



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 12:59 PM UTC

PDB ID : 1SXI / pdb_00001sxi
Title : Structure of apo transcription regulator B. megaterium
Authors : Schumacher, M.A.; Allen, G.S.; Diel, M.; Seidel, G.; Hillen, W.; Brennan, R.G.
Deposited on : 2004-03-30
Resolution : 3.00 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

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
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

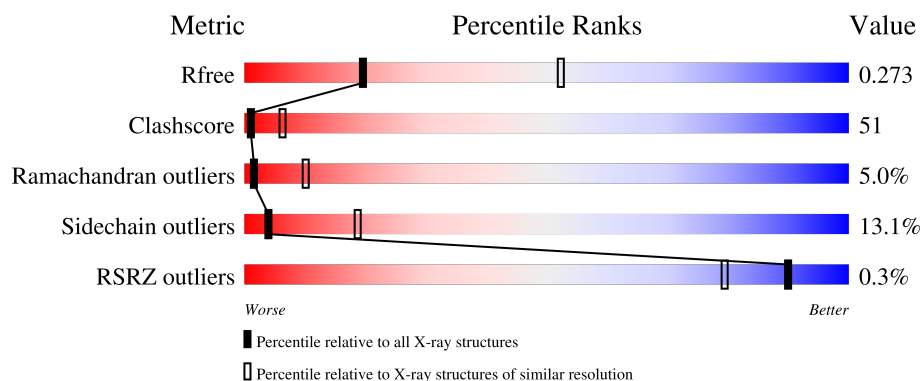
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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.



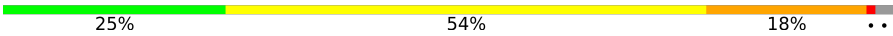
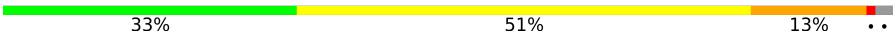
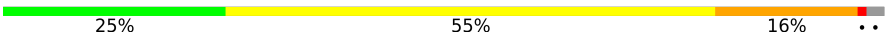
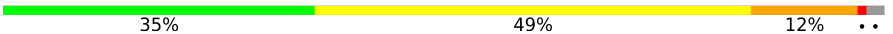
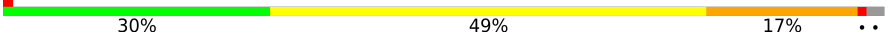
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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 $\leq 5\%$. 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	280	29% 55% 13% ..
1	B	280	31% 52% 12% ..
1	D	280	% 31% 51% 14% ..
1	G	280	26% 56% 14% ..
1	I	280	38% 48% 11% ..

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	K	280	
1	L	280	
1	M	280	
1	N	280	
1	R	280	
1	T	280	
1	W	280	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 25483 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 Glucose-resistance amylase regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	Se	0	0	0
			2123	1337	352	426	8			
1	D	273	Total	C	N	O	Se	0	0	0
			2123	1337	352	426	8			
1	B	273	Total	C	N	O	Se	0	0	0
			2123	1337	352	426	8			
1	I	273	Total	C	N	O	Se	0	0	0
			2123	1337	352	426	8			
1	R	273	Total	C	N	O	Se	0	0	0
			2123	1337	352	426	8			
1	T	273	Total	C	N	O	Se	0	0	0
			2123	1337	352	426	8			
1	L	273	Total	C	N	O	Se	0	0	0
			2123	1337	352	426	8			
1	K	273	Total	C	N	O	Se	0	0	0
			2123	1337	352	426	8			
1	W	273	Total	C	N	O	Se	0	0	0
			2123	1337	352	426	8			
1	G	273	Total	C	N	O	Se	0	0	0
			2123	1337	352	426	8			
1	N	273	Total	C	N	O	Se	0	0	0
			2123	1337	352	426	8			
1	M	273	Total	C	N	O	Se	0	0	0
			2123	1337	352	426	8			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	88	MSE	MET	modified residue	UNP P46828
A	112	MSE	MET	modified residue	UNP P46828
A	123	MSE	MET	modified residue	UNP P46828
A	250	MSE	MET	modified residue	UNP P46828
A	282	MSE	MET	modified residue	UNP P46828

