



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 20, 2020 – 10:01 PM BST

PDB ID : 3WJQ
Title : Crystal structure of the HypE CN form
Authors : Tominaga, T.; Watanabe, S.; Miki, K.
Deposited on : 2013-10-14
Resolution : 1.65 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 : 2.13.1
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) : 2.13.1

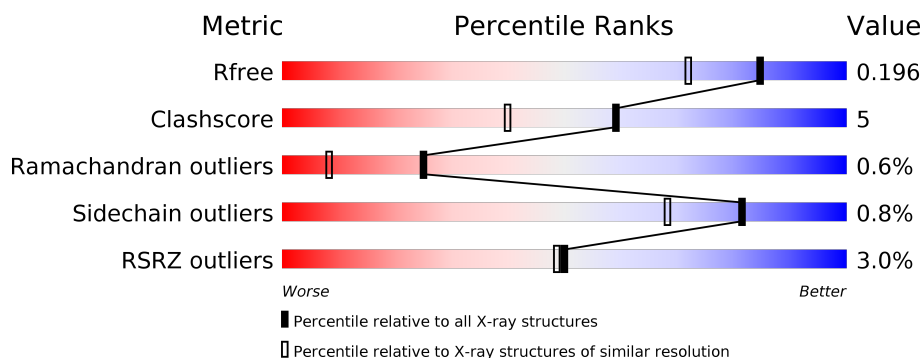
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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	3122 (1.66-1.62)
Clashscore	141614	3268 (1.66-1.62)
Ramachandran outliers	138981	3215 (1.66-1.62)
Sidechain outliers	138945	3215 (1.66-1.62)
RSRZ outliers	127900	3079 (1.66-1.62)

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	338	<div> <div>3%</div> <div>89%</div> <div>10%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
9	PO4	A	419[B]	-	X	-	-

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 3099 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.

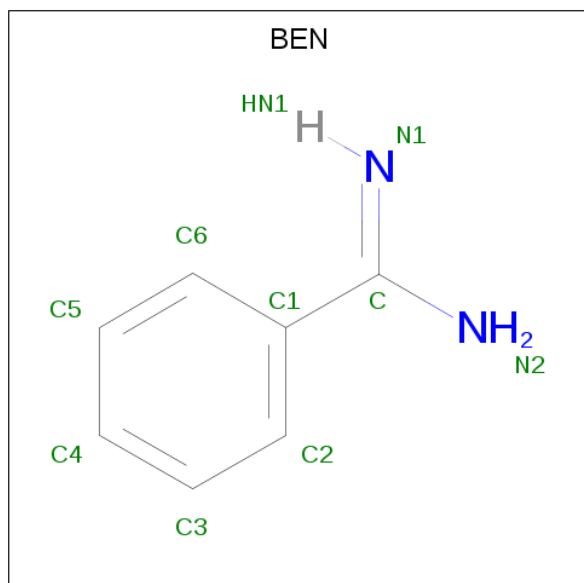
- Molecule 1 is a protein called Hydrogenase expression/formation protein HypE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	336	Total	C	N	O	S	0	15	0
			2613	1645	447	507	14			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	Mg	0	0
			4	4		

- Molecule 3 is BENZAMIDINE (three-letter code: BEN) (formula: C₇H₈N₂).



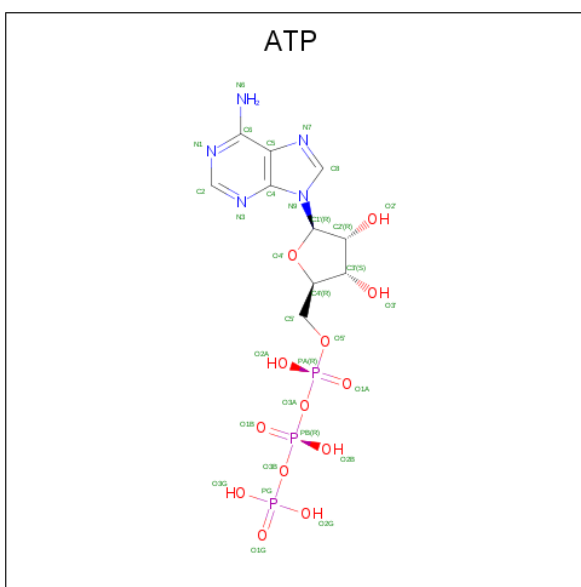
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			9	7	2		
3	A	1	Total	C	N	0	0
			9	7	2		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



