



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 07:56 PM GMT

PDB ID : 1HWZ  
Title : BOVINE GLUTAMATE DEHYDROGENASE COMPLEXED WITH  
NADPH, GLUTAMATE, AND GTP  
Authors : Smith, T.J.; Peterson, P.E.; Schmidt, T.; Fang, J.; Stanley, C.A.  
Deposited on : 2001-01-10  
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

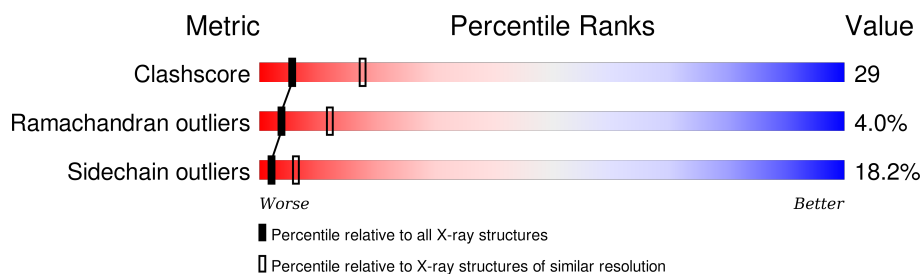
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 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))
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	501	
1	B	501	
1	C	501	
1	D	501	
1	E	501	
1	F	501	

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	GLU	A	550	-	-	X	-
2	GLU	B	554	-	-	X	-
2	GLU	C	555	-	-	X	-
2	GLU	D	556	-	-	X	-
2	GLU	E	557	-	-	X	-
2	GLU	F	558	-	-	X	-
3	NDP	A	551	-	-	X	-
3	NDP	B	559	-	-	X	-
3	NDP	C	560	-	-	X	-
3	NDP	D	561	-	-	X	-
3	NDP	E	562	-	-	X	-
3	NDP	F	563	-	-	X	-

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 24000 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 GLUTAMATE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	501	Total	C	N	O	S	0	0	0
			3910	2468	690	733	19			
1	B	501	Total	C	N	O	S	0	0	0
			3910	2468	690	733	19			
1	C	501	Total	C	N	O	S	0	0	0
			3910	2468	690	733	19			
1	D	501	Total	C	N	O	S	0	0	0
			3910	2468	690	733	19			
1	E	501	Total	C	N	O	S	0	0	0
			3910	2468	690	733	19			
1	F	501	Total	C	N	O	S	0	0	0
			3910	2468	690	733	19			

There are 30 discrepancies between the modelled and reference sequences:

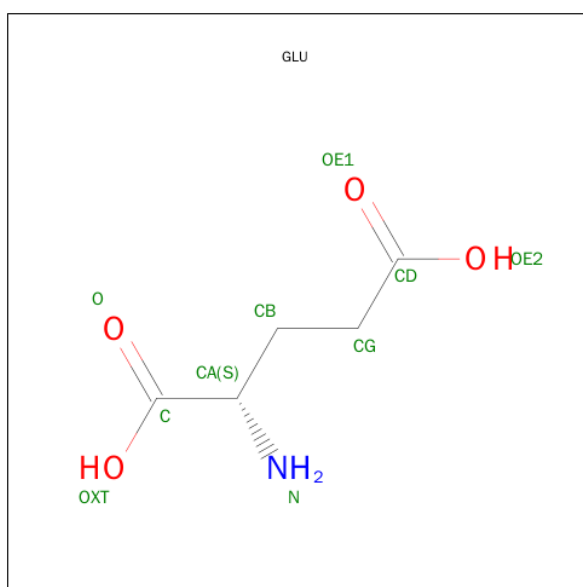
Chain	Residue	Modelled	Actual	Comment	Reference
A	200	GLY	LYS	SEE REMARK 999	UNP P00366
A	201	LYS	PRO	SEE REMARK 999	UNP P00366
A	202	PRO	GLY	SEE REMARK 999	UNP P00366
A	221	HIS	GLY	SEE REMARK 999	UNP P00366
A	222	GLY	HIS	SEE REMARK 999	UNP P00366
B	200	GLY	LYS	SEE REMARK 999	UNP P00366
B	201	LYS	PRO	SEE REMARK 999	UNP P00366
B	202	PRO	GLY	SEE REMARK 999	UNP P00366
B	221	HIS	GLY	SEE REMARK 999	UNP P00366
B	222	GLY	HIS	SEE REMARK 999	UNP P00366
C	200	GLY	LYS	SEE REMARK 999	UNP P00366
C	201	LYS	PRO	SEE REMARK 999	UNP P00366
C	202	PRO	GLY	SEE REMARK 999	UNP P00366
C	221	HIS	GLY	SEE REMARK 999	UNP P00366
C	222	GLY	HIS	SEE REMARK 999	UNP P00366
D	200	GLY	LYS	SEE REMARK 999	UNP P00366
D	201	LYS	PRO	SEE REMARK 999	UNP P00366

