



BERRY&ASSOCIATES

Dedicated to providing
specialty chemicals
to advance the **life sciences**

2012 NEW PRODUCT LIST

General Information

Welcome to our 2012 New Releases listing from Berry & Associates. Since our 2009 Full Catalog, we have added over 100 compounds to our lineup. Our product offerings have expanded further to include our Click-mates™ collection for biomolecule conjugation as well as our Click-easy™ collection of alkynes for copper-free click ligation. In addition, we also now offer a family of 5-hydroxymethyl cytidine tools for epigenetic research. This supplement includes new products in the following areas:

Blackberry® 650 Quenchers

- Non-fluorescent, stable quenchers of long-wavelength fluorophores
- Excellent performance for both FRET-mode and contact-mode quenching

BerrySelect® Reagents for DNA and RNA Synthesis

- Specialty Phosphoramidites
- Derivatized CPGs
- Functional Modifiers
- Fluorescent Building Blocks
- Quenchers of Fluorescence
- Spacers
- Carrier and coenzyme labels

Nucleosides

- One of the most comprehensive collections of nucleosides available
- Many rare nucleosides in stock in gram quantities

Specialty Chemicals

- Aminoxy Linkers and Spacers
- TEG-Azides for click chemistry
- Cyclooctynes for copper-free Click Chemistry

At Berry & Associates, we also have extensive experience with R & D collaborations in the nucleoside and phosphoramidite arena as well as with general Life Science applications. Please feel free to contact us with any of your research needs.

Ordering Information

Orders may be placed by phone, fax or e-mail:

Phone:	800.357.1145 734.426.3787
Fax:	734.426.9077
Email:	orders@berryassoc.com techhelp@berryassoc.com
Web:	www.berryassoc.com
Production, Research and Development:	Berry & Associates, Inc. 2434 Bishop Circle East Dexter, MI 48130 USA

Berry & Associates, Inc. was established in 1988 to provide the best and most comprehensive source of nucleoside analogs for researchers in the life sciences. The company is located outside of Ann Arbor, Michigan (USA) and features state-of-the-art laboratories, facilitating the production of high quality products. The credentials of our highly trained staff of chemists include over 400 publications and 80 patents in synthetic organic and medicinal chemistry. High quality chemicals, timeliness, and personalized service are the hallmarks of Berry & Associates.

Click-easy™ Alkynes for Cu-Free Click Conjugation

Since its introduction in 2002, the click reaction has become an immensely valuable tool spanning many facets of research from surface science to biomolecules. This cycloaddition reaction is remarkably efficient and reliable even in the presence of the diverse array of functional groups found in DNA and the variety of scientifically intriguing ligation partners. For seminal reading on the evolution of the use of this 1,3-dipolar cycloaddition in bioconjugation and bioorthogonal labeling, see references 1–5. Development of the click methodology has led to the production and study of many modified DNA oligonucleotides (ODNs) for biological as well as nanotechnological and surface oriented applications. The beauty of the click cycloaddition lies in the efficient formation of a non-toxic triazole on biological building blocks that have been modified with non-perturbing azides and unactivated alkynes. However, the requisite Cu(I) catalyst can degrade oligonucleotides⁶ and can compromise cell function, thereby limiting the utility of the Cu-catalysed click reaction for these purposes. As a result, a large number of cyclooctynes have been developed to capitalize upon the unique nature of a ring-bound alkyne for strain-promoted alkyne-azide cycloaddition (SPAAC).⁷ New to our catalog are the bicyclo[6.1.0]nonyne derivatives (BCNs) for copper-free click reactions.⁸ (Figure 1)

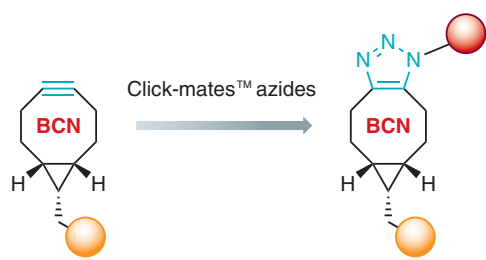


Figure 1. BCNs for copper-free clicking

In one illustrative example, 5'-Click-easy™ BCN CEP II (BA 0373) is incorporated at the 5'-terminus of an oligo with > 99% efficiency. Following a diethylamine wash, standard cleavage, and salt switch, a triethylamine salt of the T6 oligo was taken up in a buffer/acetonitrile solution and treated with PQQ-TEG azide with nearly complete consumption of oligo within 26 hours (Figure 2). When the same BCN-oligo was treated with Desthiobiotin-TEG azide (BT 1075) in approximately 100-fold molar excess in the same reaction volume, the reaction was nearly complete in just 30 minutes (Figure 3).

We have endeavored to prepare a wide variety of click ligation tools to facilitate your research. In addition to the Click-easy™ alkynes for copper free click cycloadditions, we also offer a variety of standard alkynes in our Click-mates™ collection. Naturally, the alkynes are used in conjunction with biologically relevant azides, so we also offer a large selection of Click-mates™ azides

Please consult our website (www.berryassoc.com) for the most current listing of products and prices for our entire collection of click ligation tools.

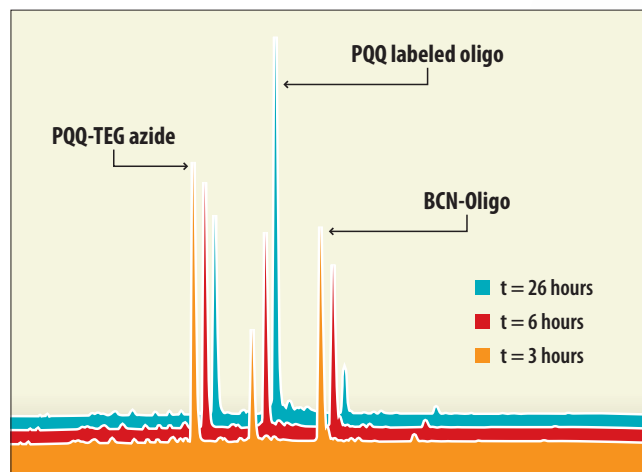


Figure 2. BCN-Oligo with PQQ-TEG azide (FC 8170)

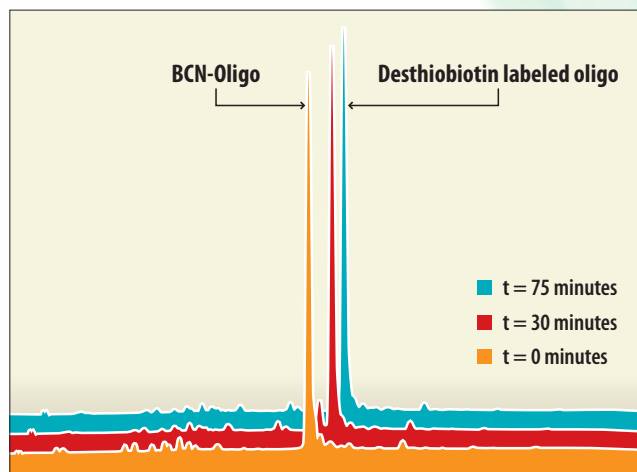


Figure 3. BCN-Oligo with excess Desthiobiotin-TEG azide (BT 1075)

1. V. V. Rostovtsev, L. G. Green, V. V. Fokin and K. B. Sharpless, *Angew. Chem., Int. Ed.*, **2002**, *41*, 2596–2599.
2. C.W. Tornøe, C. Christensen and M. Meldal, *J. Org. Chem.*, **2002**, *67*, 3057–3064.
3. H. C. Kolb, M. G. Finn and K. B. Sharpless, *Angew. Chem., Int. Ed.*, **2001**, *40*, 2004–2021.
4. P. M. E. Gramlich, C. T. Wirges, A. Manetto and T. Carell, *Angew. Chem., Int. Ed.*, **2008**, *47*, 8350–8358.
5. (a) R. Huisgen, *Proc. Chem. Soc., London*, **1961**, 357–369. (b) R. Huisgen, *Angew. Chem., Int. Ed. Engl.*, **1963**, *2*, 565–632. (c) R. Huisgen, *1,3-Dipolar Cycloadditional Chemistry*, Wiley, New York, 1984.
6. Gierlich, J.; Burley, G. A.; Gramlich, P. M. E.; Hammond, D. M.; Carell, T. *Org. Lett.* **2006**, *8*, 3639–3642.
7. Agard, N.J.; Prescher, J.A.; Bertozzi, C.R. *J. Am. Chem. Soc.* **2004**, *126*, 15046–15047.
8. a) Dommerholt, J.; Schmidt, S.; Temming, R.; Hendricks, L.J.A.; Rutjes, F.P.J.T.; van Hest, J.C.M.; Lefebvre, D.J.; Friedl, P.; van Delft, F.L.; Debets, M.F.; van Berkel, S.S.; Dommerholt, J.; Dirks, A.J.; Rutjes, F.P.J.T.; van Delft, F.L. *Accounts of Chem. Res.* **2011**, *44*, 805–815.

5-Hydroxymethyl-2'-deoxycytidine Tools for Epigenetic Research

The 5-methylation of 2'-deoxycytidine is an important and well studied modification of DNA. This methylation does not impact base pairing, but it alters DNA in ways that affect the binding of transcription factors and subsequent gene expression. Therefore, cytosine methylation is an important epigenetic marker.¹ The exact function of this new "sixth base" is the subject of much current research. As shown in Figure 1, DNA methyltransferases (DNMT) catalyze the *in vivo* conversion of dC to mC and subsequent oxidation by ten-to-eleven translocation oxidases (TET) affords hmC. It has been further postulated that a complete cycle involving further oxidation to fC and then to caC followed by decarboxylation to dC could exist *in vivo*.² Recently, fC has been discovered in embryonic stem cells with concentrations decreasing as cell differentiation proceeds.³

To facilitate research in this area, Berry & Associates offers a collection of 5hmC derivatives including the nucleoside standards 5-Hydroxymethylcytidine (PY 7596) and 5-Hydroxymethyl-2'-deoxycytidine (PY 7588), and a variety of tools for investigation of the proposed methylation cycle.

Please consult our website (www.berryassoc.com) for the most current listing of products and prices.

