



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 22, 2024 – 09:31 PM EDT

PDB ID : 6J0P  
Title : Crystal structure of HypX from Aquifex aeolicus (Crystal Form I)  
Authors : Muraki, N.; Aono, S.  
Deposited on : 2018-12-25  
Resolution : 1.79 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](mailto: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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
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)	:	2.37.1

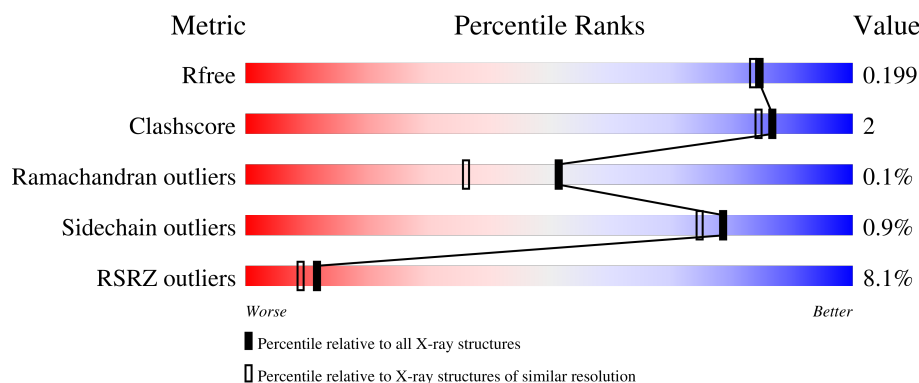
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.79 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	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	582	<div> <div>4%</div> <div>91%</div> <div>5%</div> </div>
1	B	582	<div> <div>12%</div> <div>89%</div> <div>8%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9806 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 Hydrogenase regulation HoxX.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	564	Total	C	N	O	S	0	6	0
			4633	2979	797	839	18			
1	B	563	Total	C	N	O	S	0	25	0
			4673	3004	808	842	19			

There are 40 discrepancies between the modelled and reference sequences:

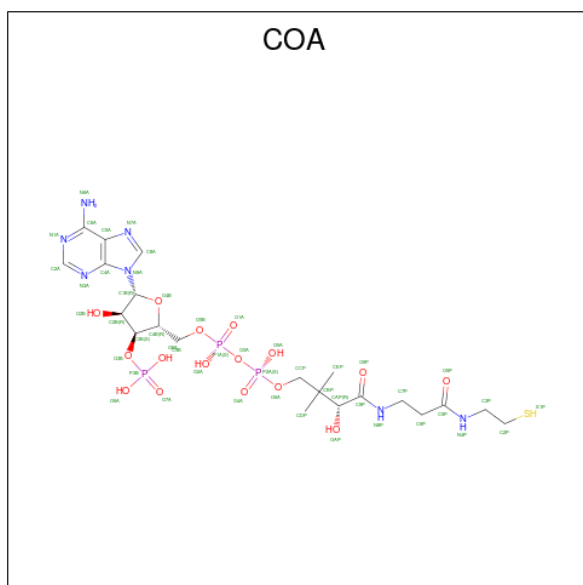
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP O67224
A	-18	GLY	-	expression tag	UNP O67224
A	-17	SER	-	expression tag	UNP O67224
A	-16	SER	-	expression tag	UNP O67224
A	-15	HIS	-	expression tag	UNP O67224
A	-14	HIS	-	expression tag	UNP O67224
A	-13	HIS	-	expression tag	UNP O67224
A	-12	HIS	-	expression tag	UNP O67224
A	-11	HIS	-	expression tag	UNP O67224
A	-10	HIS	-	expression tag	UNP O67224
A	-9	SER	-	expression tag	UNP O67224
A	-8	SER	-	expression tag	UNP O67224
A	-7	GLY	-	expression tag	UNP O67224
A	-6	LEU	-	expression tag	UNP O67224
A	-5	VAL	-	expression tag	UNP O67224
A	-4	PRO	-	expression tag	UNP O67224
A	-3	ARG	-	expression tag	UNP O67224
A	-2	GLY	-	expression tag	UNP O67224
A	-1	SER	-	expression tag	UNP O67224
A	0	HIS	-	expression tag	UNP O67224
B	-19	MET	-	initiating methionine	UNP O67224
B	-18	GLY	-	expression tag	UNP O67224
B	-17	SER	-	expression tag	UNP O67224
B	-16	SER	-	expression tag	UNP O67224
B	-15	HIS	-	expression tag	UNP O67224

