



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:43 PM UTC

PDB ID : 4XJM / pdb_00004xjm
Title : Crystal structure of 7,8-diaminopelargonic acid synthase (BioA) from Mycobacterium tuberculosis, complexed with a HTS lead compound
Authors : Finzel, B.C.; Dai, R.
Deposited on : 2015-01-08
Resolution : 1.60 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

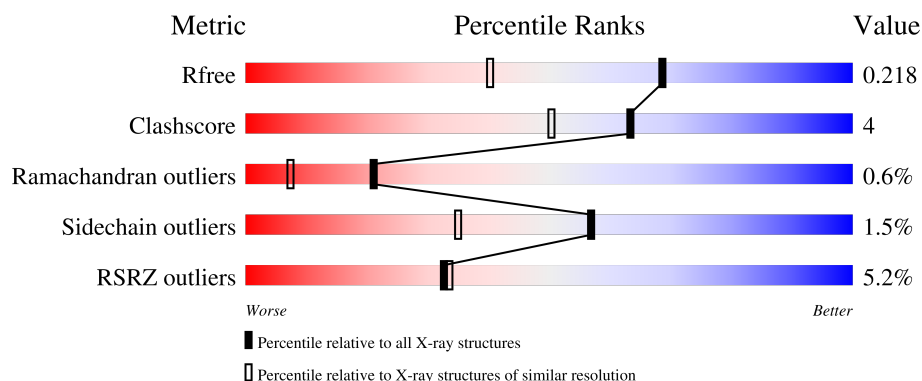
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

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	457	
1	B	457	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6871 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 Adenosylmethionine-8-amino-7-oxononanoate aminotransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	429	Total	C	N	O	S	0	8	0
			3253	2071	572	588	22			
1	B	429	Total	C	N	O	S	0	6	0
			3232	2061	566	584	21			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP P9WQ81
A	-18	GLY	-	expression tag	UNP P9WQ81
A	-17	SER	-	expression tag	UNP P9WQ81
A	-16	SER	-	expression tag	UNP P9WQ81
A	-15	HIS	-	expression tag	UNP P9WQ81
A	-14	HIS	-	expression tag	UNP P9WQ81
A	-13	HIS	-	expression tag	UNP P9WQ81
A	-12	HIS	-	expression tag	UNP P9WQ81
A	-11	HIS	-	expression tag	UNP P9WQ81
A	-10	HIS	-	expression tag	UNP P9WQ81
A	-9	SER	-	expression tag	UNP P9WQ81
A	-8	SER	-	expression tag	UNP P9WQ81
A	-7	GLY	-	expression tag	UNP P9WQ81
A	-6	LEU	-	expression tag	UNP P9WQ81
A	-5	VAL	-	expression tag	UNP P9WQ81
A	-4	PRO	-	expression tag	UNP P9WQ81
A	-3	ARG	-	expression tag	UNP P9WQ81
A	-2	GLY	-	expression tag	UNP P9WQ81
A	-1	SER	-	expression tag	UNP P9WQ81
A	0	HIS	-	expression tag	UNP P9WQ81
B	-19	MET	-	initiating methionine	UNP P9WQ81
B	-18	GLY	-	expression tag	UNP P9WQ81
B	-17	SER	-	expression tag	UNP P9WQ81
B	-16	SER	-	expression tag	UNP P9WQ81

