



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 09:12 PM GMT

PDB ID : 1NQT  
Title : Crystal structure of bovine Glutamate dehydrogenase-ADP complex  
Authors : Banerjee, S.; Schmidt, T.; Fang, J.; Stanley, C.A.; Smith, T.J.  
Deposited on : 2003-01-23  
Resolution : 3.50 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



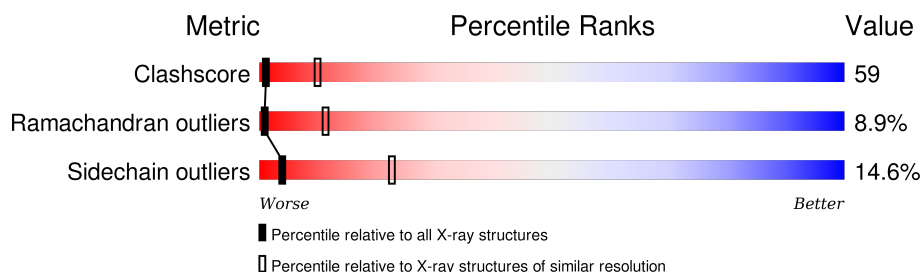
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.50 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	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)

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	496	
1	B	496	
1	C	496	
1	D	496	
1	E	496	
1	F	496	
1	G	496	

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Mol	Chain	Length	Quality of chain
1	H	496	<div><div></div><div>30%55%14%</div><div></div></div>
1	I	496	<div><div></div><div>31%52%16%</div><div></div></div>
1	J	496	<div><div></div><div>34%49%15%</div><div></div></div>
1	K	496	<div><div></div><div>24%60%15%</div><div></div></div>
1	L	496	<div><div></div><div>30%54%15%</div><div></div></div>



## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 46812 atoms, of which 0 are hydrogens and 0 are deuteriums.

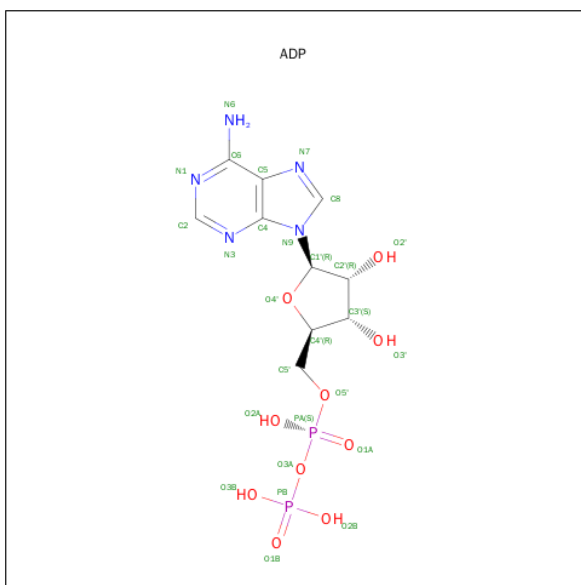
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Glutamate dehydrogenase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	496	Total	C	N	O	S	0	0	0
			3874	2450	679	726	19			
1	B	496	Total	C	N	O	S	0	0	0
			3874	2450	679	726	19			
1	C	496	Total	C	N	O	S	0	0	0
			3874	2450	679	726	19			
1	D	496	Total	C	N	O	S	0	0	0
			3874	2450	679	726	19			
1	E	496	Total	C	N	O	S	0	0	0
			3874	2450	679	726	19			
1	F	496	Total	C	N	O	S	0	0	0
			3874	2450	679	726	19			
1	G	496	Total	C	N	O	S	0	0	0
			3874	2450	679	726	19			
1	H	496	Total	C	N	O	S	0	0	0
			3874	2450	679	726	19			
1	I	496	Total	C	N	O	S	0	0	0
			3874	2450	679	726	19			
1	J	496	Total	C	N	O	S	0	0	0
			3874	2450	679	726	19			
1	K	496	Total	C	N	O	S	0	0	0
			3874	2450	679	726	19			
1	L	496	Total	C	N	O	S	0	0	0
			3874	2450	679	726	19			

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).





Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 27	C 10	N 5	O 10	P 2	0	0
2	B	1	Total 27	C 10	N 5	O 10	P 2	0	0
2	C	1	Total 27	C 10	N 5	O 10	P 2	0	0
2	D	1	Total 27	C 10	N 5	O 10	P 2	0	0
2	E	1	Total 27	C 10	N 5	O 10	P 2	0	0
2	F	1	Total 27	C 10	N 5	O 10	P 2	0	0
2	G	1	Total 27	C 10	N 5	O 10	P 2	0	0
2	H	1	Total 27	C 10	N 5	O 10	P 2	0	0
2	I	1	Total 27	C 10	N 5	O 10	P 2	0	0
2	J	1	Total 27	C 10	N 5	O 10	P 2	0	0
2	K	1	Total 27	C 10	N 5	O 10	P 2	0	0
2	L	1	Total 27	C 10	N 5	O 10	P 2	0	0