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	294	MSE	MET	modified residue	UNP P46828
A	302	MSE	MET	modified residue	UNP P46828
A	309	MSE	MET	modified residue	UNP P46828
B	88	MSE	MET	modified residue	UNP P46828
B	112	MSE	MET	modified residue	UNP P46828
B	123	MSE	MET	modified residue	UNP P46828
B	250	MSE	MET	modified residue	UNP P46828
B	282	MSE	MET	modified residue	UNP P46828
B	294	MSE	MET	modified residue	UNP P46828
B	302	MSE	MET	modified residue	UNP P46828
B	309	MSE	MET	modified residue	UNP P46828
R	88	MSE	MET	modified residue	UNP P46828
R	112	MSE	MET	modified residue	UNP P46828
R	123	MSE	MET	modified residue	UNP P46828
R	250	MSE	MET	modified residue	UNP P46828
R	282	MSE	MET	modified residue	UNP P46828
R	294	MSE	MET	modified residue	UNP P46828
R	302	MSE	MET	modified residue	UNP P46828
R	309	MSE	MET	modified residue	UNP P46828
L	88	MSE	MET	modified residue	UNP P46828
L	112	MSE	MET	modified residue	UNP P46828
L	123	MSE	MET	modified residue	UNP P46828
L	250	MSE	MET	modified residue	UNP P46828
L	282	MSE	MET	modified residue	UNP P46828
L	294	MSE	MET	modified residue	UNP P46828
L	302	MSE	MET	modified residue	UNP P46828
L	309	MSE	MET	modified residue	UNP P46828
W	88	MSE	MET	modified residue	UNP P46828
W	112	MSE	MET	modified residue	UNP P46828
W	123	MSE	MET	modified residue	UNP P46828
W	250	MSE	MET	modified residue	UNP P46828
W	282	MSE	MET	modified residue	UNP P46828
W	294	MSE	MET	modified residue	UNP P46828
W	302	MSE	MET	modified residue	UNP P46828
W	309	MSE	MET	modified residue	UNP P46828
M	88	MSE	MET	modified residue	UNP P46828
M	112	MSE	MET	modified residue	UNP P46828
M	123	MSE	MET	modified residue	UNP P46828
M	250	MSE	MET	modified residue	UNP P46828
M	282	MSE	MET	modified residue	UNP P46828
M	294	MSE	MET	modified residue	UNP P46828
M	302	MSE	MET	modified residue	UNP P46828

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
M	309	MSE	MET	modified residue	UNP P46828
D	88	MSE	MET	modified residue	UNP P46828
D	112	MSE	MET	modified residue	UNP P46828
D	123	MSE	MET	modified residue	UNP P46828
D	250	MSE	MET	modified residue	UNP P46828
D	282	MSE	MET	modified residue	UNP P46828
D	294	MSE	MET	modified residue	UNP P46828
D	302	MSE	MET	modified residue	UNP P46828
D	309	MSE	MET	modified residue	UNP P46828
I	88	MSE	MET	modified residue	UNP P46828
I	112	MSE	MET	modified residue	UNP P46828
I	123	MSE	MET	modified residue	UNP P46828
I	250	MSE	MET	modified residue	UNP P46828
I	282	MSE	MET	modified residue	UNP P46828
I	294	MSE	MET	modified residue	UNP P46828
I	302	MSE	MET	modified residue	UNP P46828
I	309	MSE	MET	modified residue	UNP P46828
T	88	MSE	MET	modified residue	UNP P46828
T	112	MSE	MET	modified residue	UNP P46828
T	123	MSE	MET	modified residue	UNP P46828
T	250	MSE	MET	modified residue	UNP P46828
T	282	MSE	MET	modified residue	UNP P46828
T	294	MSE	MET	modified residue	UNP P46828
T	302	MSE	MET	modified residue	UNP P46828
T	309	MSE	MET	modified residue	UNP P46828
K	88	MSE	MET	modified residue	UNP P46828
K	112	MSE	MET	modified residue	UNP P46828
K	123	MSE	MET	modified residue	UNP P46828
K	250	MSE	MET	modified residue	UNP P46828
K	282	MSE	MET	modified residue	UNP P46828
K	294	MSE	MET	modified residue	UNP P46828
K	302	MSE	MET	modified residue	UNP P46828
K	309	MSE	MET	modified residue	UNP P46828
G	88	MSE	MET	modified residue	UNP P46828
G	112	MSE	MET	modified residue	UNP P46828
G	123	MSE	MET	modified residue	UNP P46828
G	250	MSE	MET	modified residue	UNP P46828
G	282	MSE	MET	modified residue	UNP P46828
G	294	MSE	MET	modified residue	UNP P46828
G	302	MSE	MET	modified residue	UNP P46828
G	309	MSE	MET	modified residue	UNP P46828
N	88	MSE	MET	modified residue	UNP P46828