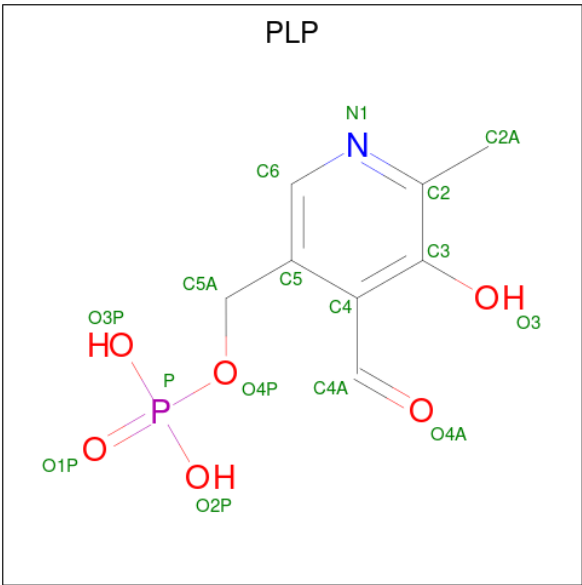
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-15	HIS	-	expression tag	UNP P9WQ81
B	-14	HIS	-	expression tag	UNP P9WQ81
B	-13	HIS	-	expression tag	UNP P9WQ81
B	-12	HIS	-	expression tag	UNP P9WQ81
B	-11	HIS	-	expression tag	UNP P9WQ81
B	-10	HIS	-	expression tag	UNP P9WQ81
B	-9	SER	-	expression tag	UNP P9WQ81
B	-8	SER	-	expression tag	UNP P9WQ81
B	-7	GLY	-	expression tag	UNP P9WQ81
B	-6	LEU	-	expression tag	UNP P9WQ81
B	-5	VAL	-	expression tag	UNP P9WQ81
B	-4	PRO	-	expression tag	UNP P9WQ81
B	-3	ARG	-	expression tag	UNP P9WQ81
B	-2	GLY	-	expression tag	UNP P9WQ81
B	-1	SER	-	expression tag	UNP P9WQ81
B	0	HIS	-	expression tag	UNP P9WQ81

- 41E
-
- Chemical structure 41E is a complex molecule. It features a central piperidine ring. The nitrogen atom of the piperidine ring is bonded to a carbonyl group (C=O), which is further connected to a thiazole ring. The thiazole ring has a methyl group (CH₃) and a carboxylate group (COO⁻). The piperidine ring is also substituted with a carboxylate group (COO⁻) and a side chain containing a carboxylate group (COO⁻) and a methyl group (CH₃). The side chain is further substituted with a carboxylate group (COO⁻) and a methyl group (CH₃). The piperidine ring is also substituted with a carboxylate group (COO⁻) and a methyl group (CH₃). The piperidine ring is also substituted with a carboxylate group (COO⁻) and a methyl group (CH₃).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 29	C 22	N 2	O 4	S 1	0	0
2	B	1	Total 29	C 22	N 2	O 4	S 1	0	0

- 
- WORLD WIDE
PDB
PROTEIN DATA BANK



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

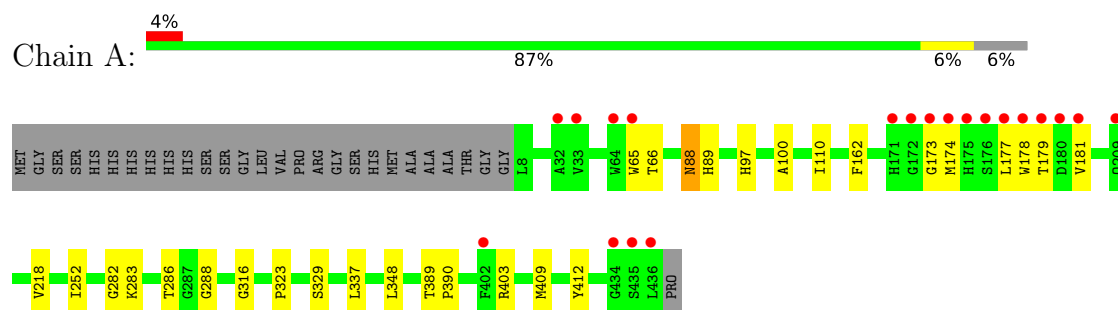
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	165	Total	O	0	0
			165	165		
4	B	133	Total	O	0	0
			133	133		

