



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 15, 2026 – 02:41 AM UTC

PDB ID : 6RH9 / pdb_00006rh9
Title : Crystal Structure of Two-Domain Laccase mutant I170F from *Streptomyces griseoflavus*
Authors : Gabdulkhakov, A.G.; Tishchenko, T.V.; Kolyadenko, I.A.
Deposited on : 2019-04-19
Resolution : 1.85 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

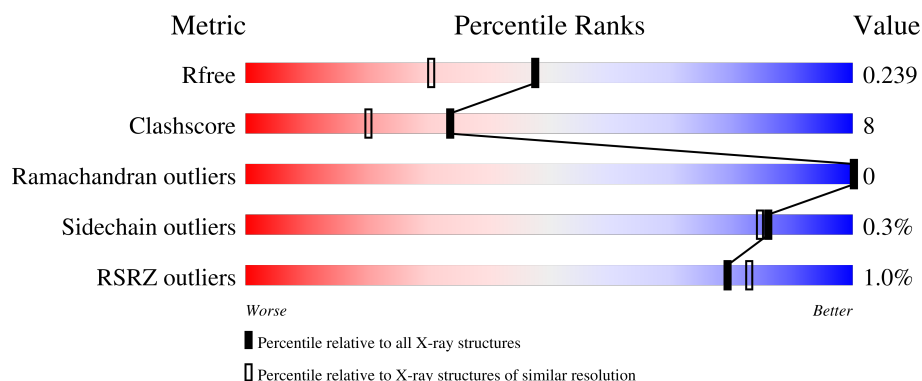
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	322	% 75% 11% 14%
1	B	322	% 76% 10% 14%
1	C	322	69% 17% 14%
1	D	322	70% 16% 14%
1	E	322	% 69% 17% 14%

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Mol	Chain	Length	Quality of chain
1	F	322	
1	G	322	
1	H	322	
1	I	322	
1	J	322	
1	K	322	
1	L	322	

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
3	EDO	D	405	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 26209 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 Two-domain laccase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	278	Total	C	N	O	S	0	5	0
			2171	1353	396	409	13			
1	B	278	Total	C	N	O	S	0	0	0
			2133	1331	389	401	12			
1	C	277	Total	C	N	O	S	0	0	0
			2128	1328	388	400	12			
1	D	278	Total	C	N	O	S	0	0	0
			2133	1331	389	401	12			
1	E	277	Total	C	N	O	S	0	1	0
			2136	1333	389	401	13			
1	F	276	Total	C	N	O	S	0	2	0
			2132	1331	388	400	13			
1	G	279	Total	C	N	O	S	0	1	0
			2151	1341	392	406	12			
1	H	278	Total	C	N	O	S	0	1	0
			2141	1336	390	402	13			
1	I	275	Total	C	N	O	S	0	2	0
			2123	1326	387	397	13			
1	J	278	Total	C	N	O	S	0	0	0
			2133	1331	389	401	12			
1	K	278	Total	C	N	O	S	0	3	0
			2155	1344	394	404	13			
1	L	275	Total	C	N	O	S	0	0	0
			2113	1319	386	396	12			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	21	PHE	LEU	conflict	UNP A0A0M4FJ81
A	170	PHE	ILE	engineered mutation	UNP A0A0M4FJ81
B	21	PHE	LEU	conflict	UNP A0A0M4FJ81
B	170	PHE	ILE	engineered mutation	UNP A0A0M4FJ81
C	21	PHE	LEU	conflict	UNP A0A0M4FJ81

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Chain	Residue	Modelled	Actual	Comment	Reference
C	170	PHE	ILE	engineered mutation	UNP A0A0M4FJ81
D	21	PHE	LEU	conflict	UNP A0A0M4FJ81
D	170	PHE	ILE	engineered mutation	UNP A0A0M4FJ81
E	21	PHE	LEU	conflict	UNP A0A0M4FJ81
E	170	PHE	ILE	engineered mutation	UNP A0A0M4FJ81
F	21	PHE	LEU	conflict	UNP A0A0M4FJ81
F	170	PHE	ILE	engineered mutation	UNP A0A0M4FJ81
G	21	PHE	LEU	conflict	UNP A0A0M4FJ81
G	170	PHE	ILE	engineered mutation	UNP A0A0M4FJ81
H	21	PHE	LEU	conflict	UNP A0A0M4FJ81
H	170	PHE	ILE	engineered mutation	UNP A0A0M4FJ81
I	21	PHE	LEU	conflict	UNP A0A0M4FJ81
I	170	PHE	ILE	engineered mutation	UNP A0A0M4FJ81
J	21	PHE	LEU	conflict	UNP A0A0M4FJ81
J	170	PHE	ILE	engineered mutation	UNP A0A0M4FJ81
K	21	PHE	LEU	conflict	UNP A0A0M4FJ81
K	170	PHE	ILE	engineered mutation	UNP A0A0M4FJ81
L	21	PHE	LEU	conflict	UNP A0A0M4FJ81
L	170	PHE	ILE	engineered mutation	UNP A0A0M4FJ81

