



wwPDB X-ray Structure Validation Summary Report

Feb 27, 2014 – 07:06 AM GMT

PDB ID : 2E9F
Title : Crystal Structure of T.th.HB8 Argininosuccinate lyase complexed with L-Arginine
Authors : Goto, M.
Deposited on : 2007-01-25
Resolution : 2.80 Å(reported)

This is a wwPDB validation summary report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

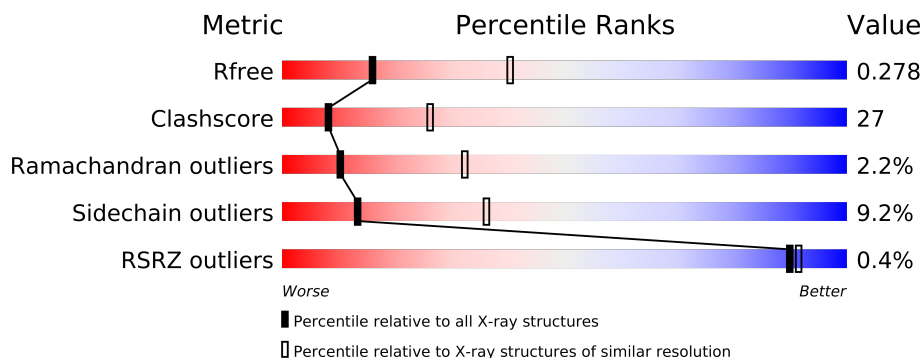
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.80 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}	66092	1799 (2.80-2.80)
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (2.80-2.80)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	462	<div><div></div><div><div></div><div></div><div></div><div></div><div></div></div></div>
1	B	462	<div><div></div><div><div></div><div></div><div></div><div></div><div></div></div></div>
1	C	462	<div><div></div><div><div></div><div></div><div></div><div></div><div></div></div></div>
1	D	462	<div><div></div><div><div></div><div></div><div></div><div></div><div></div></div></div>

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	ARG	B	491	-	X
2	ARG	B	493	-	X

2 Entry composition i

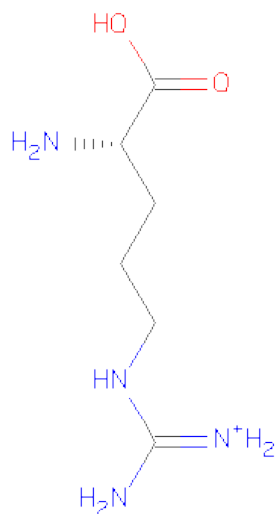
There are 3 unique types of molecules in this entry. The entry contains 14218 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Argininosuccinate lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	449	Total	C	N	O	S	0	0	0
			3530	2243	636	644	7			
1	B	450	Total	C	N	O	S	0	0	0
			3528	2241	637	643	7			
1	C	448	Total	C	N	O	S	0	0	0
			3498	2227	626	638	7			
1	D	449	Total	C	N	O	S	0	0	0
			3479	2216	625	631	7			

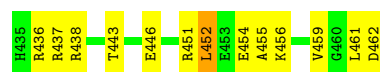
- Molecule 2 is ARGinine (three-letter code: ARG) (formula: C₆H₁₅N₄O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	4	Total	C	N	O	0	0
			48	24	16	8		