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
N	112	MSE	MET	modified residue	UNP P46828
N	123	MSE	MET	modified residue	UNP P46828
N	250	MSE	MET	modified residue	UNP P46828
N	282	MSE	MET	modified residue	UNP P46828
N	294	MSE	MET	modified residue	UNP P46828
N	302	MSE	MET	modified residue	UNP P46828
N	309	MSE	MET	modified residue	UNP P46828

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	T	2	Total Mg 2 2	0	0

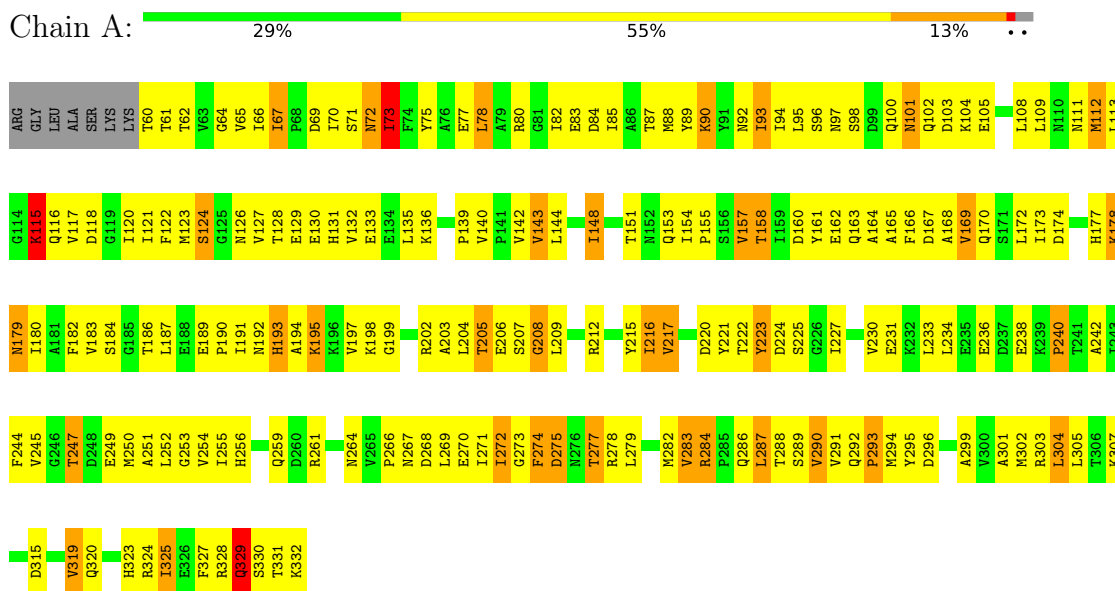
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O 1 1	0	0
3	D	3	Total O 3 3	0	0
3	T	1	Total O 1 1	0	0

