



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 15, 2023 – 11:49 PM EDT

PDB ID : 8EIM
Title : Structure of ALAS with C-terminal truncation from *S. cerevisiae*
Authors : Tran, J.U.; Brown, B.L.
Deposited on : 2022-09-15
Resolution : 2.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

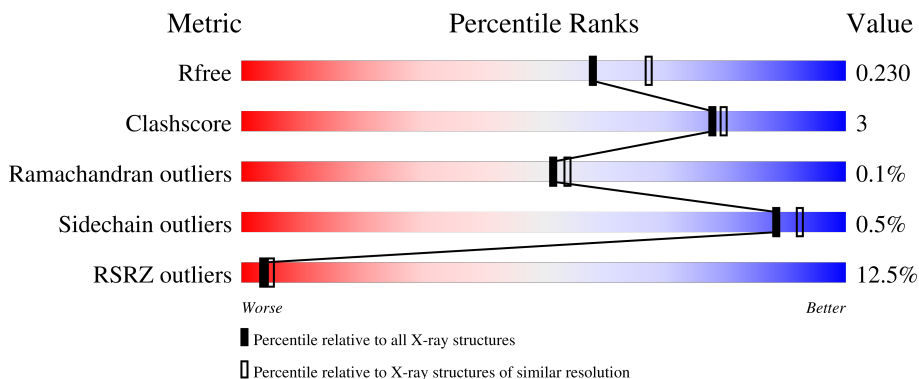
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	477	
1	B	477	

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	PLP	A	601	-	-	X	-

2 Entry composition [i](#)

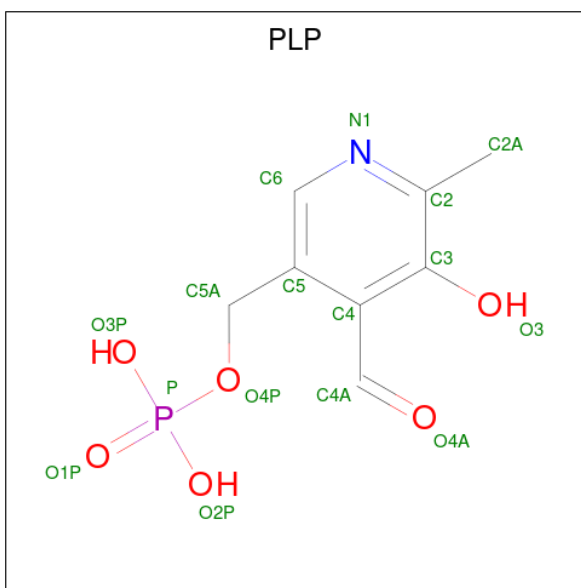
There are 5 unique types of molecules in this entry. The entry contains 7019 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 5-aminolevulinate synthase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	456	Total 3541	C 2241	N 614	O 672	S 14	0	0	0
1	B	417	Total 3226	C 2044	N 559	O 609	S 14	0	0	0

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C₈H₁₀NO₆P) (labeled as "Ligand of Interest" by depositor).



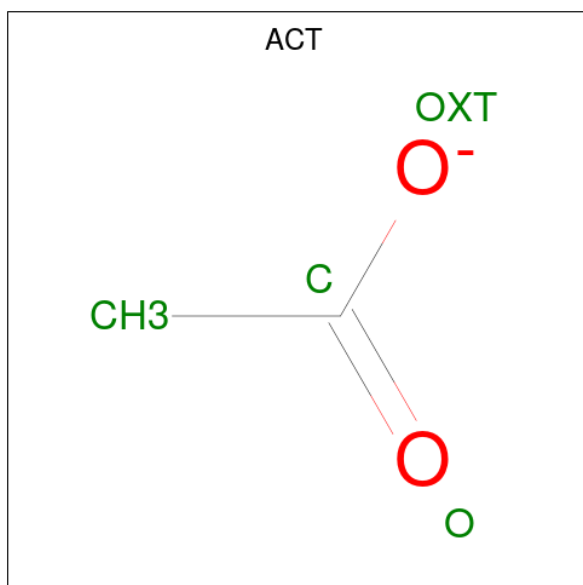
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total 16	C 8	N 1	O 6	P 1	0	0
2	B	1	Total 15	C 8	N 1	O 5	P 1	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2^-$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		