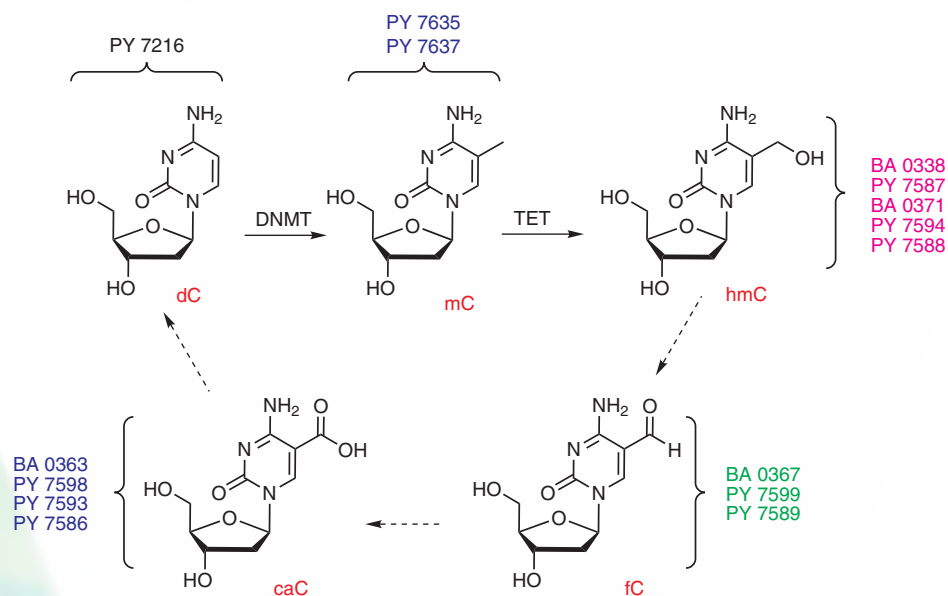
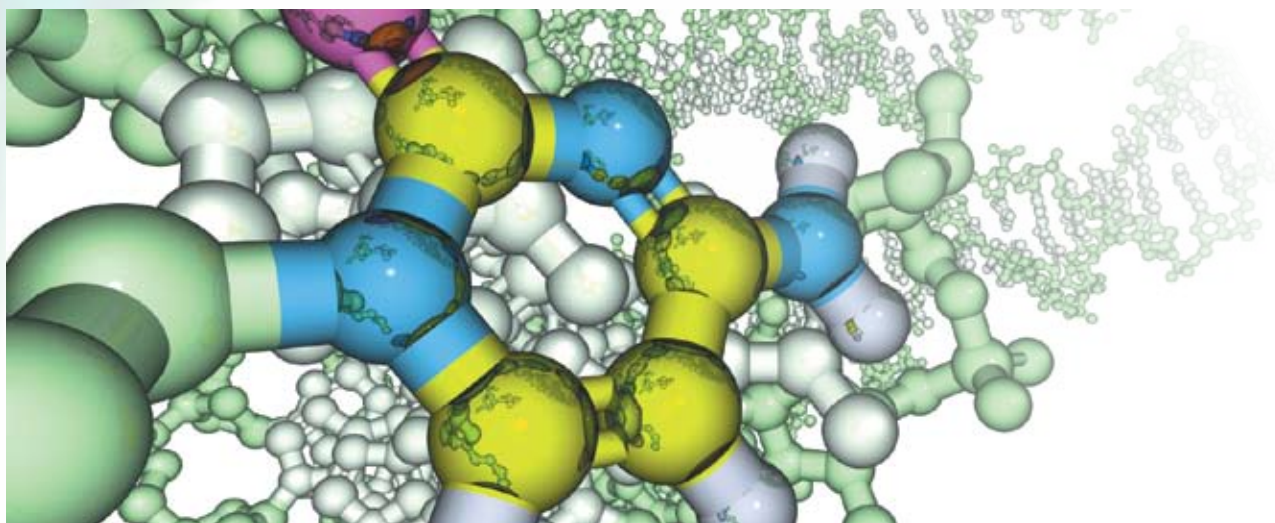
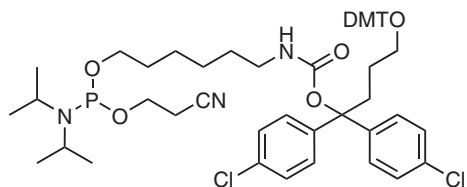


Figure 1. Cytosine methylation cycle and related Berry & Associates Products

1. a) Kriaucionis, S.; Heintz, N. *Science*, **2009**, 324, 929-930. b) Tahiliani, M.; Koh, K.P.; Shen, Y.H.; Pastor, W.A.; Bandukwala, H.; Brundo, Y.; Agrawal, S.; Iyer, I.M.; Liu, D.R.; Aravind, L.; Rao, A. *Science*, **2009**, 324, 930-935. c) Munzel, M.; Globisch, D.; Bruckl, T.; Wagner, M.; Welzmler, V.; Michalak, S.; Muller, M.; Biel, M.; Carell, T. *Angew. Chem. Int. Ed.*, **2010**, 49, 5375-5377. d) Szwagierczak, A.; Bultmann, S.; Schmidt, C.S.; Spada, F.; Leonhardt, H. *Nucleic Acids Res.* **2010**, 38, e181.
2. Munzel, M.; Globisch, D.; Carell, T. *Angew. Chem. Int. Ed.* **2010**, 50, 6460-6468.
3. Pfaffeneder, T.; Hackner, B.; Trub, M.; Munzel, M.; Muller, M.; Deiml, C.A.; Hagemeyer, C.; Carell, T. *Angew. Chem. Int. Ed.* **2011**, 50, 7008-7012.



BA 0324

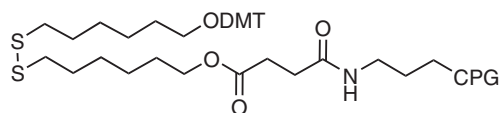


Cmoc-5'-amino-modifier-C6 CEP

MF:	C₅₃H₆₄Cl₂N₃O₇P	100 μmol	\$420
MW:	956.97	0.25 g	\$685
CAS No:	1246448-29-2		

Novel acid labile *N*-protected 5'-amino modifier.

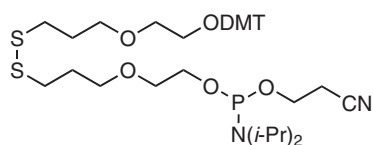
BA 0330



3'-Thiol-modifier-C6-S-S CPG (1000 Å)

MF:	N/A	100 mg	\$85
MW:	N/A	1 g	\$600
CAS No:	None Assigned		

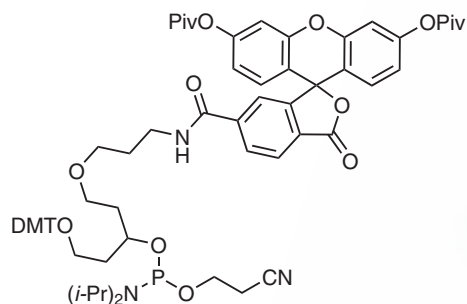
BA 0332



Thiol-modifier-oxa-C6-S-S CEP

MF:	C₄₀H₅₇N₂O₇PS₂	100 μmol	\$145
MW:	772.99	250 mg	\$360
CAS No:	1159976-43-8		

BA 0334

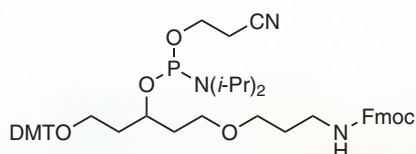


Fluorescein III CEP

MF:	C₆₉H₈₀N₃O₁₄P	100 μmol	\$380
MW:	1206.33	0.25 g	\$605
CAS No:	1306615-53-1		

Installs a 6-carboxyfluorescein internally or at the 5'-terminus of an oligonucleotide using a DMT-bearing phosphoramidite with a 1,3-diol framework.

BA 0335

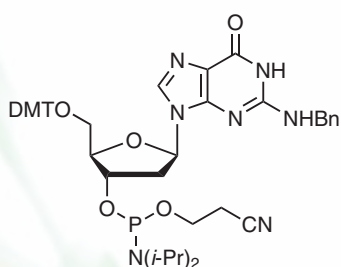


Fmoc-amino-modifier III CEP

MF:	$C_{53}H_{64}N_3O_8P$	100 μ mol	\$420
MW:	902.06	0.25 g	\$685
CAS No:	1306615-50-8		

For installation of an Fmoc protected amino group internally or at the 5'-end of an oligonucleotide.

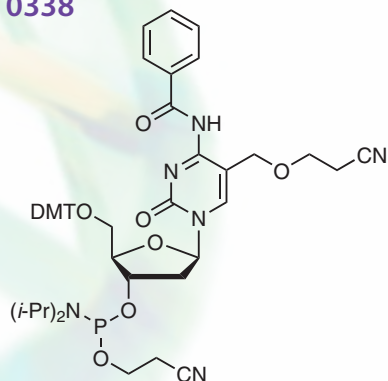
BA 0337

 N^2 -Benzyl-2-deoxyguanosine CEP

MF:	$C_{47}H_{54}N_7O_7P$	100 μ mol	\$425
MW:	859.95	250 mg	\$975
CAS No:	209785-74-0		

Useful for probing the steric requirements at N2 of dG in various applications.

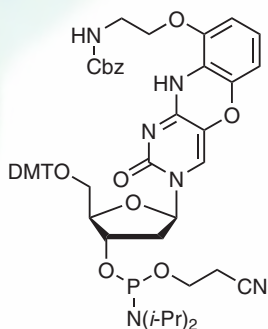
BA 0338



N-Benzoyl-5-((2-cyanoethoxy)methyl)-2'-deoxycytidine CEP

MF:	$C_{50}H_{57}N_6O_9P$	15 g minimum order	
MW:	917.00	Call for pricing	
CAS No:	188411-06-5		

BA 0339

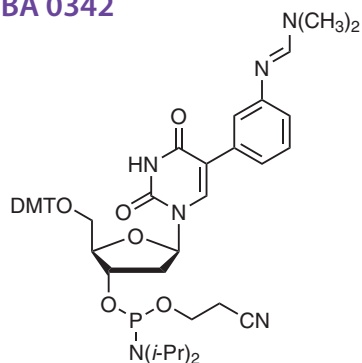


8-OxoG clamp CEP

MF:	$C_{55}H_{61}N_6O_{11}P$	100 μ mol	\$505
MW:	1013.08	0.25 g	\$958
CAS No:	1134373-47-9		

Fluorescent molecule for the selective recognition of 8-oxoG.

