Full wwPDB X-ray Structure Validation Report (i)

Nov 14, 2023 – 09:14 PM JST

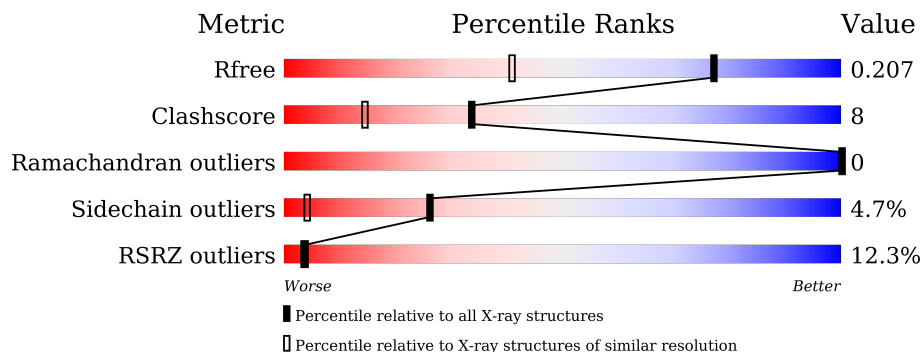|		
PDB ID	:	6AAN
Title	:	Crystal structure of Methanosarcina mazei PylRS(Y306A/Y384F) complexed with mEtZLys
Authors	:	Yanagisawa, T.; Kuratani, M.; Yokoyama, S.
Deposited on	:	2018-07-18
Resolution	:	1.51 Å(reported)

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 1.51 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	Percentile Ranks	Value
Rfree		0.207
Clashscore		8
Ramachandran outliers		0
Sidechain outliers		4.7%
RSRZ outliers		12.3%

Worse ← → Better

▮ Percentile relative to all X-ray structures
▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4009 (1.54-1.50)
Clashscore	141614	4249 (1.54-1.50)
Ramachandran outliers	138981	4148 (1.54-1.50)
Sidechain outliers	138945	4146 (1.54-1.50)
RSRZ outliers	127900	3943 (1.54-1.50)

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 <=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	274	12% / 81% 12% · 5%

2 Entry composition ⓘ

There are 7 unique types of molecules in this entry. The entry contains 2481 atoms, of which 19 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 Pyrrolysine–tRNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	261	Total	C	N	O	S	0	11	0
			2168	1383	369	404	12			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	181	GLY	-	expression tag	UNP Q8PWY1
A	182	SER	-	expression tag	UNP Q8PWY1
A	183	HIS	-	expression tag	UNP Q8PWY1
A	184	MET	-	expression tag	UNP Q8PWY1
A	306	ALA	TYR	engineered mutation	UNP Q8PWY1
A	384	PHE	TYR	engineered mutation	UNP Q8PWY1
A	444	GLY	GLU	see sequence details	UNP Q8PWY1

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Mg	0	0
			3	3		

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).

ATP

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 4 is N6-{[(3-ethynylphenyl)methoxy]carbonyl}-L-lysine (three-letter code: 9TX) (formula: $C_{16}H_{20}N_2O_4$).

9TX
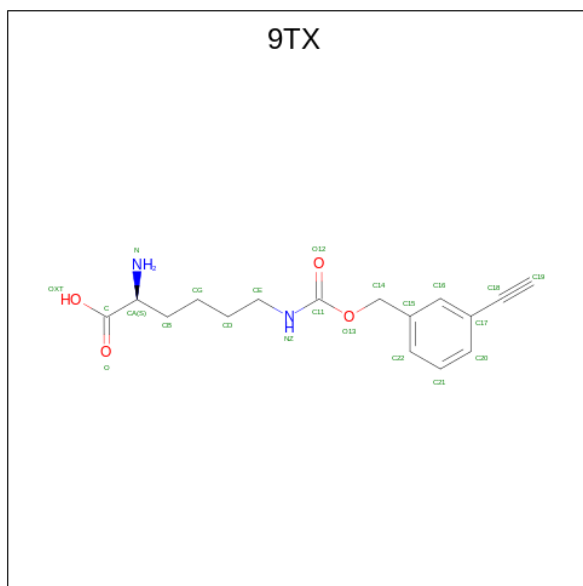
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	H	N	O	0	0
			41	16	19	2	4		

- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total 2	K 2	0	0

• Molecule 6 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).
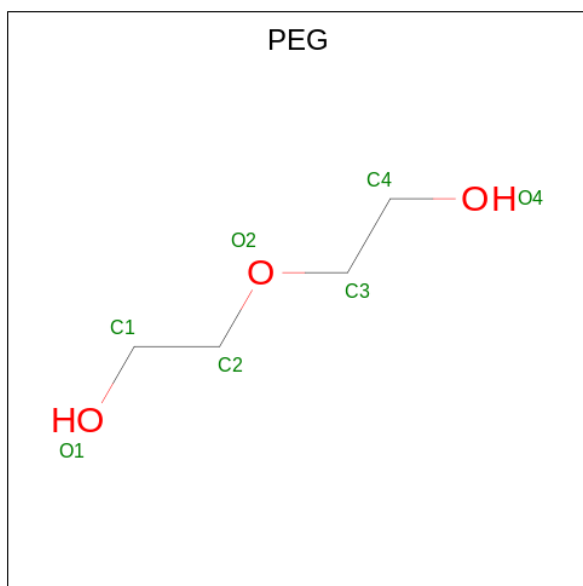
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total 7	C 4	O 3	0	0

• Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	229	Total 229	O 229	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: Pyrrolysine–tRNA ligase
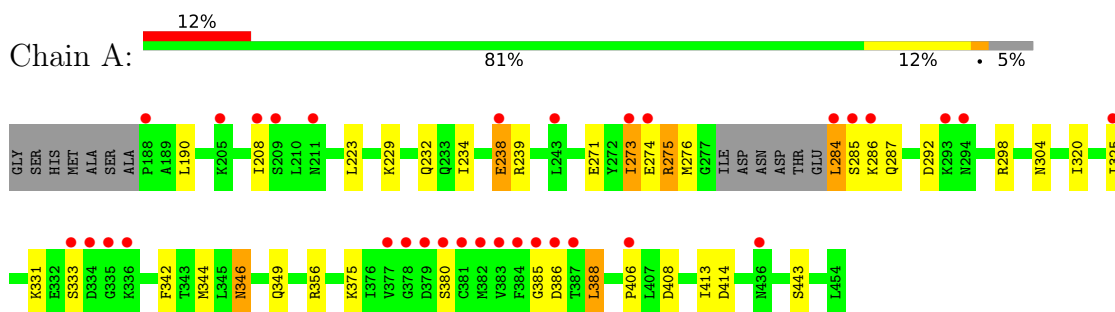
4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	101.48Å　44.07Å　72.47Å 90.00°　118.62°　90.00°	Depositor
Resolution (Å)	35.00　–　1.51 35.00　–　1.51	Depositor EDS
% Data completeness (in resolution range)	96.9 (35.00-1.51) 96.9 (35.00-1.51)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.90 (at 1.51Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.183　,　0.206 0.183　,　0.207	Depositor DCC
R_{free} test set	2000 reflections (4.64%)	wwPDB-VP
Wilson B-factor (Å^2)	23.6	Xtriage
Anisotropy	0.194	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.42 , 59.5	EDS
L-test for twinning[2]	$< |L| > = 0.49, < L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	2481	wwPDB-VP
Average B, all atoms (Å^2)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.21% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: PEG, ATP, 9TX, K, 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	0.48	0/2238	0.69	1/3005 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	298	ARG	NE-CZ-NH2	-6.88	116.86	120.30

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	2168	0	2212	37	0
2	A	3	0	0	0	0
3	A	31	0	12	1	0
4	A	22	19	0	0	0
5	A	2	0	0	0	0
6	A	7	0	10	1	0
7	A	229	0	0	9	3
All	All	2462	19	2234	38	3

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.