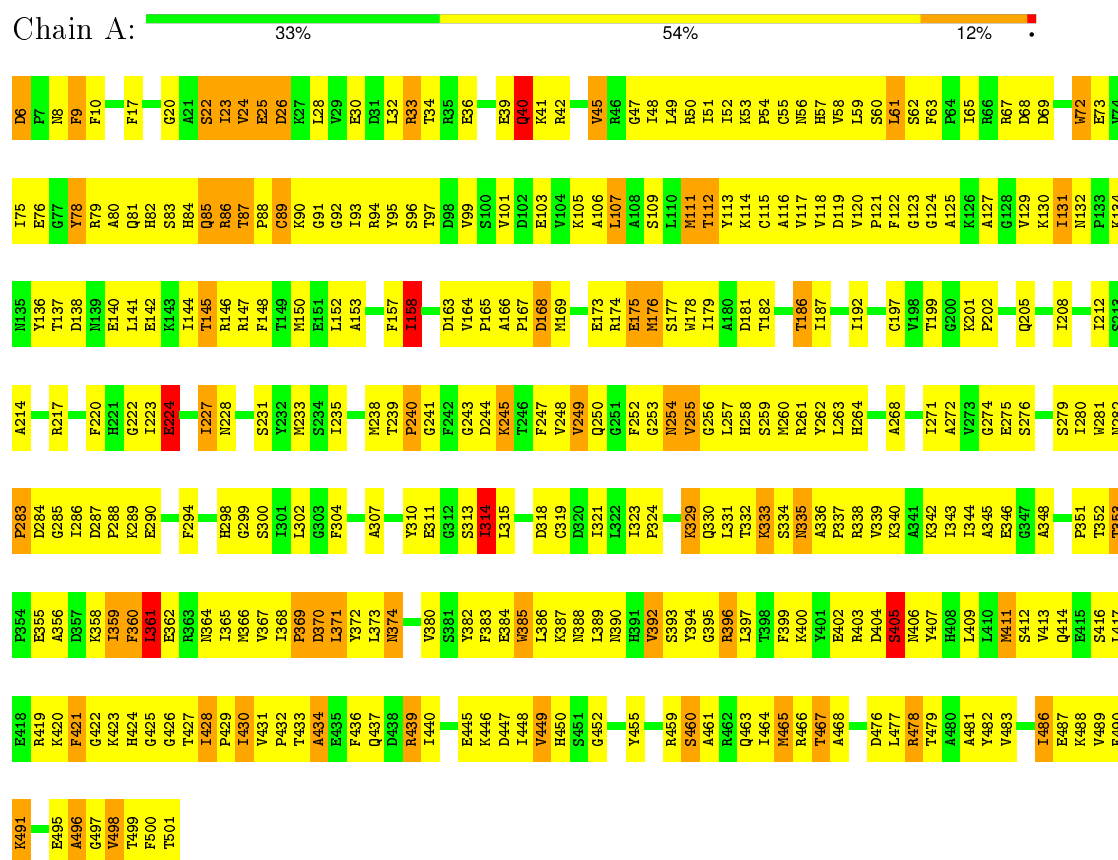


### 3 Residue-property plots

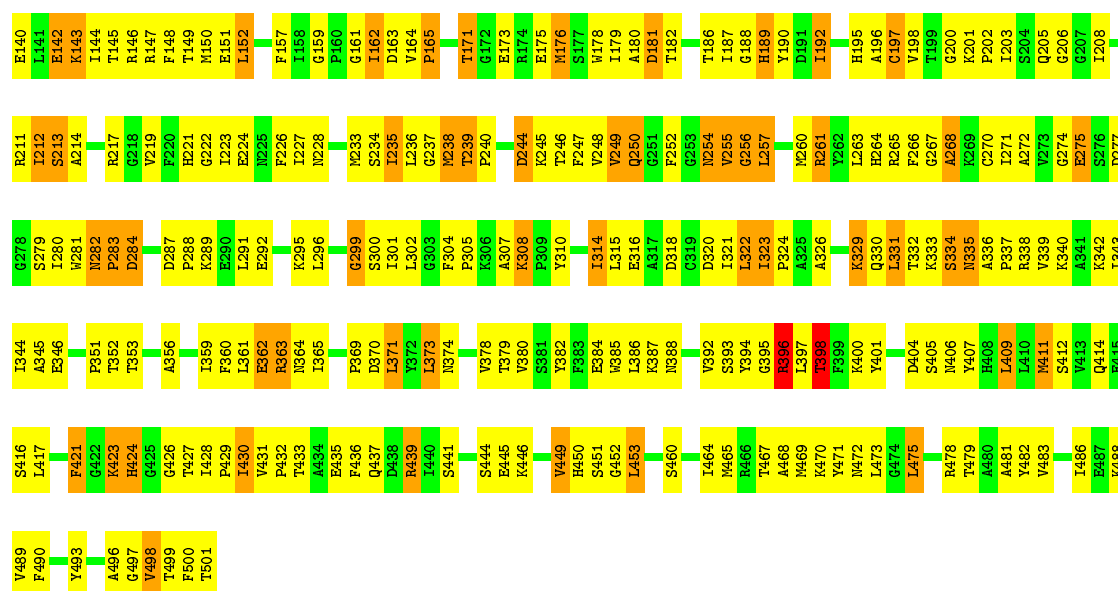
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 1