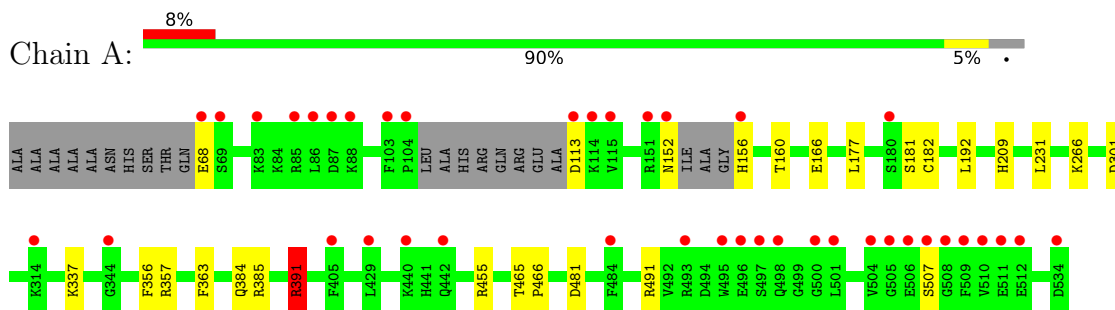
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	87	Total	O	0	0
			87	87		
5	B	102	Total	O	0	0
			102	102		

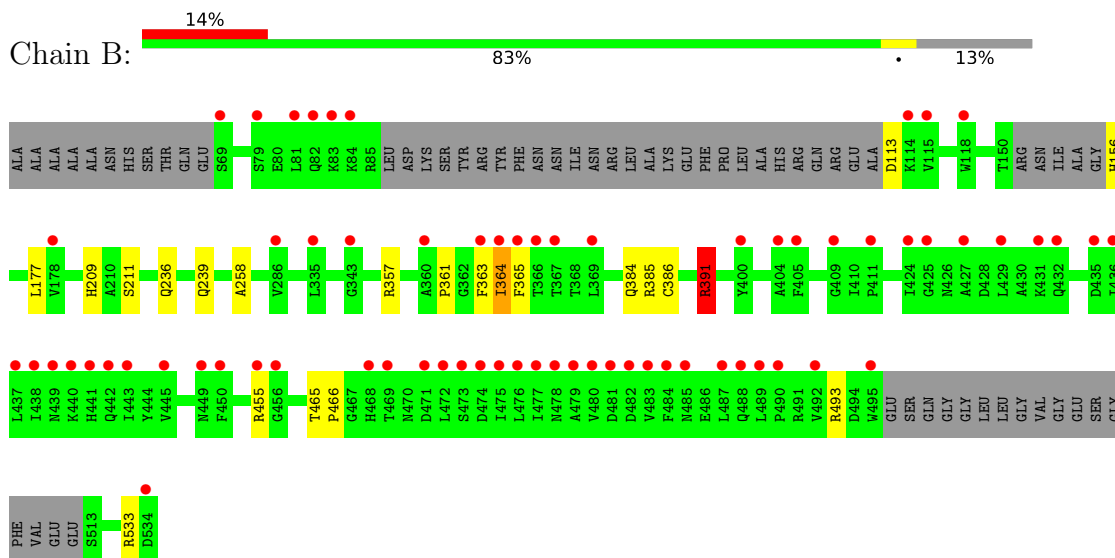
3 Residue-property plots [i](#)

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: 5-aminolevulinate synthase, mitochondrial



