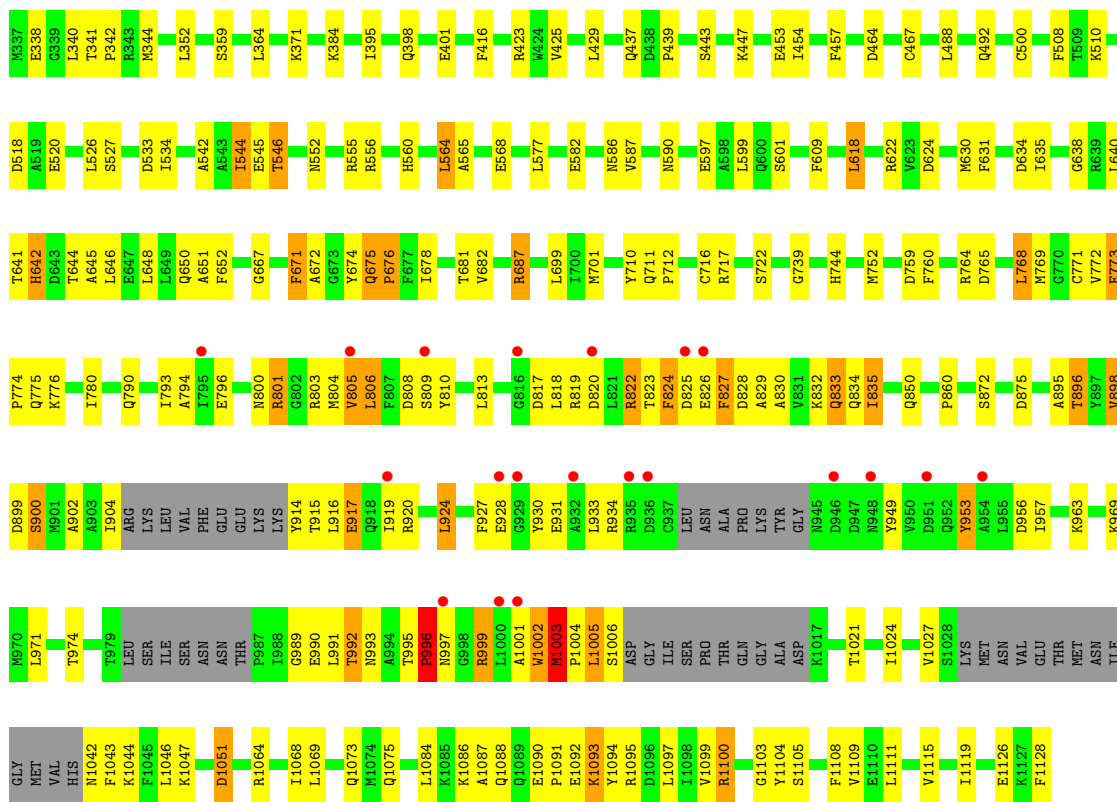• Molecule 1: CHOLINE TRIMETHYLAMINE LYASE
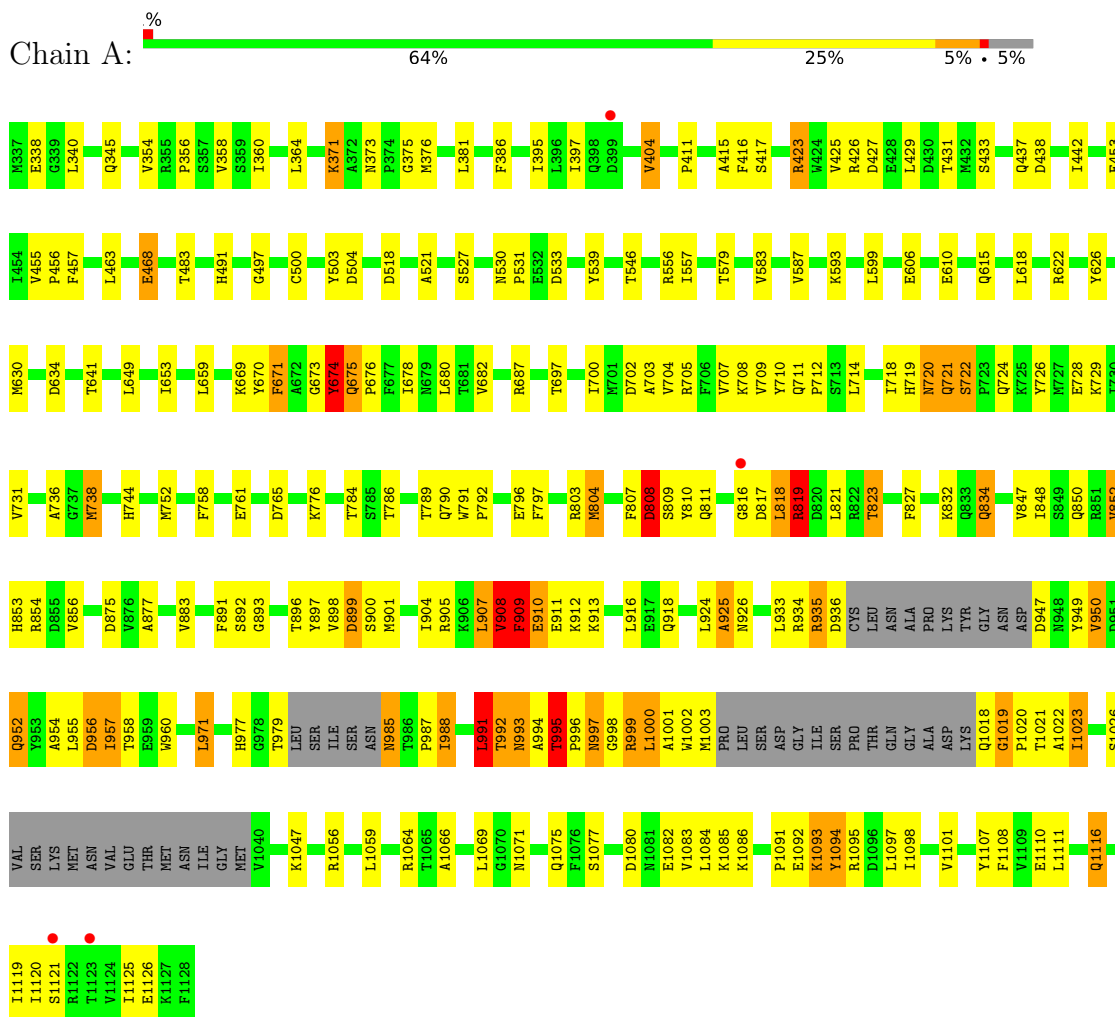
• Molecule 1: CHOLINE TRIMETHYLAMINE LYASE