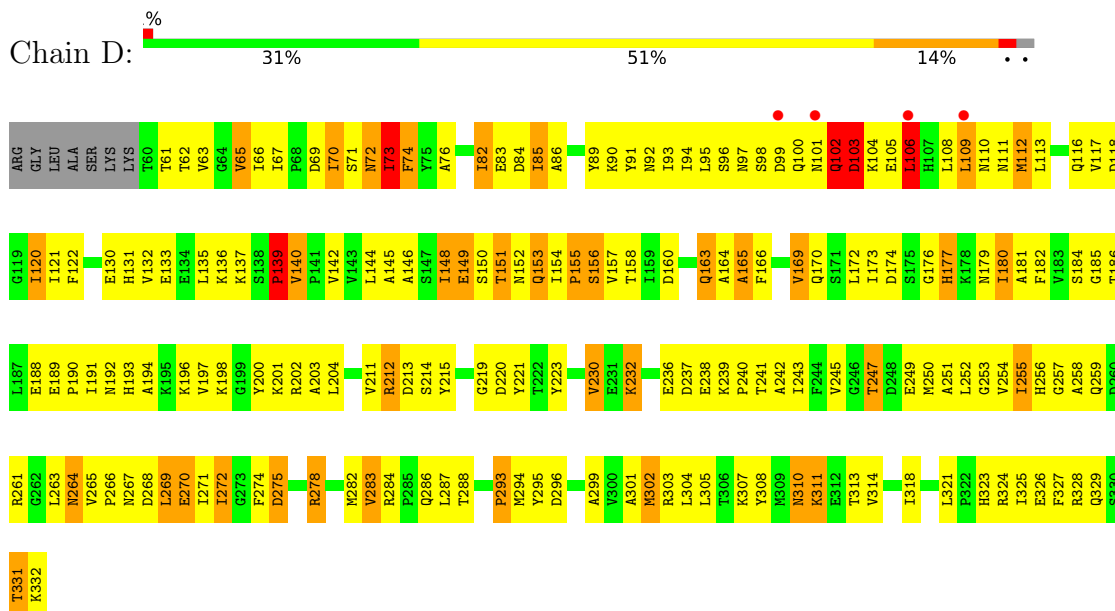
3 Residue-property plots

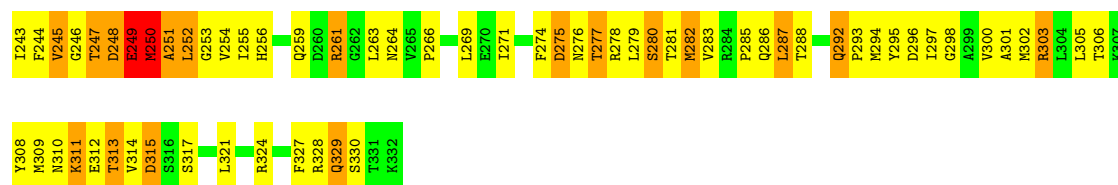
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Glucose-resistance amylase regulator



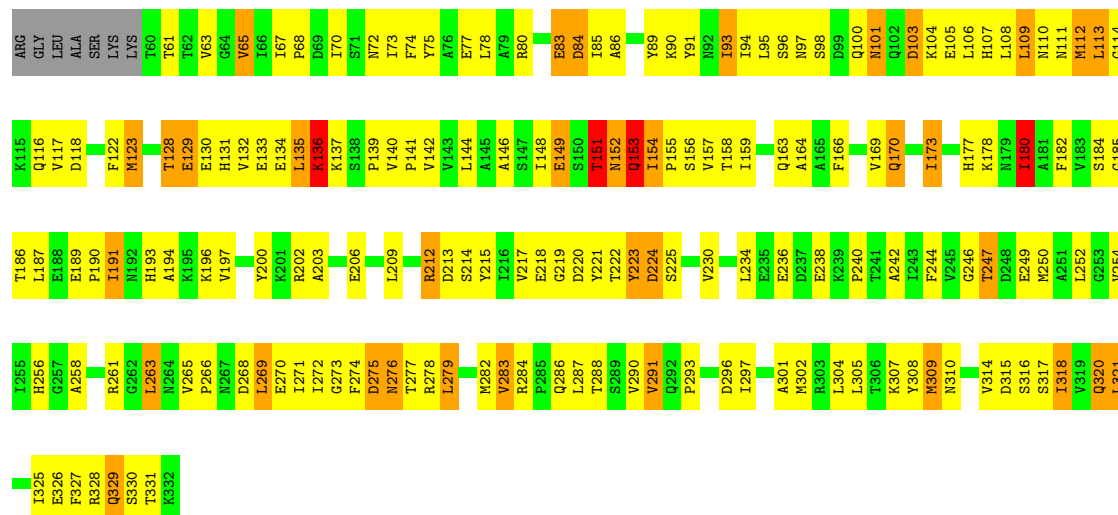
- Molecule 1: Glucose-resistance amylase regulator