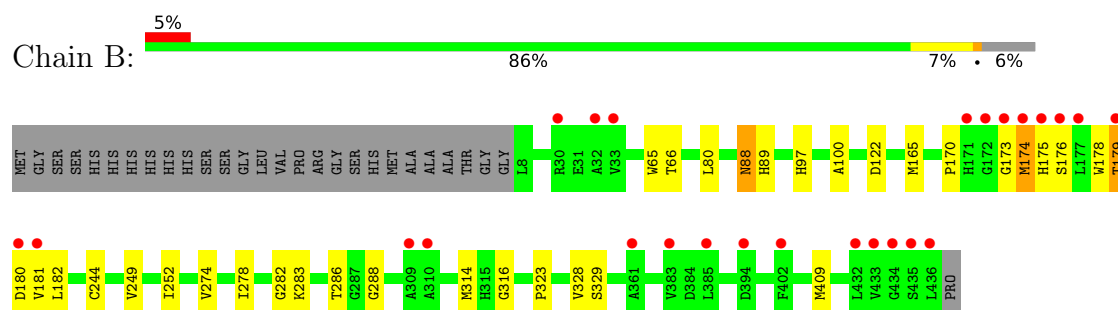
3 Residue-property plots [i](#)

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: Adenosylmethionine-8-amino-7-oxononanoate aminotransferase



- Molecule 1: Adenosylmethionine-8-amino-7-oxononanoate aminotransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.06Å 66.05Å 203.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	101.94 – 1.60 101.94 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (101.94-1.60) 99.9 (101.94-1.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.46 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.203 , 0.221 0.201 , 0.218	Depositor DCC
R_{free} test set	5669 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	17.5	Xtriage
Anisotropy	0.074	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 27.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.028 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6871	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.54% 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: PLP, 41E

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	0.67	0/3348	0.74	1/4576 (0.0%)
1	B	0.67	0/3330	0.74	0/4553
All	All	0.67	0/6678	0.74	1/9129 (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	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	162	PHE	N-CA-C	5.64	117.43	111.28

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	403[A]	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	3253	0	3242	20	0
1	B	3232	0	3230	24	0
2	A	29	0	26	8	0
2	B	29	0	26	4	0
3	A	15	0	6	4	0
3	B	15	0	6	4	0
4	A	165	0	0	0	0
4	B	133	0	0	0	0
All	All	6871	0	6536	50	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 4.