• Molecule 1: CHOLINE TRIMETHYLAMINE LYASE

Chain C:

3%

66%   24%   •   6%

• Molecule 1: CHOLINE TRIMETHYLAMINE LYASE

Chain D:

4%

62%   22%   5%   10%

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	103.60Å 154.55Å 120.78Å 90.00° 96.85° 90.00°	Depositor
Resolution (Å)	119.92 – 3.00 56.60 – 3.00	Depositor EDS
% Data completeness (in resolution range)	88.5 (119.92-3.00) 88.5 (56.60-3.00)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.05 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.8.0107	Depositor
R, R_{free}	0.232 , 0.286 0.234 , 0.284	Depositor DCC
R_{free} test set	3342 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	38.3	Xtriage
Anisotropy	0.245	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 51.2	EDS
L-test for twinning[2]	$< |L| > = 0.43, < L^2 > = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.84	EDS
Total number of atoms	23177	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.09% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ 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)

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	2/6041 (0.0%)	0.87	13/8170 (0.2%)
1	B	0.73	1/5987 (0.0%)	0.85	5/8099 (0.1%)
1	C	0.73	2/5977 (0.0%)	0.84	10/8089 (0.1%)
1	D	0.63	3/5649 (0.1%)	0.81	6/7652 (0.1%)
All	All	0.72	8/23654 (0.0%)	0.84	34/32010 (0.1%)