- Molecule 1: 5-aminolevulinate synthase, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.28Å 99.55Å 120.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.91 – 2.10 47.91 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.8 (47.91-2.10) 99.9 (47.91-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.25 (at 2.10Å)	Xtrriage
Refinement program	REFMAC 5.8.0352	Depositor
R, R_{free}	0.188 , 0.224 0.194 , 0.230	Depositor DCC
R_{free} test set	2809 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	39.5	Xtrriage
Anisotropy	0.721	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 53.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7019	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% 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: GOL, PLP, ACT

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.40	0/3615	0.68	1/4895 (0.0%)
1	B	0.43	0/3293	0.69	2/4462 (0.0%)
All	All	0.42	0/6908	0.69	3/9357 (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	A	0	2
1	B	0	3
All	All	0	5

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	391	ARG	NE-CZ-NH2	-11.10	114.75	120.30
1	A	391	ARG	NE-CZ-NH2	-10.73	114.94	120.30
1	B	391	ARG	CG-CD-NE	-5.25	100.79	111.80

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	391	ARG	Sidechain
1	A	455	ARG	Sidechain
1	B	391	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	B	455	ARG	Sidechain
1	B	493	ARG	Sidechain

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	3541	0	3508	20	0
1	B	3226	0	3208	19	0
2	A	16	0	8	6	0
2	B	15	0	7	0	0
3	A	18	0	24	2	0
3	B	6	0	8	1	0
4	B	8	0	6	0	0
5	A	87	0	0	2	0
5	B	102	0	0	1	0
All	All	7019	0	6769	39	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 3.

All (39) 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:156:HIS:NE2	1:A:357:ARG:NH1	2.29	0.80
1:A:301:ASP:OD1	3:A:602:GOL:H12	1.82	0.80
1:B:533:ARG:HH11	1:B:533:ARG:HG3	1.51	0.75
1:A:152:ASN:O	1:A:152:ASN:ND2	2.26	0.68
1:A:166:GLU:OE2	1:A:385:ARG:NH1	2.28	0.65
1:A:301:ASP:OD1	3:A:602:GOL:C1	2.45	0.65
1:B:361:PRO:HA	1:B:364:ILE:HG12	1.77	0.65
1:A:231:LEU:HD13	1:A:266:LYS:HG2	1.80	0.63
1:A:337:LYS:NZ	2:A:601:PLP:C4A	2.64	0.61
1:B:385:ARG:HG3	1:B:385:ARG:HH11	1.70	0.56
1:B:361:PRO:HA	1:B:364:ILE:CD1	2.37	0.54
1:A:481:ASP:OD1	1:A:491:ARG:NH2	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:465:THR:HB	1:B:466:PRO:HD2	1.91	0.53
1:B:357:ARG:NH1	1:B:357:ARG:O	2.40	0.52
1:B:364:ILE:HG13	1:B:365:PHE:CD2	2.45	0.52
1:A:337:LYS:HZ1	2:A:601:PLP:C4A	2.21	0.52
1:B:361:PRO:HA	1:B:364:ILE:CG1	2.40	0.51
1:B:357:ARG:CD	1:B:363:PHE:CE2	2.95	0.50
1:A:231:LEU:CD1	1:A:266:LYS:HG2	2.42	0.50
1:B:209:HIS:CD2	1:B:211:SER:H	2.30	0.49
1:A:182:CYS:HB3	2:A:601:PLP:C5A	2.43	0.49
1:A:182:CYS:HB3	2:A:601:PLP:O4P	2.13	0.49
1:A:384:GLN:OE1	1:A:391:ARG:NH2	2.47	0.47
1:B:384:GLN:OE1	1:B:391:ARG:NH2	2.48	0.47
1:A:209:HIS:HB2	2:A:601:PLP:H2A3	1.98	0.46
1:B:209:HIS:HD2	1:B:211:SER:OG	2.00	0.45
5:A:785:HOH:O	1:B:209:HIS:HE1	2.00	0.45
1:A:192:LEU:HD23	1:A:356:PHE:HD1	1.83	0.44
1:B:236:GLN:HA	1:B:239:GLN:NE2	2.32	0.44
1:A:177:LEU:HD11	1:A:363:PHE:CE1	2.52	0.44
1:B:177:LEU:HD11	1:B:363:PHE:CE1	2.53	0.43
1:B:385:ARG:HH11	1:B:385:ARG:CG	2.30	0.43
1:B:258:ALA:HB3	5:B:734:HOH:O	2.18	0.43
1:A:181:SER:HB2	2:A:601:PLP:O2P	2.19	0.43
1:A:68:GLU:O	5:A:701:HOH:O	2.21	0.43
1:A:465:THR:HB	1:A:466:PRO:HD2	2.01	0.42
1:B:357:ARG:HH12	1:B:364:ILE:HG22	1.84	0.41
1:B:386:CYS:SG	3:B:602:GOL:O2	2.80	0.40
1:A:156:HIS:ND1	1:A:160:THR:HG21	2.37	0.40