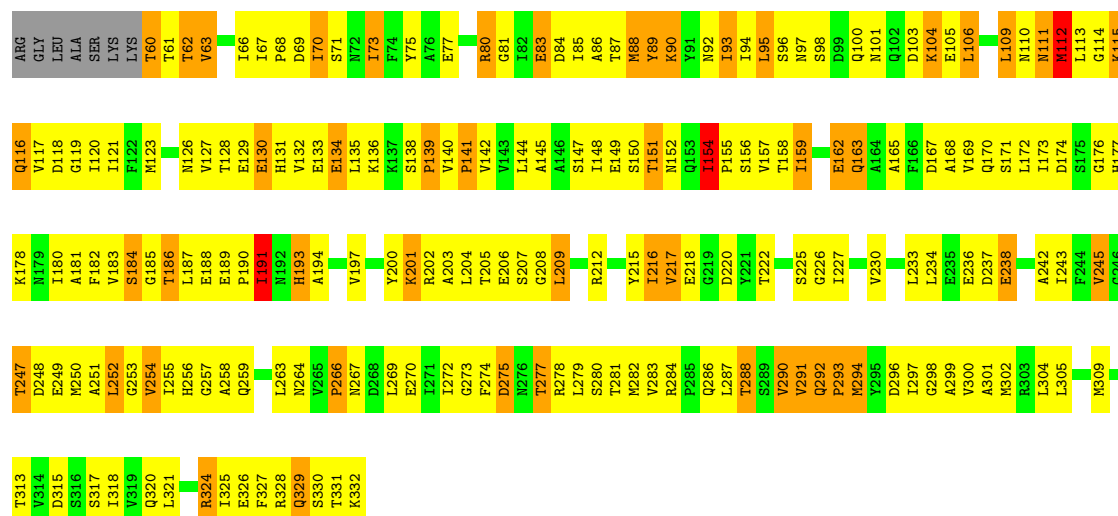
• Molecule 1: Glucose-resistance amylase regulator

Chain T: 35% 49% 12% ..



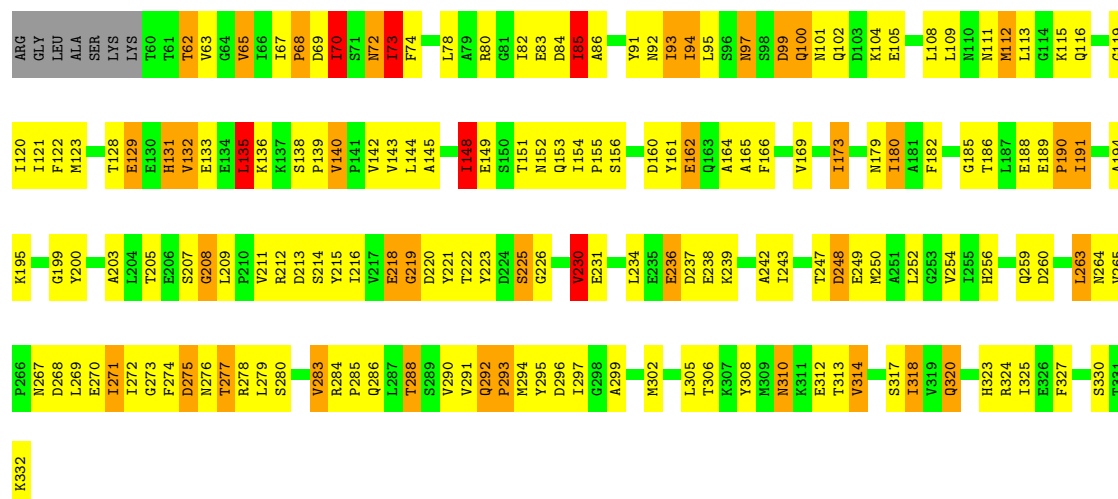
• Molecule 1: Glucose-resistance amylase regulator

Chain L: 25% 54% 18% ..