BA 0342

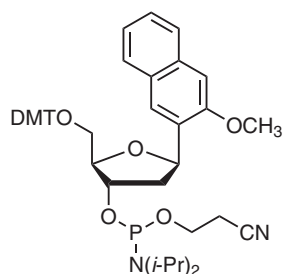


5-(3-Aminophenyl)-2'-dU CEP

MF: **C₄₈H₅₇N₆O₈P** 100 μ mol \$375
 MW: **876.98**
 CAS No: **None Assigned**

Simple organic 2'-deoxyribonucleoside derivatives for use as electroactive DNA markers.

BA 0343

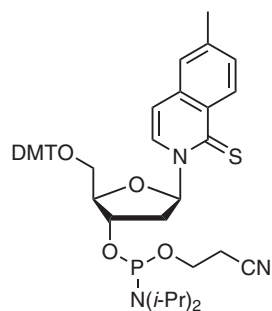


dNaM CEP

MF: **C₄₆H₅₃N₂O₇P** 100 μ mol \$415
 MW: **776.98** 0.25 g \$995
 CAS No: **1117893-10-3**

Half of a novel base pair that achieves pair recognition through hydrophobic interactions. See also d5SICS CEP (BA 0344).

BA 0344

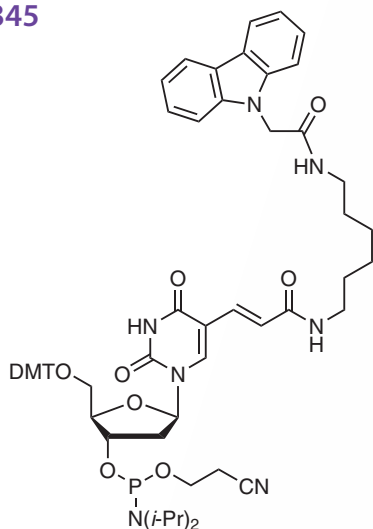


d5SICS CEP

MF: **C₄₅H₅₂N₃O₆PS** 100 μ mol \$918
 MW: **793.95** 0.25 g \$2,225
 CAS No: **1010689-04-9**

Half of a novel base pair that achieves pair recognition through hydrophobic interactions. See also dNaM CEP (BA 0343).

BA 0345

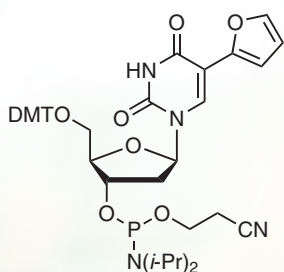


Carbazole dT CEP

MF: **C₆₂H₇₂N₇O₁₀P** 100 μ mol \$195
 MW: **1106.26** 0.25 g \$335
 CAS No: **None Assigned**

For use as a light-controlled, reversible DNA photoligation tool.

BA 0346

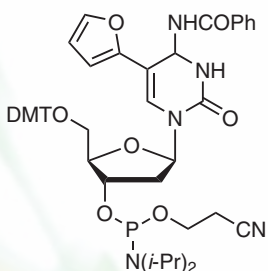


5-(2-Furyl)-dU CEP

MF:	$C_{43}H_{49}N_4O_9P$	100 μ mol	\$245
MW:	796.86	0.25 g	\$580
CAS No:	863713-51-3		

Small, fluorescent natural base mimic that can signal the presence of a basic sites in hybridized DNA oligonucleotides.

BA 0347

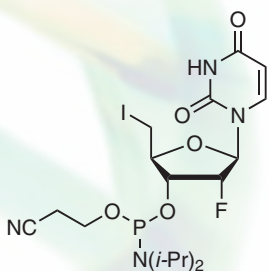


5-(Furan-2-yl)-dC CEP

MF:	$C_{50}H_{54}N_5O_9P$	100 μ mol	\$245
MW:	899.97	0.25 g	\$580
CAS No:	1119734-65-4		

Small, fluorescent natural base mimic that can be used as an *in vitro* signal the presence of G, 8-oxoG, or T on a complementary strand in hybridized DNA oligonucleotides.

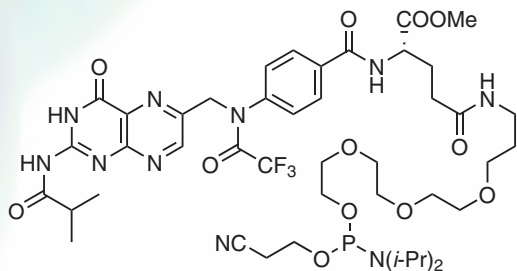
BA 0348



2'-Fluoro-5'-iodo-deoxyuridine CEP

MF:	$C_{18}H_{27}FIN_4O_5P$	100 μ mol	\$240
MW:	556.31	0.25 g	\$625
CAS No:	None Assigned		

BA 0349

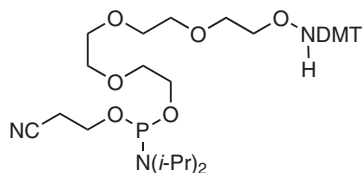


5'-Folate-TEG CEP

MF:	$C_{44}H_{62}F_3N_{10}O_{12}P$	50 μ mol	\$548
MW:	1010.99	100 μ mol	\$975
CAS No:	1356964-22-1		

Useful for the installation of a folate tag for recognition by overexpressed folate receptors in cancer cells.

BA 0350

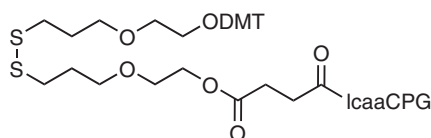


5'-Aminoxy-modifier-11 CEP

MF:	C₃₈H₅₄N₃O₈P	100 μmol	\$235
MW:	711.82	0.25 g	\$785
CAS No:	None Assigned	1 g	\$2,500

For the synthesis of oligonucleotides bearing a 5'-aminoxy group.

BA 0351

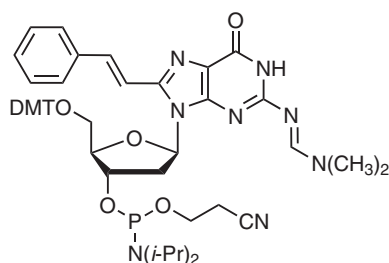


3'-Thiol modifier-oxa-C6-S-S CPG (1000 Å)

MF:	Unknown	200 nmol columns (pkg of 4)	\$69
MW:	Unknown	100 mg	\$70
CAS No:	None Assigned	1 μmol columns (pkg of 4)	\$115
		1 g	\$500

Superior thiol modifier giving higher yields and longer oligos.

BA 0352

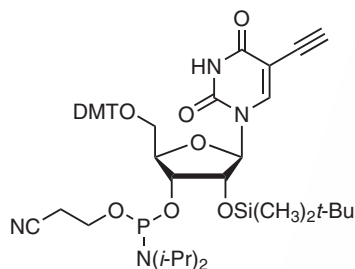


8-Styryl-dG CEP

MF:	C₅₁H₅₉N₈O₇P	100 μmol	\$752
MW:	927.04	0.25 g	\$1,560
CAS No:	1101864-12-3		

Photochromic nucleoside (PCN) that provides reversible duplex regulation via a light induced trans-cis isomerization.

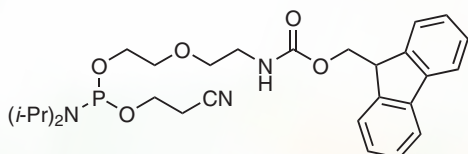
BA 0353



5-Ethynyl uridine CEP

MF:	C₄₇H₆₁N₄O₉PSi	100 μmol	\$445
MW:	885.07	0.25 g	\$960
CAS No:	1193451-06-7		

BA 0354

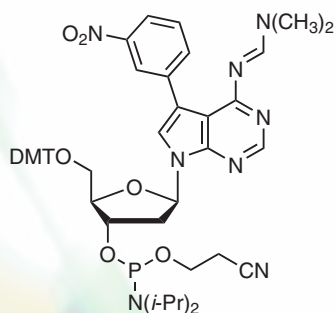


Fmoc-5'-amino-modifier-5 CEP

MF:	$C_{28}H_{38}N_3O_5P$	100 μ mol	\$95
MW:	527.59	0.25 g	\$190
CAS No:	1306615-52-0		

For 5'-incorporation of an amine that can be deprotected prior to cleaving the oligonucleotide from the solid support.

BA 0355

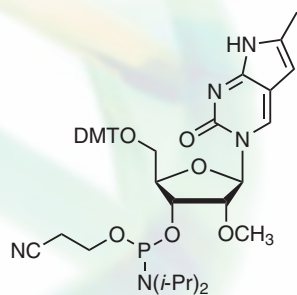


7-(3-Nitrophenyl)-7-deaza-2'-dA CEP

MF:	$C_{50}H_{57}N_8O_8P$	100 μ mol	\$315
MW:	929.01	0.25 g	\$595.55
CAS No:	None Assigned		

Simple organic 2'-deoxyribonucleoside derivatives for use as electroactive DNA markers.

BA 0356

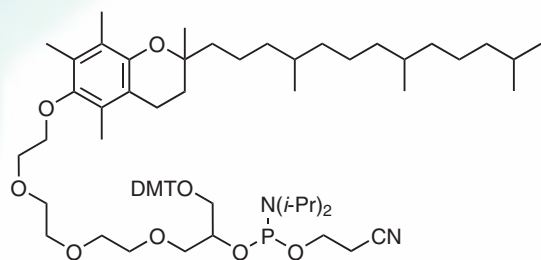


2'-O-Methyl-pyrrolo C CEP

MF:	$C_{43}H_{52}N_5O_8P$	100 μ mol	\$395
MW:	797.88	0.25 g	\$875
CAS No:	644962-95-8		

Pyrrolo-C (PC) is a fluorescent analog of cytidine.

BA 0357

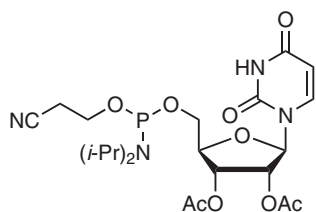


Tocopherol-TEG CEP

MF:	$C_{68}H_{103}N_2O_{10}P$	10 g minimum order	
MW:	1139.56	Call for pricing	
CAS No:	None Assigned		

Lipophilic carrier tag for internal or 5'-terminus installation in oligonucleotides.