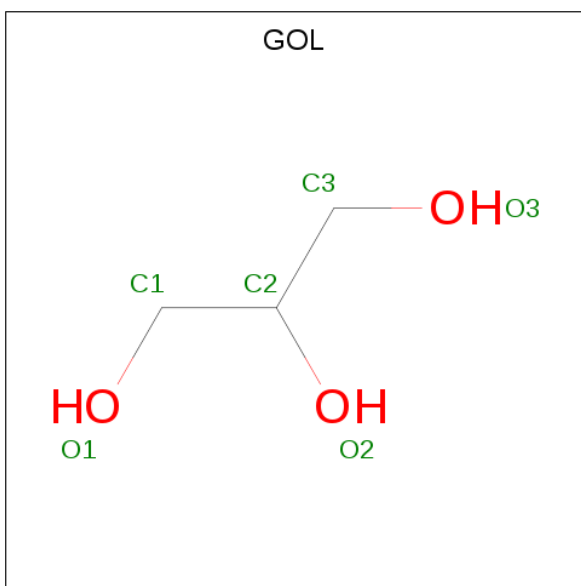
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



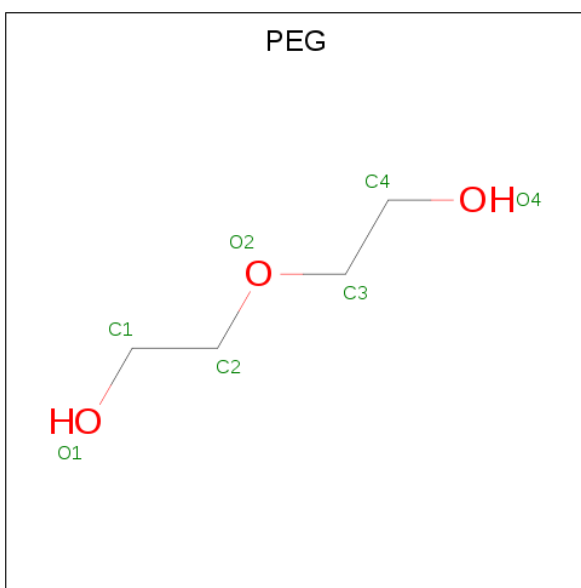
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	1
			31	10	5	13	3		

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $\text{C}_3\text{H}_8\text{O}_3$).



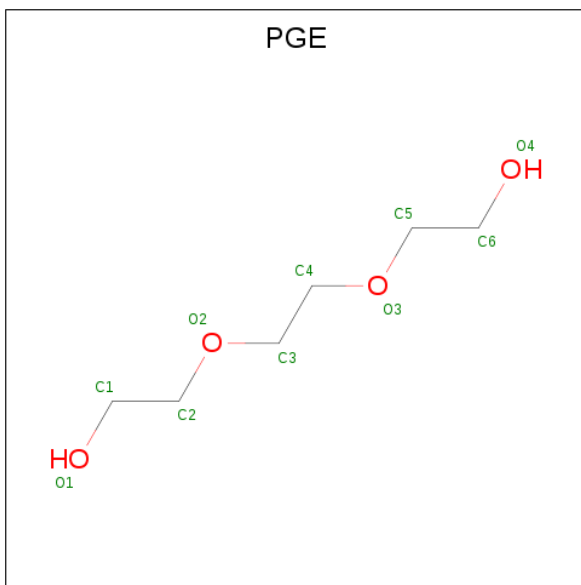
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total 6	C 3	O 3	0	0
6	A	1	Total 6	C 3	O 3	0	0
6	A	1	Total 6	C 3	O 3	0	0

- Molecule 7 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



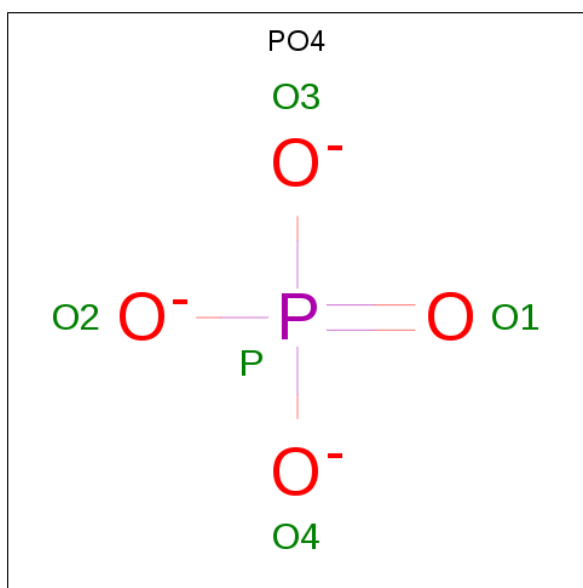
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 8 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			10	6	4		
8	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 9 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	O	P	0	1
			5	4	1		

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	339	Total	O	0	19
			358	358		

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.