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Chain	Residue	Modelled	Actual	Comment	Reference
D	202	PRO	GLY	SEE REMARK 999	UNP P00366
D	221	HIS	GLY	SEE REMARK 999	UNP P00366
D	222	GLY	HIS	SEE REMARK 999	UNP P00366
E	200	GLY	LYS	SEE REMARK 999	UNP P00366
E	201	LYS	PRO	SEE REMARK 999	UNP P00366
E	202	PRO	GLY	SEE REMARK 999	UNP P00366
E	221	HIS	GLY	SEE REMARK 999	UNP P00366
E	222	GLY	HIS	SEE REMARK 999	UNP P00366
F	200	GLY	LYS	SEE REMARK 999	UNP P00366
F	201	LYS	PRO	SEE REMARK 999	UNP P00366
F	202	PRO	GLY	SEE REMARK 999	UNP P00366
F	221	HIS	GLY	SEE REMARK 999	UNP P00366
F	222	GLY	HIS	SEE REMARK 999	UNP P00366

- Molecule 2 is GLUTAMIC ACID (three-letter code: GLU) (formula: C<sub>5</sub>H<sub>9</sub>NO<sub>4</sub>).



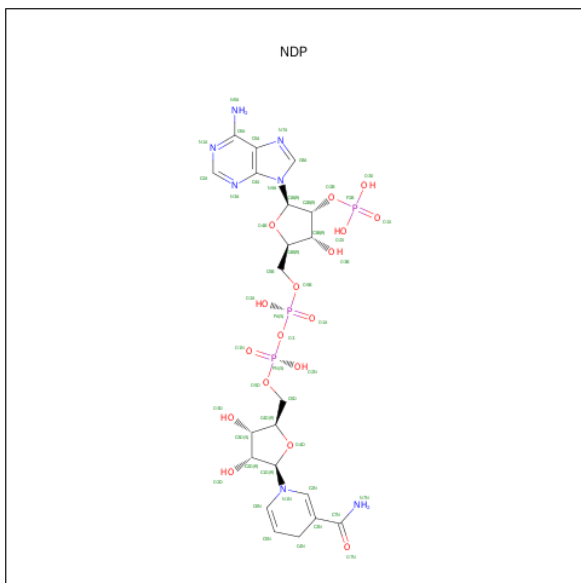
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			10	5	1	4		
2	B	1	Total	C	N	O	0	0
			10	5	1	4		
2	C	1	Total	C	N	O	0	0
			10	5	1	4		
2	D	1	Total	C	N	O	0	0
			10	5	1	4		
2	E	1	Total	C	N	O	0	0
			10	5	1	4		

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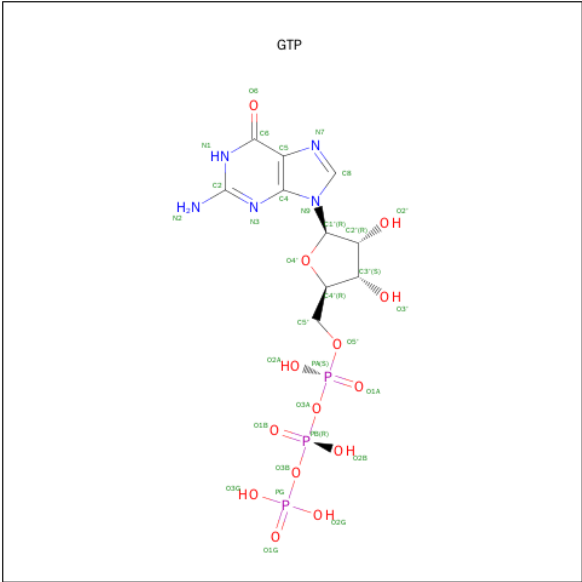
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	F	1	Total	C	N	O	0	0
			10	5	1	4		

- Molecule 3 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula:  $C_{21}H_{30}N_7O_{17}P_3$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	D	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	E	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	F	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 4 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula:  $C_{10}H_{16}N_5O_{14}P_3$ ).