BA 0358

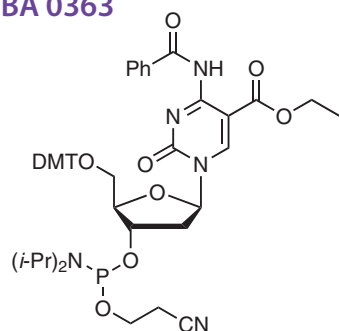


2',3'-Di-O-acetyl-U-5'-CEP

MF:	C₂₂H₃₃N₄O₉P	100 μmol	\$155
MW:	528.49	0.25 g	\$465
CAS No:	208655-84-9		

Used for the preparation of uridylated amino acids. See also BX 00004.

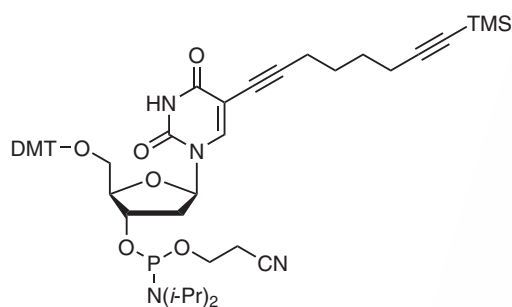
BA 0363



5-Carboethoxy dC CEP

MF:	C₄₉H₅₆N₅O₁₀P	5 g minimum order
MW:	905.97	Call for pricing
CAS No:	None Assigned	

BA 0364



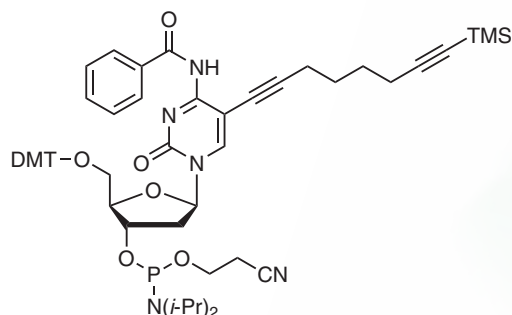
5-Octadiynyl-TMS-dU CEP

MF:	C₅₀H₆₃N₄O₈PSi	2 g minimum order
MW:	907.12	Call for pricing
CAS No:	None Assigned	

A useful diyne from our Click-mates™ collection for allowing multiple varied click reactions in a single oligo.

This compound is sold under license from baseclick GmbH, and the purchase of these products for use in applications relating to copper catalyzed azide-alkyne cycloaddition chemistry ("Click Chemistry") includes a limited, nontransferable license to intellectual property owned by TSRI to use this product solely for internal non-commercial research activities and specifically excludes clinical, therapeutic, or diagnostic use in humans or animals. Information regarding a license for commercial use in Click Chemistry may be obtained directly from The Scripps Research Institute, 10550 N. Torrey Pines Rd., La Jolla, CA 92037, or by contacting 858-784-8140 or click@scripps.edu.

BA 0365



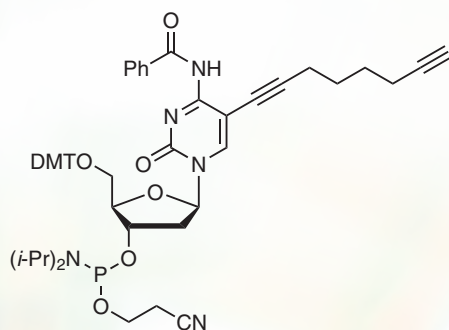
5-Octadiynyl-TMS-dC CEP

MF:	C₅₇H₆₈N₅O₈PSi	2 g minimum order
MW:	1010.24	Call for pricing
CAS No:	1021301-02-9	

A useful diyne from our Click-mates™ collection for multiple, varied click reactions in a single oligo.

This compound is sold under license from baseclick GmbH, and the purchase of these products for use in applications relating to copper catalyzed azide-alkyne cycloaddition chemistry ("Click Chemistry") includes a limited, nontransferable license to intellectual property owned by TSRI to use this product solely for internal non-commercial research activities and specifically excludes clinical, therapeutic, or diagnostic use in humans or animals. Information regarding a license for commercial use in Click Chemistry may be obtained directly from The Scripps Research Institute, 10550 N. Torrey Pines Rd., La Jolla, CA 92037, or by contacting 858-784-8140 or click@scripps.edu.

BA 0366



5-Octadiynyl-dC CEP

MF: $C_{54}H_{60}N_5O_8P$

MW: 938.06

CAS No: 1021300-97-9

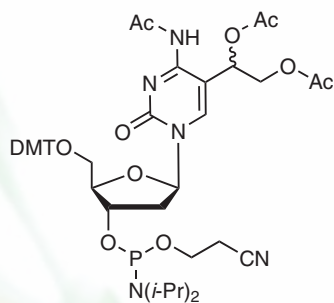
2 g minimum order

Call for pricing

A useful diyne from our Click-mates™ collection for multiple, varied click reactions in a single oligo.

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BA 0367



Masked 5-formyl-dC CEP

MF: $C_{47}H_{58}N_5O_{12}P$

MW: 915.96

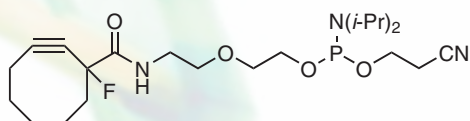
CAS No: 364613-34-3

50 μmol \$640

100 μmol \$995

Useful for investigating the role of cytidine methylation in control of gene expression.

BA 0368



5'-Click-easy™ MFCO CEP

MF: $C_{22}H_{37}FN_3O_4P$

MW: 457.52

CAS No: 1306615-49-5

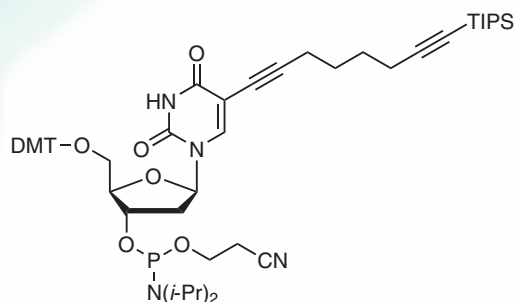
50 μmol \$325

100 μmol \$590

BA 0368 from our Click-mates™ collection is for the installation of a monofluoro-substituted cyclooctyne (MFCO) into an oligonucleotide for subsequent copper-free click elaboration.

BA 0368 is protected by US 7,807,619, the rights of which are assigned to The Regents of the University of California. Additional patents covering methods for their use in the modification of biomolecules are pending. It may be used for research purposes only. It is not licensed for resale and may only be used by the buyer. This product may not be used and are not licensed for clinical assays, where the results of such assays are provided as a diagnostic service. If a diagnostic or therapeutic use is anticipated, then a license must be requested from the University of California. The availability of such diagnostic and therapeutic use license(s) cannot be guaranteed from the University of California.

BA 0369



5-Octadiynyl-TIPS-dU CEP

MF: $C_{56}H_{75}N_4O_8PSi$

MW: 991.28

CAS No: None Assigned

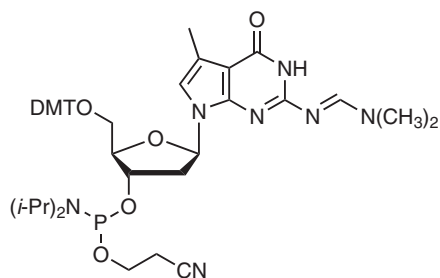
2 g minimum order

Call for pricing

A useful diyne from our Click-mates™ collection for multiple, varied click reactions in a single oligo.

This compound is sold under license from baseclick GmbH, and the purchase of these products for use in applications relating to copper catalyzed azide-alkyne cycloaddition chemistry ("Click Chemistry") includes a limited, nontransferable license to intellectual property owned by TSRI to use this product solely for internal non-commercial research activities and specifically excludes clinical, therapeutic, or diagnostic use in humans or animals. Information regarding a license for commercial use in Click Chemistry may be obtained directly from The Scripps Research Institute, 10550 N. Torrey Pines Rd., La Jolla, CA 92037, or by contacting 858-784-8140 or click@scripps.edu.

BA 0370

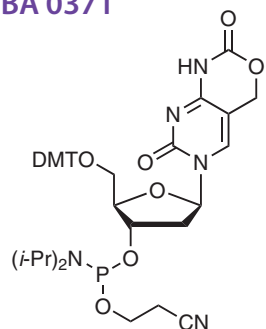


7-Deaza-7-methyl-2'-dG CEP

MF:	C₄₅H₅₆N₇O₇P	100 μmol	\$725
MW:	837.94	250 mg	\$2120
CAS No:	None Assigned		

BA 0370 is a useful tool for investigations of duplex stabilization as a result of substituents on the purine scaffold.

BA 0371

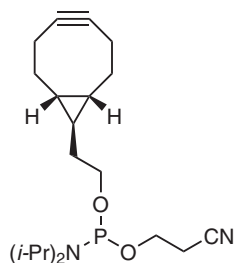


5-Hydroxymethyl-dC cyclic carbamate CEP

MF:	C₄₁H₄₈N₅O₉P	50 μmol	\$540
MW:	785.82	100 μmol	\$825
CAS No:	1257247-42-9		

Useful for investigating the role of hydroxymethyl cytidine in control of gene expression.

BA 0372



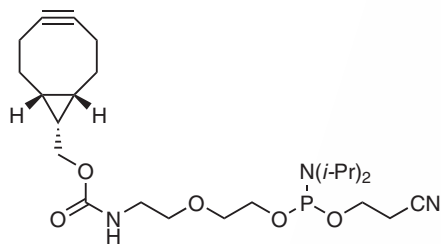
5'-Click-easy™ BCN CEP I

MF:	C₂₂H₃₃N₂O₂P	100 μmol	\$450
MW:	364.46	250 μmol	\$990
CAS No:	None Assigned		

BA 0372 is for the installation of a strained cyclooctyne (BCN) into an oligonucleotide for subsequent copper-free click elaboration.

Patent Pending.

BA 0373



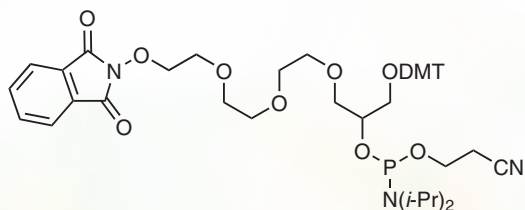
5'-Click-easy™ BCN CEP II

MF:	C₂₄H₄₀N₃O₅P	100 μmol	\$499
MW:	481.57	250 μmol	\$1100
CAS No:	1352811-59-6		

BA 0373 is for the installation of a strained cyclooctyne (BCN) into an oligonucleotide for subsequent copper-free click elaboration.