All (50) 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:283:LYS:HZ1	3:A:502:PLP:C4A	1.44	1.26
1:B:283:LYS:HZ1	3:B:502:PLP:C4A	1.56	1.14
1:B:283:LYS:HZ2	3:B:502:PLP:C4A	1.60	1.03
1:A:283:LYS:HZ2	3:A:502:PLP:C4A	1.58	0.98
2:A:501:41E:OAD	2:A:501:41E:H20	1.76	0.84
1:A:389:THR:HB	1:A:390:PRO:HD3	1.69	0.73
1:B:65:TRP:HB2	1:B:283:LYS:HD3	1.74	0.69
1:A:65:TRP:HB2	1:A:283:LYS:HD3	1.75	0.68
1:B:178:TRP:HE3	1:B:181:VAL:HG21	1.67	0.58
2:A:501:41E:CAJ	2:A:501:41E:H7	2.32	0.57
1:A:286:THR:HG22	1:A:329[A]:SER:OG	2.06	0.56
1:B:66:THR:HB	1:B:288:GLY:HA2	1.88	0.56
1:B:179:THR:O	1:B:180:ASP:HB2	2.06	0.54
2:B:501:41E:H7	2:B:501:41E:CAJ	2.37	0.54
1:B:286:THR:HG22	1:B:329[A]:SER:OG	2.08	0.53
1:B:178:TRP:CE3	1:B:181:VAL:HG21	2.43	0.53
2:A:501:41E:H7	2:A:501:41E:H5	1.92	0.52
1:A:66:THR:HB	1:A:288:GLY:HA2	1.91	0.51
1:A:174:MET:HB3	1:A:177:LEU:HD13	1.92	0.51
1:A:178:TRP:HE1	2:A:501:41E:H23	1.75	0.51
1:A:88:ASN:HD22	1:A:89:HIS:H	1.58	0.49
1:B:88:ASN:HD22	1:B:89:HIS:H	1.60	0.49
1:A:348:LEU:HD11	1:A:412:TYR:HA	1.97	0.47
1:A:88:ASN:ND2	1:A:89:HIS:H	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:HIS:HA	1:B:323:PRO:HD2	1.97	0.46
1:A:89:HIS:HA	1:A:323:PRO:HD2	1.97	0.46
1:B:170:PRO:HA	1:B:175:HIS:HD2	1.82	0.45
1:B:173:GLY:HA2	2:B:501:41E:CAX	2.47	0.45
1:B:244:CYS:HB3	1:B:249:VAL:O	2.18	0.44
1:B:80:LEU:HD13	1:B:328[B]:VAL:CG2	2.46	0.44
1:A:110:ILE:HG21	1:A:337:LEU:HD11	2.00	0.43
1:B:88:ASN:ND2	1:B:89:HIS:H	2.14	0.43
2:A:501:41E:H17	2:A:501:41E:H13	1.66	0.43
2:A:501:41E:OAD	2:A:501:41E:CAK	2.47	0.43
1:B:88:ASN:HD22	1:B:88:ASN:N	2.16	0.43
1:A:283:LYS:NZ	3:A:502:PLP:C4	2.70	0.42
1:B:283:LYS:HZ2	3:B:502:PLP:C4	2.25	0.42
1:B:97:HIS:CE1	1:B:100:ALA:HB2	2.55	0.42
1:B:165:MET:HE3	2:B:501:41E:H24	2.02	0.41
1:A:283:LYS:HZ2	3:A:502:PLP:C4	2.26	0.41
1:B:283:LYS:NZ	3:B:502:PLP:C4	2.73	0.41
1:A:173:GLY:HA2	2:A:501:41E:CAX	2.51	0.41
1:B:88:ASN:ND2	1:B:89:HIS:N	2.69	0.41
1:A:173:GLY:O	1:A:174:MET:C	2.63	0.41
1:A:218:VAL:HG22	1:A:252:ILE:HD12	2.03	0.41
1:A:97:HIS:CE1	1:A:100:ALA:HB2	2.56	0.40
1:B:252:ILE:HG12	1:B:278:ILE:HB	2.02	0.40
2:A:501:41E:H26	1:B:314:MET:O	2.21	0.40
1:B:174:MET:HB2	2:B:501:41E:CAF	2.51	0.40
1:A:178:TRP:HE3	1:A:181:VAL:HG21	1.87	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	435/457 (95%)	422 (97%)	10 (2%)	3 (1%)	18	6
1	B	433/457 (95%)	420 (97%)	11 (2%)	2 (0%)	24	10
All	All	868/914 (95%)	842 (97%)	21 (2%)	5 (1%)	21	7

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	179	THR
1	A	282	GLY
1	B	282	GLY
1	A	316	GLY
1	B	316	GLY

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	334/346 (96%)	332 (99%)	2 (1%)	78	66
1	B	332/346 (96%)	324 (98%)	8 (2%)	43	19
All	All	666/692 (96%)	656 (98%)	10 (2%)	57	35

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	88	ASN
1	A	409	MET
1	B	88	ASN
1	B	122	ASP
1	B	174	MET
1	B	176	SER
1	B	179	THR
1	B	182	LEU
1	B	274	VAL
1	B	409	MET

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	88	ASN
1	A	209	GLN
1	B	42	HIS
1	B	78	GLN
1	B	88	ASN
1	B	175	HIS
1	B	209	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](#)

4 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	41E	A	501	-	31,31,31	2.65	5 (16%)	41,42,42	1.56	7 (17%)
3	PLP	A	502	1	15,15,16	2.83	3 (20%)	21,22,23	1.44	4 (19%)
2	41E	B	501	-	31,31,31	2.44	5 (16%)	41,42,42	1.38	5 (12%)
3	PLP	B	502	1	15,15,16	2.77	3 (20%)	21,22,23	1.44	3 (14%)