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	450/477 (94%)	437 (97%)	12 (3%)	1 (0%)	47	49
1	B	409/477 (86%)	399 (98%)	10 (2%)	0	100	100
All	All	859/954 (90%)	836 (97%)	22 (3%)	1 (0%)	51	54

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	507	SER

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	386/398 (97%)	385 (100%)	1 (0%)	92	95
1	B	353/398 (89%)	350 (99%)	3 (1%)	81	86
All	All	739/796 (93%)	735 (100%)	4 (0%)	88	92

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	ASP
1	B	113	ASP
1	B	156	HIS
1	B	364	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	195	GLN
1	A	449	ASN
1	A	498	GLN
1	A	514	ASN
1	B	209	HIS
1	B	232	ASN

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Mol	Chain	Res	Type
1	B	478	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	PLP	B	601	1	15,15,16	0.82	1 (6%)	20,22,23	1.16	2 (10%)
3	GOL	A	604	-	5,5,5	0.39	0	5,5,5	0.61	0
4	ACT	B	603	-	3,3,3	0.72	0	3,3,3	1.04	0
3	GOL	A	602	-	5,5,5	0.16	0	5,5,5	0.37	0
2	PLP	A	601	-	16,16,16	0.58	0	20,23,23	0.85	1 (5%)
3	GOL	A	603	-	5,5,5	0.24	0	5,5,5	0.62	0
4	ACT	B	604	-	3,3,3	1.15	0	3,3,3	0.66	0
3	GOL	B	602	-	5,5,5	0.11	0	5,5,5	0.35	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PLP	B	601	1	-	1/6/6/8	0/1/1/1
3	GOL	A	604	-	-	4/4/4/4	-
3	GOL	A	602	-	-	2/4/4/4	-
2	PLP	A	601	-	-	5/8/8/8	0/1/1/1
3	GOL	A	603	-	-	0/4/4/4	-
3	GOL	B	602	-	-	3/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	PLP	C3-C2	-2.21	1.38	1.40

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	PLP	O3-C3-C2	2.60	123.17	117.49
2	B	601	PLP	O3P-P-O1P	2.47	120.34	110.68
2	B	601	PLP	O3-C3-C4	2.00	123.37	118.10

There are no chirality outliers.