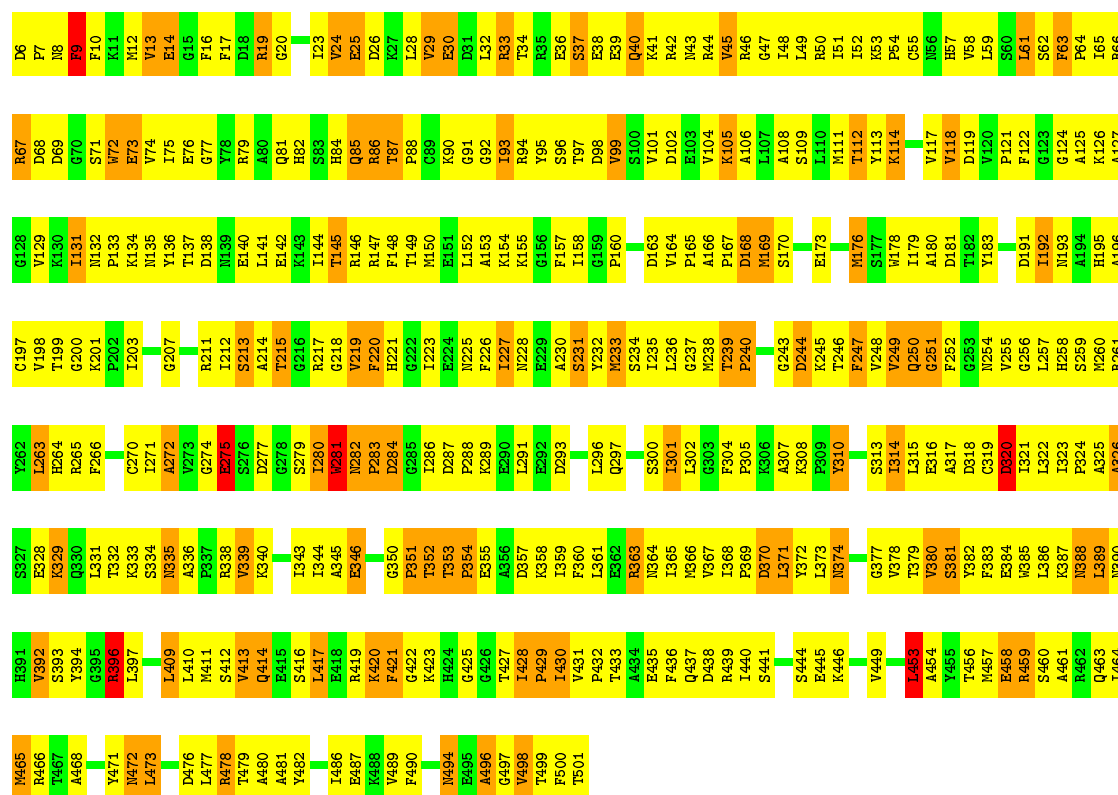






• Molecule 1: Glutamate dehydrogenase 1

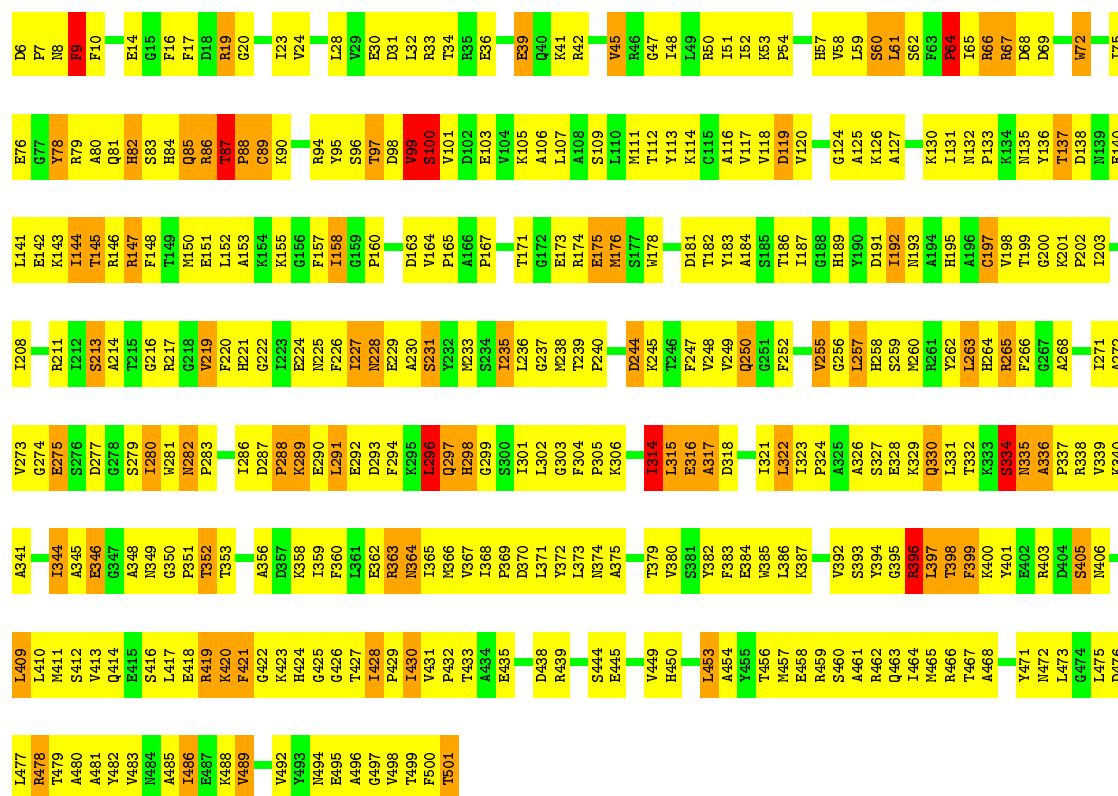
Chain C: 26% 54% 18%



• Molecule 1: Glutamate dehydrogenase 1

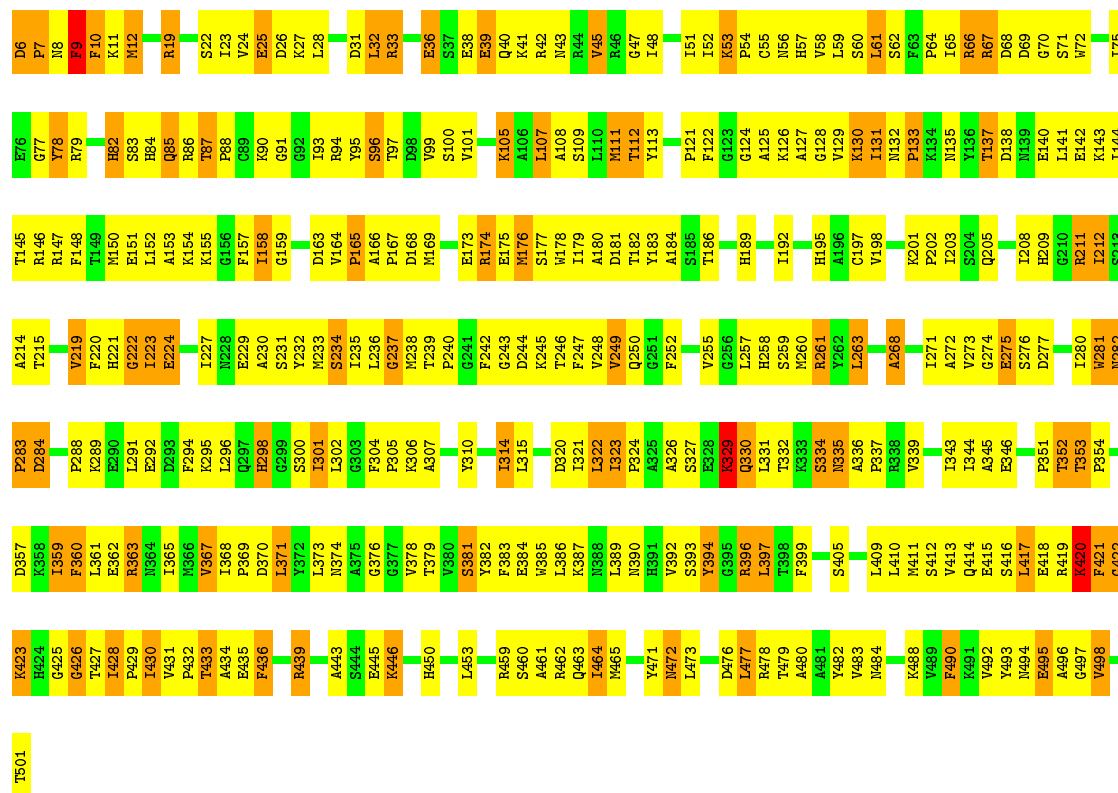
Chain D: 29% 55% 15%