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	41E	A	501	-	-	5/23/33/33	0/3/3/3
3	PLP	A	502	1	-	0/6/6/8	0/1/1/1
2	41E	B	501	-	-	3/23/33/33	0/3/3/3
3	PLP	B	502	1	-	0/6/6/8	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	41E	CAU-CAZ	-9.32	1.34	1.47
2	B	501	41E	CAU-CAZ	-8.93	1.35	1.47
2	A	501	41E	CAW-CBA	-7.89	1.35	1.48
2	B	501	41E	CAW-CBA	-7.55	1.35	1.48
3	A	502	PLP	C3-C2	7.05	1.48	1.41
3	A	502	PLP	C5-C4	6.57	1.47	1.40
3	B	502	PLP	C3-C2	6.55	1.47	1.41
3	B	502	PLP	C5-C4	6.51	1.47	1.40
3	B	502	PLP	C3-C4	4.93	1.49	1.40
2	A	501	41E	CAZ-SAT	-4.72	1.63	1.73
3	A	502	PLP	C3-C4	4.63	1.48	1.40
2	A	501	41E	CBA-SAT	-4.37	1.64	1.73
2	B	501	41E	CBA-SAT	-3.61	1.66	1.73
2	A	501	41E	CAX-NAR	-3.47	1.34	1.41
2	B	501	41E	CAX-NAR	-3.34	1.34	1.41
2	B	501	41E	CAZ-SAT	-2.99	1.67	1.73

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	41E	CAX-NAR-CAV	-4.30	119.91	127.52
2	A	501	41E	CBA-CAW-NBC	3.74	123.68	119.10
2	B	501	41E	CBA-CAW-NBC	3.40	123.26	119.10
2	A	501	41E	CAP-NBC-CAQ	3.29	119.40	112.68
2	B	501	41E	CAU-CAZ-SAT	3.29	124.14	118.85
2	A	501	41E	CBA-SAT-CAZ	3.13	96.19	91.95
2	A	501	41E	CAU-CAZ-SAT	2.87	123.46	118.85
2	B	501	41E	CBA-SAT-CAZ	2.69	95.59	91.95
2	B	501	41E	CAP-NBC-CAQ	2.64	118.07	112.68
3	B	502	PLP	C4A-C4-C5	-2.60	118.26	120.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	PLP	C6-N1-C2	2.54	123.80	119.20
3	B	502	PLP	C6-N1-C2	2.36	123.48	119.20
2	A	501	41E	CAA-OAS-CAY	-2.25	112.67	117.50
3	A	502	PLP	O4P-C5A-C5	2.24	113.55	109.36
3	B	502	PLP	O3-C3-C4	2.23	123.91	118.10
3	A	502	PLP	O3P-P-O2P	2.22	116.13	107.80
2	A	501	41E	CAM-CAL-CAV	-2.15	108.26	113.06
3	A	502	PLP	C4A-C4-C5	-2.09	118.79	120.94
2	B	501	41E	CAX-NAR-CAV	-2.02	123.95	127.52

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	41E	CAK-CAY-OAS-CAA
2	A	501	41E	CAH-CAY-OAS-CAA
2	A	501	41E	CAV-CAL-CAM-CBB
2	B	501	41E	CAV-CAL-CAM-CBB
2	A	501	41E	CAL-CAM-CBB-CAN
2	A	501	41E	CAL-CAM-CBB-CAO
2	B	501	41E	CAM-CAL-CAV-OAD
2	B	501	41E	CAM-CAL-CAV-NAR

