



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 08:03 pm BST

PDB ID : 3FPZ
Title : Saccharomyces cerevisiae THI4p is a suicide thiamin thiazole synthase
Authors : Bale, S.; Chatterjee, A.; Dorrestein, P.C.; Begley, T.P.; Ealick, S.E.
Deposited on : 2009-01-06
Resolution : 1.82 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 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 : 2.11
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) : 2.11

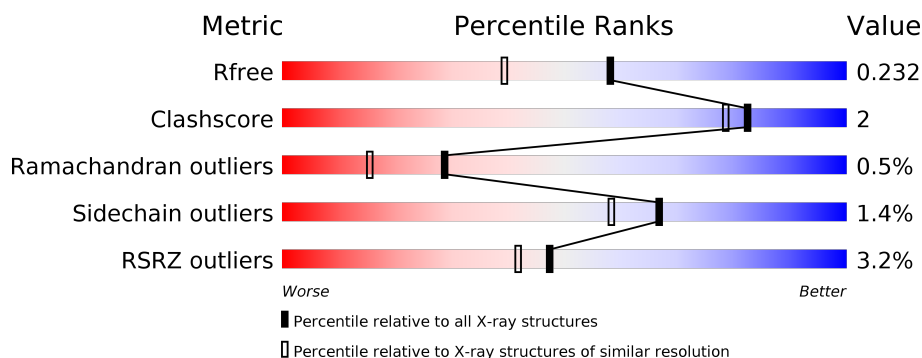
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.82 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	7484 (1.84-1.80)
Clashscore	141614	8401 (1.84-1.80)
Ramachandran outliers	138981	8290 (1.84-1.80)
Sidechain outliers	138945	8290 (1.84-1.80)
RSRZ outliers	127900	7371 (1.84-1.80)

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 on 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	326	<div> <div>3%</div> <div> <div></div> <div>90%</div> <div>6%</div> <div>5%</div> </div> </div>
1	B	326	<div> <div>3%</div> <div> <div></div> <div>88%</div> <div>6%</div> <div>6%</div> </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5237 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.

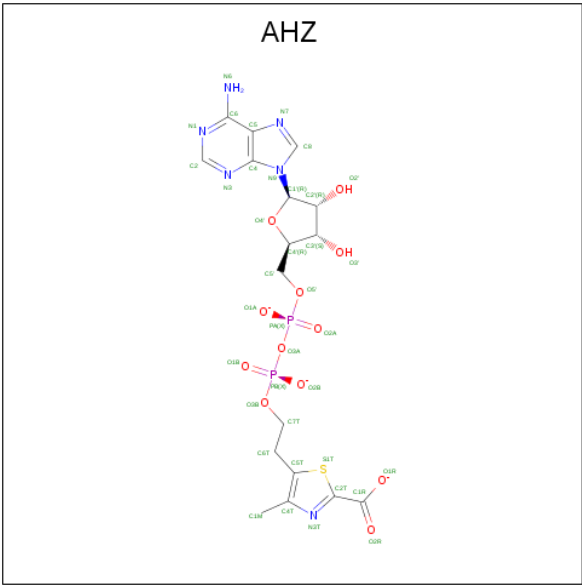
- Molecule 1 is a protein called Thiazole biosynthetic enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	311	Total	C	N	O	S	0	2	0
			2342	1486	398	442	16			
1	B	307	Total	C	N	O	S	0	2	0
			2313	1468	393	436	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	205	DHA	CYS	SEE REMARK 999	UNP P32318
B	205	DHA	CYS	SEE REMARK 999	UNP P32318

- Molecule 2 is ADENOSINE DIPHOSPHATE 5-(BETA-ETHYL)-4-METHYL-THIAZOLE-2-CARBOXYLIC ACID (three-letter code: AHZ) (formula: C₁₇H₁₉N₆O₁₂P₂S).