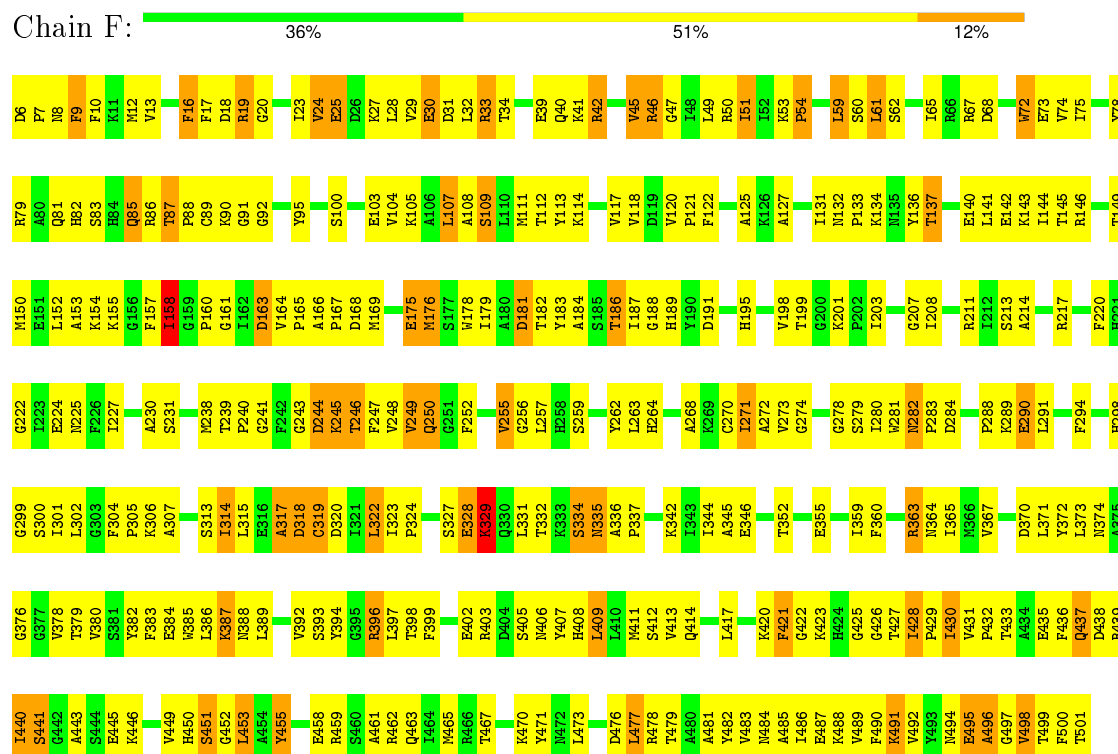
# Molecule 1: Glutamate dehydrogenase 1

Chain E: 33% 49% 17%

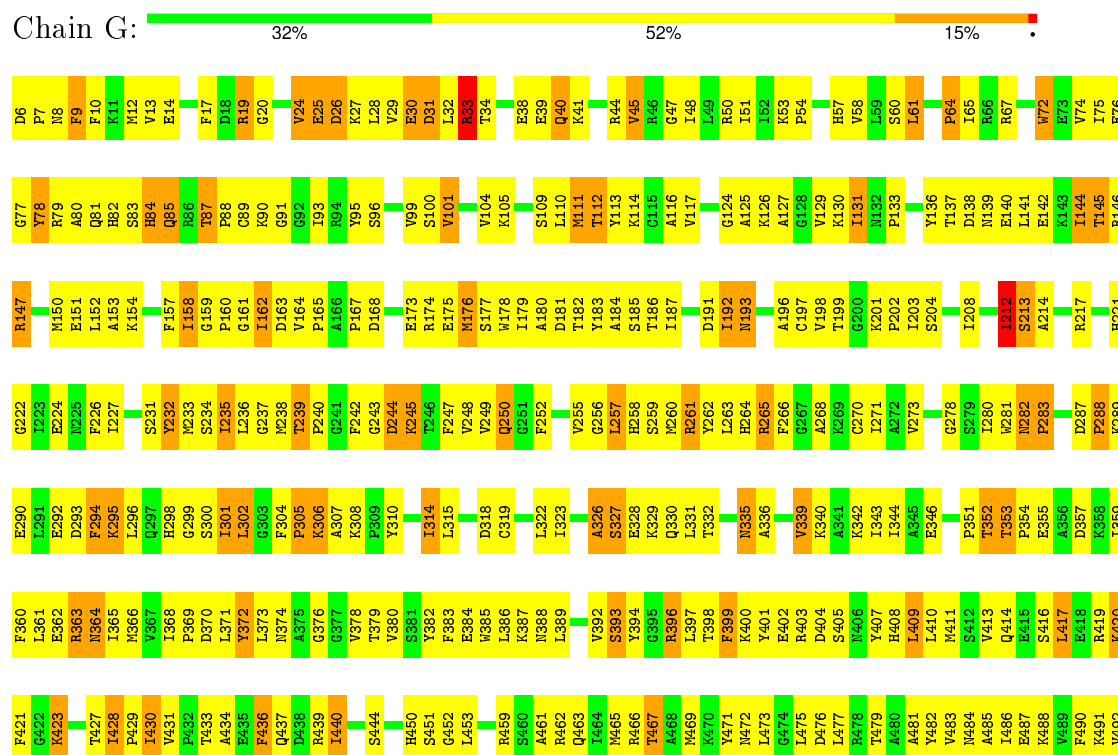




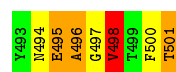
● Molecule 1: Glutamate dehydrogenase 1