*Continued on next page...*

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP O67224
B	-13	HIS	-	expression tag	UNP O67224
B	-12	HIS	-	expression tag	UNP O67224
B	-11	HIS	-	expression tag	UNP O67224
B	-10	HIS	-	expression tag	UNP O67224
B	-9	SER	-	expression tag	UNP O67224
B	-8	SER	-	expression tag	UNP O67224
B	-7	GLY	-	expression tag	UNP O67224
B	-6	LEU	-	expression tag	UNP O67224
B	-5	VAL	-	expression tag	UNP O67224
B	-4	PRO	-	expression tag	UNP O67224
B	-3	ARG	-	expression tag	UNP O67224
B	-2	GLY	-	expression tag	UNP O67224
B	-1	SER	-	expression tag	UNP O67224
B	0	HIS	-	expression tag	UNP O67224

- Molecule 2 is COENZYME A (three-letter code: COA) (formula:  $C_{21}H_{36}N_7O_{16}P_3S$ ).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		
2	B	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



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

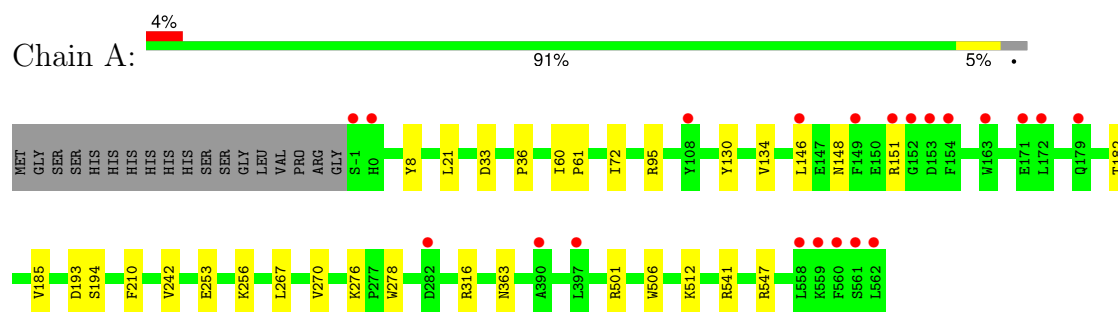
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	247	Total	O	0	0
			247	247		
4	B	145	Total	O	0	0
			145	145		

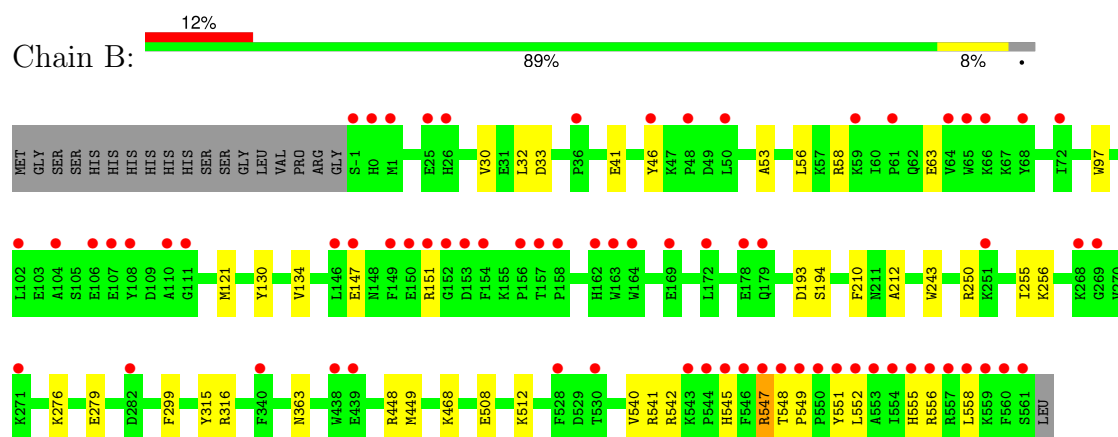
### 3 Residue-property plots [i](#)