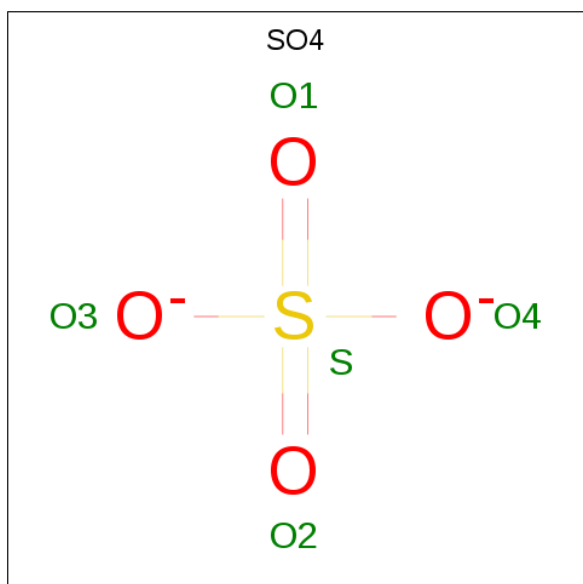
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P S	0	0
			38	17	6	12	2 1		

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	S	0	0
			38	17	6	12	2	1		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

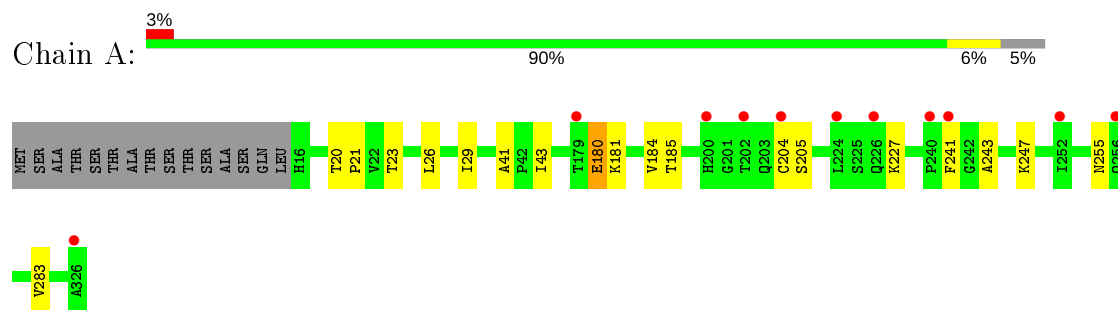
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	241	Total	O	0	0
			241	241		
4	B	250	Total	O	0	0
			250	250		

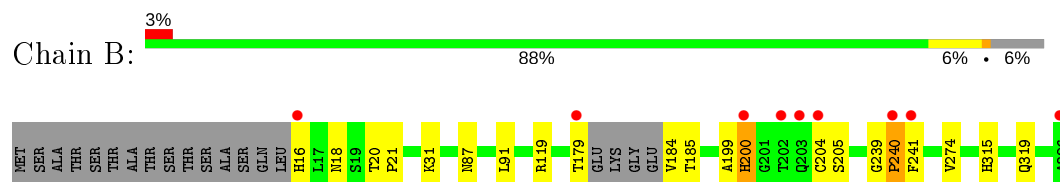
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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 ($\text{RSRZ} > 2$). 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: Thiazole biosynthetic enzyme



• Molecule 1: Thiazole biosynthetic enzyme



4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	140.70 Å 140.70 Å 73.33 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.75 – 1.82 49.74 – 1.82	Depositor EDS
% Data completeness (in resolution range)	98.8 (49.75-1.82) 98.7 (49.74-1.82)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 1.82 Å)	Xtriage
Refinement program	REFMAC 5.4.0066	Depositor
R, R_{free}	0.195 , 0.236 0.194 , 0.232	Depositor DCC
R_{free} test set	3212 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	24.2	Xtriage
Anisotropy	0.295	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.479 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5237	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.01% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: AHZ, SO4, DHA

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	0.63	0/2387	0.62	0/3236
1	B	0.65	0/2357	0.63	0/3193
All	All	0.64	0/4744	0.62	0/6429

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
1	B	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	205	DHA	Peptide
1	B	205	DHA	Peptide

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	2342	0	2301	12	0
1	B	2313	0	2272	14	0
2	A	38	0	19	1	0
2	B	38	0	19	1	0
3	A	10	0	0	0	0
3	B	5	0	0	0	0
4	A	241	0	0	2	0
4	B	250	0	0	0	0
All	All	5237	0	4611	23	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 2.