- Molecule 2 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	4	Total Cu 4 4	0	0
2	B	4	Total Cu 4 4	0	0
2	C	4	Total Cu 4 4	0	0
2	D	4	Total Cu 4 4	0	0
2	E	4	Total Cu 4 4	0	0
2	F	4	Total Cu 4 4	0	0
2	G	5	Total Cu 5 5	0	0
2	H	3	Total Cu 3 3	0	0
2	I	4	Total Cu 4 4	0	0
2	J	4	Total Cu 4 4	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	K	4	Total	Cu	0	0
			4	4		
2	L	4	Total	Cu	0	0
			4	4		

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



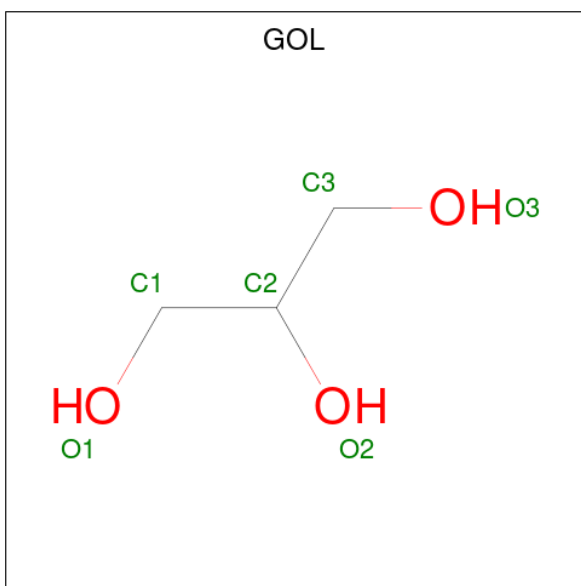
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	G	1	Total	C	O	0	0
			4	2	2		
3	G	1	Total	C	O	0	0
			4	2	2		
3	J	1	Total	C	O	0	0
			4	2	2		
3	L	1	Total	C	O	0	0
			4	2	2		
3	L	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



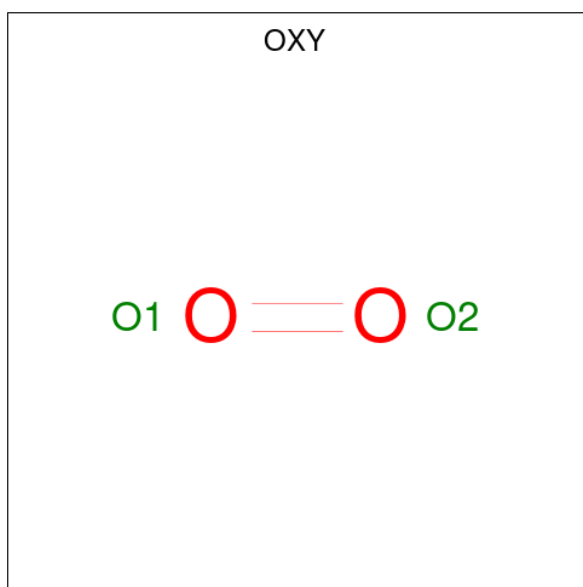
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	G	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is OXYGEN MOLECULE (CCD ID: OXY) (formula: O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	I	1	Total O 2 2	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	40	Total O 40 40	0	0
7	B	43	Total O 43 43	0	0
7	C	43	Total O 43 43	0	0
7	D	45	Total O 45 45	0	0
7	E	41	Total O 41 41	0	0
7	F	23	Total O 23 23	0	0
7	G	51	Total O 51 51	0	0
7	H	51	Total O 51 51	0	0
7	I	34	Total O 34 34	0	0
7	J	38	Total O 38 38	0	0
7	K	32	Total O 32 32	0	0