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: Hydrogenase regulation HoxX



#### • Molecule 1: Hydrogenase regulation HoxX



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.84Å 124.30Å 290.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.47 – 1.79 48.48 – 1.79	Depositor EDS
% Data completeness (in resolution range)	98.7 (42.47-1.79) 98.7 (48.48-1.79)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.08 (at 1.79Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, $R_{free}$	0.173 , 0.196 0.177 , 0.199	Depositor DCC
$R_{free}$ test set	6672 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.9	Xtriage
Anisotropy	0.162	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 50.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	9806	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.55% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: COA, GOL

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.41	0/4769	0.55	0/6441
1	B	8.66	2/4812 (0.0%)	0.72	4/6498 (0.1%)
All	All	6.14	2/9581 (0.0%)	0.64	4/12939 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	547[A]	ARG	CB-CG	424.38	12.98	1.52
1	B	547[B]	ARG	CB-CG	424.38	12.98	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	547[A]	ARG	CA-CB-CG	-28.09	51.60	113.40
1	B	547[B]	ARG	CA-CB-CG	-28.09	51.60	113.40
1	B	547[A]	ARG	CB-CG-CD	9.01	135.02	111.60
1	B	547[B]	ARG	CB-CG-CD	9.01	135.02	111.60

There are no chirality outliers.

There are no planarity outliers.

### 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	4633	0	4583	18	0
1	B	4673	0	4602	22	0
2	A	48	0	32	0	0
2	B	48	0	32	0	0
3	A	6	0	8	0	0
3	B	6	0	8	1	0
4	A	247	0	0	3	0
4	B	145	0	0	1	0
All	All	9806	0	9265	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 2.

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:148:ASN:OD1	1:A:151:ARG:NH1	2.25	0.70
1:B:30:VAL:HB	1:B:552[A]:LEU:HB3	1.81	0.62
1:B:508:GLU:O	1:B:512[B]:LYS:HE2	2.03	0.58
1:B:33:ASP:O	1:B:58:ARG:NH1	2.38	0.56
4:A:737:HOH:O	1:B:512[B]:LYS:HE3	2.05	0.56
1:A:316:ARG:NH2	4:A:701:HOH:O	2.41	0.54
1:A:267:LEU:O	1:A:270:VAL:HG12	2.07	0.53
1:B:41:GLU:OE2	1:B:556[A]:ARG:NH1	2.33	0.53
1:A:253:GLU:O	1:A:541:ARG:NH1	2.42	0.52
1:B:97:TRP:HB2	1:B:121:MET:HG2	1.91	0.52
1:B:193:ASP:OD1	1:B:194:SER:HA	2.11	0.51
1:A:36:PRO:HB3	1:A:61:PRO:HG3	1.93	0.50
1:A:21:LEU:HD22	1:A:146:LEU:HD11	1.93	0.50
1:B:210:PHE:CE1	1:B:256:LYS:HE2	2.49	0.48
1:A:501:ARG:HD3	1:A:506:TRP:CZ2	2.49	0.48
1:B:316:ARG:NH2	4:B:702:HOH:O	2.47	0.48
1:A:193:ASP:OD1	1:A:194:SER:HA	2.14	0.47
1:B:448:ARG:HD3	1:B:549[B]:PRO:HB3	1.95	0.47
1:A:185:VAL:HG11	1:A:242[B]:VAL:HG21	1.96	0.47
1:A:95:ARG:NH1	1:B:315[A]:TYR:OH	2.48	0.46
1:A:8:TYR:HA	1:A:33:ASP:HB3	1.98	0.45
1:A:182:THR:HG23	1:A:242[A]:VAL:HG22	1.97	0.45
1:B:449:MET:SD	1:B:549[B]:PRO:HG3	2.56	0.45
1:B:32:LEU:HD11	1:B:548[B]:THR:HG21	1.99	0.45
1:A:276:LYS:HD3	1:A:278:TRP:CZ2	2.51	0.45
1:B:130:TYR:HA	1:B:134:VAL:HB	2.00	0.44