Patent Pending.

BA 0374

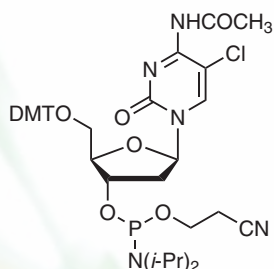


Aminoxy-modifier CEP

MF:	$C_{47}H_{58}N_3O_{11}P$	100 μ mol	\$260
MW:	871.95	0.25 g	\$865
CAS No:	None Assigned		

For the synthesis of oligonucleotides bearing an aminoxy moiety.

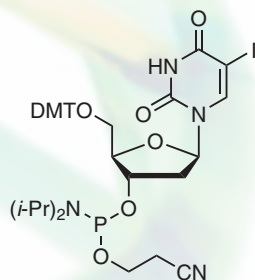
BA 0375



5-Chloro-dC CEP

MF:	$C_{41}H_{49}ClN_5O_8P$	100 μ mol	\$285
MW:	806.28	0.25 g	\$670
CAS No:	1356140-67-4		

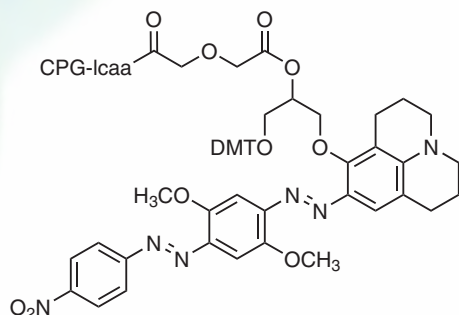
BA 0376



5-Iodo-dU CEP

MF:	$C_{39}H_{46}IN_4O_8P$	0.25 g	\$160
MW:	856.68	1 g	\$485
CAS No:	178925-48-9		

BL 2020

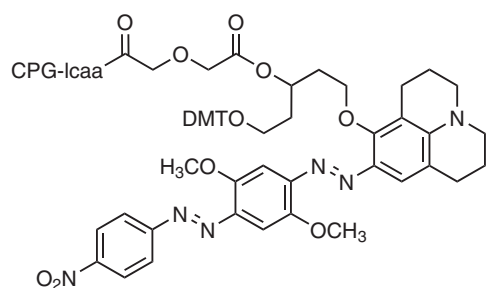


3'-BBQ-650® CPG II (1000 Å)

MF:	N/A	200 nmol columns (pkg of 4)	\$75
MW:	N/A	100 mg	\$135
CAS No:	None Assigned	1 μ mol columns (pkg of 4)	\$210
		1 g	\$1,050

BlackBerry® Quencher 650 CPG II (3'-BBQ-650® CPG II) is used to install BBQ-650® at the 3' end of an oligonucleotide, and is more rapidly cleaved from the solid support than its predecessor BL 2010.

BL 2030

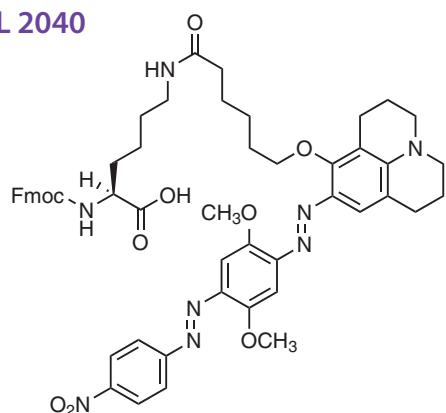


3'-BBQ-650® CPG III (1000 Å)

MF:	N/A	200 nmol columns (pkg of 4)	\$80
MW:	N/A	100 mg	\$147
CAS No:	None Assigned	1 μmol columns (pkg of 4)	\$230
		1 g	\$1,140

BlackBerry® Quencher 650 CPG III (3'-BBQ-650® CPG III) is used to install BBQ-650® at the 3' end of an oligonucleotide, and is more efficiently cleaved from the solid support than its predecessor BL 2020.

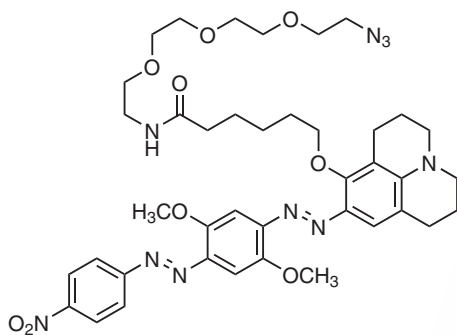
BL 2040



Fmoc-Lysine (BBQ-650™)-OH

MF:	C ₅₃ H ₅₈ N ₈ O ₁₀	50 mg	\$465
MW:	967.08	100 mg	\$710
CAS No:	None Assigned		

BL 3030



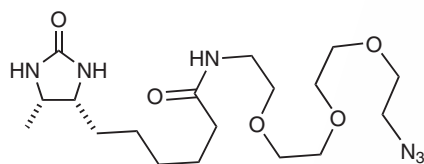
BBQ-650™-TEG azide

MF:	C ₄₀ H ₅₂ N ₁₀ O ₉	5 mg	\$250
MW:	816.90	10 mg	\$425
CAS No:	1333148-80-3		

A useful tool from our Click-mates™ collection for adding our long wavelength quencher, BBQ-650®, to a variety of substrates.

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BT 1075



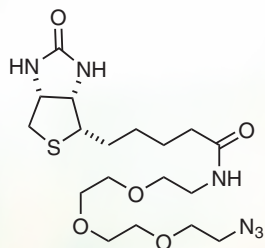
Desthiobiotin-TEG azide

MF:	C ₁₈ H ₃₄ N ₆ O ₅	25 mg	\$175
MW:	414.50	100 mg	\$595
CAS No:	1306615-47-3		

A useful tool from our Click-mates™ collection for ligation of desthiobiotin via Saudinger ligation or click chemistry.

The purchase of these products for use in applications relating to copper catalyzed azide-alkyne cycloaddition chemistry ("Click Chemistry") includes a limited, nontransferable license to intellectual property owned by TSRI to use this product solely for internal non-commercial research activities and specifically excludes clinical, therapeutic, or diagnostic use in humans or animals. Information regarding a license for commercial use in Click Chemistry may be obtained directly from The Scripps Research Institute, 10550 N. Torrey Pines Rd., La Jolla, CA 92037, or by contacting 858-784-8140 or click@scripps.edu

BT 1085



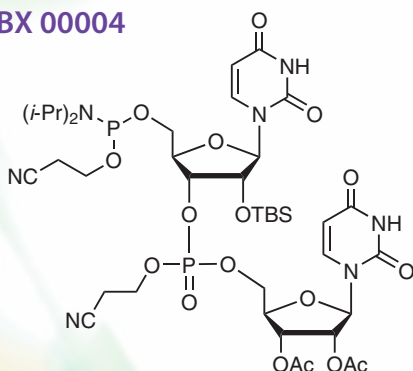
Biotin-TEG azide

MF:	C₁₈H₃₂N₆O₅S	25 mg	\$190
MW:	444.55	100 mg	\$490
CAS No:	875770-34-6		

A useful tool from our Click-mates™ collection for ligation of biotin *via* Saudinger ligation or click chemistry.

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BX 00004

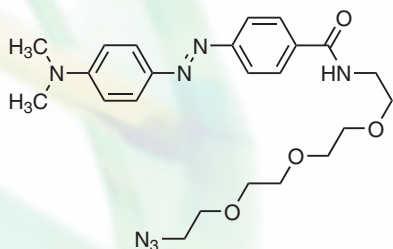


Protected U-U 5'-CEP

MF:	C₄₀H₆₁N₇O₁₇P₂Si	50 mg	\$495
MW:	1001.98	100 mg	\$980
CAS No:	915098-72-5		

Valuable nucleoside diphosphate tool for glycobiology. See also BA 0358.

DB 8010

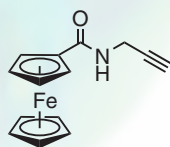


DabcyI-TEG azide

MF:	C₂₃H₃₁N₇O₄	5 mg	\$95
MW:	469.54	25 mg	\$210
CAS No:	None Assigned		

The purchase of these products for use in applications relating to copper catalyzed azide-alkyne cycloaddition chemistry ("Click Chemistry") includes a limited, nontransferable license to intellectual property owned by TSRI to use this product solely for internal non-commercial research activities and specifically excludes clinical, therapeutic, or diagnostic use in humans or animals. Information regarding a license for commercial use in Click Chemistry may be obtained directly from The Scripps Research Institute, 10550 N. Torrey Pines Rd., La Jolla, CA 92037, or by contacting 858-784-8140 or click@scripps.edu.

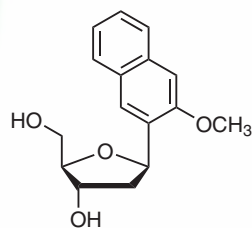
FC 8100



Ferrocenoyl propargylamide

MF:	C₁₄H₁₃FeNO	10 mg	\$98
MW:	267.10	100 mg	\$175
CAS No:	260444-13-1		

FC 8110

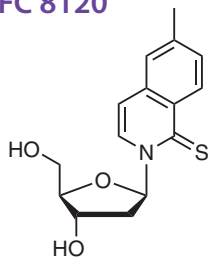


dNaM

MF:	C₁₆H₁₈O₇	25 mg	\$165
MW:	274.31	100 mg	\$495
CAS No:	1117893-19-2		

Unnatural base for expansion of the genetic alphabet.

FC 8120

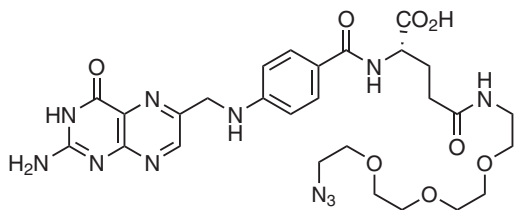


d5SICS

MF:	C₁₆H₁₇NO₃S	25 mg	\$390
MW:	291.37	100 mg	\$995
CAS No:	None Assigned		

Unnatural base for expansion of the genetic alphabet.