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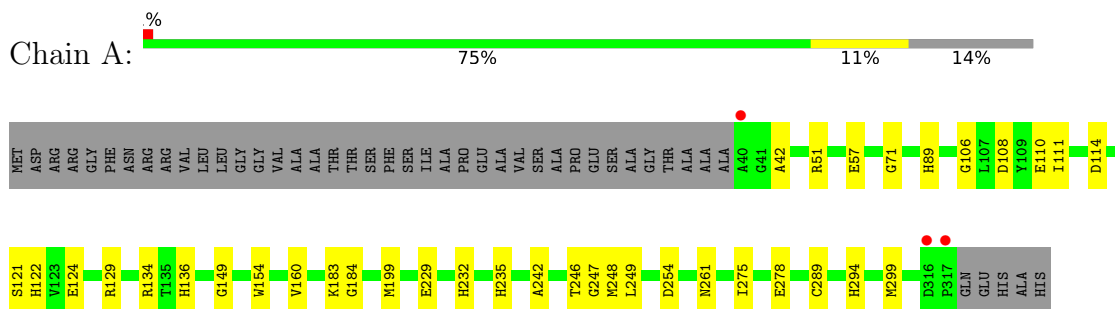
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	L	30	Total	O	0	0
			30	30		

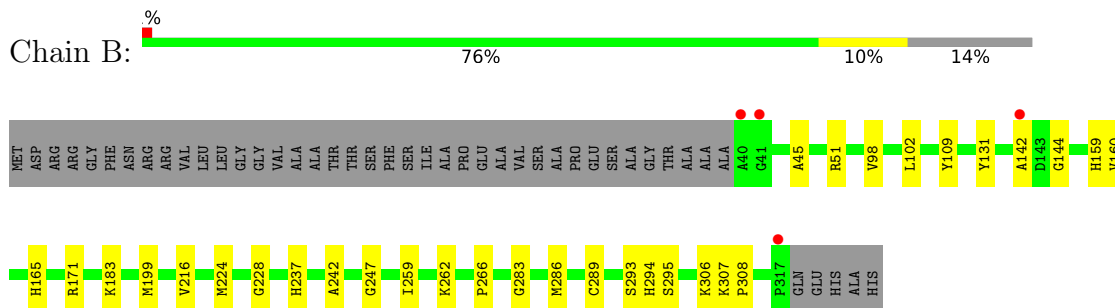
3 Residue-property plots [i](#)

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.