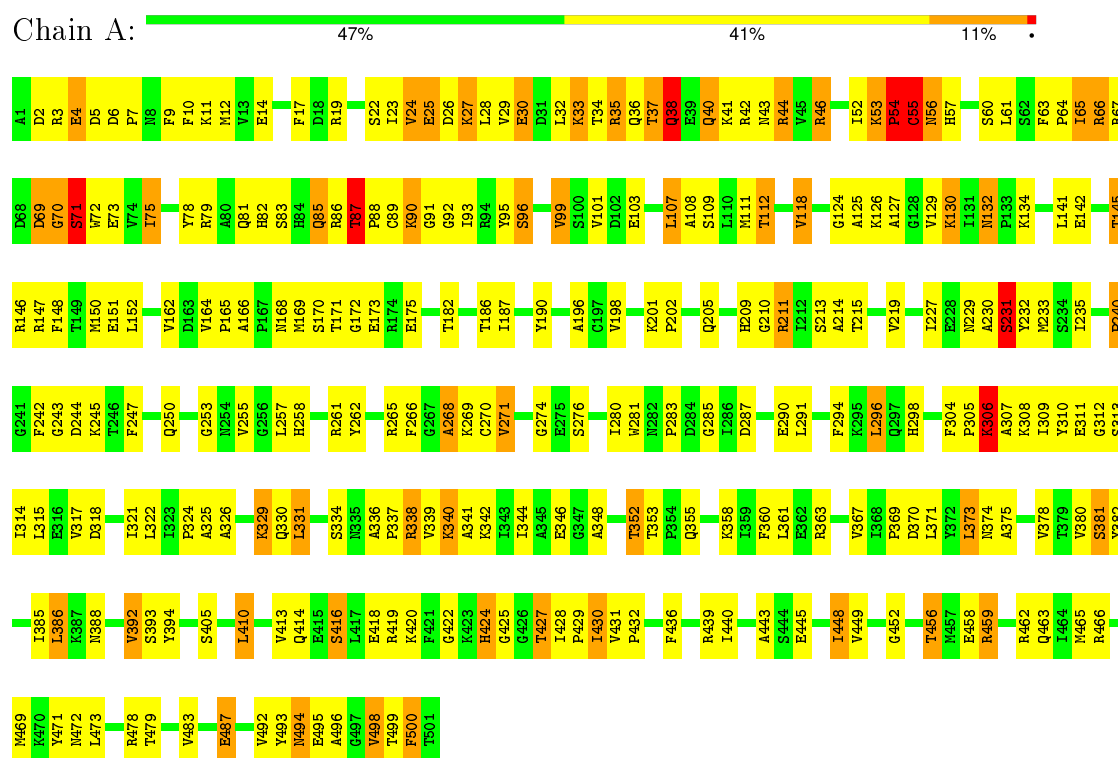


### 3 Residue-property plots

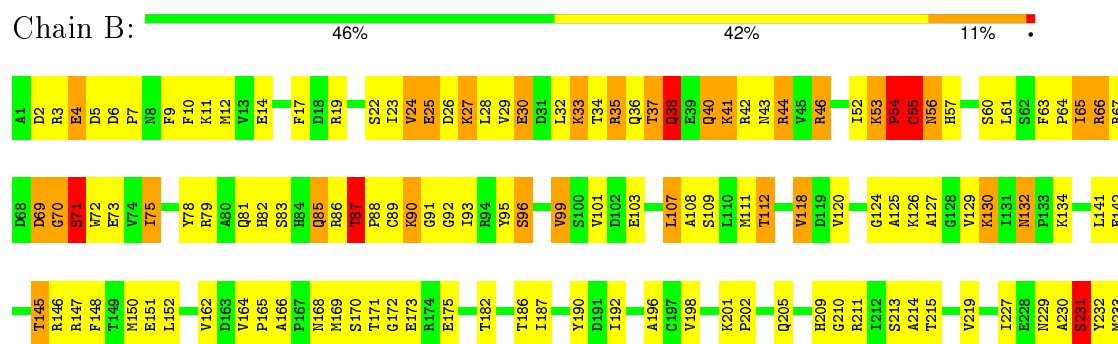
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