- Chain A:
-
- 3% 89% 10%
- MET GLY E3 K4 G12 M16 R21 L38 D89 F49 K52 V55 F67 A90 E91 P92 I93 D106 R112 D118 E141 H154 L193 D196 K220 T223 R236 E246 R253 E254 Y255 R256 P266 E272 P282
- Y283 R298 D309 Y310 L316 E317 T318 G319 I320 G321 G322 K323 R324 F325 M326 C338

4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	102.85Å 102.85Å 104.31Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.38 – 1.65 46.52 – 1.64	Depositor EDS
% Data completeness (in resolution range)	99.8 (42.38-1.65) 99.8 (46.52-1.64)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.07 (at 1.64Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.175 , 0.196 0.175 , 0.196	Depositor DCC
R_{free} test set	3427 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	18.6	Xtriage
Anisotropy	0.571	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 53.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.011 for -h,l,k 0.005 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3099	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.76% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, MG, PGE, BEN, PO4, XCN, SO4, ATP, PEG

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	0.33	0/2634	0.54	0/3566

There are no bond length outliers.

There are no bond angle outliers.

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	2613	0	2635	24	1
2	A	4	0	0	0	0
3	A	18	0	14	0	0
4	A	25	0	0	1	0
5	A	31	0	12	1	0
6	A	18	0	24	3	0
7	A	7	0	10	2	0
8	A	20	0	28	4	0
9	A	5	0	0	0	0
10	A	358	0	0	6	0
All	All	3099	0	2723	26	1

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 5.

All (26) 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:21:ARG:HH21	1:A:236:ARG:HH22	1.42	0.67
1:A:220[B]:LYS:NZ	10:A:833:HOH:O	2.30	0.63
1:A:106[A]:ASP:OD2	10:A:741:HOH:O	2.16	0.59
1:A:67:PHE:HB3	7:A:416:PEG:H12	1.92	0.52
1:A:118:ASP:OD2	10:A:836:HOH:O	2.18	0.51
1:A:38:LEU:HD11	5:A:411[A]:ATP:N3	2.26	0.51
1:A:193:LEU:HD21	1:A:254:GLU:HB3	1.93	0.50
1:A:39:ASP:HB2	6:A:413:GOL:H11	1.93	0.49
8:A:417:PGE:H62	8:A:417:PGE:H22	1.95	0.48
1:A:193:LEU:HD21	1:A:254:GLU:CB	2.46	0.46
1:A:282:GLU:HG2	1:A:283:TYR:CD2	2.50	0.46
1:A:55:VAL:HG21	1:A:90:ALA:HB2	1.96	0.46
1:A:266:PRO:HD3	8:A:418:PGE:H22	1.97	0.45
1:A:272:GLU:OE2	1:A:338[A]:XCN:CS	2.65	0.45
1:A:252:ARG:HH21	6:A:415:GOL:H11	1.81	0.45
1:A:196:ASP:HB3	6:A:415:GOL:H12	1.99	0.45
1:A:92:PRO:HG3	10:A:835:HOH:O	2.17	0.45
1:A:298:ARG:NH2	4:A:406:SO4:O1	2.51	0.43
1:A:154[B]:HIS:HE1	10:A:590:HOH:O	2.01	0.43
1:A:316:LEU:HD21	1:A:326:MET:HE3	2.01	0.43
1:A:112[A]:ARG:HH22	7:A:416:PEG:H21	1.84	0.43
1:A:246:GLU:HB2	1:A:316:LEU:HD22	2.00	0.43
1:A:12:GLY:HA2	1:A:16:MET:CE	2.50	0.41
1:A:52:LYS:NZ	10:A:676[A]:HOH:O	2.39	0.41
8:A:418:PGE:O2	8:A:418:PGE:H52	2.21	0.41
1:A:256:ARG:HD2	8:A:417:PGE:H42	2.03	0.40

All (1) 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)
1:A:4:LYS:NZ	1:A:106[B]:ASP:OD1[7_555]	2.15	0.05