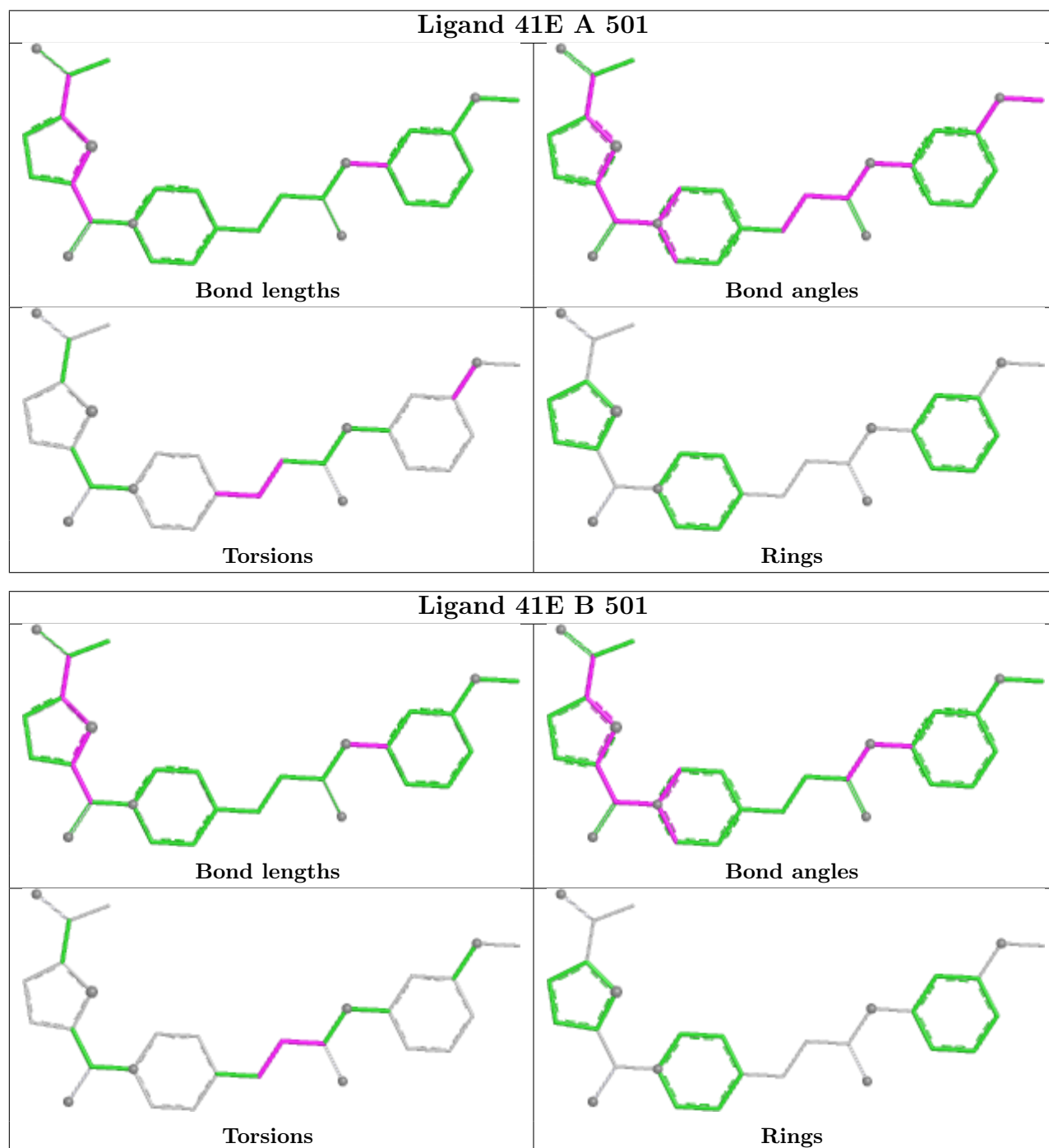
There are no ring outliers.

4 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	41E	8	0
3	A	502	PLP	4	0
2	B	501	41E	4	0
3	B	502	PLP	4	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 ⓘ

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	429/457 (93%)	0.22	20 (4%) 36 38	7, 15, 27, 48	11 (2%)
1	B	429/457 (93%)	0.38	25 (5%) 29 29	7, 17, 32, 57	10 (2%)
All	All	858/914 (93%)	0.30	45 (5%) 33 33	7, 16, 31, 57	21 (2%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	436	LEU	7.2
1	B	175	HIS	7.1
1	A	173	GLY	6.9
1	A	175	HIS	6.4
1	A	436	LEU	6.3
1	B	173	GLY	6.0
1	A	177	LEU	6.0
1	A	174	MET	5.8
1	B	171	HIS	5.5
1	B	177	LEU	5.5
1	B	179	THR	5.3
1	A	32	ALA	5.2
1	B	176	SER	5.0
1	B	172	GLY	5.0
1	B	174	MET	4.8
1	A	179	THR	4.8
1	A	178	TRP	4.7
1	A	176	SER	4.5
1	A	65	TRP	4.0
1	A	180	ASP	3.8
1	A	171	HIS	3.5
1	B	180	ASP	3.5
1	A	33	VAL	3.5
1	B	434	GLY	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	181	VAL	3.3
1	A	181	VAL	3.2
1	A	64	TRP	3.0
1	B	385	LEU	2.9
1	B	310	ALA	2.8
1	B	309	ALA	2.7
1	B	361	ALA	2.7
1	A	402	PHE	2.7
1	A	172	GLY	2.7
1	B	435	SER	2.6
1	B	402	PHE	2.5
1	A	435	SER	2.5
1	B	432	LEU	2.4
1	B	383	VAL	2.4
1	A	434	GLY	2.3
1	B	394	ASP	2.3
1	B	33	VAL	2.2
1	A	209	GLN	2.2
1	B	433	VAL	2.1
1	B	32	ALA	2.1
1	B	30	ARG	2.1