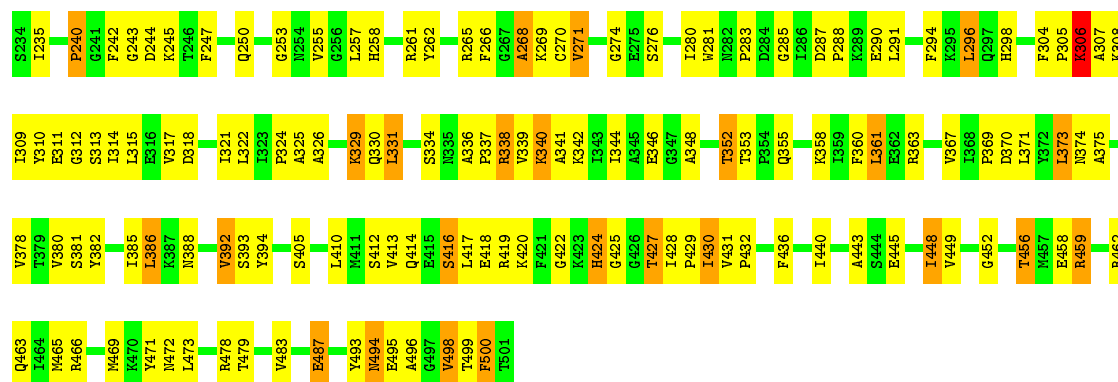
#### • Molecule 1: GLUTAMATE DEHYDROGENASE