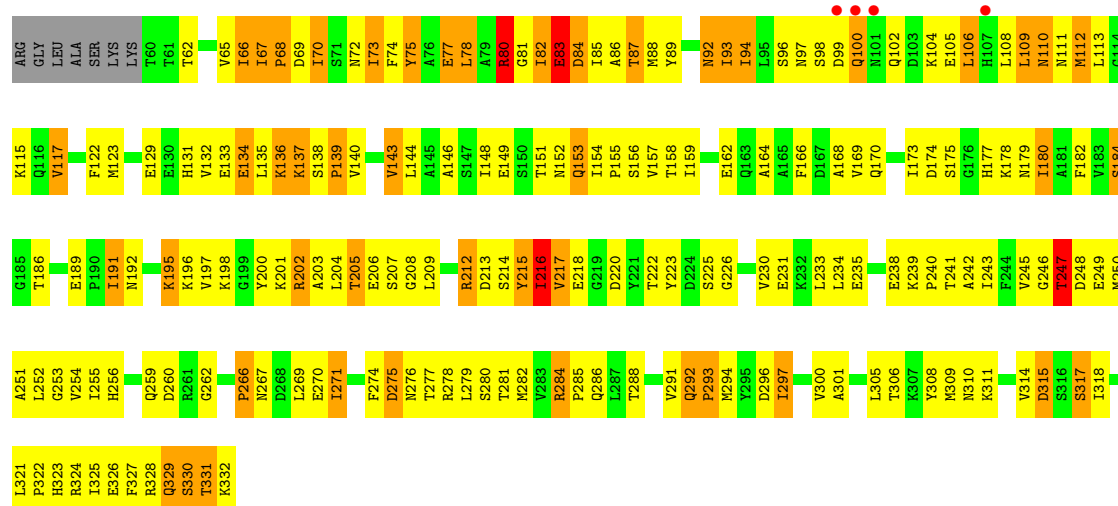


• Molecule 1: Glucose-resistance amylase regulator

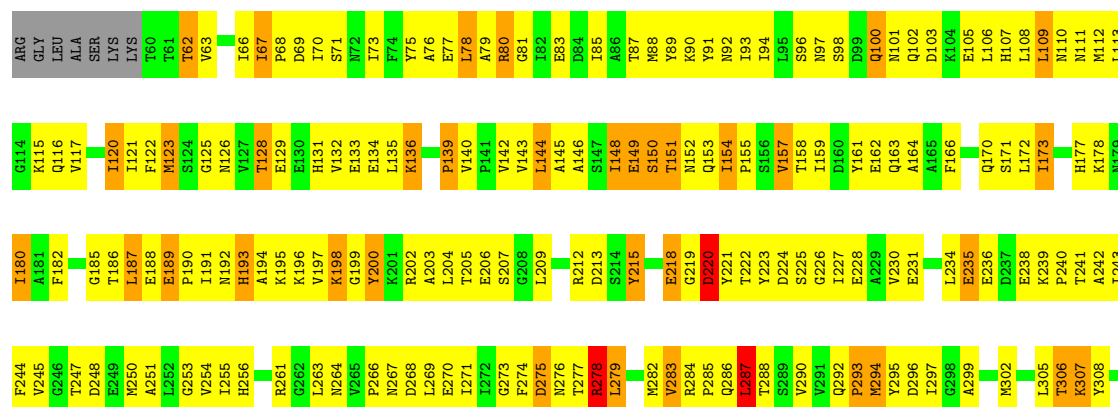
Chain K: 37% 45% 13% ..



● Molecule 1: Glucose-resistance amylase regulator



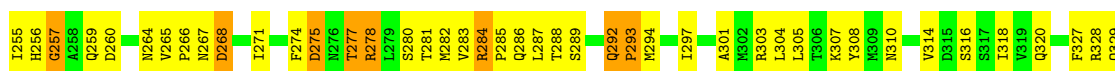
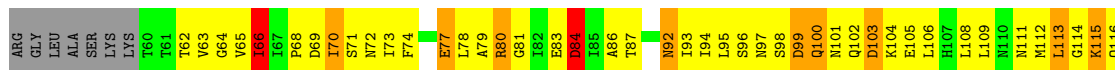
● Molecule 1: Glucose-resistance amylase regulator





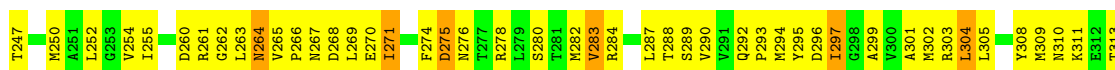
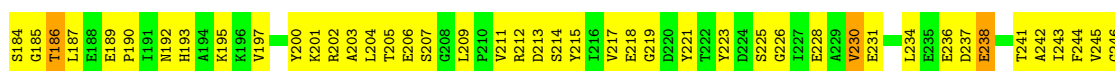
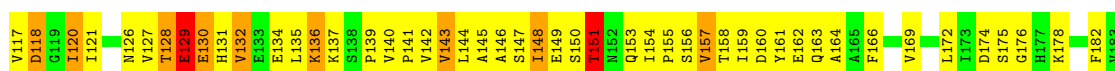
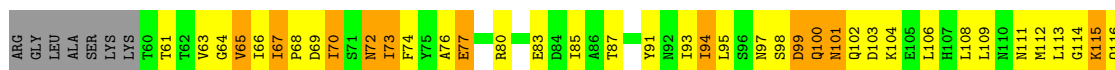
• Molecule 1: Glucose-resistance amylase regulator

Chain N: 33% 51% 13% ..