All (38) 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:234:ILE:HG23	7:A:605:HOH:O	1.71	0.90
1:A:238:GLU:HB2	7:A:605:HOH:O	1.77	0.85
1:A:273:ILE:HG23	1:A:276:MET:HB2	1.73	0.70
1:A:239:ARG:N	7:A:605:HOH:O	2.24	0.69
1:A:292:ASP:OD1	7:A:601:HOH:O	2.08	0.69
1:A:273:ILE:HG22	1:A:274:GLU:HA	1.72	0.69
1:A:273:ILE:HA	1:A:275:ARG:N	2.12	0.65
1:A:356:ARG:NH2	1:A:388:LEU:HD22	2.14	0.62
1:A:304:ASN:HB2	7:A:701:HOH:O	2.01	0.60
6:A:508:PEG:H12	7:A:759:HOH:O	2.04	0.56
1:A:342:PHE:HE2	1:A:344[B]:MET:CE	2.20	0.54
1:A:208:ILE:HD12	1:A:223:LEU:HD21	1.90	0.52
1:A:273:ILE:CG2	1:A:274:GLU:HA	2.40	0.51
1:A:414:ASP:OD2	7:A:603:HOH:O	2.19	0.51
1:A:344[B]:MET:HE1	3:A:504:ATP:O4'	2.11	0.50
1:A:271:GLU:OE1	1:A:275:ARG:HD2	2.11	0.50
1:A:273:ILE:HA	1:A:275:ARG:H	1.77	0.48
1:A:406:PRO:HD2	7:A:708:HOH:O	2.12	0.48
1:A:284:LEU:O	1:A:285:SER:OG	2.24	0.48
1:A:275:ARG:NH1	7:A:602:HOH:O	2.10	0.46
1:A:346:ASN:HD22	1:A:346:ASN:C	2.18	0.46
1:A:229[B]:LYS:HD3	1:A:232:GLN:OE1	2.16	0.45
1:A:356:ARG:CZ	1:A:388:LEU:HD22	2.45	0.45
1:A:273:ILE:CB	1:A:274:GLU:HA	2.45	0.45
1:A:271:GLU:OE1	1:A:275:ARG:CD	2.65	0.45
1:A:287:GLN:HA	1:A:331:LYS:HB2	1.99	0.45
1:A:284:LEU:HD12	1:A:284:LEU:HA	1.78	0.44
1:A:342:PHE:HE2	1:A:344[B]:MET:HE3	1.82	0.44
1:A:344[B]:MET:HE3	1:A:344[B]:MET:HB2	1.35	0.43
1:A:285:SER:OG	1:A:286:LYS:N	2.51	0.43
1:A:325:ILE:HA	1:A:344[A]:MET:O	2.18	0.42
1:A:320:ILE:O	1:A:349:GLN:HA	2.20	0.42
1:A:342:PHE:HE2	1:A:344[B]:MET:HE1	1.85	0.42
1:A:229[B]:LYS:HD3	1:A:229[B]:LYS:HA	1.82	0.41
1:A:271:GLU:C	1:A:273:ILE:N	2.73	0.41
1:A:273:ILE:CA	1:A:275:ARG:H	2.34	0.41
1:A:380:SER:HA	1:A:385:GLY:O	2.20	0.41
1:A:408:ASP:HB3	1:A:413:ILE:O	2.20	0.41

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:748:HOH:O	7:A:758:HOH:O[2_556]	2.07	0.13
7:A:653:HOH:O	7:A:813:HOH:O[4_455]	2.08	0.12
7:A:807:HOH:O	7:A:820:HOH:O[1_556]	2.17	0.03

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	270/274 (98%)	260 (96%)	10 (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	246/243 (101%)	235 (96%)	11 (4%)	27 5

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	190	LEU
1	A	238	GLU
1	A	273	ILE
1	A	275	ARG

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Mol	Chain	Res	Type
1	A	284	LEU
1	A	333	SER
1	A	346	ASN
1	A	375	LYS
1	A	386	ASP
1	A	388	LEU
1	A	443	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	346	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 5 are monoatomic - leaving 3 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	ATP	A	504	2	26,33,33	0.97	1 (3%)	31,52,52	1.49	3 (9%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	#|Z| > 2	Counts	RMSZ	#|Z| > 2
6	PEG	A	508	-	6,6,6	0.30	0	5,5,5	0.80	0
4	9TX	A	505	-	21,22,22	1.77	3 (14%)	26,27,27	1.91	5 (19%)

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	ATP	A	504	2	-	4/18/38/38	0/3/3/3
6	PEG	A	508	-	-	3/4/4/4	-
4	9TX	A	505	-	-	8/19/19/19	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	505	9TX	O-C	4.96	1.37	1.22
4	A	505	9TX	C11-NZ	4.90	1.45	1.34
3	A	504	ATP	C5-C4	2.55	1.47	1.40
4	A	505	9TX	O13-C11	2.49	1.39	1.35