● Molecule 1: Glutamate dehydrogenase 1

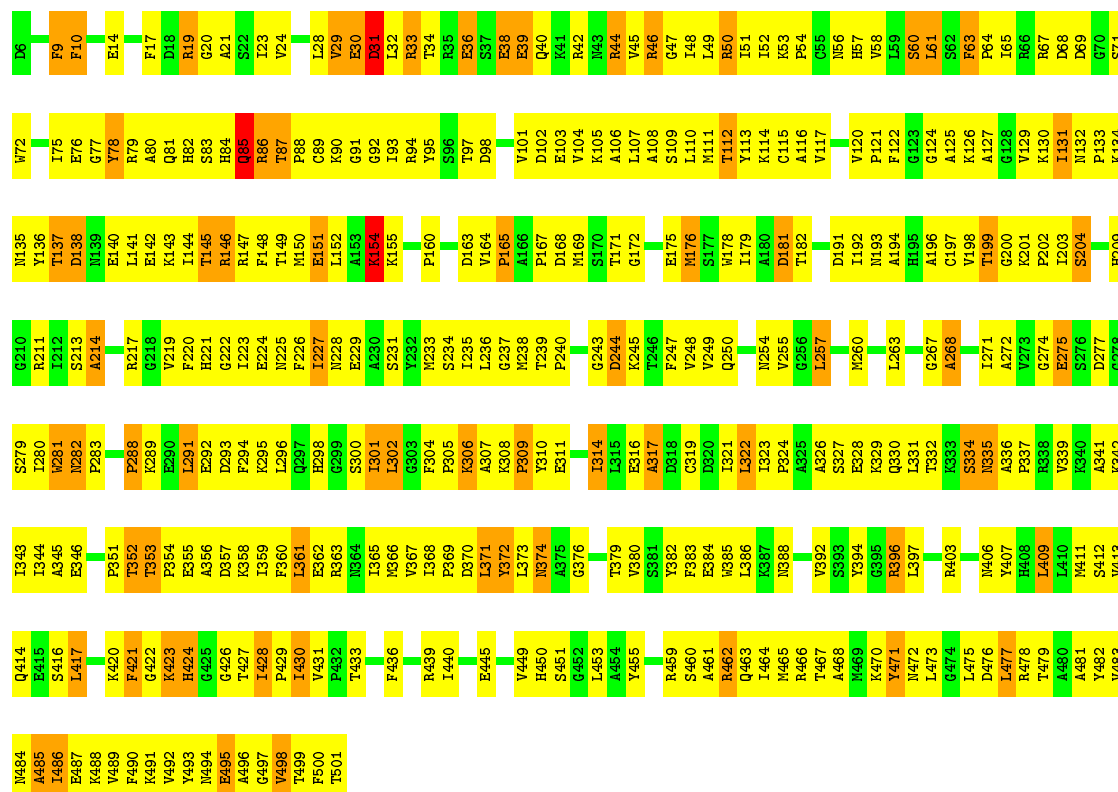






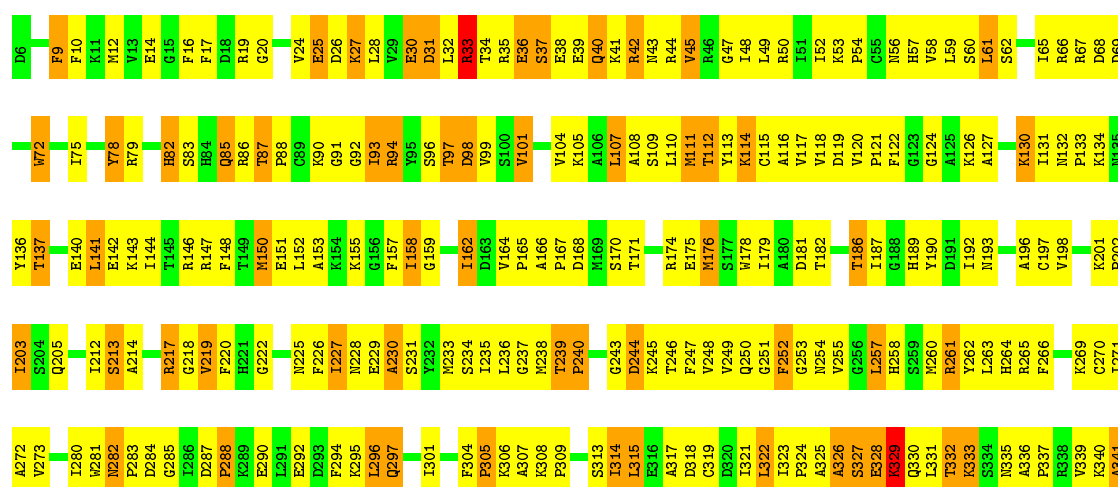
• Molecule 1: Glutamate dehydrogenase 1

Chain H: 30% 55% 14% .



• Molecule 1: Glutamate dehydrogenase 1

Chain I: 31% 52% 16% .