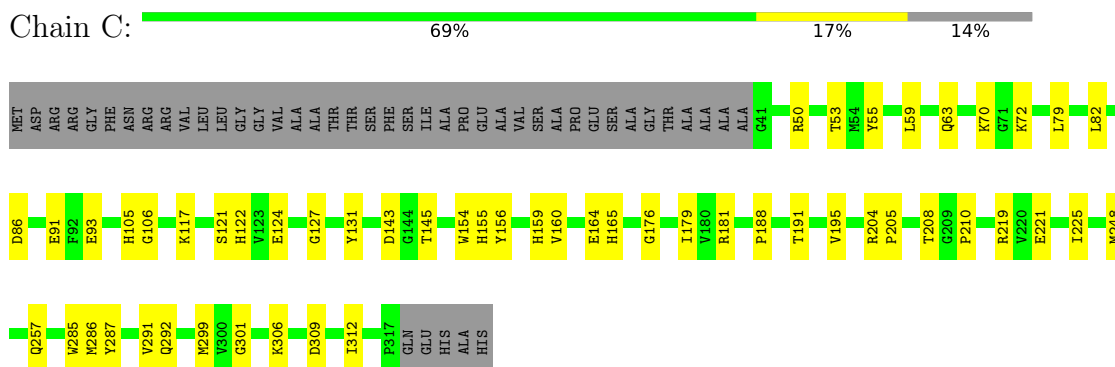
- Molecule 1: Two-domain laccase



- Molecule 1: Two-domain laccase

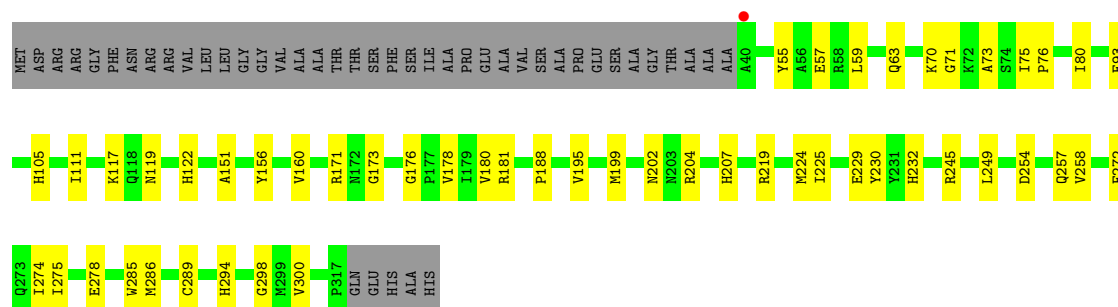


- Molecule 1: Two-domain laccase

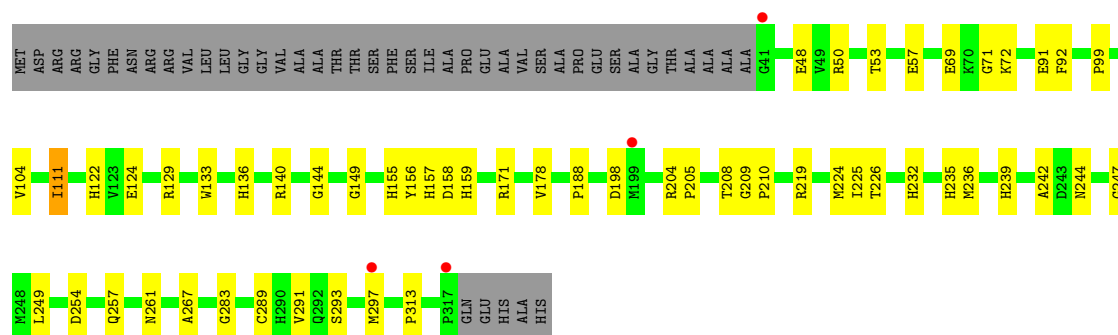


- Molecule 1: Two-domain laccase

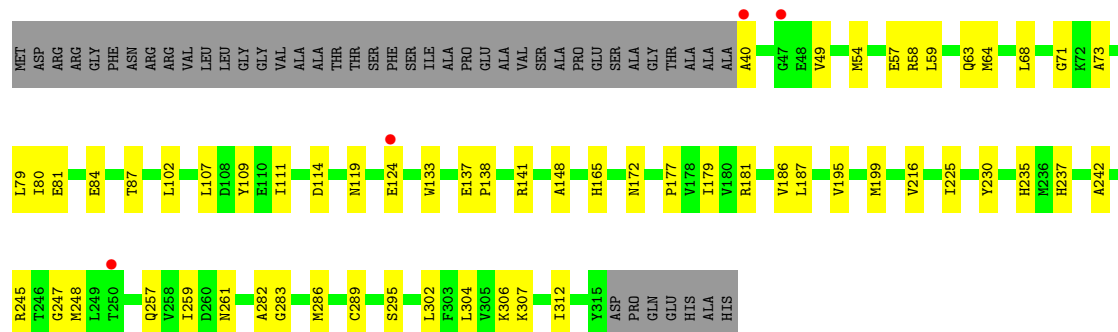




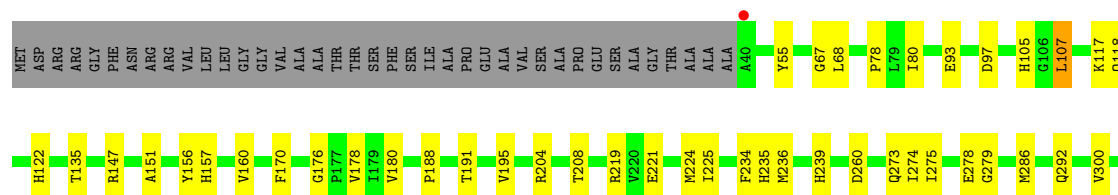
• Molecule 1: Two-domain laccase

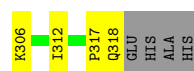


• Molecule 1: Two-domain laccase

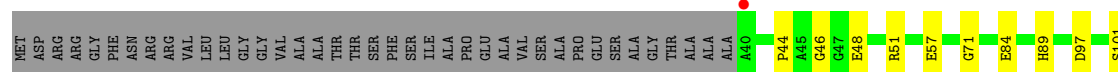
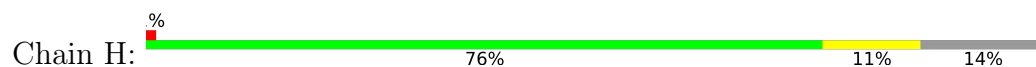