#### • Molecule 1: GLUTAMATE DEHYDROGENASE

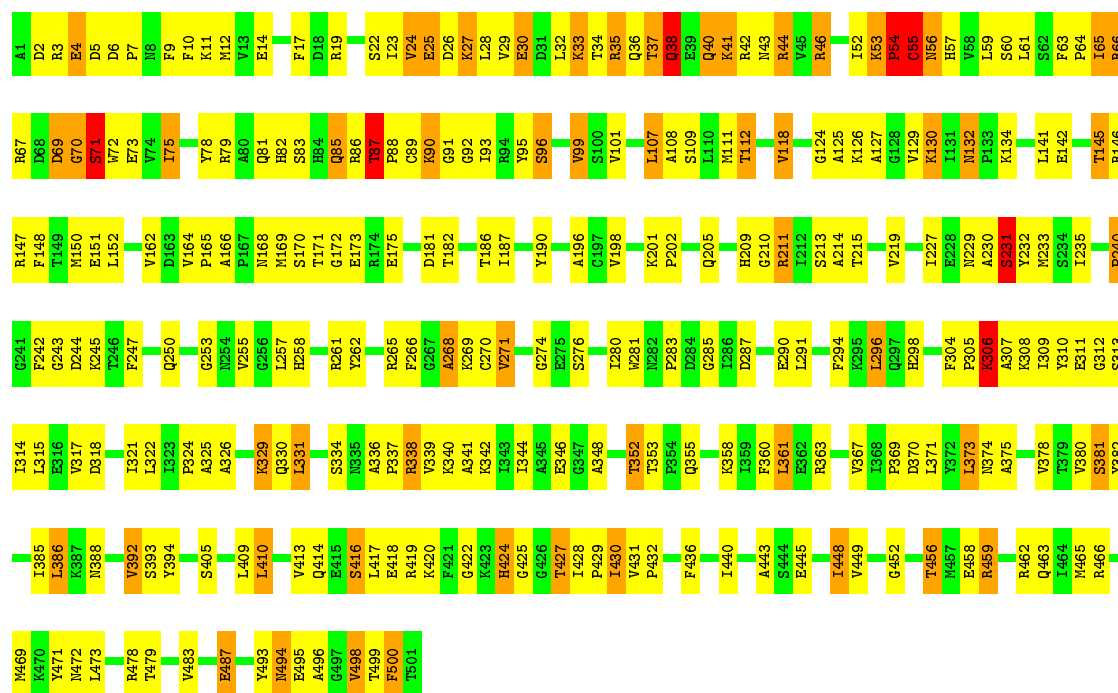






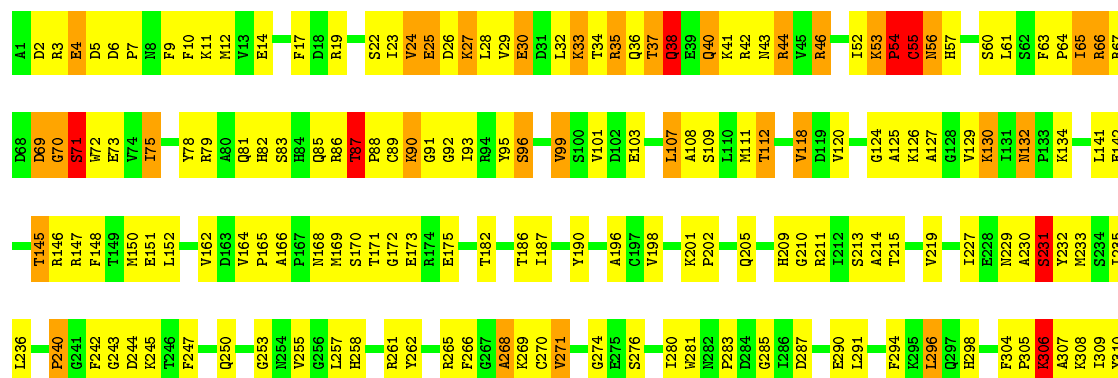
### • Molecule 1: GLUTAMATE DEHYDROGENASE