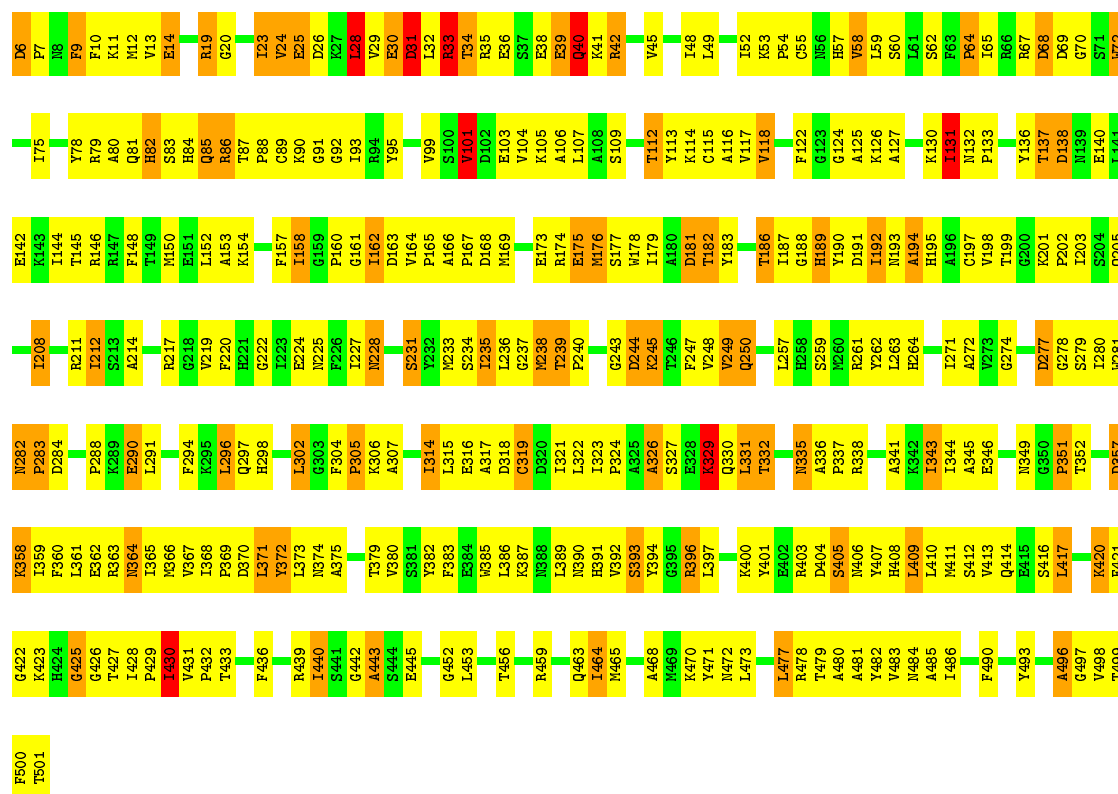






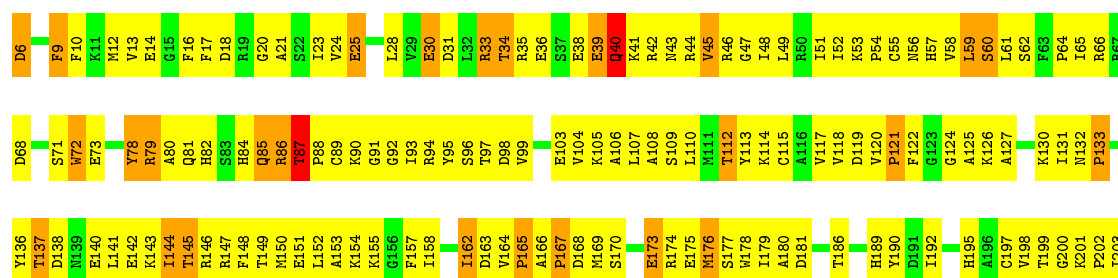
• Molecule 1: Glutamate dehydrogenase 1

Chain J: 34% 49% 15% .