• Molecule 1: Two-domain laccase

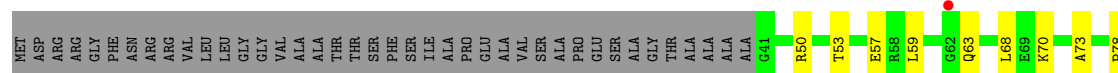




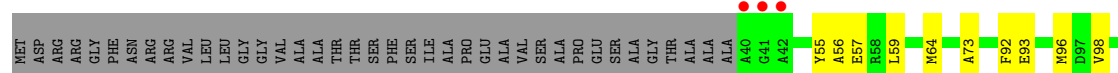
- Molecule 1: Two-domain laccase



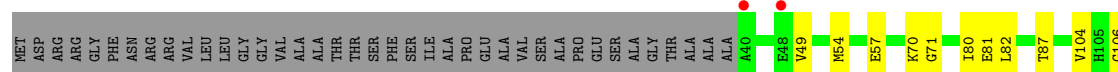
- Molecule 1: Two-domain laccase

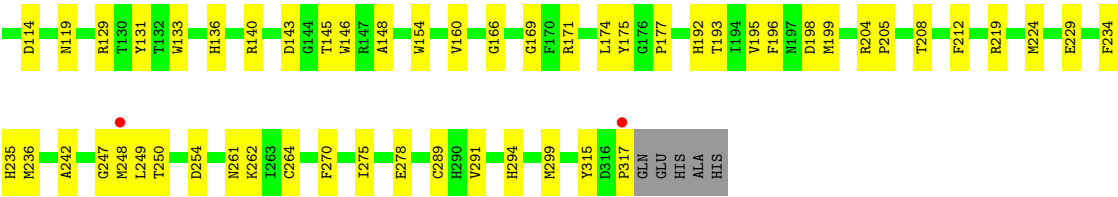


- Molecule 1: Two-domain laccase

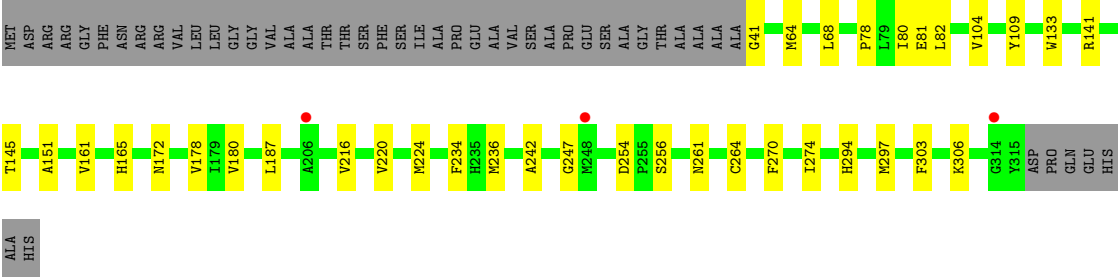


- Molecule 1: Two-domain laccase





● Molecule 1: Two-domain laccase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	77.25Å 94.77Å 116.18Å 90.13° 89.95° 92.56°	Depositor
Resolution (Å)	47.34 – 1.85 47.34 – 1.85	Depositor EDS
% Data completeness (in resolution range)	91.3 (47.34-1.85) 91.4 (47.34-1.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.23 (at 1.86Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.178 , 0.237 0.180 , 0.239	Depositor DCC
R_{free} test set	12841 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	30.3	Xtriage
Anisotropy	0.193	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.036 for h,-k,-l 0.097 for -h,k,-l 0.065 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	26209	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.86% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: OXY, SO4, CU, EDO, GOL

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	0.34	0/2235	0.54	0/3036
1	B	0.44	2/2196 (0.1%)	0.52	1/2983 (0.0%)
1	C	0.33	0/2191	0.50	0/2976
1	D	0.37	0/2196	0.51	0/2983
1	E	0.47	1/2199 (0.0%)	0.54	0/2986
1	F	0.38	2/2197 (0.1%)	0.53	0/2983
1	G	0.42	1/2214 (0.0%)	0.54	0/3008
1	H	0.37	1/2204 (0.0%)	0.52	0/2993
1	I	0.31	0/2188	0.49	0/2971
1	J	0.33	0/2196	0.50	0/2983
1	K	0.35	0/2219	0.52	0/3013
1	L	0.35	0/2175	0.54	0/2953
All	All	0.37	7/26410 (0.0%)	0.52	1/35868 (0.0%)