FC 8150



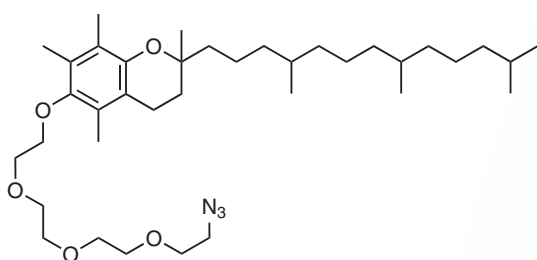
Folate TEG azide

MF:	C₂₇H₃₅N₁₁O₈	1 mg	\$295
MW:	641.64	5 mg	\$750
CAS No:	1313026-32-2		

Useful azide from our Click-mates™ collection for the installation of a folate tag for recognition by overexpressed folate receptors in cancer cells.

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FC 8160



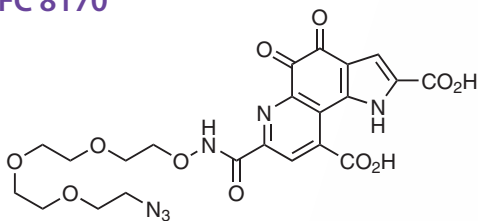
Tocopherol TEG azide

MF:	C₃₇H₆₅N₃O₅	5 mg	\$725
MW:	631.93	10 mg	\$975
CAS No:	None Assigned		

Lipophilic carrier tag from our Click-mates™ collection.

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FC 8170



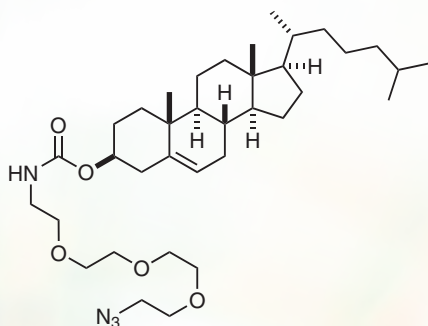
PQQ-TEG azide

MF:	C₂₂H₂₂N₆O₁₁	1 mg	\$295
MW:	546.44		
CAS No:	1333145-23-5		

Useful azide from our Click-mates™ collection for the installation of a methoxatin tag.

The purchase of these products for use in applications relating to copper catalyzed azide-alkyne cycloaddition chemistry ("Click Chemistry") includes a limited, nontransferable license to intellectual property owned by TSRI to use this product solely for internal non-commercial research activities and specifically excludes clinical, therapeutic, or diagnostic use in humans or animals. Information regarding a license for commercial use in Click Chemistry may be obtained directly from The Scripps Research Institute, 10550 N. Torrey Pines Rd., La Jolla, CA 92037, or by contacting 858-784-8140 or click@scripps.edu.

FC 8180



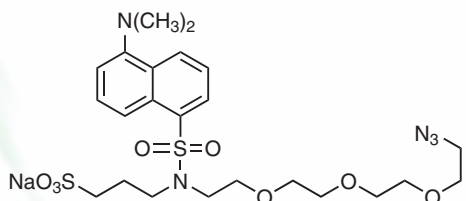
Cholesteryl-TEG azide

MF:	C₃₆H₆₂N₄O₅	25 mg	\$175
MW:	630.90	100 mg	\$595
CAS No:	None Assigned		

Useful azide from our Click-mates™ collection for the installation of a cholesteryl tag.

The purchase of these products for use in applications relating to copper catalyzed azide-alkyne cycloaddition chemistry ("Click Chemistry") includes a limited, nontransferable license to intellectual property owned by TSRI to use this product solely for internal non-commercial research activities and specifically excludes clinical, therapeutic, or diagnostic use in humans or animals. Information regarding a license for commercial use in Click Chemistry may be obtained directly from The Scripps Research Institute, 10550 N. Torrey Pines Rd., La Jolla, CA 92037, or by contacting 858-784-8140 or click@scripps.edu.

FD 13005



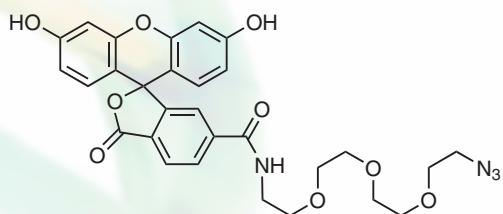
Water Soluble Dansyl-TEG azide

MF:	C₂₃H₃₄N₅NaO₈S₂	25 mg	\$95
MW:	595.67	100 mg	\$325
CAS No:	None Assigned		

Useful tool from our Click-mates™ collection for ligation of a water soluble dansyl group *via* Saudinger ligation or click chemistry.

The purchase of these products for use in applications relating to copper catalyzed azide-alkyne cycloaddition chemistry ("Click Chemistry") includes a limited, nontransferable license to intellectual property owned by TSRI to use this product solely for internal non-commercial research activities and specifically excludes clinical, therapeutic, or diagnostic use in humans or animals. Information regarding a license for commercial use in Click Chemistry may be obtained directly from The Scripps Research Institute, 10550 N. Torrey Pines Rd., La Jolla, CA 92037, or by contacting 858-784-8140 or click@scripps.edu.

FF 6110



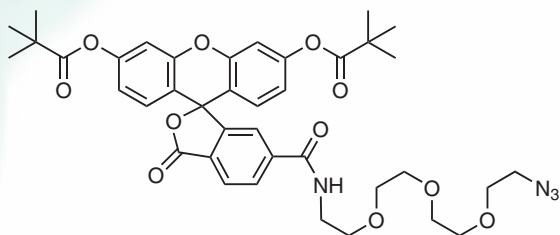
6-Carboxyfluorescein-TEG azide

MF:	C₂₉H₂₈N₄O₉	25 mg	\$175
MW:	576.55	100 mg	\$595
CAS No:	412319-45-0		

Useful fluorescein azide from our Click-mates™ collection for fluorescein incorporation *via* Staudinger ligation or click chemistry.

The purchase of these products for use in applications relating to copper catalyzed azide-alkyne cycloaddition chemistry ("Click Chemistry") includes a limited, nontransferable license to intellectual property owned by TSRI to use this product solely for internal non-commercial research activities and specifically excludes clinical, therapeutic, or diagnostic use in humans or animals. Information regarding a license for commercial use in Click Chemistry may be obtained directly from The Scripps Research Institute, 10550 N. Torrey Pines Rd., La Jolla, CA 92037, or by contacting 858-784-8140 or click@scripps.edu.

FF 6120



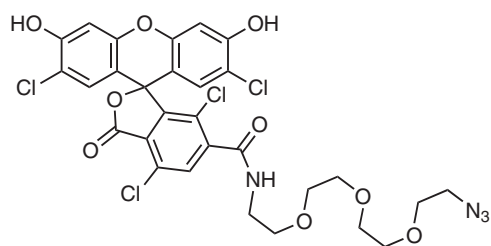
6-Carboxyfluorescein-dipivalate TEG azide

MF:	C₃₉H₄₄N₄O₁₁	25 mg	\$175
MW:	744.79	100 mg	\$595
CAS No:	1254709-11-9		

Useful fluorescein azide from our Click-mates™ collection for fluorescein incorporation *via* Staudinger ligation or click chemistry.

The purchase of these products for use in applications relating to copper catalyzed azide-alkyne cycloaddition chemistry ("Click Chemistry") includes a limited, nontransferable license to intellectual property owned by TSRI to use this product solely for internal non-commercial research activities and specifically excludes clinical, therapeutic, or diagnostic use in humans or animals. Information regarding a license for commercial use in Click Chemistry may be obtained directly from The Scripps Research Institute, 10550 N. Torrey Pines Rd., La Jolla, CA 92037, or by contacting 858-784-8140 or click@scripps.edu.

FF 6130



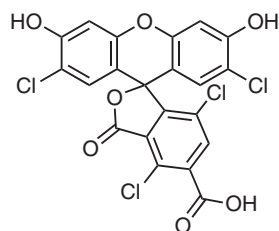
6-TET-TEG azide

MF: $C_{29}H_{24}Cl_4N_4O_9$ 1 mg \$295
 MW: 714.33
 CAS No: 1333145-22-4

Useful fluorescein azide from our Click-mates™ collection for tetrachloro-fluorescein incorporation *via* Staudinger ligation or click chemistry.

The purchase of these products for use in applications relating to copper catalyzed azide-alkyne cycloaddition chemistry ("Click Chemistry") includes a limited, nontransferable license to intellectual property owned by TSRI to use this product solely for internal non-commercial research activities and specifically excludes clinical, therapeutic, or diagnostic use in humans or animals. Information regarding a license for commercial use in Click Chemistry may be obtained directly from The Scripps Research Institute, 10550 N. Torrey Pines Rd., La Jolla, CA 92037, or by contacting 858-784-8140 or click@scripps.edu.

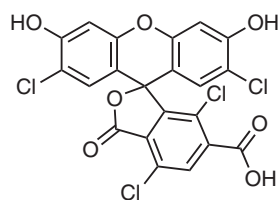
FF 6150



5-Carboxy-TET

MF: $C_{21}H_8Cl_4O_7$ 50 mg \$200
 MW: 514.10 250 mg \$850
 CAS No: 155911-13-0

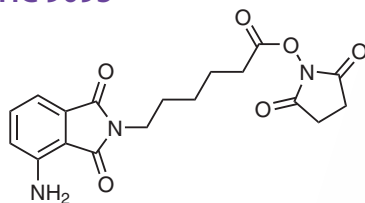
FF 6160



6-Carboxy-TET

MF: $C_{21}H_8Cl_4O_7$ 50 mg \$200
 MW: 514.10 250 mg \$850
 CAS No: 155911-14-1

HC 9095

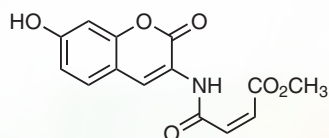


Luminol synthon N-hydroxysuccinimide ester

MF: $C_{18}H_{19}N_3O_6$ 10 mg \$75
 MW: 373.36 50 mg \$210
 CAS No: 1111822-73-1

Luminol precursor.

HC 9097

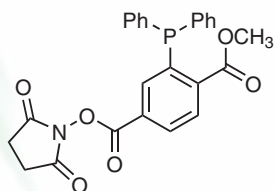


Glutathione reductase probe

MF:	C₁₄H₁₁NO₆	10 mg	\$125
MW:	289.24	50 mg	\$365
CAS No:	1173888-41-9		

Fast thiol detection with a probe utilizing coumarin as the fluorophore and maleimide as the thiol acceptor.