All (23) 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:B:199:ALA:HA	1:B:200:HIS:CB	1.90	1.02
1:A:180:GLU:N	1:A:181:LYS:HA	1.85	0.88
1:B:199:ALA:CA	1:B:200:HIS:CB	2.66	0.72
1:B:20:THR:HB	1:B:21:PRO:HD2	1.76	0.67
1:A:180:GLU:H	1:A:181:LYS:HA	1.57	0.67
1:A:26:LEU:HD22	1:A:29:ILE:HD11	1.74	0.67
1:B:16:HIS:HA	1:B:18:ASN:OD1	1.96	0.65
4:A:390:HOH:O	1:B:315:HIS:HE1	1.84	0.60
1:B:241:PHE:HB3	2:B:1101:AHZ:N7	2.18	0.58
1:B:20:THR:HB	1:B:21:PRO:CD	2.38	0.54
1:A:184:VAL:HB	1:A:283:VAL:HG22	1.93	0.51
1:A:23:THR:H	1:B:87:ASN:HD21	1.58	0.51
1:B:184:VAL:CB	1:B:185:THR:HA	2.41	0.51
1:A:185:THR:HA	4:A:693:HOH:O	2.11	0.50
1:B:274:VAL:O	1:B:315:HIS:HD2	1.95	0.50
1:A:43:ILE:HB	1:B:119:ARG:HD3	1.94	0.48
1:A:180:GLU:H	1:A:181:LYS:CA	2.24	0.48
1:B:315:HIS:CE1	1:B:319:GLN:HE21	2.32	0.47
1:A:241:PHE:HB2	2:A:1100:AHZ:N7	2.29	0.47
1:A:243:ALA:HB3	1:A:247:LYS:HE3	1.97	0.46
1:A:41:ALA:O	1:B:119:ARG:HD2	2.19	0.42
1:B:239:GLY:HA3	1:B:240:PRO:HD3	1.83	0.41
1:A:20:THR:HB	1:A:21:PRO:HD2	2.02	0.41

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	310/326 (95%)	297 (96%)	12 (4%)	1 (0%)	41	27
1	B	304/326 (93%)	293 (96%)	9 (3%)	2 (1%)	22	10
All	All	614/652 (94%)	590 (96%)	21 (3%)	3 (0%)	29	15

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	200	HIS
1	B	240	PRO
1	A	180	GLU

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	248/265 (94%)	245 (99%)	3 (1%)	71	64
1	B	245/265 (92%)	241 (98%)	4 (2%)	62	53
All	All	493/530 (93%)	486 (99%)	7 (1%)	67	58

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	204	CYS
1	A	227	LYS
1	A	255	ASN

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Mol	Chain	Res	Type
1	B	31	LYS
1	B	91	LEU
1	B	179	THR
1	B	204	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	87	ASN
1	B	203	GLN
1	B	285	ASN
1	B	315	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	DHA	B	205	1	4,4,5	1.34	1 (25%)	2,4,6	2.90	2 (100%)
1	DHA	A	205	1	4,4,5	1.41	1 (25%)	2,4,6	3.32	2 (100%)

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
1	DHA	B	205	1	-	0/0/2/4	-
1	DHA	A	205	1	-	0/0/2/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	205	DHA	C-CA	2.75	1.49	1.45
1	B	205	DHA	C-CA	2.59	1.49	1.45

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	205	DHA	CB-CA-N	-3.80	116.82	125.81
1	B	205	DHA	CB-CA-N	-3.33	117.94	125.81
1	A	205	DHA	O-C-CA	-2.77	120.38	125.54
1	B	205	DHA	O-C-CA	-2.39	121.09	125.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

5 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	$\# Z > 2$
3	SO4	A	328	-	4,4,4	0.14	0	6,6,6	0.07	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	AHZ	A	1100	-	32,41,41	1.43	3 (9%)	28,62,62	1.38	3 (10%)
3	SO4	B	327	-	4,4,4	0.14	0	6,6,6	0.12	0
3	SO4	A	327	-	4,4,4	0.14	0	6,6,6	0.07	0
2	AHZ	B	1101	-	32,41,41	1.50	5 (15%)	28,62,62	1.33	4 (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	AHZ	A	1100	-	-	5/18/43/43	0/4/4/4
2	AHZ	B	1101	-	-	5/18/43/43	0/4/4/4

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1101	AHZ	O4'-C1'	4.75	1.47	1.41
2	A	1100	AHZ	O4'-C1'	3.98	1.46	1.41
2	A	1100	AHZ	C6T-C5T	3.55	1.52	1.50
2	B	1101	AHZ	C6T-C5T	3.09	1.52	1.50
2	B	1101	AHZ	C2T-S1T	-2.76	1.69	1.73
2	A	1100	AHZ	C2T-S1T	-2.59	1.70	1.73
2	B	1101	AHZ	C5T-S1T	-2.02	1.70	1.74
2	B	1101	AHZ	C1M-C4T	2.02	1.53	1.50

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1100	AHZ	O4'-C1'-C2'	-4.30	100.64	106.93
2	B	1101	AHZ	O4'-C1'-C2'	-3.91	101.20	106.93
2	A	1100	AHZ	N3-C2-N1	-3.87	122.63	128.68
2	B	1101	AHZ	N3-C2-N1	-3.81	122.72	128.68
2	A	1100	AHZ	C6T-C5T-C4T	2.33	129.30	127.43
2	B	1101	AHZ	C4-C5-N7	-2.07	107.24	109.40
2	B	1101	AHZ	C6T-C5T-C4T	2.04	129.07	127.43

There are no chirality outliers.