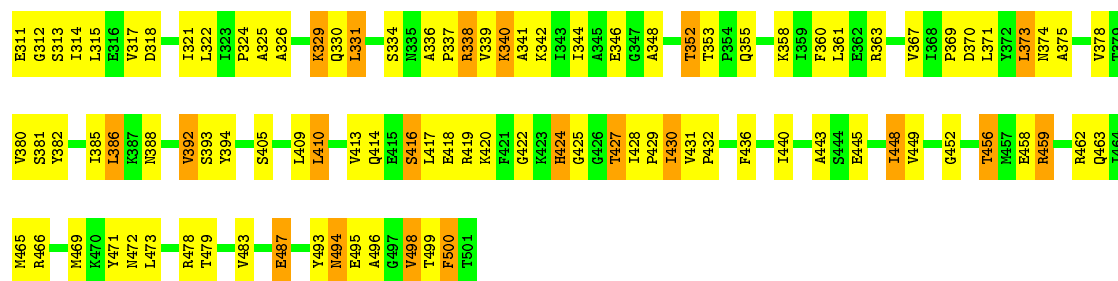
Chain C: 47% 41% 11%



### • Molecule 1: GLUTAMATE DEHYDROGENASE

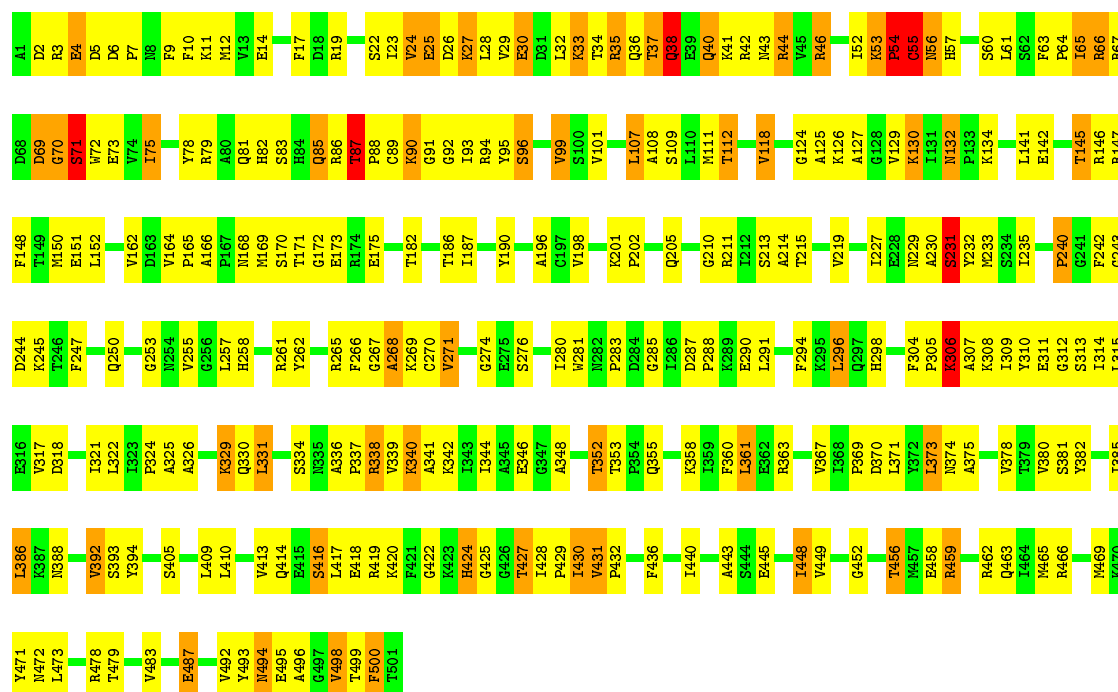
Chain D: 47% 42% 10%





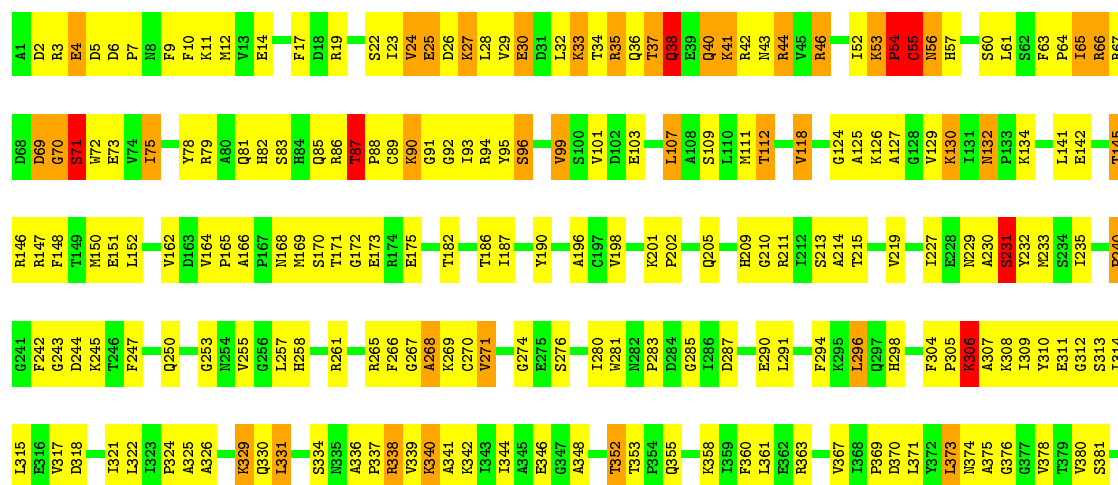
• Molecule 1: GLUTAMATE DEHYDROGENASE

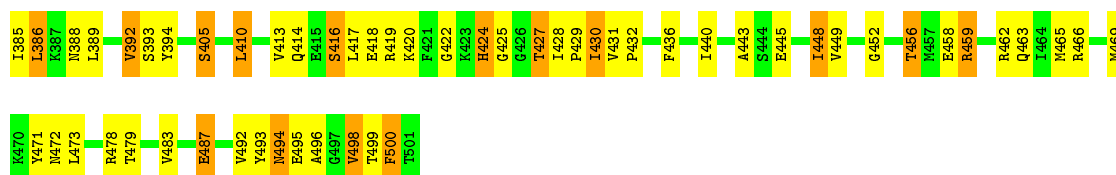
Chain E: 47% 42% 11%



• Molecule 1: GLUTAMATE DEHYDROGENASE

Chain F: 47% 41% 11%





## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	124.30Å 102.50Å 169.20Å 90.00° 102.20° 90.00°	Depositor
Resolution (Å)	8.00 – 2.80	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.80)	Depositor
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.250 , 0.300	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	24000	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

## 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: GTP, NDP

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	0.58	0/3991	0.84	8/5384 (0.1%)
1	B	0.58	0/3991	0.84	8/5384 (0.1%)
1	C	0.58	0/3991	0.84	8/5384 (0.1%)
1	D	0.58	0/3991	0.84	8/5384 (0.1%)
1	E	0.58	0/3991	0.84	8/5384 (0.1%)
1	F	0.58	0/3991	0.84	8/5384 (0.1%)
All	All	0.58	0/23946	0.84	48/32304 (0.1%)

There are no bond length outliers.