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	505	9TX	O13-C11-NZ	6.74	121.10	110.61
4	A	505	9TX	O12-C11-NZ	-4.37	118.26	124.96
3	A	504	ATP	N3-C2-N1	-4.19	122.12	128.68
3	A	504	ATP	PB-O3B-PG	-3.40	121.16	132.83
3	A	504	ATP	C2-N1-C6	2.86	123.65	118.75
4	A	505	9TX	OXT-C-O	-2.28	118.90	124.09
4	A	505	9TX	C16-C17-C18	2.07	122.74	120.09
4	A	505	9TX	CB-CA-C	2.06	115.21	110.30

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	504	ATP	C5'-O5'-PA-O3A
4	A	505	9TX	N-CA-CB-CG
4	A	505	9TX	NZ-C11-O13-C14

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Mol	Chain	Res	Type	Atoms
4	A	505	9TX	O12-C11-O13-C14
6	A	508	PEG	O1-C1-C2-O2
4	A	505	9TX	CG-CD-CE-NZ
4	A	505	9TX	C-CA-CB-CG
3	A	504	ATP	C5'-O5'-PA-O2A
6	A	508	PEG	C1-C2-O2-C3
4	A	505	9TX	OXT-C-CA-CB
6	A	508	PEG	C4-C3-O2-C2
4	A	505	9TX	O-C-CA-CB
3	A	504	ATP	PA-O3A-PB-O1B
4	A	505	9TX	OXT-C-CA-N
3	A	504	ATP	PA-O3A-PB-O2B

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	504	ATP	1	0
6	A	508	PEG	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 then 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.