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	348/338 (103%)	342 (98%)	4 (1%)	2 (1%)	25 8

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	223	THR
1	A	141	GLU

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	272/264 (103%)	270 (99%)	2 (1%)	84 71

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	PHE
1	A	309	ASP

Some 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 ⓘ

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	XCN	A	338[A]	1	4,8,8	1.05	0	2,9,9	0.69	0
1	XCN	A	338[B]	1	4,6,8	5.38	1 (25%)	2,7,9	9.70	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	XCN	A	338[A]	1	-	0/3/8/8	-
1	XCN	A	338[B]	1	-	0/3/6/8	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	338[B]	XCN	CS-SG	10.59	2.29	1.68

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	338[B]	XCN	CB-SG-CS	-10.14	75.27	100.03
1	A	338[B]	XCN	SG-CS-NC	-9.24	134.32	175.93

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	338[A]	XCN	1	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 4 are monoatomic - leaving 15 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
4	SO4	A	409	-	4,4,4	0.14	0	6,6,6	0.07	0
3	BEN	A	412	-	9,9,9	1.44	1 (11%)	7,11,11	0.76	0
5	ATP	A	411[A]	2	26,33,33	0.90	1 (3%)	31,52,52	1.27	4 (12%)
4	SO4	A	406	-	4,4,4	0.12	0	6,6,6	0.06	0
6	GOL	A	414	-	5,5,5	0.48	0	5,5,5	0.23	0
6	GOL	A	413	-	5,5,5	0.52	0	5,5,5	0.67	0
7	PEG	A	416	-	6,6,6	0.53	0	5,5,5	0.71	0
6	GOL	A	415	-	5,5,5	0.41	0	5,5,5	0.55	0
4	SO4	A	408	-	4,4,4	0.14	0	6,6,6	0.07	0
8	PGE	A	418	-	9,9,9	0.65	0	8,8,8	0.83	0
8	PGE	A	417	-	9,9,9	0.59	0	8,8,8	0.74	0
4	SO4	A	410	-	4,4,4	0.13	0	6,6,6	0.06	0
4	SO4	A	407	-	4,4,4	0.13	0	6,6,6	0.07	0
9	PO4	A	419[B]	2	4,4,4	2.96	4 (100%)	6,6,6	0.42	0
3	BEN	A	405	-	9,9,9	1.34	1 (11%)	7,11,11	0.70	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
3	BEN	A	412	-	-	0/4/4/4	0/1/1/1
5	ATP	A	411[A]	2	-	1/18/38/38	0/3/3/3
3	BEN	A	405	-	-	0/4/4/4	0/1/1/1
6	GOL	A	414	-	-	4/4/4/4	-
6	GOL	A	413	-	-	4/4/4/4	-
7	PEG	A	416	-	-	2/4/4/4	-
6	GOL	A	415	-	-	2/4/4/4	-
8	PGE	A	418	-	-	3/7/7/7	-
8	PGE	A	417	-	-	2/7/7/7	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	419[B]	PO4	P-O1	4.50	1.61	1.50
3	A	412	BEN	C1-C	-3.96	1.40	1.47
3	A	405	BEN	C1-C	-3.69	1.40	1.47
9	A	419[B]	PO4	P-O3	2.31	1.61	1.54
9	A	419[B]	PO4	P-O2	2.26	1.61	1.54
5	A	411[A]	ATP	C5-C4	2.16	1.46	1.40
9	A	419[B]	PO4	P-O4	-2.06	1.48	1.54

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	411[A]	ATP	N3-C2-N1	-3.18	123.71	128.68
5	A	411[A]	ATP	PB-O3B-PG	-2.72	123.48	132.83
5	A	411[A]	ATP	C4-C5-N7	-2.40	106.90	109.40
5	A	411[A]	ATP	O2A-PA-O1A	2.21	123.17	112.24

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	414	GOL	O1-C1-C2-C3
6	A	414	GOL	C1-C2-C3-O3
6	A	413	GOL	O1-C1-C2-C3
6	A	413	GOL	C1-C2-C3-O3
8	A	418	PGE	O3-C5-C6-O4
6	A	414	GOL	O2-C2-C3-O3
8	A	417	PGE	O2-C3-C4-O3
6	A	414	GOL	O1-C1-C2-O2