All (15) torsion outliers are listed below:

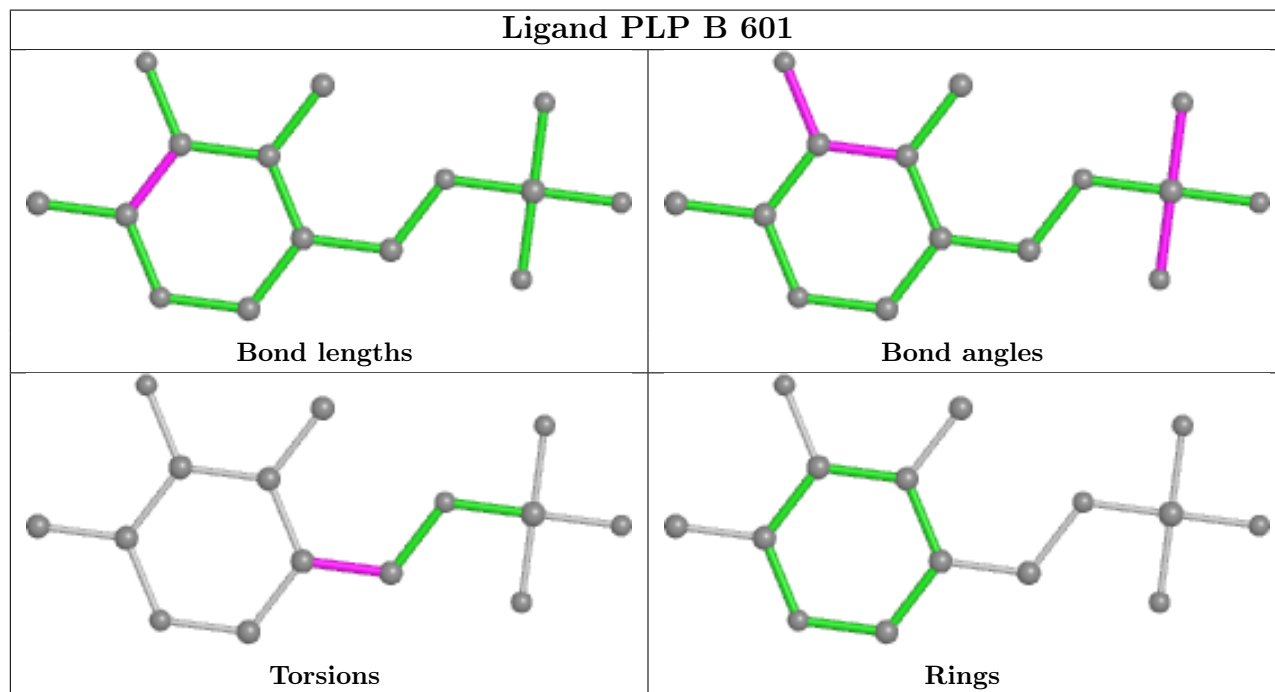
Mol	Chain	Res	Type	Atoms
2	A	601	PLP	C5A-O4P-P-O2P
2	A	601	PLP	C5A-O4P-P-O3P
3	A	604	GOL	C1-C2-C3-O3
3	B	602	GOL	C1-C2-C3-O3
3	B	602	GOL	O2-C2-C3-O3
2	A	601	PLP	C3-C4-C4A-O4A
3	A	602	GOL	C1-C2-C3-O3
3	A	604	GOL	O1-C1-C2-C3
3	A	602	GOL	O2-C2-C3-O3
3	A	604	GOL	O1-C1-C2-O2
3	A	604	GOL	O2-C2-C3-O3
2	A	601	PLP	C5A-O4P-P-O1P
2	B	601	PLP	C6-C5-C5A-O4P
2	A	601	PLP	C5-C4-C4A-O4A
3	B	602	GOL	O1-C1-C2-O2