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:255:ILE:HG21	1:B:540:VAL:HG12	2.00	0.44
1:A:512:LYS:HD3	1:A:512:LYS:HA	1.77	0.43
1:A:547:ARG:NH2	4:A:703:HOH:O	2.46	0.43
1:B:276:LYS:HB2	1:B:279:GLU:HG3	2.00	0.43
1:A:210:PHE:CE1	1:A:256:LYS:HE2	2.53	0.43
1:B:147:GLU:O	1:B:151:ARG:HG2	2.19	0.42
1:A:130:TYR:HA	1:A:134:VAL:HB	2.01	0.42
1:B:212:ALA:HA	1:B:243:TRP:O	2.20	0.41
1:B:551[A]:TYR:CE1	1:B:552[A]:LEU:HG	2.55	0.41
1:B:299:PHE:CG	3:B:602:GOL:H12	2.56	0.40
1:B:53:ALA:HB1	1:B:56:LEU:HD13	2.03	0.40
1:A:60:ILE:HD12	1:A:72:ILE:HD11	2.04	0.40
1:B:46:TYR:HB2	1:B:555[A]:HIS:HB3	2.02	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	568/582 (98%)	558 (98%)	10 (2%)	0	100	100
1	B	574/582 (99%)	553 (96%)	20 (4%)	1 (0%)	47	33
All	All	1142/1164 (98%)	1111 (97%)	30 (3%)	1 (0%)	51	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	558[A]	LEU

### 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	491/511 (96%)	490 (100%)	1 (0%)	93	92
1	B	492/511 (96%)	481 (98%)	11 (2%)	52	39
All	All	983/1022 (96%)	971 (99%)	12 (1%)	78	65

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	363	ASN
1	B	63	GLU
1	B	250	ARG
1	B	363	ASN
1	B	468	LYS
1	B	541	ARG
1	B	542[A]	ARG
1	B	542[B]	ARG
1	B	545[A]	HIS
1	B	545[B]	HIS
1	B	547[A]	ARG
1	B	547[B]	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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](#)

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	COA	A	601	-	43,50,50	3.62	13 (30%)	56,75,75	1.83	11 (19%)
3	GOL	B	602	-	5,5,5	0.93	0	5,5,5	1.13	0
2	COA	B	601	-	43,50,50	3.66	13 (30%)	56,75,75	1.79	11 (19%)
3	GOL	A	602	-	5,5,5	1.06	0	5,5,5	1.09	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	COA	A	601	-	-	4/44/64/64	0/3/3/3
3	GOL	B	602	-	-	2/4/4/4	-
2	COA	B	601	-	-	3/44/64/64	0/3/3/3
3	GOL	A	602	-	-	2/4/4/4	-

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	COA	O4B-C1B	16.76	1.62	1.40
2	A	601	COA	O4B-C1B	15.71	1.61	1.40
2	A	601	COA	P2A-O3A	9.23	1.69	1.59
2	B	601	COA	P2A-O3A	8.42	1.68	1.59
2	A	601	COA	C9P-N8P	7.91	1.52	1.33
2	B	601	COA	C9P-N8P	7.67	1.51	1.33

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	COA	P3B-O3B	5.79	1.69	1.59
2	A	601	COA	P3B-O3B	5.56	1.69	1.59
2	B	601	COA	O4B-C4B	-5.21	1.33	1.45
2	A	601	COA	O4B-C4B	-5.15	1.33	1.45
2	A	601	COA	C2A-N3A	4.46	1.39	1.32
2	A	601	COA	C5P-N4P	4.04	1.43	1.33
2	B	601	COA	P1A-O3A	3.94	1.63	1.59
2	B	601	COA	C5P-N4P	3.59	1.41	1.33
2	A	601	COA	P1A-O3A	3.46	1.63	1.59
2	B	601	COA	C2A-N3A	3.39	1.37	1.32
2	B	601	COA	O3B-C3B	-2.90	1.34	1.44
2	B	601	COA	OAP-CAP	-2.85	1.37	1.42
2	A	601	COA	C7P-N8P	2.72	1.52	1.46
2	A	601	COA	O3B-C3B	-2.60	1.35	1.44
2	B	601	COA	C7P-N8P	2.56	1.51	1.46
2	A	601	COA	C6A-N6A	2.48	1.43	1.34
2	B	601	COA	C6A-N6A	2.33	1.42	1.34
2	A	601	COA	OAP-CAP	-2.30	1.38	1.42
2	B	601	COA	C1B-N9A	-2.12	1.44	1.49
2	A	601	COA	C1B-N9A	-2.10	1.44	1.49