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	5
1	B	0	4
1	C	0	4
1	D	0	2
All	All	0	15

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	726	TYR	CE1-CZ	-5.62	1.31	1.38
1	B	468	GLU	CD-OE2	-5.40	1.19	1.25
1	D	774	PRO	N-CD	5.26	1.55	1.47
1	C	342	PRO	N-CD	5.22	1.55	1.47
1	A	674	TYR	CE1-CZ	-5.09	1.31	1.38

The worst 5 of 34 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	996	PRO	CA-N-CD	-7.46	101.05	111.50
1	C	675	GLN	C-N-CD	6.72	142.51	128.40
1	B	1019	GLY	C-N-CD	6.40	141.85	128.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	675	GLN	C-N-CD	6.35	141.74	128.40
1	A	1019	GLY	C-N-CD	6.31	141.65	128.40

There are no chirality outliers.

5 of 15 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	891	PHE	Peptide
1	A	909	PHE	Peptide
1	A	925	ALA	Peptide
1	A	950	VAL	Peptide
1	A	992	THR	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	5919	0	5803	173	0
1	B	5868	0	5753	166	0
1	C	5856	0	5720	214	0
1	D	5534	0	5388	211	0
All	All	23177	0	22664	762	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 17.

The worst 5 of 762 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:C:989:GLY:CA	1:C:1005:LEU:CD2	1.78	1.60
1:D:1099:VAL:HG11	1:D:1119:ILE:CG2	1.28	1.58
1:D:949:TYR:O	1:D:953:TYR:CE1	1.69	1.43
1:C:989:GLY:CA	1:C:1005:LEU:HD23	1.43	1.36
1:C:989:GLY:HA3	1:C:1005:LEU:CD2	1.44	1.36

There are no symmetry-related clashes.

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5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	740/792 (93%)	656 (89%)	76 (10%)	8 (1%)	14	50
1	B	734/792 (93%)	664 (90%)	66 (9%)	4 (0%)	29	68
1	C	734/792 (93%)	661 (90%)	69 (9%)	4 (0%)	29	68
1	D	700/792 (88%)	635 (91%)	60 (9%)	5 (1%)	22	60
All	All	2908/3168 (92%)	2616 (90%)	271 (9%)	21 (1%)	22	60

5 of 21 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	825	ASP
1	C	996	PRO
1	C	1003	MET
1	D	825	ASP
1	A	810	TYR

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	624/668 (93%)	544 (87%)	80 (13%)	4	19
1	B	618/668 (92%)	544 (88%)	74 (12%)	5	22
1	C	617/668 (92%)	542 (88%)	75 (12%)	5	21
1	D	580/668 (87%)	496 (86%)	84 (14%)	3	15
All	All	2439/2672 (91%)	2126 (87%)	313 (13%)	4	19

5 of 313 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	370	VAL
1	D	955	LEU
1	D	448	LYS
1	D	708	LYS
1	D	1075	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 59 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	345	GLN
1	D	833	GLN
1	C	642	HIS
1	D	790	GLN
1	D	569	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 monosaccharides in this entry.

5.6   Ligand geometry ⓘ

There are no ligands in this entry.

5.7   Other polymers ⓘ

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	750/792 (94%)	-0.35	4 (0%)	91	75	6, 27, 56, 87	0
1	B	746/792 (94%)	-0.34	3 (0%)	92	79	3, 26, 65, 91	0
1	C	746/792 (94%)	-0.17	20 (2%)	54	26	14, 34, 82, 146	0
1	D	710/792 (89%)	0.11	30 (4%)	36	14	17, 51, 91, 115	0
All	All	2952/3168 (93%)	-0.19	57 (1%)	66	37	3, 33, 78, 146	0

The worst 5 of 57 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	827	PHE	5.0
1	C	1001	ALA	4.5
1	C	816	GLY	4.4
1	D	992	THR	4.3
1	D	822	ARG	4.3

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

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

6.3   Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4   Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.