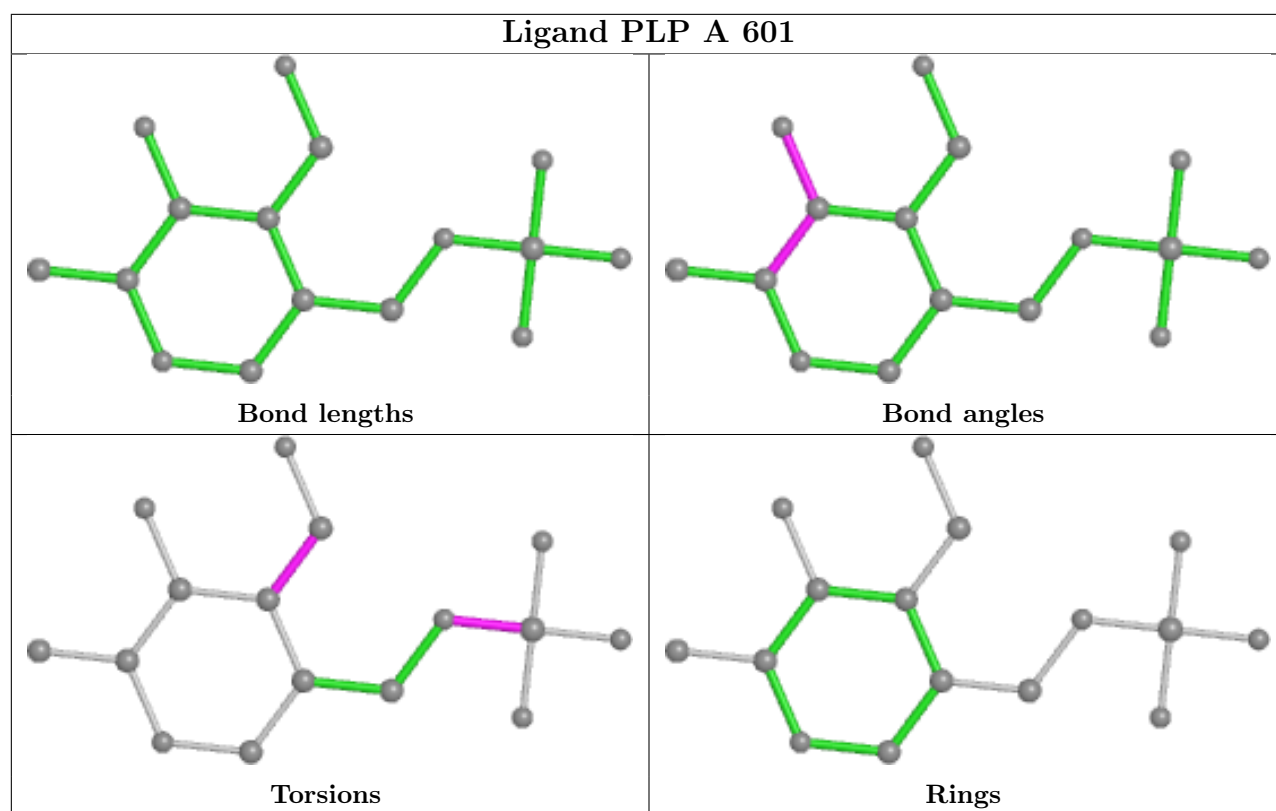
There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	GOL	2	0
2	A	601	PLP	6	0
3	B	602	GOL	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	456/477 (95%)	0.40	40 (8%) 10 12	35, 56, 93, 127	0
1	B	417/477 (87%)	0.77	69 (16%) 1 2	31, 54, 122, 173	0
All	All	873/954 (91%)	0.58	109 (12%) 3 5	31, 56, 110, 173	0

All (109) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	510	VAL	9.6
1	A	507	SER	9.4
1	B	478	ASN	9.3
1	B	436	ILE	8.8
1	A	509	PHE	7.2
1	B	482	ASP	6.4
1	B	438	ILE	6.2
1	B	443	ILE	6.2
1	A	104	PRO	5.8
1	B	439	ASN	5.5
1	A	103	PHE	5.5
1	A	85	ARG	5.1
1	B	488	GLN	5.0
1	B	485	ASN	4.9
1	A	314	LYS	4.8
1	B	440	LYS	4.8
1	A	87	ASP	4.8
1	B	114	LYS	4.6
1	B	481	ASP	4.5
1	B	363	PHE	4.5
1	B	490	PRO	4.4
1	B	442	GLN	4.2
1	B	456	GLY	4.2
1	B	441	HIS	4.2