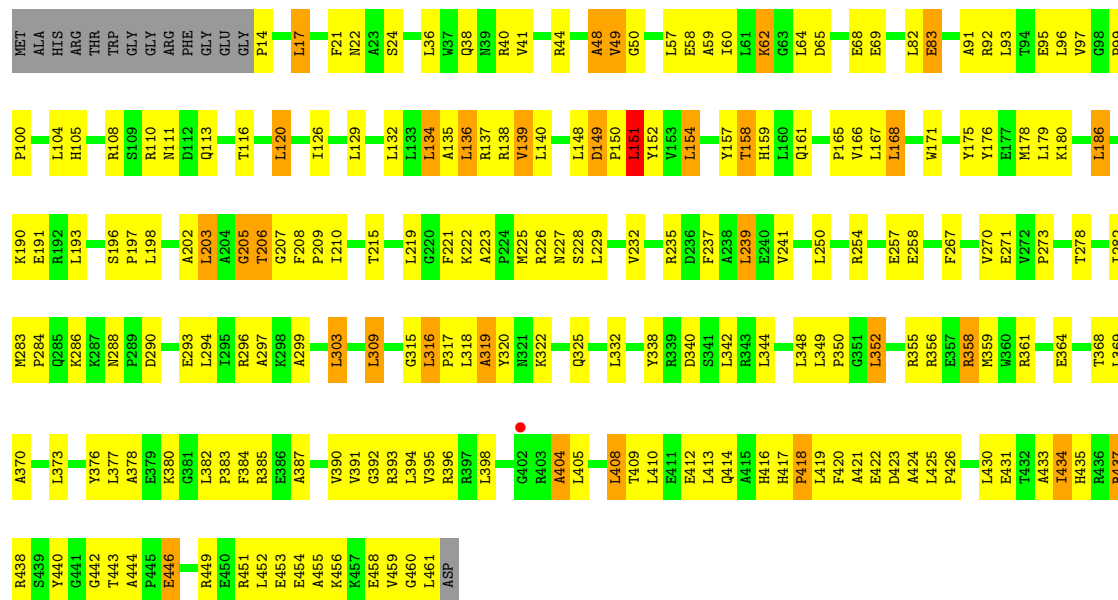
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	34	Total 34	O 34	0	0
3	B	41	Total 41	O 41	0	0
3	C	30	Total 30	O 30	0	0
3	D	30	Total 30	O 30	0	0



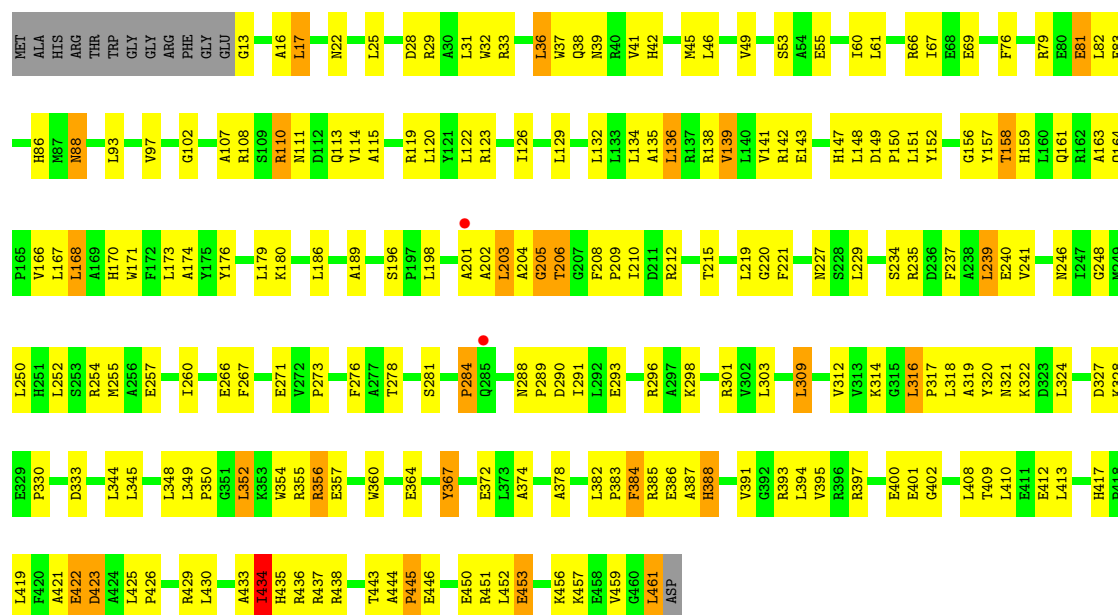
• Molecule 1: Argininosuccinate lyase

Chain C:



• Molecule 1: Argininosuccinate lyase

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.34Å 119.79Å 257.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.29 – 2.80 34.97 – 2.80	Depositor EDS
% Data completeness (in resolution range)	85.6 (34.29-2.80) 85.7 (34.97-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.95 (at 2.81Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.205 , 0.271 0.216 , 0.278	Depositor DCC
R_{free} test set	5234 reflections (10.08%)	DCC
Wilson B-factor (Å ²)	50.7	Xtriage
Anisotropy	0.404	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 26.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 51944 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	14218	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.08% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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	0.40	0/3600	0.65	0/4877
1	B	0.40	0/3597	0.65	2/4873 (0.0%)
1	C	0.38	0/3568	0.64	0/4839
1	D	0.38	0/3548	0.64	3/4814 (0.1%)
All	All	0.39	0/14313	0.64	5/19403 (0.0%)