• Molecule 1: Glucose-resistance amylase regulator

Chain M: 30% 55% 12% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	110.24Å 106.22Å 157.74Å 90.00° 108.20° 90.00°	Depositor
Resolution (Å)	74.58 – 3.00 74.58 – 3.00	Depositor EDS
% Data completeness (in resolution range)	95.5 (74.58-3.00) 95.5 (74.58-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 3.01Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.220 , 0.282 0.215 , 0.273	Depositor DCC
R_{free} test set	6726 reflections (10.09%)	wwPDB-VP
Wilson B-factor (Å ²)	61.6	Xtriage
Anisotropy	0.084	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 60.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.020 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	25483	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% of the height of the origin peak. No significant pseudotranslation is detected.*

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 ⓘ

5.1 Standard geometry ⓘ

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

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	0.59	1/2148 (0.0%)	1.08	9/2899 (0.3%)
1	B	0.54	0/2148	1.10	15/2899 (0.5%)
1	D	0.59	1/2148 (0.0%)	1.08	17/2899 (0.6%)
1	G	0.52	0/2148	1.02	11/2899 (0.4%)
1	I	0.60	2/2148 (0.1%)	1.16	20/2899 (0.7%)
1	K	0.60	1/2148 (0.0%)	1.10	20/2899 (0.7%)
1	L	0.53	0/2148	1.07	13/2899 (0.4%)
1	M	0.53	0/2148	1.05	12/2899 (0.4%)
1	N	0.49	0/2148	1.02	9/2899 (0.3%)
1	R	0.52	0/2148	1.06	15/2899 (0.5%)
1	T	0.58	1/2148 (0.0%)	1.14	16/2899 (0.6%)
1	W	0.61	1/2147 (0.0%)	1.13	19/2896 (0.7%)
All	All	0.56	7/25775 (0.0%)	1.09	176/34785 (0.5%)

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	112	MSE	SE-CE	-7.21	1.73	1.95
1	A	112	MSE	SE-CE	-6.37	1.76	1.95
1	T	151	THR	CA-CB	-6.08	1.43	1.53
1	K	112	MSE	SE-CE	-5.46	1.79	1.95
1	I	112	MSE	CG-SE	-5.27	1.79	1.95

The worst 5 of 176 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	T	151	THR	N-CA-C	13.75	125.96	110.97
1	K	70	ILE	N-CA-C	-10.41	102.96	113.47
1	I	117	VAL	N-CA-C	-9.37	99.83	110.05
1	L	109	LEU	N-CA-C	-9.20	101.23	111.07
1	K	140	VAL	N-CA-C	9.02	116.39	108.63

There are no chirality outliers.

There are no planarity outliers.

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	2123	0	2121	247	0
1	B	2123	0	2121	197	0
1	D	2123	0	2121	205	0
1	G	2123	0	2121	249	0
1	I	2123	0	2121	176	0
1	K	2123	0	2121	209	0
1	L	2123	0	2121	251	0
1	M	2123	0	2121	213	0
1	N	2123	0	2121	257	0
1	R	2123	0	2121	247	0
1	T	2123	0	2121	200	0
1	W	2123	0	2120	258	0
2	T	2	0	0	0	0
3	A	1	0	0	0	0
3	D	3	0	0	0	0
3	T	1	0	0	0	0
All	All	25483	0	25451	2587	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 51.