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Mol	Chain	Res	Type	RSRZ
1	B	479	ALA	4.1
1	B	480	VAL	4.1
1	B	483	VAL	4.1
1	A	534	ASP	4.0
1	B	455	ARG	4.0
1	A	115	VAL	4.0
1	A	505	GLY	4.0
1	B	492	VAL	4.0
1	A	508	GLY	3.9
1	A	114	LYS	3.9
1	B	409	GLY	3.9
1	A	113	ASP	3.8
1	B	469	THR	3.8
1	B	495	TRP	3.7
1	B	400	TYR	3.6
1	B	468	HIS	3.6
1	B	471	ASP	3.6
1	B	366	THR	3.5
1	B	475	ILE	3.4
1	A	86	LEU	3.3
1	A	504	VAL	3.3
1	B	445	VAL	3.3
1	A	442	GLN	3.2
1	A	506	GLU	3.2
1	A	495	TRP	3.2
1	B	364	ILE	3.1
1	B	435	ASP	3.1
1	B	424	ILE	3.1
1	A	484	PHE	3.1
1	A	429	LEU	3.0
1	B	84	LYS	3.0
1	A	498	GLN	3.0
1	B	487	LEU	3.0
1	A	88	LYS	2.9
1	A	151	ARG	2.9
1	B	79	SER	2.9
1	B	474	ASP	2.8
1	B	365	PHE	2.8
1	B	450	PHE	2.8
1	A	69	SER	2.8
1	B	449	ASN	2.7
1	B	432	GLN	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	411	PRO	2.7
1	B	425	GLY	2.7
1	B	360	ALA	2.7
1	B	367	THR	2.7
1	B	115	VAL	2.6
1	B	427	ALA	2.6
1	B	477	ILE	2.6
1	A	440	LYS	2.6
1	B	437	LEU	2.6
1	B	534	ASP	2.5
1	B	404	ALA	2.5
1	B	405	PHE	2.5
1	B	472	LEU	2.5
1	A	405	PHE	2.5
1	B	489	LEU	2.4
1	A	511	GLU	2.4
1	B	484	PHE	2.4
1	A	496	GLU	2.4
1	A	493	ARG	2.4
1	B	343	GLY	2.4
1	A	180	SER	2.3
1	B	83	LYS	2.3
1	A	152	ASN	2.3
1	A	83	LYS	2.3
1	A	497	SER	2.3
1	B	429	LEU	2.3
1	A	156	HIS	2.3
1	A	500	GLY	2.3
1	B	82	GLN	2.2
1	A	512	GLU	2.2
1	B	69	SER	2.2
1	B	431	LYS	2.2
1	A	501	LEU	2.1
1	B	118	TRP	2.1
1	B	369	LEU	2.1
1	B	473	SER	2.1
1	A	68	GLU	2.1
1	B	286	VAL	2.1
1	A	344	GLY	2.1
1	B	81	LEU	2.1
1	B	178	VAL	2.1
1	B	476	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	335	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