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	198	ASP	C-O	-7.17	1.18	1.24
1	F	124[A]	GLU	CA-C	5.62	1.59	1.52
1	F	124[B]	GLU	CA-C	5.62	1.59	1.52
1	H	183	LYS	C-O	-5.50	1.17	1.23
1	B	159	HIS	C-O	-5.41	1.16	1.24

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	160	VAL	N-CA-C	5.30	120.38	109.34

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	2171	0	2029	25	0
1	B	2133	0	1999	23	0
1	C	2128	0	1994	40	0
1	D	2133	0	1999	39	0
1	E	2136	0	2002	34	0
1	F	2132	0	2002	49	0
1	G	2151	0	2013	29	0
1	H	2141	0	2007	25	0
1	I	2123	0	1996	30	0
1	J	2133	0	1999	38	0
1	K	2155	0	2016	51	0
1	L	2113	0	1983	31	0
2	A	4	0	0	0	0
2	B	4	0	0	0	0
2	C	4	0	0	0	0
2	D	4	0	0	0	0
2	E	4	0	0	0	0
2	F	4	0	0	0	0
2	G	5	0	0	0	0
2	H	3	0	0	0	0
2	I	4	0	0	0	0
2	J	4	0	0	0	0
2	K	4	0	0	0	0
2	L	4	0	0	0	0
3	C	4	0	6	1	0
3	D	4	0	6	6	0
3	G	8	0	12	0	0
3	J	4	0	6	1	0
3	L	8	0	12	1	0
4	D	5	0	0	0	0
5	G	6	0	8	1	0
6	I	2	0	0	0	0
7	A	40	0	0	0	0
7	B	43	0	0	0	0
7	C	43	0	0	7	0
7	D	45	0	0	1	0
7	E	41	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	F	23	0	0	4	0
7	G	51	0	0	4	0
7	H	51	0	0	1	0
7	I	34	0	0	1	0
7	J	38	0	0	4	0
7	K	32	0	0	2	0
7	L	30	0	0	1	0
All	All	26209	0	24089	384	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 8.

The worst 5 of 384 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:F:282:ALA:HB2	1:F:307:LYS:HE3	1.40	1.02
1:F:282:ALA:CB	1:F:307:LYS:HE3	1.92	0.98
1:J:141:ARG:HH12	1:K:248:MET:HE1	1.41	0.84
1:K:294:HIS:HB3	1:K:299:MET:HE2	1.63	0.80
1:L:220:VAL:HG21	1:L:303:PHE:HZ	1.45	0.80

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	281/322 (87%)	270 (96%)	11 (4%)	0	100	100
1	B	276/322 (86%)	266 (96%)	10 (4%)	0	100	100
1	C	275/322 (85%)	265 (96%)	10 (4%)	0	100	100
1	D	276/322 (86%)	269 (98%)	7 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	276/322 (86%)	266 (96%)	10 (4%)	0	100	100
1	F	276/322 (86%)	264 (96%)	12 (4%)	0	100	100
1	G	278/322 (86%)	272 (98%)	6 (2%)	0	100	100
1	H	277/322 (86%)	267 (96%)	10 (4%)	0	100	100
1	I	275/322 (85%)	263 (96%)	12 (4%)	0	100	100
1	J	276/322 (86%)	265 (96%)	11 (4%)	0	100	100
1	K	279/322 (87%)	271 (97%)	8 (3%)	0	100	100
1	L	273/322 (85%)	259 (95%)	14 (5%)	0	100	100
All	All	3318/3864 (86%)	3197 (96%)	121 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	223/249 (90%)	223 (100%)	0	100	100
1	B	219/249 (88%)	218 (100%)	1 (0%)	81	77
1	C	219/249 (88%)	219 (100%)	0	100	100
1	D	219/249 (88%)	217 (99%)	2 (1%)	70	64
1	E	220/249 (88%)	216 (98%)	4 (2%)	51	40
1	F	219/249 (88%)	218 (100%)	1 (0%)	81	77
1	G	221/249 (89%)	221 (100%)	0	100	100
1	H	220/249 (88%)	220 (100%)	0	100	100
1	I	219/249 (88%)	218 (100%)	1 (0%)	81	77
1	J	219/249 (88%)	219 (100%)	0	100	100
1	K	221/249 (89%)	221 (100%)	0	100	100
1	L	217/249 (87%)	217 (100%)	0	100	100
All	All	2636/2988 (88%)	2627 (100%)	9 (0%)	86	85