The worst 5 of 48 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	33	LYS	N-CA-C	8.29	133.39	111.00
1	F	33	LYS	N-CA-C	8.29	133.38	111.00
1	E	33	LYS	N-CA-C	8.28	133.37	111.00
1	D	33	LYS	N-CA-C	8.28	133.36	111.00
1	C	33	LYS	N-CA-C	8.28	133.35	111.00

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	3910	0	3889	233	0
1	B	3910	0	3889	239	0
1	C	3910	0	3889	239	0
1	D	3910	0	3889	240	0
1	E	3910	0	3889	245	0
1	F	3910	0	3889	241	0
2	A	10	0	5	7	0
2	B	10	0	5	7	0
2	C	10	0	5	7	0
2	D	10	0	5	7	0
2	E	10	0	5	7	0
2	F	10	0	5	7	0
3	A	48	0	26	24	0
3	B	48	0	26	22	0
3	C	48	0	26	24	0
3	D	48	0	26	23	0
3	E	48	0	26	25	0
3	F	48	0	26	23	0
4	A	32	0	12	2	0
4	B	32	0	12	2	0
4	C	32	0	12	2	0
4	D	32	0	12	2	0
4	E	32	0	12	2	0
4	F	32	0	12	1	0
All	All	24000	0	23592	1401	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 29.

The worst 5 of 1401 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:215:THR:HG21	3:A:551:NDP:H42N	1.15	1.15
1:B:374:ASN:HB2	3:B:559:NDP:H41N	1.16	1.14
1:A:374:ASN:HB2	3:A:551:NDP:H41N	1.16	1.13
1:F:215:THR:HG21	3:F:563:NDP:H42N	1.15	1.12
1:C:374:ASN:HB2	3:C:560:NDP:H41N	1.16	1.12

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	499/501 (100%)	437 (88%)	42 (8%)	20 (4%)	4	12
1	B	499/501 (100%)	436 (87%)	43 (9%)	20 (4%)	4	12
1	C	499/501 (100%)	437 (88%)	42 (8%)	20 (4%)	4	12
1	D	499/501 (100%)	437 (88%)	42 (8%)	20 (4%)	4	12
1	E	499/501 (100%)	437 (88%)	42 (8%)	20 (4%)	4	12
1	F	499/501 (100%)	437 (88%)	42 (8%)	20 (4%)	4	12
All	All	2994/3006 (100%)	2621 (88%)	253 (8%)	120 (4%)	4	12

5 of 120 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	5	ASP
1	A	35	ARG
1	A	38	GLN
1	A	54	PRO
1	A	55	CYS

### 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	417/417 (100%)	341 (82%)	76 (18%)	2	6
1	B	417/417 (100%)	341 (82%)	76 (18%)	2	6
1	C	417/417 (100%)	341 (82%)	76 (18%)	2	6

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	417/417 (100%)	341 (82%)	76 (18%)	2	6
1	E	417/417 (100%)	341 (82%)	76 (18%)	2	6
1	F	417/417 (100%)	341 (82%)	76 (18%)	2	6
All	All	2502/2502 (100%)	2046 (82%)	456 (18%)	2	6

5 of 456 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	385	ILE
1	D	107	LEU
1	F	313	SER
1	C	416	SER
1	D	9	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 93 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	437	GLN
1	D	209	HIS
1	F	258	HIS
1	C	450	HIS
1	D	82	HIS