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	COA	C7P-C6P-C5P	-6.81	101.05	112.39
2	B	601	COA	N3A-C2A-N1A	-6.75	119.51	128.67
2	A	601	COA	N3A-C2A-N1A	-6.22	120.22	128.67
2	B	601	COA	C7P-C6P-C5P	-6.17	102.11	112.39
2	B	601	COA	C5A-C6A-N6A	4.83	127.67	120.31
2	A	601	COA	C5A-C6A-N6A	4.57	127.27	120.31
2	A	601	COA	O4B-C1B-N9A	-3.61	103.96	108.75
2	B	601	COA	N6A-C6A-N1A	-3.11	111.70	118.33
2	B	601	COA	C4B-O4B-C1B	-2.80	107.36	109.92
2	B	601	COA	O6A-CCP-CBP	-2.71	106.18	110.55
2	A	601	COA	O6A-CCP-CBP	-2.66	106.28	110.55
2	A	601	COA	O5P-C5P-C6P	-2.65	117.21	122.02
2	A	601	COA	N6A-C6A-N1A	-2.59	112.81	118.33
2	A	601	COA	C1B-N9A-C4A	-2.45	122.33	126.64
2	A	601	COA	CDP-CBP-CAP	2.43	112.91	108.77
2	A	601	COA	C4B-O4B-C1B	-2.37	107.75	109.92
2	B	601	COA	C6P-C5P-N4P	2.33	120.59	116.34
2	A	601	COA	C6P-C5P-N4P	2.28	120.49	116.34

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	COA	C1B-N9A-C4A	-2.21	122.76	126.64
2	B	601	COA	O5P-C5P-C6P	-2.21	118.02	122.02
2	B	601	COA	C7P-N8P-C9P	-2.11	118.75	122.55
2	B	601	COA	O3A-P2A-O4A	-2.01	104.67	110.70

There are no chirality outliers.

All (11) torsion outliers are listed below:

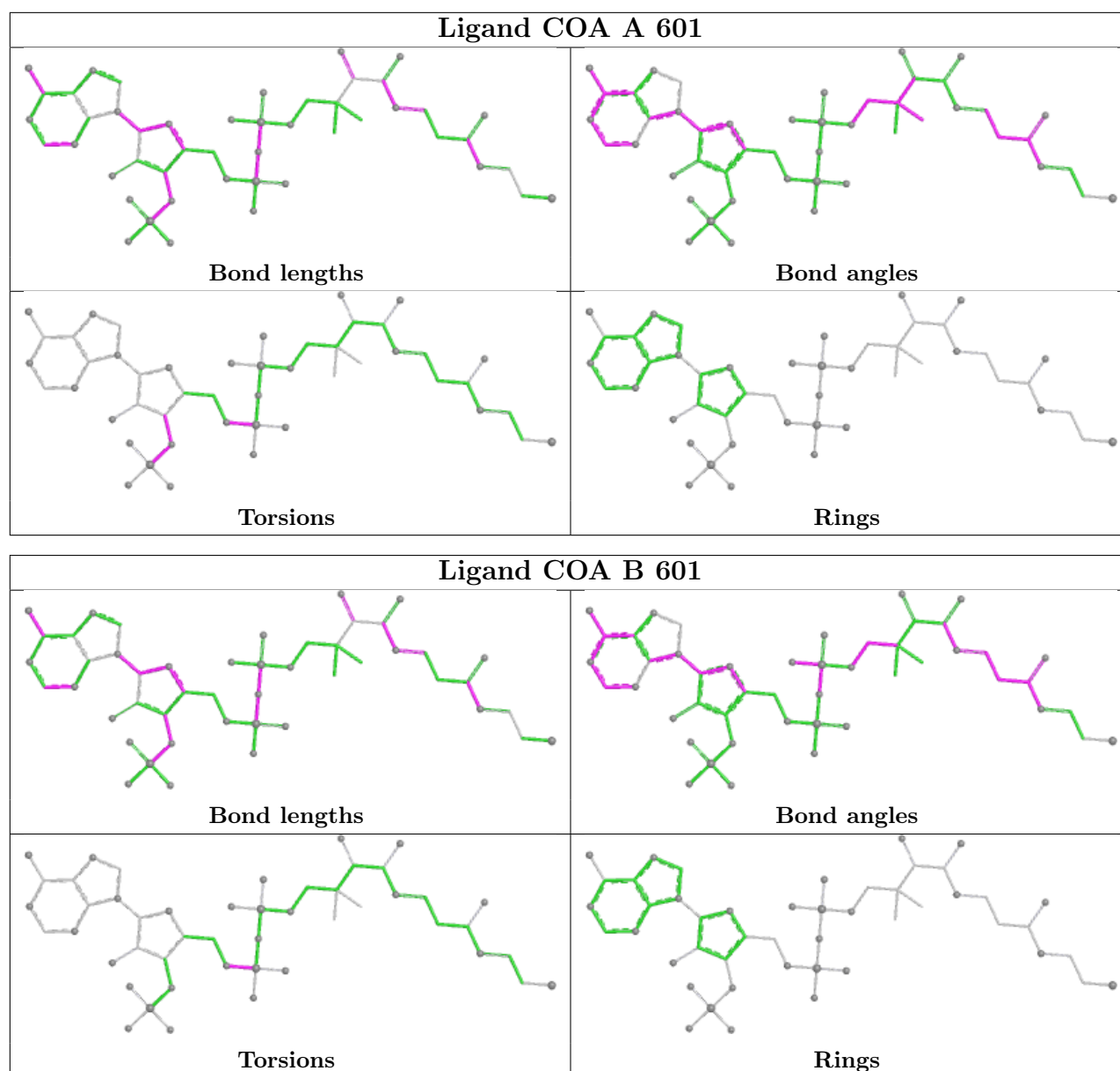
Mol	Chain	Res	Type	Atoms
2	A	601	COA	C5B-O5B-P1A-O2A
2	B	601	COA	C5B-O5B-P1A-O2A
3	A	602	GOL	O1-C1-C2-O2
3	A	602	GOL	O1-C1-C2-C3
3	B	602	GOL	C1-C2-C3-O3
3	B	602	GOL	O2-C2-C3-O3
2	A	601	COA	C5B-O5B-P1A-O1A
2	B	601	COA	C5B-O5B-P1A-O1A
2	B	601	COA	C5B-O5B-P1A-O3A
2	A	601	COA	C2B-C3B-O3B-P3B
2	A	601	COA	C3B-O3B-P3B-O8A

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	564/582 (96%)	0.21	21 (3%) 41 36	18, 28, 49, 78	0
1	B	563/582 (96%)	0.72	70 (12%) 4 3	21, 36, 67, 99	13 (2%)
All	All	1127/1164 (96%)	0.47	91 (8%) 12 9	18, 31, 62, 99	13 (1%)

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	108	TYR	9.8
1	B	-1	SER	7.7
1	B	152	GLY	6.7
1	B	558[A]	LEU	6.6
1	B	528	PHE	6.6
1	B	149	PHE	6.1
1	A	558	LEU	5.9
1	B	560[A]	PHE	5.8
1	B	546[A]	PHE	5.7
1	B	557[A]	ARG	5.6
1	B	553[A]	ALA	5.3
1	B	154	PHE	5.1
1	B	544[A]	PRO	4.9
1	B	146	LEU	4.9
1	B	151	ARG	4.9
1	A	153	ASP	4.8
1	B	547[A]	ARG	4.7
1	B	0	HIS	4.6
1	A	562	LEU	4.5
1	B	548[A]	THR	4.5
1	B	163	TRP	4.5
1	A	151	ARG	4.4
1	B	554[A]	ILE	4.4
1	A	108	TYR	4.4