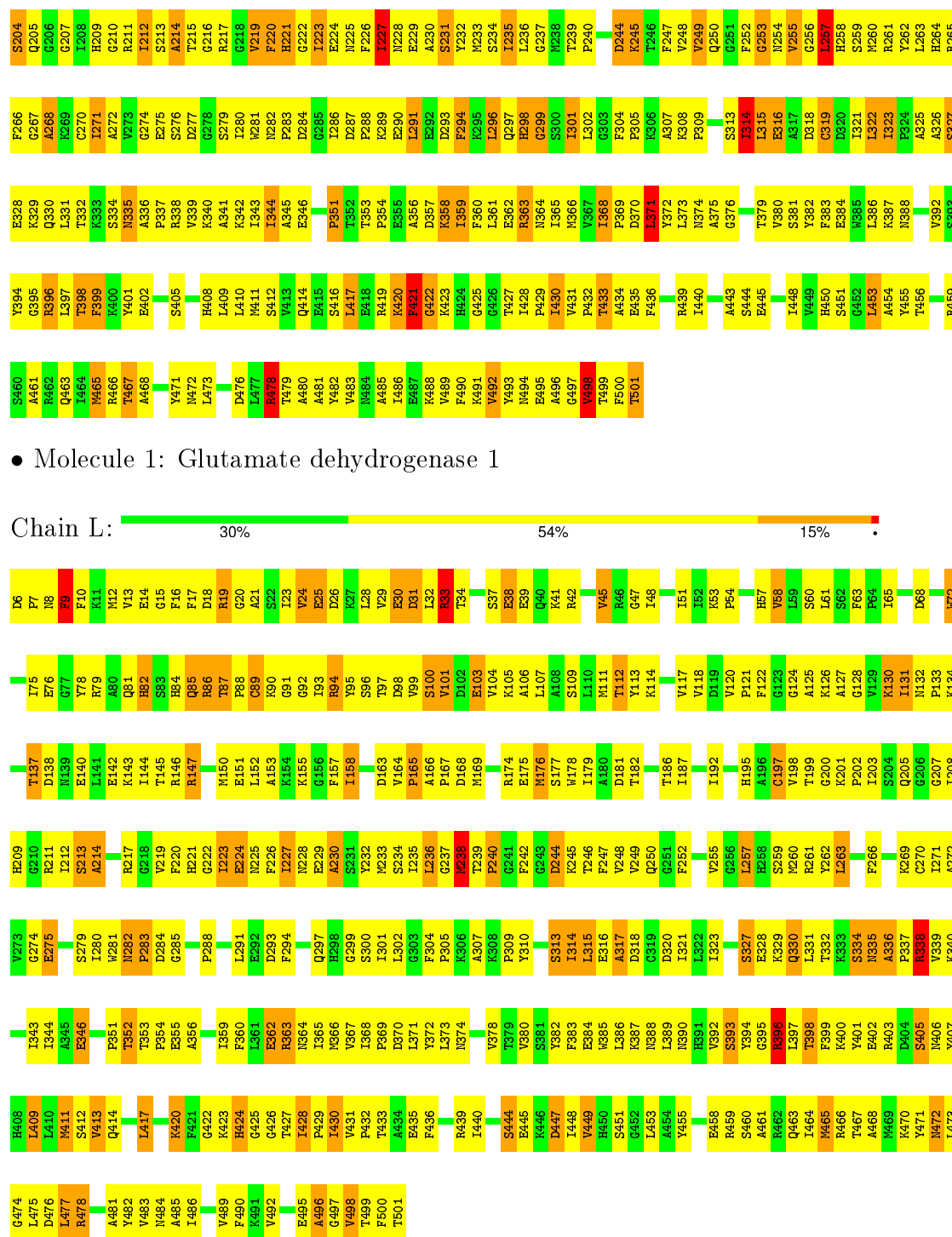


• Molecule 1: Glutamate dehydrogenase 1

Chain K: 24% 60% 15% .







• Molecule 1: Glutamate dehydrogenase 1



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	96.46 Å   171.93 Å   439.68 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	19.97 – 3.50	Depositor
% Data completeness (in resolution range)	94.0 (19.97-3.50)	Depositor
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.207 , 0.276	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	46812	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP



## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

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

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	0.70	1/3958 (0.0%)	0.89	2/5340 (0.0%)
1	B	0.76	1/3958 (0.0%)	0.91	2/5340 (0.0%)
1	C	0.72	1/3958 (0.0%)	0.91	2/5340 (0.0%)
1	D	0.69	2/3958 (0.1%)	0.89	4/5340 (0.1%)
1	E	0.71	0/3958	0.91	2/5340 (0.0%)
1	F	0.73	0/3958	0.89	0/5340
1	G	0.79	0/3958	0.95	2/5340 (0.0%)
1	H	0.78	1/3958 (0.0%)	0.89	1/5340 (0.0%)
1	I	0.71	1/3958 (0.0%)	0.88	1/5340 (0.0%)
1	J	0.74	1/3958 (0.0%)	0.91	1/5340 (0.0%)
1	K	0.70	0/3958	0.89	2/5340 (0.0%)
1	L	0.73	2/3958 (0.1%)	0.88	0/5340
All	All	0.73	10/47496 (0.0%)	0.90	19/64080 (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	E	0	2
1	F	0	1
1	G	0	1
1	H	0	1
All	All	0	5

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	197	CYS	CB-SG	-7.54	1.69	1.82
1	B	197	CYS	CB-SG	-6.96	1.70	1.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	89	CYS	CB-SG	-6.15	1.71	1.82
1	C	55	CYS	CB-SG	-6.08	1.72	1.82
1	L	89	CYS	CB-SG	-5.47	1.73	1.81

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	245	LYS	CD-CE-NZ	7.41	128.75	111.70
1	J	296	LEU	CA-CB-CG	6.82	130.98	115.30
1	B	61	LEU	CA-CB-CG	6.37	129.94	115.30
1	D	87	THR	N-CA-C	6.27	127.92	111.00
1	H	146	ARG	NE-CZ-NH2	-6.19	117.20	120.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	310	TYR	Sidechain
1	E	490	PHE	Sidechain
1	F	455	TYR	Sidechain
1	G	113	TYR	Sidechain
1	H	310	TYR	Sidechain

## 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	3874	0	3841	456	0
1	B	3874	0	3841	464	0
1	C	3874	0	3841	527	0
1	D	3874	0	3841	499	0
1	E	3874	0	3841	437	0
1	F	3874	0	3841	420	0
1	G	3874	0	3841	453	0
1	H	3874	0	3841	481	0
1	I	3874	0	3841	509	0
1	J	3874	0	3841	446	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	K	3874	0	3841	569	0
1	L	3874	0	3841	508	0
2	A	27	0	12	6	0
2	B	27	0	12	5	0
2	C	27	0	12	6	0
2	D	27	0	12	5	0
2	E	27	0	12	6	0
2	F	27	0	12	3	0
2	G	27	0	12	4	0
2	H	27	0	12	2	0
2	I	27	0	12	5	0
2	J	27	0	12	3	0
2	K	27	0	12	2	0
2	L	27	0	12	4	0
All	All	46812	0	46236	5447	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 59.