All (10) torsion outliers are listed below:

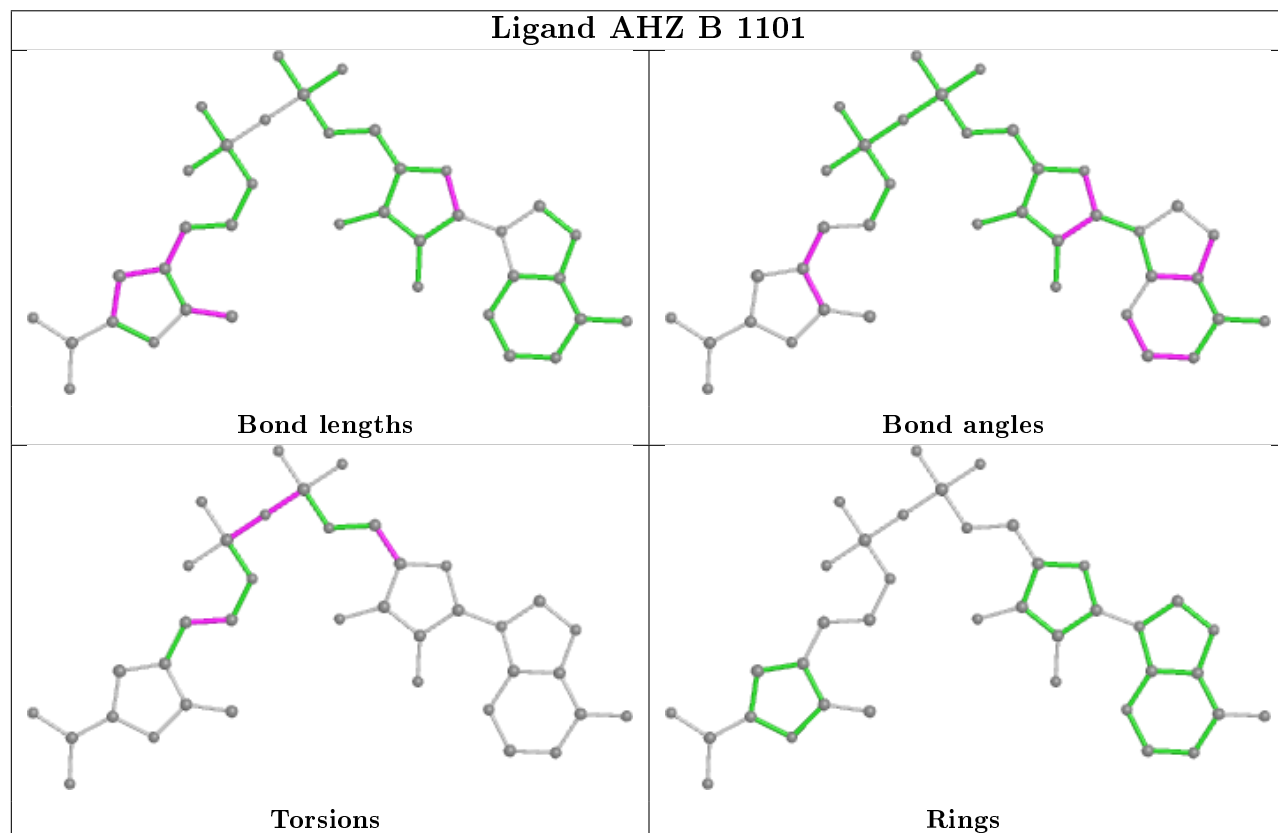
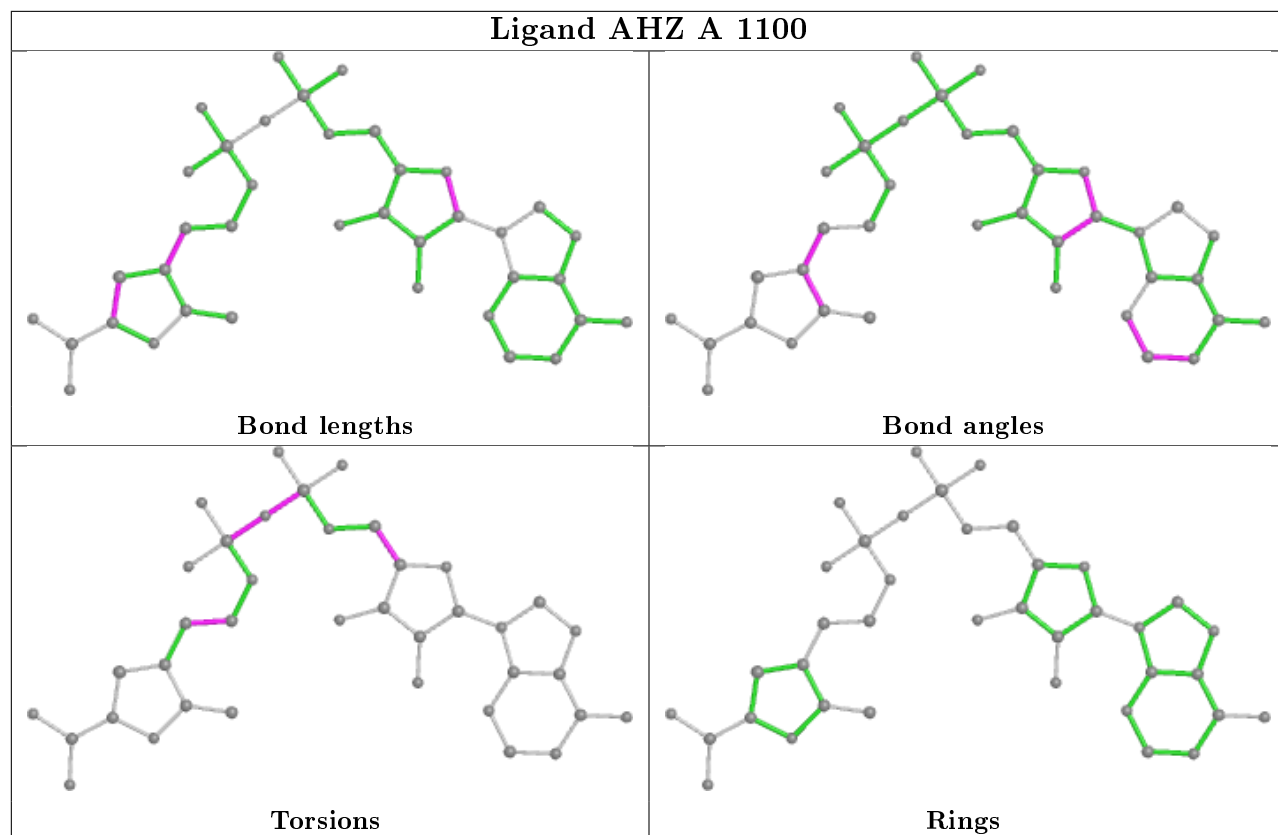
Mol	Chain	Res	Type	Atoms
2	A	1100	AHZ	PA-O3A-PB-O3B
2	A	1100	AHZ	C5T-C6T-C7T-O3B
2	B	1101	AHZ	PA-O3A-PB-O3B
2	B	1101	AHZ	C5T-C6T-C7T-O3B
2	A	1100	AHZ	PB-O3A-PA-O2A
2	A	1100	AHZ	O4'-C4'-C5'-O5'
2	B	1101	AHZ	O4'-C4'-C5'-O5'
2	A	1100	AHZ	PB-O3A-PA-O1A
2	B	1101	AHZ	PB-O3A-PA-O2A
2	B	1101	AHZ	PB-O3A-PA-O1A