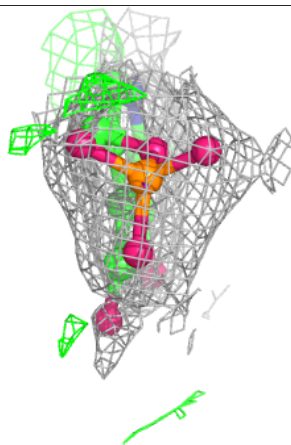
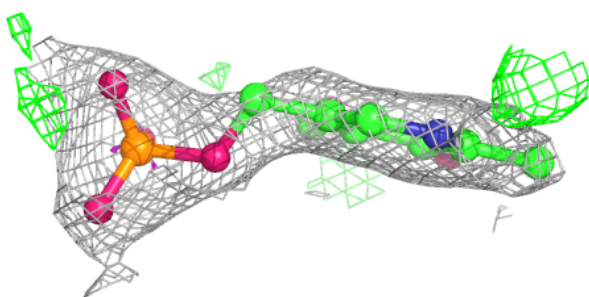
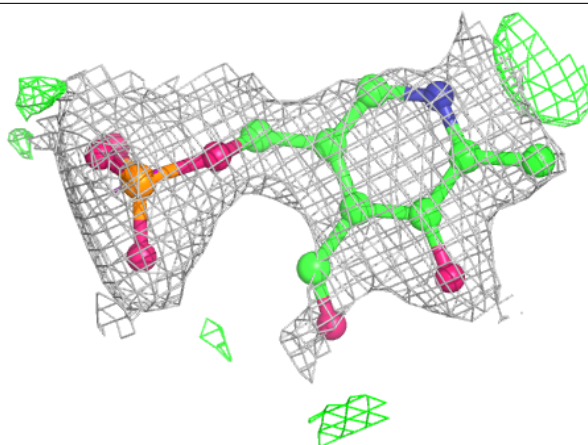
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GOL	A	603	6/6	0.78	0.25	62,71,77,79	0
3	GOL	A	604	6/6	0.84	0.25	50,56,74,77	0
4	ACT	B	603	4/4	0.84	0.15	68,70,72,78	0
4	ACT	B	604	4/4	0.84	0.16	70,72,73,78	0
3	GOL	B	602	6/6	0.90	0.19	47,60,67,74	0
3	GOL	A	602	6/6	0.91	0.17	61,66,71,72	0
2	PLP	A	601	16/16	0.94	0.16	44,54,65,68	16
2	PLP	B	601	15/16	0.97	0.14	33,40,51,51	0

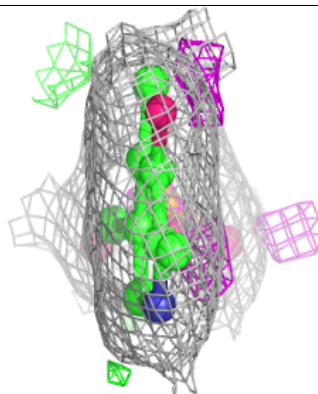
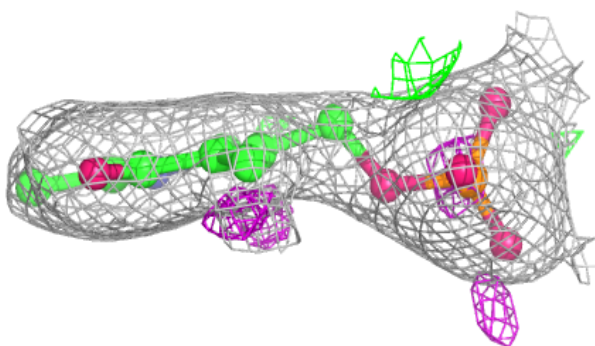
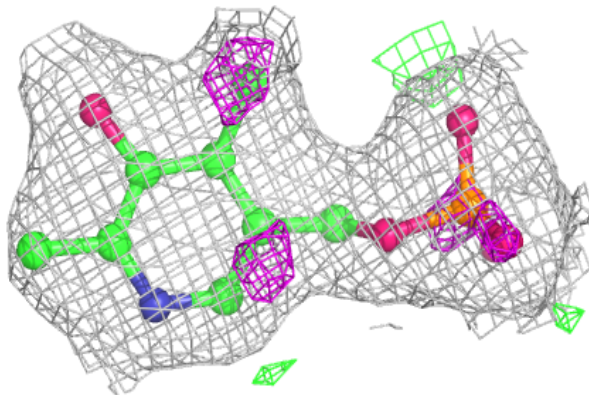
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around PLP A 601:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around PLP B 601:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.