The worst 5 of 2587 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:W:115:LYS:HE2	1:G:94:ILE:HD11	1.25	1.17
1:D:121:ILE:HD11	1:D:302:MSE:HE3	1.20	1.16
1:G:234:LEU:HD13	1:G:261:ARG:HE	1.09	1.16
1:B:288:THR:HG22	1:B:328:ARG:H	1.07	1.15
1:B:70:ILE:HD11	1:I:70:ILE:HD11	1.20	1.15

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	271/280 (97%)	217 (80%)	43 (16%)	11 (4%)	2	13
1	B	271/280 (97%)	218 (80%)	39 (14%)	14 (5%)	1	9
1	D	271/280 (97%)	224 (83%)	33 (12%)	14 (5%)	1	9
1	G	271/280 (97%)	199 (73%)	56 (21%)	16 (6%)	1	7
1	I	271/280 (97%)	226 (83%)	39 (14%)	6 (2%)	5	26
1	K	271/280 (97%)	233 (86%)	30 (11%)	8 (3%)	3	19
1	L	271/280 (97%)	212 (78%)	42 (16%)	17 (6%)	1	6
1	M	271/280 (97%)	222 (82%)	38 (14%)	11 (4%)	2	13
1	N	271/280 (97%)	204 (75%)	47 (17%)	20 (7%)	1	4
1	R	271/280 (97%)	209 (77%)	41 (15%)	21 (8%)	1	4
1	T	271/280 (97%)	231 (85%)	29 (11%)	11 (4%)	2	13
1	W	269/280 (96%)	219 (81%)	38 (14%)	12 (4%)	2	12
All	All	3250/3360 (97%)	2614 (80%)	475 (15%)	161 (5%)	1	10

5 of 161 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	208	GLY
1	D	102	GLN
1	B	275	ASP
1	R	275	ASP
1	R	313	THR

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	239/236 (101%)	209 (87%)	30 (13%)	4	20
1	B	239/236 (101%)	202 (84%)	37 (16%)	2	13
1	D	239/236 (101%)	209 (87%)	30 (13%)	4	20
1	G	239/236 (101%)	207 (87%)	32 (13%)	4	18
1	I	239/236 (101%)	211 (88%)	28 (12%)	5	23
1	K	239/236 (101%)	208 (87%)	31 (13%)	4	19
1	L	239/236 (101%)	203 (85%)	36 (15%)	3	14
1	M	239/236 (101%)	212 (89%)	27 (11%)	5	24
1	N	239/236 (101%)	219 (92%)	20 (8%)	10	37
1	R	239/236 (101%)	205 (86%)	34 (14%)	3	16
1	T	239/236 (101%)	205 (86%)	34 (14%)	3	16
1	W	239/236 (101%)	202 (84%)	37 (16%)	2	13
All	All	2868/2832 (101%)	2492 (87%)	376 (13%)	4	19

5 of 376 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	K	67	ILE
1	W	216	ILE
1	K	99	ASP
1	W	66	ILE
1	G	80	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 126 such sidechains are listed below:

Mol	Chain	Res	Type
1	T	256	HIS
1	N	264	ASN
1	K	100	GLN
1	N	179	ASN
1	M	152	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	W	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	W	98:SER	C	99:ASP	N	2.69

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	265/280 (94%)	-0.01	0 100 100	29, 57, 81, 98	0
1	B	265/280 (94%)	-0.22	0 100 100	22, 50, 73, 83	0
1	D	265/280 (94%)	-0.32	4 (1%) 72 49	17, 37, 93, 108	0
1	G	265/280 (94%)	-0.03	0 100 100	27, 57, 84, 91	0
1	I	265/280 (94%)	-0.40	1 (0%) 88 76	13, 36, 89, 102	0
1	K	265/280 (94%)	-0.31	0 100 100	12, 42, 66, 83	0
1	L	265/280 (94%)	-0.25	0 100 100	22, 49, 74, 91	0
1	M	265/280 (94%)	-0.23	0 100 100	14, 49, 82, 94	0
1	N	265/280 (94%)	0.11	0 100 100	33, 63, 103, 116	0
1	R	265/280 (94%)	-0.12	0 100 100	18, 50, 93, 100	0
1	T	265/280 (94%)	-0.31	0 100 100	16, 37, 86, 112	0
1	W	265/280 (94%)	-0.10	4 (1%) 72 49	26, 49, 71, 85	0
All	All	3180/3360 (94%)	-0.18	9 (0%) 90 79	12, 49, 85, 116	0

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	W	100	GLN	3.7
1	D	99	ASP	3.3
1	W	101	ASN	2.7
1	W	99	ASP	2.7
1	D	106	LEU	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 [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	MG	T	754	1/1	0.92	0.23	61,61,61,61	0
2	MG	T	764	1/1	0.94	0.10	30,30,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.