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1100	AHZ	1	0
2	B	1101	AHZ	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

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	310/326 (95%)	-0.02	11 (3%) 44 38	14, 24, 47, 56	0
1	B	306/326 (93%)	-0.10	9 (2%) 51 46	14, 24, 45, 59	0
All	All	616/652 (94%)	-0.06	20 (3%) 47 42	14, 24, 47, 59	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	241	PHE	6.6
1	A	241	PHE	6.1
1	A	202	THR	5.4
1	B	240	PRO	4.8
1	B	202	THR	4.8
1	A	326	ALA	4.5
1	A	240	PRO	4.1
1	B	200	HIS	3.5
1	A	252	ILE	3.5
1	A	200	HIS	3.4
1	A	179	THR	3.2
1	A	224	LEU	3.2
1	A	226	GLN	3.1
1	A	256	GLN	2.8
1	A	204	CYS	2.6
1	B	179	THR	2.6
1	B	204	CYS	2.6
1	B	16	HIS	2.2
1	B	326	ALA	2.2
1	B	203	GLN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(\AA^2)	Q<0.9
1	DHA	A	205	5/6	0.73	0.30	38,38,40,41	0
1	DHA	B	205	5/6	0.82	0.28	37,37,39,41	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

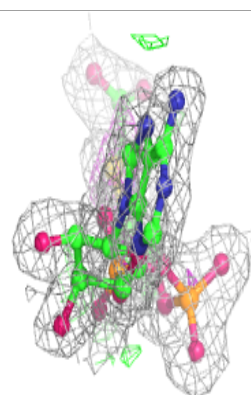
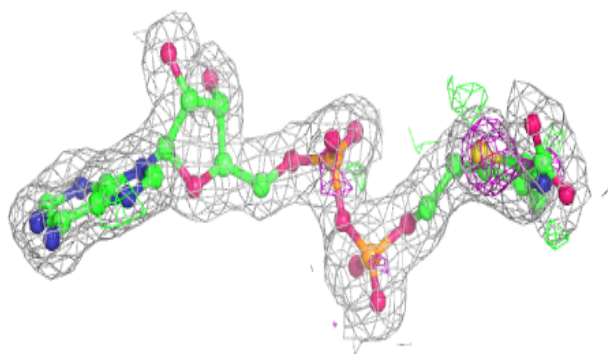
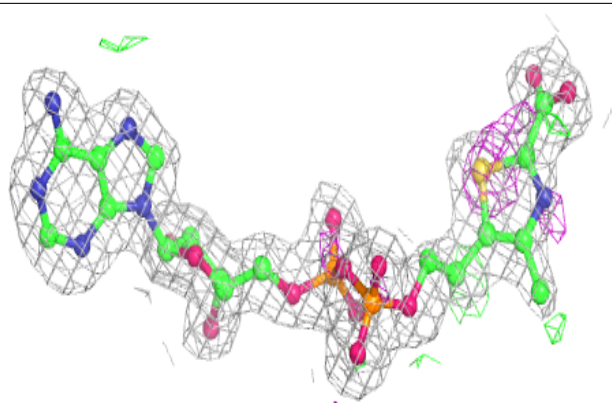
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	SO4	A	328	5/5	0.79	0.22	89,89,89,89	0
3	SO4	B	327	5/5	0.89	0.13	70,71,71,71	0
3	SO4	A	327	5/5	0.94	0.08	73,73,74,74	0
2	AHZ	A	1100	38/38	0.95	0.11	19,23,44,44	0
2	AHZ	B	1101	38/38	0.96	0.09	19,22,41,42	0

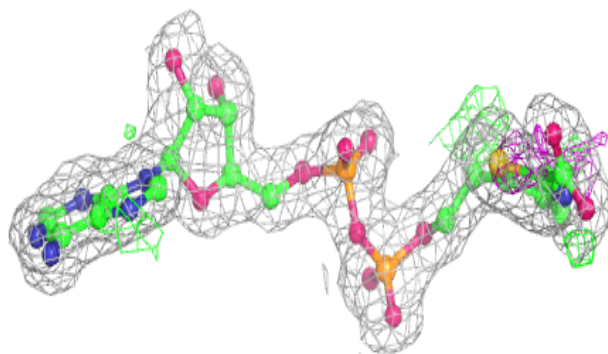
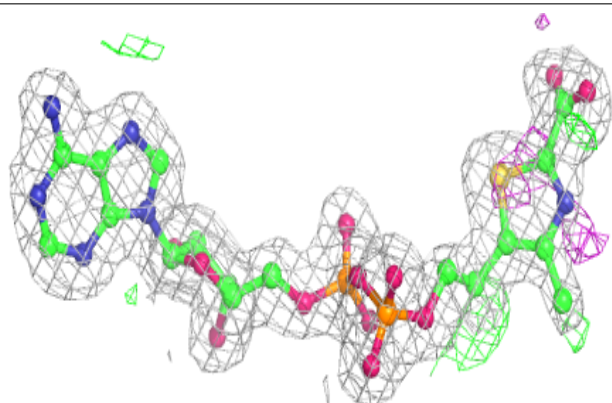
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around AHZ A 1100:

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

**Electron density around AHZ B 1101:**

$2mF_o-DF_c$ (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.