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 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	41E	A	501	29/29	0.85	0.15	20,28,36,37	0
2	41E	B	501	29/29	0.91	0.12	16,23,30,31	0

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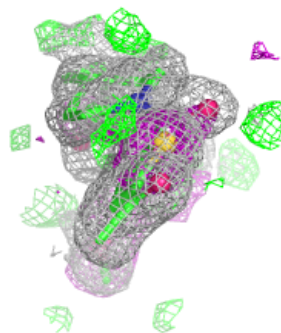
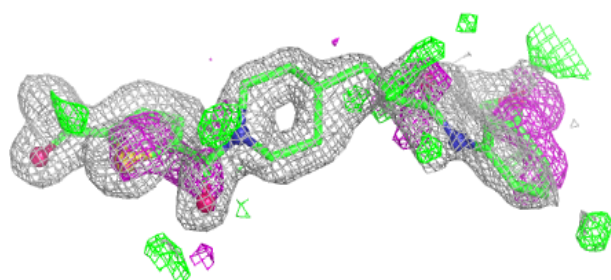
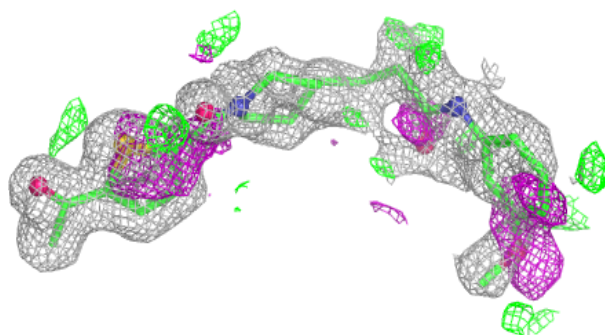
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	PLP	A	502	15/16	0.97	0.05	10,11,11,12	0
3	PLP	B	502	15/16	0.97	0.06	11,11,12,13	0

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.

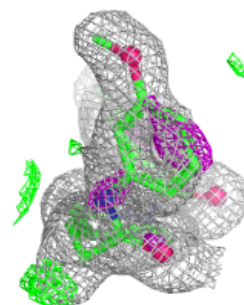
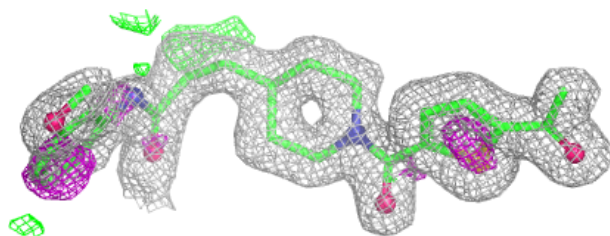
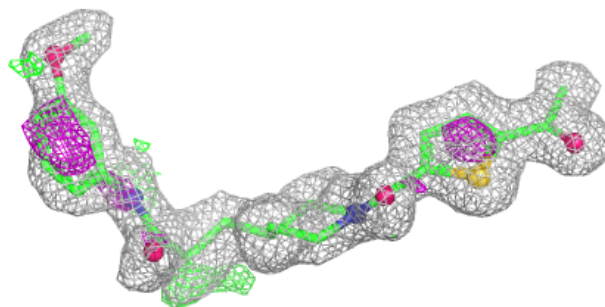
Electron density around 41E A 501:

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 41E B 501:

$2mF_o-DF_c$ (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.