Ligand ATP A 504

Bond lengths

Bond angles

Torsions

Rings

Ligand 9TX A 505

Bond lengths

Bond angles

Torsions

Rings
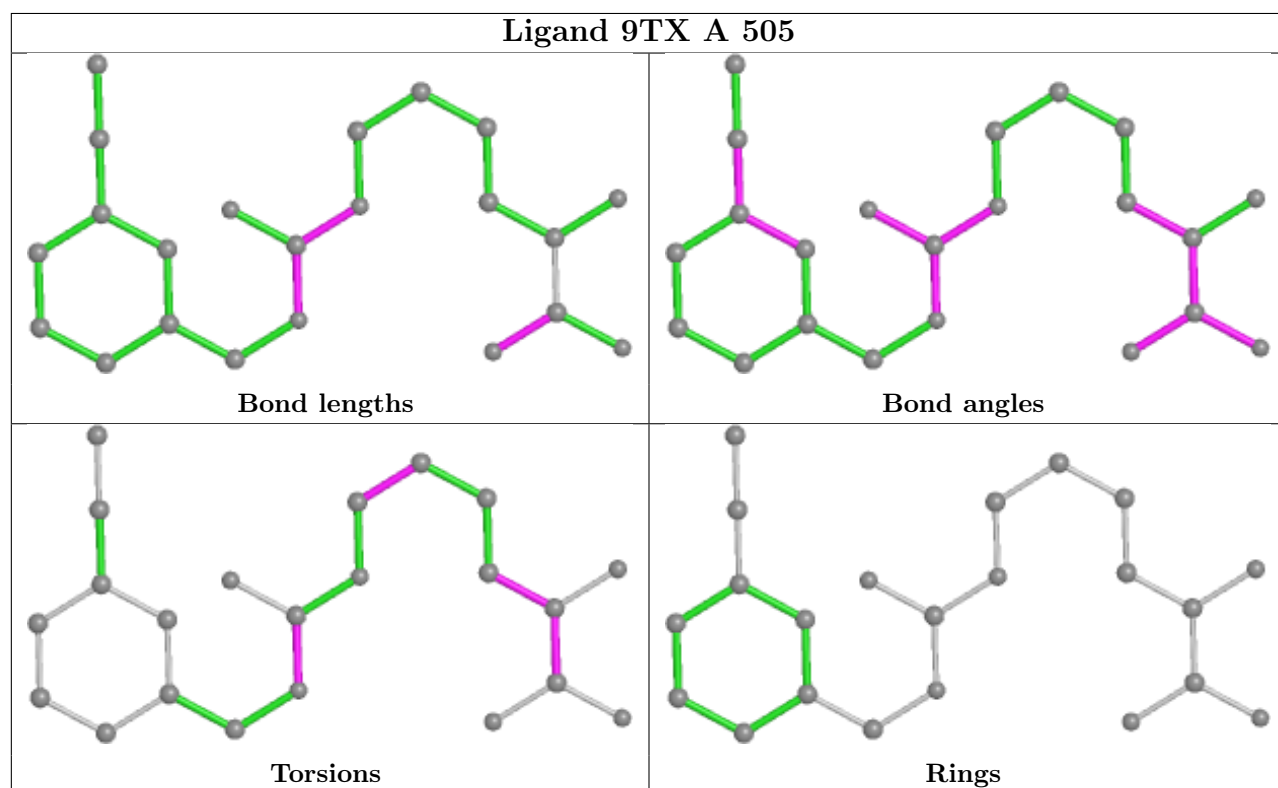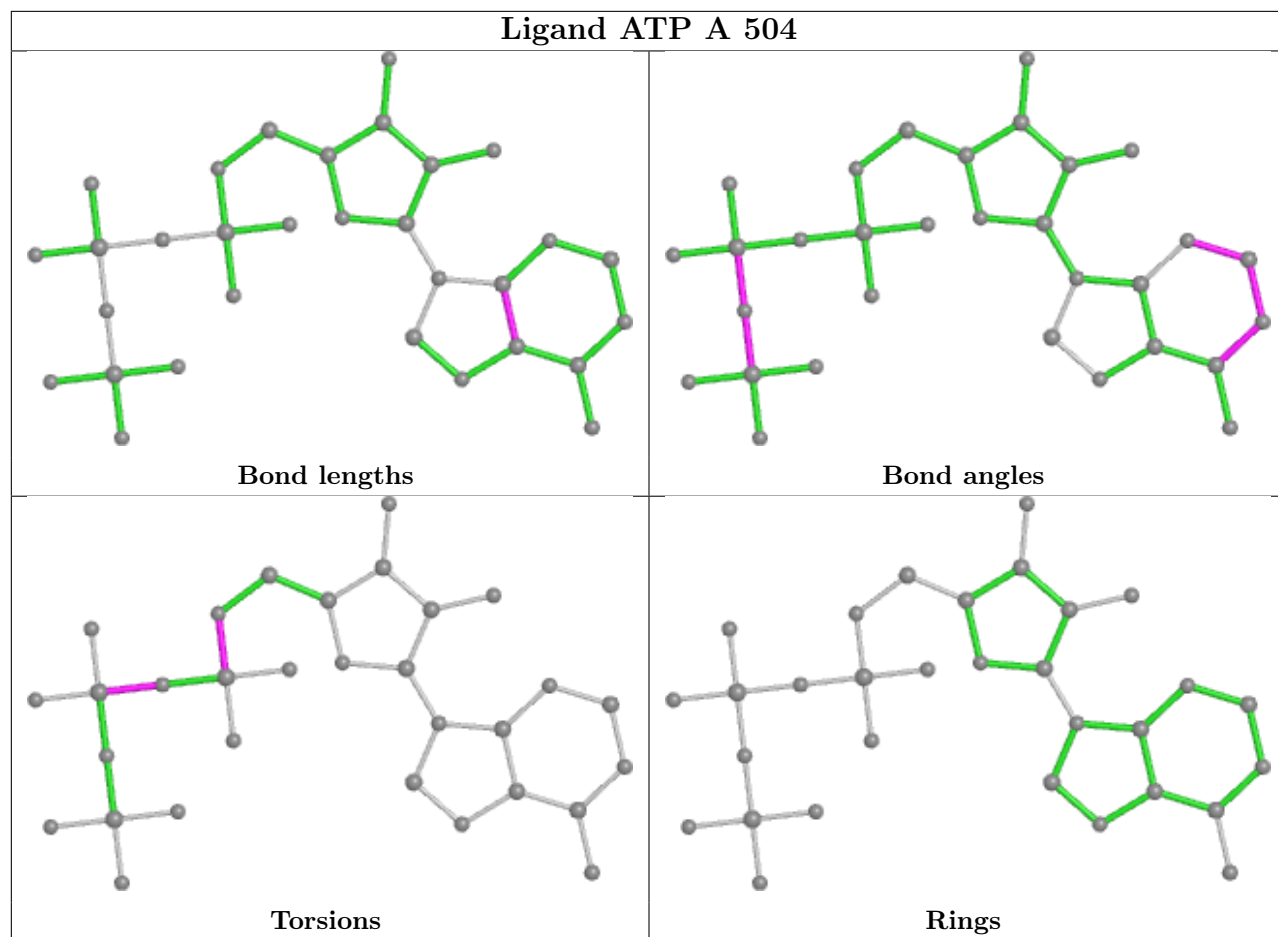
5.7 Other polymers ⓘ