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	153	ASP	4.2
1	B	552[A]	LEU	4.1
1	B	551[A]	TYR	4.0
1	B	556[A]	ARG	4.0
1	B	107	GLU	3.9
1	B	549[A]	PRO	3.9
1	B	110	ALA	3.8
1	A	152	GLY	3.8
1	B	25	GLU	3.7
1	B	156	PRO	3.6
1	B	1[A]	MET	3.6
1	B	271	LYS	3.6
1	B	545[A]	HIS	3.5
1	B	555[A]	HIS	3.5
1	A	-1	SER	3.5
1	B	65	TRP	3.4
1	B	150	GLU	3.4
1	B	164	TRP	3.4
1	A	149	PHE	3.4
1	A	561	SER	3.4
1	A	560	PHE	3.4
1	B	162	HIS	3.3
1	A	0	HIS	3.3
1	A	282	ASP	3.3
1	B	46	TYR	3.2
1	B	59	LYS	3.0
1	B	111	GLY	2.9
1	A	179	GLN	2.9
1	B	438	TRP	2.9
1	B	561[A]	SER	2.8
1	B	550[A]	PRO	2.7
1	B	269	GLY	2.7
1	B	66	LYS	2.6
1	B	178	GLU	2.6
1	B	543[A]	LYS	2.6
1	A	172	LEU	2.6
1	B	559[A]	LYS	2.6
1	B	172	LEU	2.6
1	A	146	LEU	2.5
1	A	154	PHE	2.5
1	B	106	GLU	2.5
1	A	390	ALA	2.5

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	102	LEU	2.5
1	B	61	PRO	2.4
1	B	147	GLU	2.4
1	A	163	TRP	2.4
1	B	26	HIS	2.4
1	B	179	GLN	2.4
1	B	72	ILE	2.3
1	A	171	GLU	2.3
1	B	169	GLU	2.3
1	B	157	THR	2.3
1	B	282	ASP	2.3
1	A	397	LEU	2.2
1	B	50	LEU	2.2
1	B	158	PRO	2.2
1	B	68	TYR	2.1
1	A	559	LYS	2.1
1	B	340	PHE	2.1
1	B	439	GLU	2.1
1	B	64	VAL	2.1
1	B	268	LYS	2.1
1	B	251	LYS	2.1
1	B	48	PRO	2.0
1	B	530	THR	2.0
1	B	36	PRO	2.0
1	B	104	ALA	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