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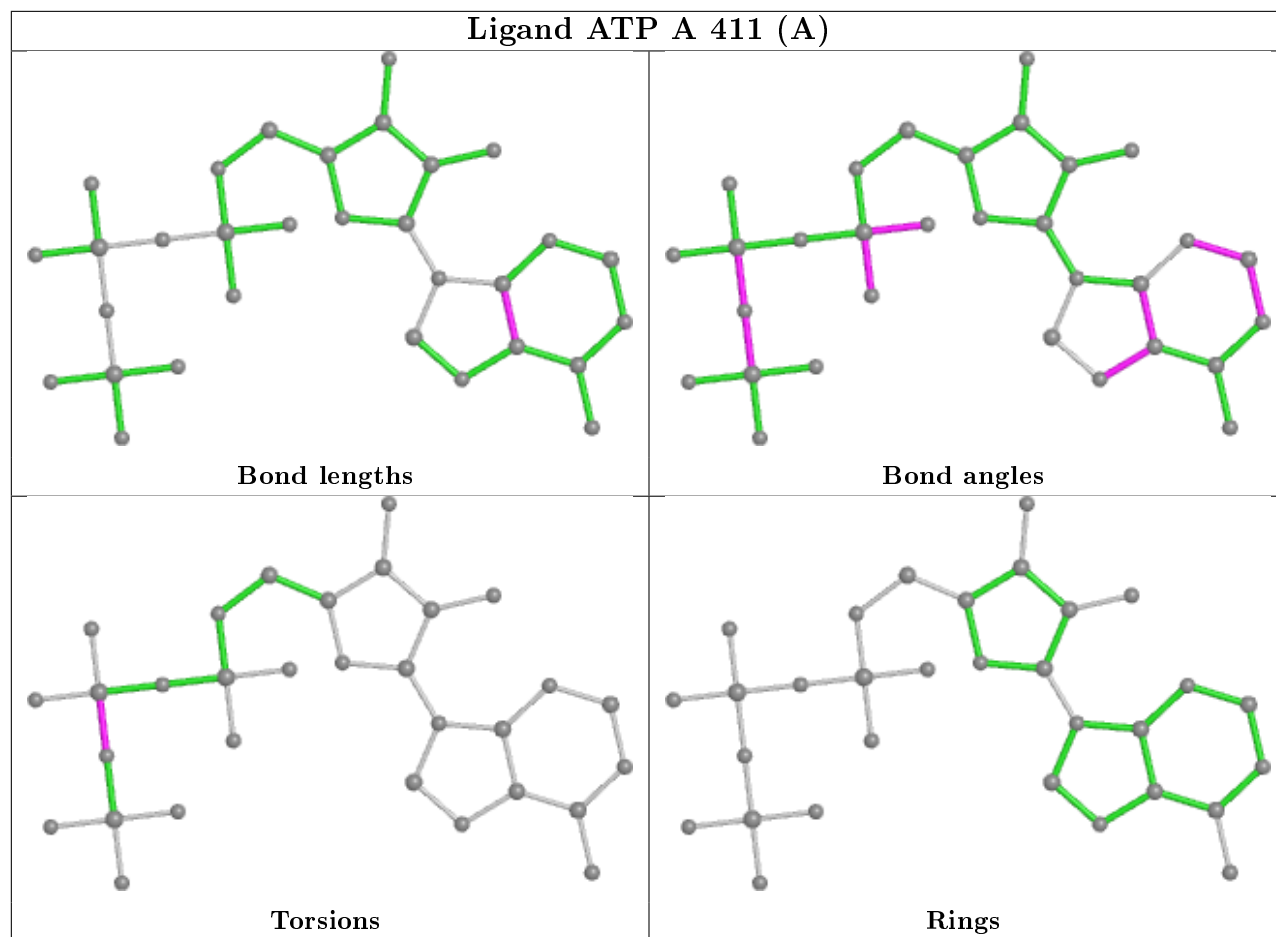
Mol	Chain	Res	Type	Atoms
6	A	413	GOL	O1-C1-C2-O2
6	A	413	GOL	O2-C2-C3-O3
8	A	417	PGE	O3-C5-C6-O4
7	A	416	PEG	C4-C3-O2-C2
6	A	415	GOL	O1-C1-C2-O2
8	A	418	PGE	O1-C1-C2-O2
5	A	411[A]	ATP	PG-O3B-PB-O1B
7	A	416	PEG	O2-C3-C4-O4
8	A	418	PGE	C1-C2-O2-C3
6	A	415	GOL	O2-C2-C3-O3

There are no ring outliers.