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	261/274 (95%)	0.45	32 (12%)	4	3	17, 28, 63, 80	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	383	VAL	6.8
1	A	382	MET	6.5
1	A	274	GLU	5.5
1	A	384	PHE	5.4
1	A	273	ILE	5.4
1	A	385	GLY	4.5
1	A	336	LYS	4.5
1	A	293	LYS	4.4
1	A	386	ASP	4.3
1	A	378	GLY	4.2
1	A	284	LEU	4.1
1	A	286	LYS	3.7
1	A	381	CYS	3.6
1	A	380	SER	3.5
1	A	379	ASP	3.4
1	A	334	ASP	3.4
1	A	243[A]	LEU	3.3
1	A	335	GLY	3.3
1	A	188	PRO	3.1
1	A	205	LYS	2.7
1	A	406	PRO	2.7
1	A	211	ASN	2.6
1	A	238	GLU	2.5
1	A	333	SER	2.5
1	A	208	ILE	2.5
1	A	325	ILE	2.4
1	A	294	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	209	SER	2.2
1	A	377	VAL	2.1
1	A	285	SER	2.1
1	A	387	THR	2.1
1	A	436[A]	ASN	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
4	9TX	A	505	22/22	0.79	0.19	34,53,64,75	0
6	PEG	A	508	7/7	0.79	0.13	35,42,49,59	0
2	MG	A	503	1/1	0.91	0.24	52,52,52,52	0
3	ATP	A	504	31/31	0.91	0.10	20,26,53,61	0
5	K	A	507	1/1	0.95	0.16	52,52,52,52	0
5	K	A	506	1/1	0.95	0.11	86,86,86,86	0
2	MG	A	501	1/1	0.97	0.12	39,39,39,39	0
2	MG	A	502	1/1	0.97	0.11	39,39,39,39	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 9TX A 505:

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

Electron density around ATP A 504:

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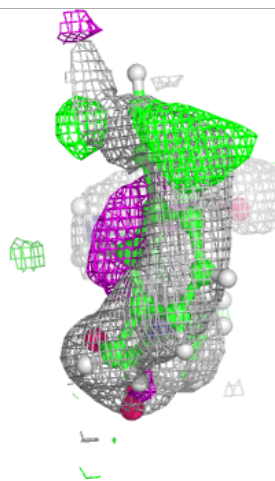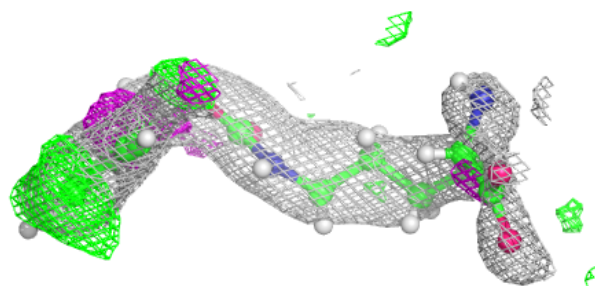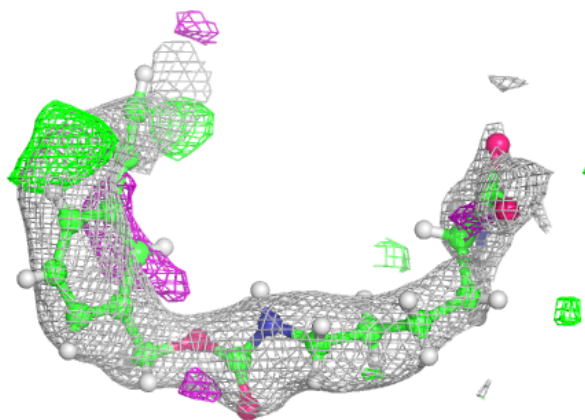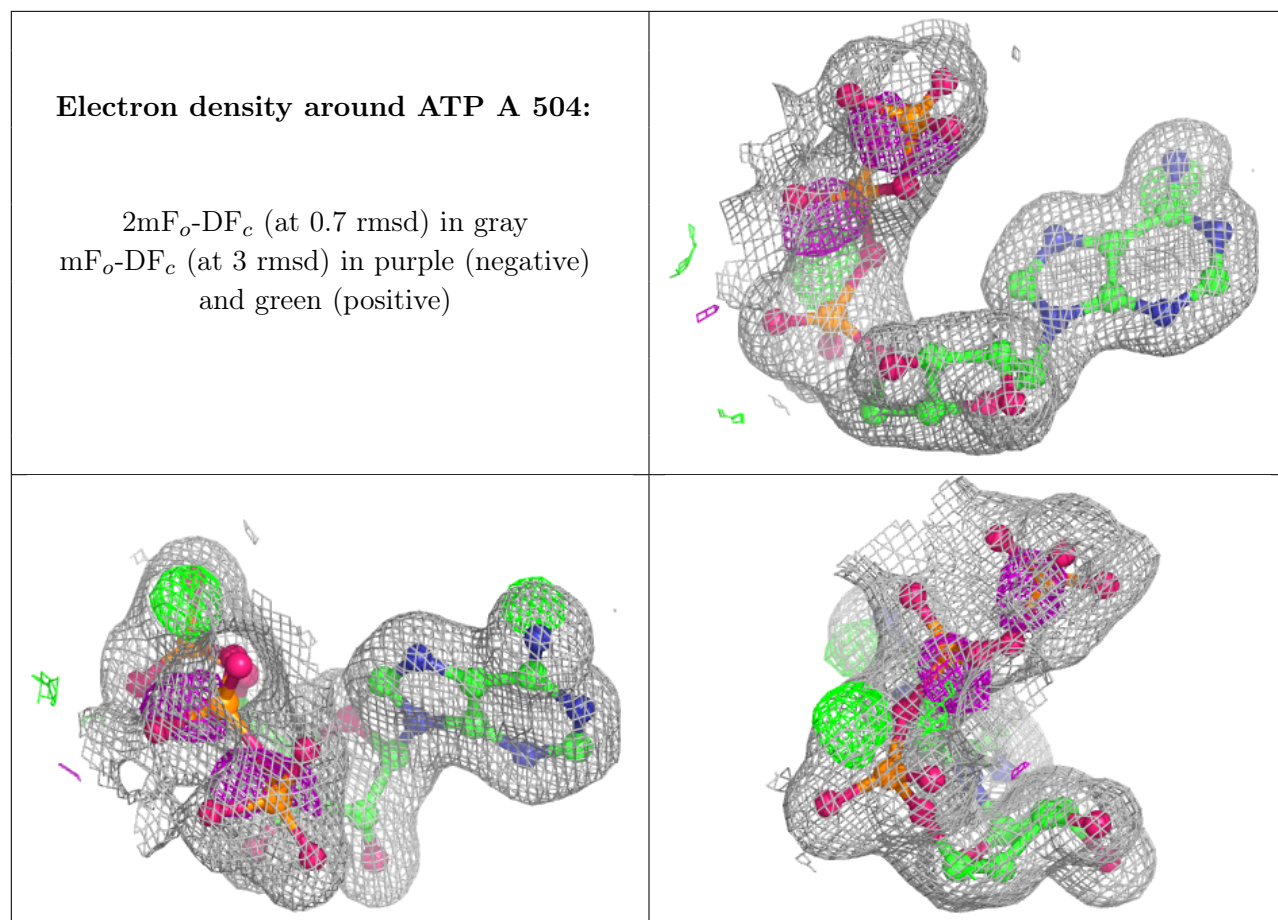
6.5 Other polymers (i)

There are no such residues in this entry.