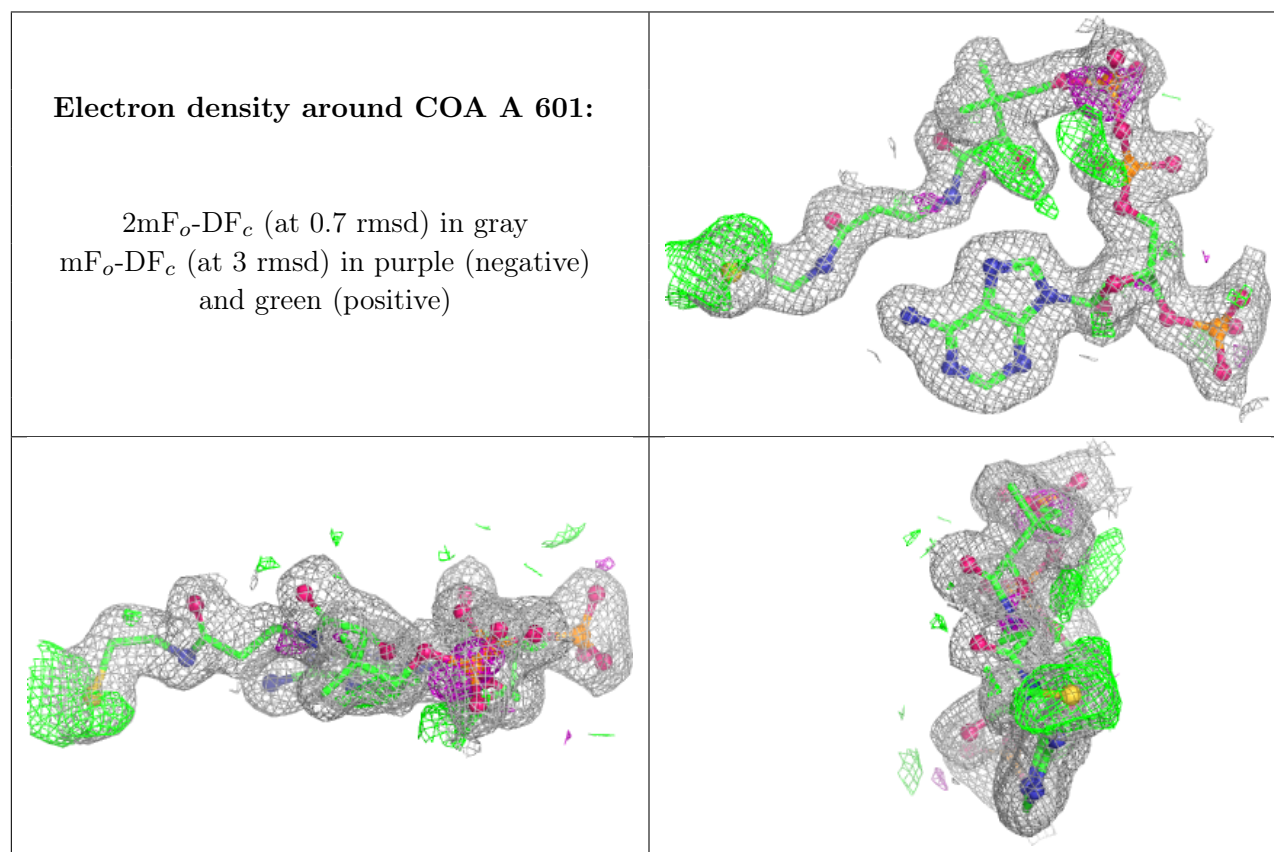
There are no monosaccharides in this entry.

## 6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> 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.

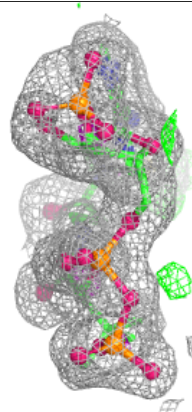
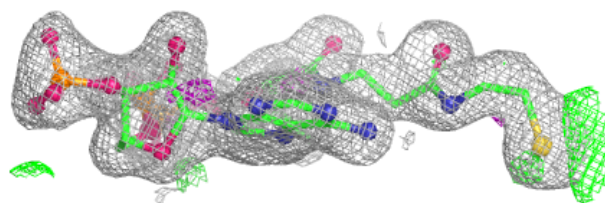
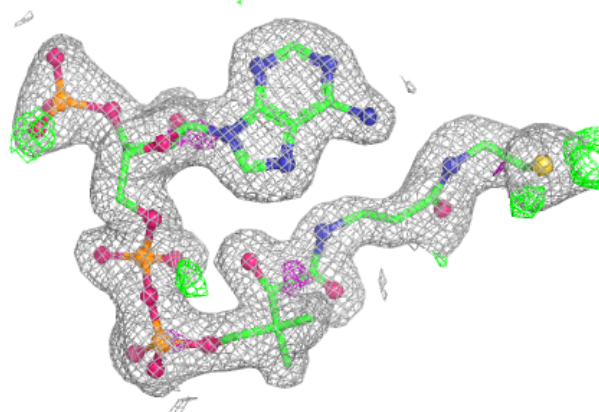
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	GOL	A	602	6/6	0.88	0.12	34,44,46,50	0
3	GOL	B	602	6/6	0.90	0.13	40,45,50,50	0
2	COA	A	601	48/48	0.94	0.13	18,26,33,35	0
2	COA	B	601	48/48	0.95	0.12	23,32,40,49	0

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 COA 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.