7 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	411[A]	ATP	1	0
4	A	406	SO4	1	0
6	A	413	GOL	1	0
7	A	416	PEG	2	0
6	A	415	GOL	2	0
8	A	418	PGE	2	0
8	A	417	PGE	2	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	335/338 (99%)	0.01	10 (2%) 50 48	10, 19, 36, 46	1 (0%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	319	GLY	6.4
1	A	320	ILE	5.9
1	A	318	THR	5.0
1	A	321	GLY	3.4
1	A	324	ARG	2.3
1	A	55	VAL	2.2
1	A	93[A]	ILE	2.2
1	A	310	TYR	2.1
1	A	323	LYS	2.1
1	A	316	LEU	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(Å ²)	Q<0.9
1	XCN	A	338[A]	9/9	0.97	0.08	15,17,25,27	9
1	XCN	A	338[B]	7/9	0.97	0.08	14,16,20,22	7

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

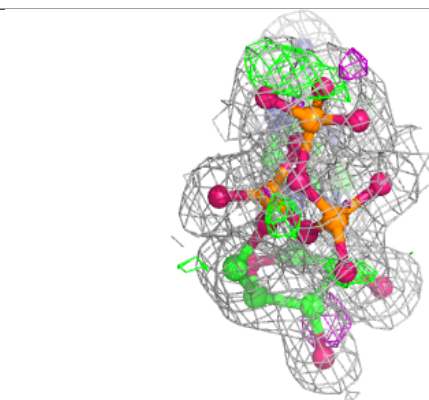
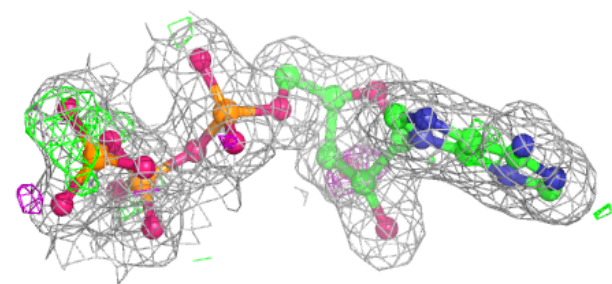
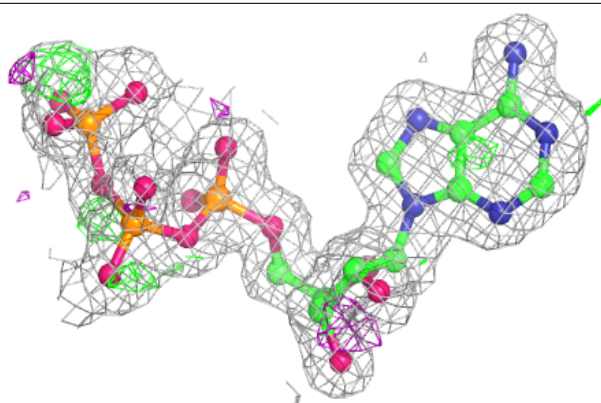
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	PGE	A	417	10/10	0.67	0.32	19,26,30,32	10
6	GOL	A	415	6/6	0.71	0.32	26,35,37,38	0
6	GOL	A	414	6/6	0.81	0.33	25,41,51,61	0
8	PGE	A	418	10/10	0.81	0.21	30,35,36,42	10
3	BEN	A	412	9/9	0.82	0.13	25,26,35,37	0
7	PEG	A	416	7/7	0.83	0.29	37,45,57,61	0
9	PO4	A	419[B]	5/5	0.87	0.25	18,19,23,27	5
4	SO4	A	408	5/5	0.89	0.40	50,58,70,77	0
4	SO4	A	407	5/5	0.90	0.28	48,53,60,73	0
4	SO4	A	409	5/5	0.91	0.20	38,58,59,76	0
6	GOL	A	413	6/6	0.91	0.21	20,33,36,45	0
2	MG	A	404	1/1	0.94	0.07	18,18,18,18	1
4	SO4	A	410	5/5	0.94	0.31	63,66,72,76	0
5	ATP	A	411[A]	31/31	0.95	0.08	11,15,20,23	8
3	BEN	A	405	9/9	0.96	0.09	15,17,20,20	0
4	SO4	A	406	5/5	0.97	0.18	47,58,59,64	0
2	MG	A	402	1/1	0.98	0.06	12,12,12,12	1
2	MG	A	403	1/1	0.98	0.04	15,15,15,15	1
2	MG	A	401	1/1	0.99	0.04	14,14,14,14	1

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 ATP A 411 (A):

$2mF_o-DF_c$ (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.