The worst 5 of 5447 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:423:LYS:HD3	1:A:426:GLY:HA3	1.23	1.17
1:B:95:TYR:OH	1:B:145:THR:HG22	1.44	1.16
1:L:57:HIS:CD2	1:L:84:HIS:HE1	1.62	1.15
1:C:235:ILE:HG22	1:C:364:ASN:HD21	1.12	1.13
1:C:396:ARG:HH11	1:C:396:ARG:HG3	1.14	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	494/496 (100%)	341 (69%)	109 (22%)	44 (9%)	1	11
1	B	494/496 (100%)	359 (73%)	94 (19%)	41 (8%)	1	13
1	C	494/496 (100%)	351 (71%)	82 (17%)	61 (12%)	0	6
1	D	494/496 (100%)	344 (70%)	105 (21%)	45 (9%)	1	11
1	E	494/496 (100%)	352 (71%)	100 (20%)	42 (8%)	1	13
1	F	494/496 (100%)	378 (76%)	85 (17%)	31 (6%)	2	21
1	G	494/496 (100%)	362 (73%)	94 (19%)	38 (8%)	1	14
1	H	494/496 (100%)	341 (69%)	116 (24%)	37 (8%)	1	15
1	I	494/496 (100%)	355 (72%)	89 (18%)	50 (10%)	1	9
1	J	494/496 (100%)	353 (72%)	99 (20%)	42 (8%)	1	13
1	K	494/496 (100%)	328 (66%)	113 (23%)	53 (11%)	0	8
1	L	494/496 (100%)	347 (70%)	102 (21%)	45 (9%)	1	11
All	All	5928/5952 (100%)	4211 (71%)	1188 (20%)	529 (9%)	1	11

5 of 529 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	9	PHE
1	A	25	GLU
1	A	26	ASP
1	A	30	GLU
1	A	33	ARG

### 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	413/413 (100%)	368 (89%)	45 (11%)	8	37
1	B	413/413 (100%)	352 (85%)	61 (15%)	4	22
1	C	413/413 (100%)	348 (84%)	65 (16%)	3	19
1	D	413/413 (100%)	354 (86%)	59 (14%)	4	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	413/413 (100%)	343 (83%)	70 (17%)	2	15
1	F	413/413 (100%)	359 (87%)	54 (13%)	5	27
1	G	413/413 (100%)	354 (86%)	59 (14%)	4	24
1	H	413/413 (100%)	348 (84%)	65 (16%)	3	19
1	I	413/413 (100%)	362 (88%)	51 (12%)	6	29
1	J	413/413 (100%)	348 (84%)	65 (16%)	3	19
1	K	413/413 (100%)	346 (84%)	67 (16%)	3	17
1	L	413/413 (100%)	351 (85%)	62 (15%)	3	21
All	All	4956/4956 (100%)	4233 (85%)	723 (15%)	4	22

5 of 723 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	238	MET
1	G	428	ILE
1	L	58	VAL
1	F	322	LEU
1	G	100	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 164 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	84	HIS
1	G	139	ASN
1	L	205	GLN
1	F	139	ASN
1	F	388	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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

12 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	ADP	A	1	-	22,29,29	1.50	3 (13%)	27,45,45	2.43	5 (18%)
2	ADP	B	2	-	22,29,29	1.20	2 (9%)	27,45,45	2.46	4 (14%)
2	ADP	C	3	-	22,29,29	1.34	3 (13%)	27,45,45	2.38	5 (18%)
2	ADP	D	4	-	22,29,29	1.85	4 (18%)	27,45,45	2.34	4 (14%)
2	ADP	E	5	-	22,29,29	1.90	3 (13%)	27,45,45	2.50	4 (14%)
2	ADP	F	502	-	22,29,29	1.25	2 (9%)	27,45,45	2.41	5 (18%)
2	ADP	G	502	-	22,29,29	1.64	4 (18%)	27,45,45	2.38	5 (18%)
2	ADP	H	502	-	22,29,29	1.17	2 (9%)	27,45,45	2.44	4 (14%)
2	ADP	I	502	-	22,29,29	1.19	2 (9%)	27,45,45	2.44	5 (18%)
2	ADP	J	502	-	22,29,29	1.18	2 (9%)	27,45,45	2.42	5 (18%)
2	ADP	K	502	-	22,29,29	1.24	2 (9%)	27,45,45	2.40	5 (18%)
2	ADP	L	502	-	22,29,29	1.26	2 (9%)	27,45,45	2.39	5 (18%)

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	ADP	A	1	-	-	0/12/32/32	0/3/3/3
2	ADP	B	2	-	-	0/12/32/32	0/3/3/3
2	ADP	C	3	-	-	0/12/32/32	0/3/3/3
2	ADP	D	4	-	-	0/12/32/32	0/3/3/3
2	ADP	E	5	-	-	0/12/32/32	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ADP	F	502	-	-	0/12/32/32	0/3/3/3
2	ADP	G	502	-	-	0/12/32/32	0/3/3/3
2	ADP	H	502	-	-	0/12/32/32	0/3/3/3
2	ADP	I	502	-	-	0/12/32/32	0/3/3/3
2	ADP	J	502	-	-	0/12/32/32	0/3/3/3
2	ADP	K	502	-	-	0/12/32/32	0/3/3/3
2	ADP	L	502	-	-	0/12/32/32	0/3/3/3

The worst 5 of 31 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	3	ADP	C5-N7	-2.12	1.32	1.39
2	A	1	ADP	C5-N7	-2.06	1.32	1.39
2	D	4	ADP	C5-N7	-2.02	1.32	1.39
2	G	502	ADP	PA-O5'	2.03	1.68	1.59
2	G	502	ADP	C2-N3	2.03	1.35	1.32

The worst 5 of 56 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	ADP	N3-C2-N1	-10.69	120.71	128.89
2	H	502	ADP	N3-C2-N1	-10.52	120.84	128.89
2	E	5	ADP	N3-C2-N1	-10.45	120.89	128.89
2	L	502	ADP	N3-C2-N1	-10.38	120.95	128.89
2	I	502	ADP	N3-C2-N1	-10.36	120.96	128.89

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

12 monomers are involved in 51 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1	ADP	6	0
2	B	2	ADP	5	0
2	C	3	ADP	6	0
2	D	4	ADP	5	0
2	E	5	ADP	6	0
2	F	502	ADP	3	0
2	G	502	ADP	4	0
2	H	502	ADP	2	0
2	I	502	ADP	5	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	J	502	ADP	3	0
2	K	502	ADP	2	0
2	L	502	ADP	4	0

## 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 ⓘ

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.