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	107	LEU
1	I	257	GLN
1	E	111	ILE
1	E	171	ARG
1	E	224[A]	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 21 such sidechains are listed below:

Mol	Chain	Res	Type
1	H	122	HIS
1	I	83	ASN
1	L	165	HIS
1	I	292	GLN
1	I	63	GLN

5.3.3 RNA [i](#)

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 58 ligands modelled in this entry, 48 are monoatomic - leaving 10 for Mogul analysis.

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
3	EDO	D	405	-	3,3,3	0.40	0	2,2,2	0.25	0
5	GOL	G	408	-	5,5,5	0.34	0	5,5,5	0.44	0
3	EDO	J	405	-	3,3,3	0.47	0	2,2,2	0.30	0
3	EDO	L	406	-	3,3,3	0.33	0	2,2,2	0.90	0
6	OXY	I	402	2	1,1,1	0.16	0	-		
3	EDO	C	405	-	3,3,3	0.44	0	2,2,2	0.39	0
3	EDO	L	405	-	3,3,3	0.41	0	2,2,2	0.41	0
4	SO4	D	406	-	4,4,4	0.29	0	6,6,6	0.14	0
3	EDO	G	406	-	3,3,3	0.42	0	2,2,2	0.37	0
3	EDO	G	407	-	3,3,3	0.41	0	2,2,2	0.52	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
3	EDO	D	405	-	-	0/1/1/1	-
5	GOL	G	408	-	-	2/4/4/4	-
3	EDO	J	405	-	-	1/1/1/1	-
3	EDO	L	406	-	-	1/1/1/1	-
3	EDO	C	405	-	-	0/1/1/1	-
3	EDO	L	405	-	-	0/1/1/1	-
3	EDO	G	406	-	-	0/1/1/1	-
3	EDO	G	407	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	G	408	GOL	O1-C1-C2-O2
5	G	408	GOL	O1-C1-C2-C3
3	J	405	EDO	O1-C1-C2-O2
3	L	406	EDO	O1-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	405	EDO	6	0
5	G	408	GOL	1	0
3	J	405	EDO	1	0
3	C	405	EDO	1	0
3	L	405	EDO	1	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 ⓘ

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	278/322 (86%)	-0.01	3 (1%) 78 81	12, 37, 53, 82	6 (2%)
1	B	278/322 (86%)	0.08	4 (1%) 73 77	24, 38, 60, 104	1 (0%)
1	C	277/322 (86%)	0.04	0 100 100	24, 39, 55, 70	2 (0%)
1	D	278/322 (86%)	0.01	1 (0%) 88 91	24, 37, 52, 66	3 (1%)
1	E	277/322 (86%)	0.06	4 (1%) 73 77	15, 38, 55, 80	4 (1%)
1	F	276/322 (85%)	0.43	4 (1%) 73 77	17, 46, 64, 76	4 (1%)
1	G	279/322 (86%)	-0.04	1 (0%) 88 91	22, 36, 53, 72	2 (0%)
1	H	278/322 (86%)	-0.01	2 (0%) 84 87	16, 36, 51, 86	4 (1%)
1	I	275/322 (85%)	0.17	4 (1%) 72 76	15, 41, 60, 69	5 (1%)
1	J	278/322 (86%)	0.18	3 (1%) 78 81	25, 42, 60, 77	3 (1%)
1	K	278/322 (86%)	0.18	4 (1%) 73 77	14, 41, 59, 86	5 (1%)
1	L	275/322 (85%)	0.29	3 (1%) 78 81	26, 42, 59, 82	1 (0%)
All	All	3327/3864 (86%)	0.11	33 (0%) 79 83	12, 39, 58, 104	40 (1%)

The worst 5 of 33 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	40	ALA	4.0
1	K	40	ALA	3.9
1	A	317	PRO	3.6
1	B	317	PRO	3.5
1	B	40	ALA	3.5