LK 4260



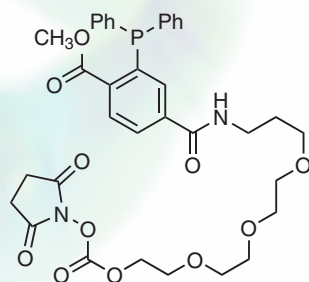
AmAz Coupler

MF:	C₂₅H₂₀NO₆P	50 mg	\$125
MW:	461.40	100 mg	\$200
CAS No:	532931-48-9		

Our AmAz Coupler allows for the linking of an amine-containing component to an azide-containing component. The amine-containing component is first reacted with the NHS ester and then the azide-containing component is reacted with the phosphine. An intramolecular trapping of the phosphorimidate by the ortho- methyl ester results in a stable aryl-1,4-diamide linkage between the two components.

LK 4260 and the use thereof in chemoselective ligation, are protected by one or more of the following US Patents, the rights of which are assigned to The Regents of the University of California: US 6,570,040, US 7,122,703, and US 7,838,665. Additional patents are pending. It may be used for research purposes only. It is not licensed for resale and may only be used by the buyer. This product may not be used and is not licensed for clinical assays, where the results of such assays are provided as a diagnostic service. If a diagnostic or therapeutic use is anticipated, then a license must be requested from the University of California. The availability of such diagnostic and therapeutic use license(s) cannot be guaranteed from the University of California.

LK 4265



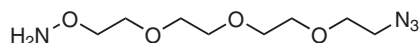
AmAz Coupler II

MF:	C₃₅H₃₉N₂O₁₁P	50 mg	\$250
MW:	694.66	100 mg	\$400
CAS No:	None Assigned		

Our AmAz Coupler II allows for the linking of an amine-containing component to an azide-containing component. The amine-containing component is first reacted with the NHS ester and then the azide-containing component is reacted with the phosphine. An intramolecular trapping of the phosphorimidate by the ortho- methyl ester results in a stable aryl-1,4-diamide linkage between the two components.

LK 4260 and the use thereof in chemoselective ligation, are protected by one or more of the following US Patents, the rights of which are assigned to The Regents of the University of California: US 6,570,040, US 7,122,703, and US 7,838,665. Additional patents are pending. It may be used for research purposes only. It is not licensed for resale and may only be used by the buyer. This product may not be used and is not licensed for clinical assays, where the results of such assays are provided as a diagnostic service. If a diagnostic or therapeutic use is anticipated, then a license must be requested from the University of California. The availability of such diagnostic and therapeutic use license(s) cannot be guaranteed from the University of California.

LK 4270



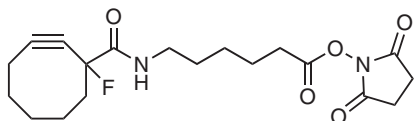
Aminoxy-TEG-Azide

MF:	C₈H₁₈N₄O₄	100 mg	\$325
MW:	234.25	1 g	\$2,600
CAS No:	1306615-51-9		

A bi-functional linker from our Click-mates™ collection with potential applications in oligonucleotide ligations.

The purchase of these products for use in applications relating to copper catalyzed azide-alkyne cycloaddition chemistry ("Click Chemistry") includes a limited, nontransferable license to intellectual property owned by TSRI to use this product solely for internal non-commercial research activities and specifically excludes clinical, therapeutic, or diagnostic use in humans or animals. Information regarding a license for commercial use in Click Chemistry may be obtained directly from The Scripps Research Institute, 10550 N. Torrey Pines Rd., La Jolla, CA 92037, or by contacting 858-784-8140 or click@scripps.edu.

LK 4300



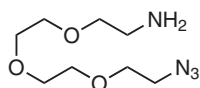
Click-easy™ MFCO-N-hydroxysuccinimide

MF:	C₁₉H₂₅FN₂O₅	25 mg	\$255
MW:	380.41	50 mg	\$470
CAS No:	1306615-48-4		

Monofluoro-substituted cyclooctyne (MFCO) from our Click-mates™ collection suitable for copper free click reactions.

LK 4300 is protected by US 7,807,619, the rights of which are assigned to The Regents of the University of California. Additional patents covering methods for their use in the modification of biomolecules are pending. It may be used for research purposes only. It is not licensed for resale and may only be used by the buyer. This product may not be used and are not licensed for clinical assays, where the results of such assays are provided as a diagnostic service. If a diagnostic or therapeutic use is anticipated, then a license must be requested from the University of California. The availability of such diagnostic and therapeutic use license(s) cannot be guaranteed from the University of California.

LK 4310



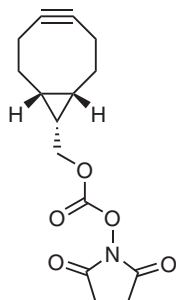
Amino-TEG azide

MF:	C₈H₁₈N₄O₃	500 mg	\$120
MW:	218.25	1 g	\$175
CAS No:	134179-38-7		

A bi-functional linker from our Click-mates™ collection with potential applications in oligonucleotide ligations.

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LK 4320



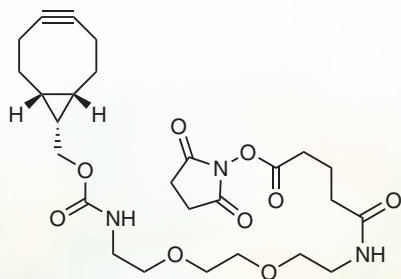
Click-easy™ BCN N-hydroxysuccinimide ester I

MF:	C₁₅H₁₇NO₅	50 mg	\$185
MW:	291.30	100 mg	\$354
CAS No:	1352642-02-4		

A state-of-the-art cycloalkyne for catalyst free strain-promoted azide-alkyne cycloadditions.

BCN reagents are sold for research use only under a licensing agreement with Synaffix, B.V.

LK 4330



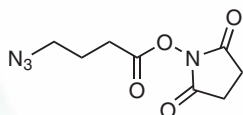
Click-easy™ BCN N-hydroxysuccinimide ester II

MF:	C₂₆H₃₇N₃O₉	10 mg	\$150
MW:	535.59	25 mg	\$220
CAS No:	1314639-16-1		

A state-of-the-art cycloalkyne for catalyst free strain-promoted azide-alkyne cycloadditions.

BCN reagents are sold for research use only under a licensing agreement with Synaffix, B.V.

LK 4340



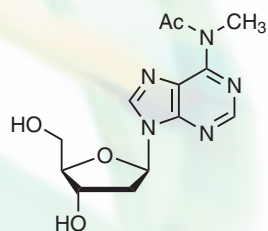
4-Azidobutyric acid N-hydroxysuccinimide ester

MF:	C₈H₁₀N₄O₄	10 mg	\$135
MW:	226.19	25 mg	\$300
CAS No:	943858-70-6		

A state-of-the-art cycloalkyne for catalyst free strain-promoted azide-alkyne cycloadditions.

The purchase of these products for use in applications relating to copper catalyzed azide-alkyne cycloaddition chemistry ("Click Chemistry") includes a limited, nontransferable license to intellectual property owned by TSRI to use this product solely for internal non-commercial research activities and specifically excludes clinical, therapeutic, or diagnostic use in humans or animals. Information regarding a license for commercial use in Click Chemistry may be obtained directly from The Scripps Research Institute, 10550 N. Torrey Pines Rd., La Jolla, CA 92037, or by contacting 858-784-8140 or click@scripps.edu.

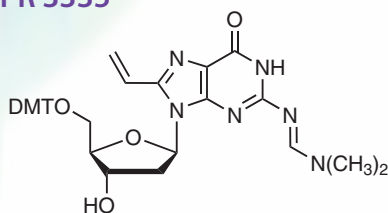
PR 3006



6-N-Acyl-6-N-methyl-2'-deoxyadenosine

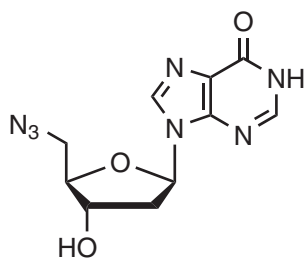
MF:	C₁₃H₁₇N₅O₄	25 mg	\$168
MW:	307.31	100 mg	\$595
CAS No:	None Assigned		

PR 3335

5'-O-Dimethoxytrityl-N²-(dimethylaminomethylidene)-8-ethenyl-2'-deoxyguanosine

MF:	C₃₆H₃₈N₆O₆	100 mg	\$335
MW:	650.72	1 g	\$2,450
CAS No:	1101864-07-6		

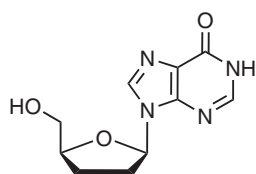
PR 3464



5'-Azido-2',5'-dideoxyinosine

MF:	$C_{10}H_{11}N_7O_3$	5 mg	\$75
MW:	277.24	25 mg	\$257
CAS No:	496809-83-7		

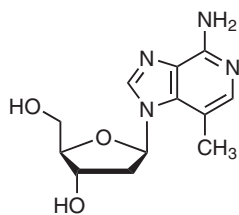
PR 3727



2',3'-Dideoxyinosine

MF:	$C_{10}H_{12}N_4O_3$	100 mg	\$75
MW:	236.23	1 g	\$225
CAS No:	69655-05-6	5 g	\$890

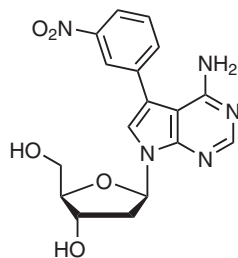
PRA 10013



3-Deaza-3-methyl-2'-deoxyadenosine

MF:	$C_{12}H_{16}N_4O_3$	10 mg	\$95
MW:	264.28	25 mg	\$215
CAS No:	515815-12-0		

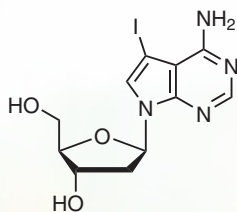
PRA 10032



7-(3-Nitrophenyl)-7-deaza-2'-deoxyadenosine

MF:	$C_{17}H_{17}N_5O_5$	10 mg	\$119
MW:	371.35	50 mg	\$495
CAS No:	None Assigned		

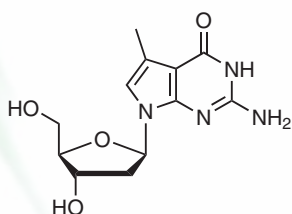
PRA 10036



5-Iodo-2'-deoxytubercidin

MF:	$C_{11}H_{13}IN_4O_3$	10 mg	\$87
MW:	376.15	50 mg	\$355
CAS No:	166247-63-8		

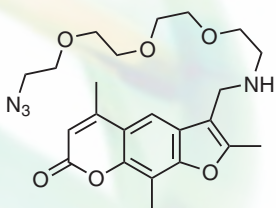
PRA 10042



7-Methyl-7-deaza-2'-deoxyguanosine

MF:	$C_{12}H_{16}N_4O_4$	10 mg	\$107
MW:	280.29	50 mg	\$375
CAS No:	90358-21-7		

PS 5030



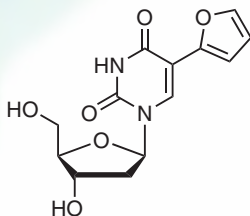
Psoralen-TEG azide

MF:	$C_{23}H_{30}N_4O_6$	1 mg	\$210
MW:	458.22	5 mg	\$450
CAS No:	1352815-11-2		

Useful azide from our Click-mates™ collection for the installation of a psoralen label.