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

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

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	284	PRO	N-CA-CB	5.68	110.11	103.30
1	D	156	GLY	N-CA-C	-5.52	99.29	113.10
1	B	104	LEU	CA-CB-CG	5.46	127.85	115.30
1	D	348	LEU	CA-CB-CG	5.35	127.61	115.30
1	B	284	PRO	N-CA-CB	5.25	109.60	103.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	320	TYR	Sidechain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3530	0	3538	217	0
1	B	3528	0	3530	214	0
1	C	3498	0	3495	196	0
1	D	3479	0	3456	197	0
2	B	48	0	48	22	0
3	A	34	0	0	1	0
3	B	41	0	0	2	0
3	C	30	0	0	1	0
3	D	30	0	0	3	0
All	All	14218	0	14067	747	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 27.

The worst 5 of 747 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:59:ALA:HB1	1:B:97:VAL:HG21	1.32	1.11
1:A:456:LYS:HB3	1:A:461:LEU:HB3	1.28	1.05
1:B:167:LEU:HD13	1:B:443:THR:HG22	1.42	1.02
1:A:167:LEU:H	1:A:443:THR:HG21	1.26	0.99
1:C:99:PRO:HG2	1:C:100:PRO:HD3	1.43	0.98

There are no symmetry-related clashes.

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	447/462 (97%)	401 (90%)	34 (8%)	12 (3%)	8	25
1	B	448/462 (97%)	410 (92%)	30 (7%)	8 (2%)	13	39
1	C	446/462 (96%)	400 (90%)	34 (8%)	12 (3%)	8	25
1	D	447/462 (97%)	395 (88%)	44 (10%)	8 (2%)	13	39
All	All	1788/1848 (97%)	1606 (90%)	142 (8%)	40 (2%)	10	32

5 of 40 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	282	ILE
1	A	319	ALA
1	B	149	ASP
1	B	205	GLY
1	C	203	LEU

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	353/369 (96%)	323 (92%)	30 (8%)	15	41
1	B	351/369 (95%)	318 (91%)	33 (9%)	13	34
1	C	348/369 (94%)	314 (90%)	34 (10%)	12	32
1	D	340/369 (92%)	309 (91%)	31 (9%)	14	37
All	All	1392/1476 (94%)	1264 (91%)	128 (9%)	13	36

5 of 128 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	389	HIS
1	C	139	VAL
1	D	357	GLU
1	B	422	GLU
1	C	104	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 33 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	325	GLN
1	C	42	HIS
1	D	170	HIS
1	B	417	HIS
1	C	22	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

4 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	ARG	B	491	-	11,11,11	0.54	0	13,13,13	0.80	0
2	ARG	B	492	-	11,11,11	0.47	0	13,13,13	1.00	0
2	ARG	B	493	-	11,11,11	0.53	0	13,13,13	1.11	1 (7%)
2	ARG	B	494	-	11,11,11	0.51	0	13,13,13	0.66	0

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	ARG	B	491	-	-	0/11/11/11	0/0/0/0
2	ARG	B	492	-	-	0/11/11/11	0/0/0/0
2	ARG	B	493	-	-	0/11/11/11	0/0/0/0
2	ARG	B	494	-	-	0/11/11/11	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	B	493	ARG	C-CA-N	2.74	113.90	109.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	449/462 (97%)	-0.56	2 (0%) 90 91	25, 42, 58, 81	0
1	B	450/462 (97%)	-0.57	3 (0%) 84 85	22, 38, 57, 79	0
1	C	448/462 (96%)	-0.55	1 (0%) 93 95	25, 42, 64, 72	0
1	D	449/462 (97%)	-0.45	2 (0%) 90 91	28, 45, 61, 77	0
All	All	1796/1848 (97%)	-0.53	8 (0%) 90 91	22, 42, 62, 81	0

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	279	GLY	3.0
1	A	462	ASP	2.7
1	D	285	GLN	2.6
1	B	280	SER	2.5
1	B	285	GLN	2.4

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 carbohydrates in this entry.

6.4 Ligands

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	ARG	B	491	12/12	0.21	2.84	33,35,38,40	0
2	ARG	B	493	12/12	0.19	2.03	33,43,50,51	0
2	ARG	B	492	12/12	0.21	1.70	45,49,50,51	0
2	ARG	B	494	12/12	0.17	1.42	42,46,48,49	0

6.5 Other polymers ⓘ

There are no such residues in this entry.