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 oligosaccharides 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. 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(Å ²)	Q<0.9
4	SO4	D	406	5/5	0.86	0.11	77,78,82,84	0
2	CU	F	404	1/1	0.89	0.14	97,97,97,97	1
2	CU	G	403	1/1	0.91	0.22	219,219,219,219	1
3	EDO	G	406	4/4	0.91	0.11	38,40,41,44	0
3	EDO	L	405	4/4	0.91	0.14	52,56,56,56	0
2	CU	D	403	1/1	0.91	0.17	207,207,207,207	1
2	CU	G	405	1/1	0.92	0.11	71,71,71,71	1
3	EDO	G	407	4/4	0.92	0.13	38,40,42,43	0
3	EDO	C	405	4/4	0.94	0.10	53,54,55,55	0
3	EDO	D	405	4/4	0.94	0.14	47,48,49,51	0
3	EDO	L	406	4/4	0.94	0.09	38,40,43,45	0
2	CU	A	403	1/1	0.94	0.14	100,100,100,100	1
5	GOL	G	408	6/6	0.94	0.08	51,52,54,57	0
2	CU	K	403	1/1	0.95	0.07	59,59,59,59	1
2	CU	L	404	1/1	0.95	0.11	116,116,116,116	1
2	CU	I	405	1/1	0.95	0.08	58,58,58,58	1
3	EDO	J	405	4/4	0.95	0.11	32,33,39,45	0
2	CU	E	403	1/1	0.96	0.11	82,82,82,82	1
2	CU	B	403	1/1	0.96	0.10	99,99,99,99	1
6	OXY	I	402	2/2	0.96	0.10	57,57,57,59	0
2	CU	C	404	1/1	0.97	0.06	38,38,38,38	1
2	CU	J	403	1/1	0.98	0.08	65,65,65,65	1
2	CU	L	402	1/1	0.99	0.03	34,34,34,34	0
2	CU	D	404	1/1	0.99	0.03	29,29,29,29	0
2	CU	G	404	1/1	0.99	0.04	29,29,29,29	0
2	CU	E	402	1/1	0.99	0.03	39,39,39,39	0
2	CU	I	403	1/1	0.99	0.02	31,31,31,31	0
2	CU	I	404	1/1	0.99	0.02	42,42,42,42	0
2	CU	A	402	1/1	0.99	0.02	36,36,36,36	0
2	CU	J	401	1/1	0.99	0.02	31,31,31,31	0
2	CU	J	402	1/1	0.99	0.02	36,36,36,36	0
2	CU	F	401	1/1	0.99	0.02	39,39,39,39	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CU	J	404	1/1	0.99	0.02	37,37,37,37	0
2	CU	C	401	1/1	0.99	0.03	39,39,39,39	0
2	CU	A	401	1/1	1.00	0.02	30,30,30,30	0
2	CU	E	401	1/1	1.00	0.02	32,32,32,32	0
2	CU	B	404	1/1	1.00	0.02	36,36,36,36	0
2	CU	A	404	1/1	1.00	0.01	30,30,30,30	0
2	CU	E	404	1/1	1.00	0.02	37,37,37,37	0
2	CU	C	402	1/1	1.00	0.01	29,29,29,29	0
2	CU	K	401	1/1	1.00	0.02	30,30,30,30	0
2	CU	K	402	1/1	1.00	0.02	37,37,37,37	0
2	CU	F	402	1/1	1.00	0.03	38,38,38,38	0
2	CU	K	404	1/1	1.00	0.02	33,33,33,33	0
2	CU	L	401	1/1	1.00	0.01	33,33,33,33	0
2	CU	F	403	1/1	1.00	0.02	38,38,38,38	0
2	CU	L	403	1/1	1.00	0.02	40,40,40,40	0
2	CU	C	403	1/1	1.00	0.02	38,38,38,38	0
2	CU	G	401	1/1	1.00	0.02	31,31,31,31	0
2	CU	G	402	1/1	1.00	0.03	37,37,37,37	1
2	CU	B	401	1/1	1.00	0.02	31,31,31,31	0
2	CU	D	401	1/1	1.00	0.02	30,30,30,30	0
2	CU	D	402	1/1	1.00	0.02	36,36,36,36	1
2	CU	H	401	1/1	1.00	0.02	31,31,31,31	0
2	CU	H	402	1/1	1.00	0.02	43,43,43,43	0
2	CU	H	403	1/1	1.00	0.01	34,34,34,34	0
2	CU	I	401	1/1	1.00	0.02	35,35,35,35	0
2	CU	B	402	1/1	1.00	0.01	40,40,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.