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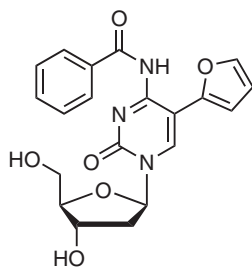
PY 7053



5-Furan-2-yl dU

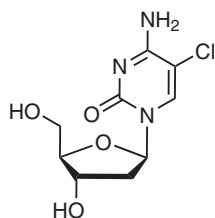
MF:	$C_{13}H_{14}N_2O_6$	25 mg	\$95
MW:	294.26	100 mg	\$345
CAS No:	92233-50-6		

PY 7054

*N*⁴-Benzoyl-5-furan-2-yl dC

MF:	C₂₀H₁₉N₃O₆	25 mg	\$125
MW:	397.38	100 mg	\$395
CAS No:	1119734-59-6		

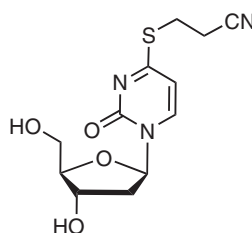
PY 7114



5-Chloro-2'-deoxycytidine

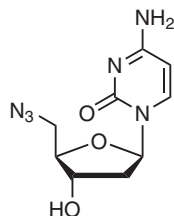
MF:	C₉H₁₂ClN₃O₅	25 mg	\$295
MW:	261.66	100 mg	\$985
CAS No:	32387-56-7		

PY 7212

*S*⁴-(2-Cyanoethyl)-4-thio-2'-deoxyuridine

MF:	C₁₂H₁₅N₃O₄S	10 mg	\$66
MW:	297.33	100 mg	\$536
CAS No:	136055-15-7		

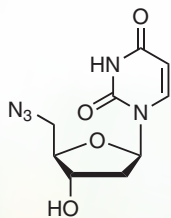
PY 7219



5'-Azido-2,5'-dideoxycytidine

MF:	C₉H₁₂N₆O₃	25 mg	\$98
MW:	252.23	100 mg	\$345
CAS No:	4803-88-7		

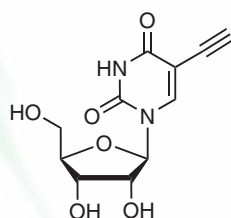
PY 7276



5'-Azido-2',5'-dideoxyuridine

MF:	$C_9H_{11}N_5O_4$	25 mg	\$98
MW:	253.21	100 mg	\$345
CAS No:	35959-37-6		

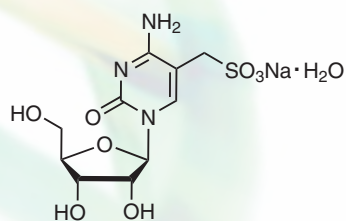
PY 7563



5-Ethynyl uridine

MF:	$C_{11}H_{12}N_2O_6$	10 mg	\$151
MW:	268.22	100 mg	\$1220
CAS No:	69075-42-9		

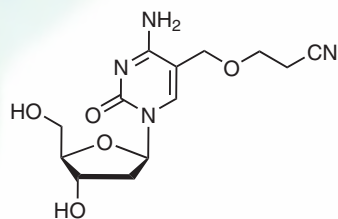
PY 7586



Cytidin-5-yl-methanesulfonate sodium salt hydrate

MF:	$C_{10}H_{16}N_3NaO_9S$	50 mg	\$525
MW:	377.30	100 mg	\$897
CAS No:	None Assigned		

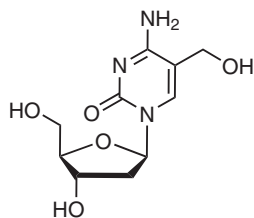
PY 7587



5-[(2-Cyanoethoxy)methyl]-2'-deoxycytidine

MF:	$C_{13}H_{18}N_4O_5$	10 mg	\$210
MW:	310.31	50 mg	\$750
CAS No:	188411-05-4		

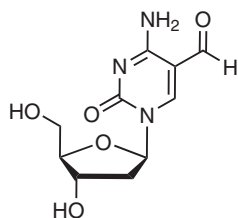
PY 7588



5-Hydroxymethyl-2'-deoxycytidine

MF:	$C_{10}H_{15}N_3O_5$	10 mg	\$245
MW:	257.24	50 mg	\$875
CAS No:	7226-77-9		

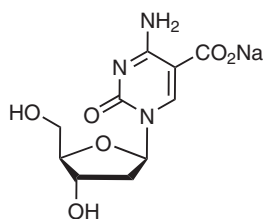
PY 7589



5-Formyl-2'-deoxycytidine

MF:	$C_{10}H_{13}N_3O_5$	5 mg	\$185
MW:	255.03	10 mg	\$315
CAS No:	137017-45-9		

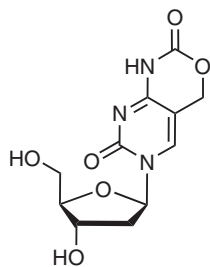
PY 7593



2'-Deoxycytidine-5-Carboxylic acid sodium salt

MF:	$C_{10}H_{12}N_3NaO_6$	25 mg	\$147
MW:	293.21	100 mg	\$335
CAS No:	1009808-6-2-1 (negative ion)		

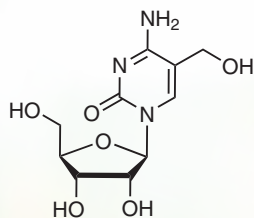
PY 7594



5-Hydroxymethyl-2'-deoxycytidine cyclic carbamate

MF:	$C_{11}H_{13}N_3O_6$	10 mg	\$150
MW:	283.24	50 mg	\$525
CAS No:	1257247-41-8		

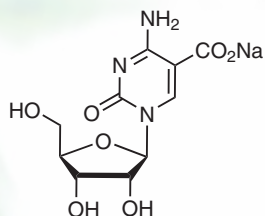
PY 7596



5-Hydroxymethylcytidine

MF:	$C_{10}H_{15}N_3O_6$	50 mg	\$295
MW:	273.24	100 mg	\$495
CAS No:	19235-17-7		

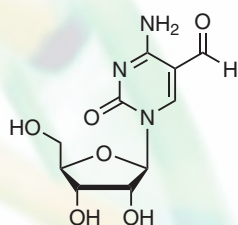
PY 7598



Cytidine-5-carboxylic acid, sodium salt

MF:	$C_{10}H_{12}N_3NaO_7$	10 mg	\$225
MW:	309.21	50 mg	\$690
CAS No:	64623-37-6 (parent acid)		

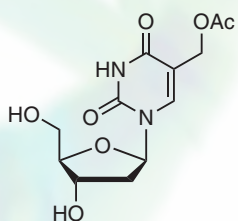
PY 7599



5-Formylcytidine

MF:	$C_{10}H_{13}N_3O_6$	5 mg	\$191.50
MW:	271.23	10 mg	\$325
CAS No:	148608-53-1		

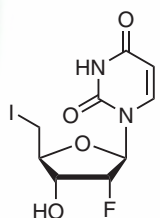
PY 7604



5-Acetoxymethyl-2'-deoxyuridine

MF:	$C_{12}H_{16}N_2O_7$	25 mg	\$315
MW:	300.26	100 mg	\$1,095
CAS No:	148380-55-6		

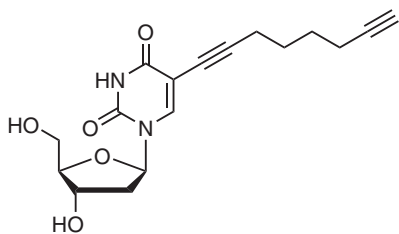
PY 7612



2'-Fluoro-5'-iodo deoxyuridine

MF:	$C_9H_{10}FIN_2O_4$	25 mg	\$88
MW:	356.09	100 mg	\$297
CAS No:	211694-25-6		

PY 7713

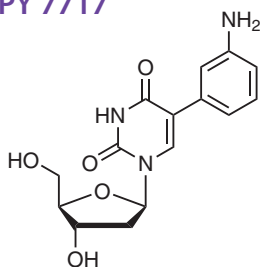


5-(1,7-Octadiyn-1-yl)-2'-deoxyuridine

MF:	C₁₇H₂₀N₂O₅	50 mg	\$68
MW:	332.35	250 mg	\$225
CAS No:	909398-18-1		

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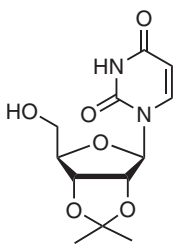
PY 7717



5-(3-Aminophenyl)-2'-deoxyuridine

MF:	C₁₅H₁₇N₃O₅	25 mg	\$235
MW:	319.31	100 mg	\$795
CAS No:	188482-60-2		

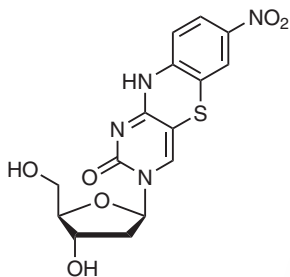
PY 7781



2'3'-O-Isopropylidene uridine

MF:	C₁₂H₁₆N₂O₆	5 g	\$95
MW:	284.27	25 g	\$345
CAS No:	362-43-6		

PYA 11115

tC_{nitro} nucleoside

MF:	C₁₅H₁₄N₄O₆S	5 mg	\$138
MW:	378.36	25 mg	\$575
CAS No:	1139889-25-0		



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