### 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 carbohydrates in this entry.



## 5.6 Ligand geometry

18 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	GLU	A	550	-	3,9,9	0.68	0	2,11,11	0.42	0
3	NDP	A	551	-	42,52,52	1.98	9 (21%)	55,80,80	2.25	19 (34%)
4	GTP	A	553	-	25,34,34	1.34	2 (8%)	34,54,54	2.34	10 (29%)
2	GLU	B	554	-	3,9,9	0.68	0	2,11,11	0.42	0
3	NDP	B	559	-	42,52,52	1.98	9 (21%)	55,80,80	2.25	19 (34%)
4	GTP	B	564	-	25,34,34	1.34	2 (8%)	34,54,54	2.34	10 (29%)
2	GLU	C	555	-	3,9,9	0.68	0	2,11,11	0.42	0
3	NDP	C	560	-	42,52,52	1.98	9 (21%)	55,80,80	2.25	19 (34%)
4	GTP	C	565	-	25,34,34	1.34	2 (8%)	34,54,54	2.34	10 (29%)
2	GLU	D	556	-	3,9,9	0.67	0	2,11,11	0.42	0
3	NDP	D	561	-	42,52,52	1.98	9 (21%)	55,80,80	2.25	19 (34%)
4	GTP	D	566	-	25,34,34	1.35	3 (12%)	34,54,54	2.34	10 (29%)
2	GLU	E	557	-	3,9,9	0.68	0	2,11,11	0.42	0
3	NDP	E	562	-	42,52,52	1.98	9 (21%)	55,80,80	2.25	19 (34%)
4	GTP	E	567	-	25,34,34	1.35	2 (8%)	34,54,54	2.34	10 (29%)
2	GLU	F	558	-	3,9,9	0.69	0	2,11,11	0.42	0
3	NDP	F	563	-	42,52,52	1.98	9 (21%)	55,80,80	2.25	19 (34%)
4	GTP	F	568	-	25,34,34	1.34	2 (8%)	34,54,54	2.33	10 (29%)

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	GLU	A	550	-	-	0/3/9/9	0/0/0/0
3	NDP	A	551	-	-	0/30/77/77	0/5/5/5
4	GTP	A	553	-	-	0/18/38/38	0/3/3/3
2	GLU	B	554	-	-	0/3/9/9	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NDP	B	559	-	-	0/30/77/77	0/5/5/5
4	GTP	B	564	-	-	0/18/38/38	0/3/3/3
2	GLU	C	555	-	-	0/3/9/9	0/0/0/0
3	NDP	C	560	-	-	0/30/77/77	0/5/5/5
4	GTP	C	565	-	-	0/18/38/38	0/3/3/3
2	GLU	D	556	-	-	0/3/9/9	0/0/0/0
3	NDP	D	561	-	-	0/30/77/77	0/5/5/5
4	GTP	D	566	-	-	0/18/38/38	0/3/3/3
2	GLU	E	557	-	-	0/3/9/9	0/0/0/0
3	NDP	E	562	-	-	0/30/77/77	0/5/5/5
4	GTP	E	567	-	-	0/18/38/38	0/3/3/3
2	GLU	F	558	-	-	0/3/9/9	0/0/0/0
3	NDP	F	563	-	-	0/30/77/77	0/5/5/5
4	GTP	F	568	-	-	0/18/38/38	0/3/3/3

The worst 5 of 67 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	551	NDP	C3B-C2B	-6.14	1.39	1.53
3	D	561	NDP	C3B-C2B	-6.14	1.39	1.53
3	C	560	NDP	C3B-C2B	-6.14	1.39	1.53
3	E	562	NDP	C3B-C2B	-6.13	1.39	1.53
3	B	559	NDP	C3B-C2B	-6.12	1.39	1.53

The worst 5 of 174 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	567	GTP	C5-C6-N1	-7.56	113.26	123.59
4	D	566	GTP	C5-C6-N1	-7.54	113.28	123.59
4	C	565	GTP	C5-C6-N1	-7.54	113.28	123.59
4	A	553	GTP	C5-C6-N1	-7.52	113.30	123.59
4	B	564	GTP	C5-C6-N1	-7.51	113.31	123.59

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

18 monomers are involved in 158 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	550	GLU	7	0
3	A	551	NDP	24	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	553	GTP	2	0
2	B	554	GLU	7	0
3	B	559	NDP	22	0
4	B	564	GTP	2	0
2	C	555	GLU	7	0
3	C	560	NDP	24	0
4	C	565	GTP	2	0
2	D	556	GLU	7	0
3	D	561	NDP	23	0
4	D	566	GTP	2	0
2	E	557	GLU	7	0
3	E	562	NDP	25	0
4	E	567	GTP	2	0
2	F	558	GLU	7	0
3	F	563	NDP	23	0
4	F	568	GTP	1	0

## 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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.