METKINEN chemistry

Reagents and building blocks for oligonucleotide synthesis



Synthetic RNA oligonucleotides, incorporating 2'-Fluoro-2'-deoxynucleosides (2'-F-RNA) find constantly increasing applications in research and new drug development. Among important areas of their application are such as design and synthesis of efficient and stable siRNAs,¹ aptamers,² anti-microRNAs (Antagomirs)³ as well as general use for affinity purification of RNAbinding proteins.⁴

Today Metkinen Chemistry offers more than 99% pure purine 2'-fluoronucleosides, prepared by bioconversion of synthetic 2'-F-uridine, using evolved recombinant E.coli strains. These strains, in turn are grown on the media, incorporation only ingredients, derived from yeast and vegetable. That means that in the production process no secondary reagents of animal source (either BSE certified or non- BSE certified) are employed at any stage of 2'-F- A and G preparation. Hence, we guarantee the complete absence of ANY BSE/TSE risk, associated with our 2'-Fnucleosides and derivatives.

In addition to purine 2'-F-nucleosides, we offer a full set of 2'-F-nucleoside CE-phosphoramidites for the chemical synthesis and a full set of 2'-F-nucleoside 5'-triphosphates for enzymatic preparation of 2'-F-RNA. Our latest addition to series of 2'-fluoronucleoside derivatives – 2'-Fnucleoside bound solid supports for 2'-F-RNA oligosynthesis, complements the range of these compounds and opens up a full spectrum of 2'-F-RNA applications.

References:

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2'-Fluoro-2'-Deoxyadenosine

9-(2-Deoxy-2-fluoro-ß-D-ribofuranosyl)-adenine



Catalogue number: 203-17

CAS number: 64183-27-3

IUPAC name: 2'-Deoxy-2'-fluoroadenosine

Purity: > 99%

Description: white to off-white crystals

Storage of dry compound: 3 years at +4°C

Note: Gram to Kilogram scale production. Please request prices for bulk!





Purity: > 95%

Description: white to off-white amorphous powder

Storage of dry compound: 3 years at -20°C



2'-Fluoro-2'-Deoxyadenosine-3'-CE-phosphoramidite



Catalogue number: 203-51

Purity: > 95%

Description: white to off-white amorphous powder

Storage of dry compound: 3 years at -20°C





Catalogue number: 203-52

Purity: > 95%

Description: white to off-white amorphous powder

Storage of dry compound: 3 years at -20°C



2'-Fluoro-2'-Deoxycytidine-3'-CE-phosphoramidite

Catalogue number: 203-53

Purity: > 95%

Description: white to off-white amorphous powder

Storage of dry compound: 3 years at -20°C

METKINEN chemistry

2'-Fluoro-2'-deoxynucleosides & Derivatives



2'-Fluoro-2'-deoxynucleoside 5'-triphosphates are popular products for various applications: PCR, DNA- and RNA-polymerase, reverse transcriptase based assays, etc. Nowadays these compounds find diverse use in siRNA and Aptamer research.

Product Note: All of our Triphosphates are lithium salts. They are analyzed by NMR, UV, HPLC and are accompanied by a Product Data Sheet



New 2'-F-Nucleoside bound Solid Supports (F-CS-supports)

The F-CS Solid Supports incorporate a new linker that allows 3-4 times faster cleavage of target oligonucleotide off the solid phase under the the regular cleavage/deprotection conditions. This feature make them indispensable, when UltraFAST Deprotection ULtraMILD Deprotection is necessary in order to insure the convenient isolation of modified and dye labelled oligonucleotides that appear not stable to deprotection with ammonium hydroxide or AMA at regular conditions (e.g. oligonucleotides, labled with Cy5 or Cy 5.5 dyes)

The novel F-CS supports appear absolutely similar, if not better than the conventional nucleoside bound supports (incorporating the traditional succinate linker) in terms of performance as solid phases for RNA Oligonucleotide synthesis. Noteworthy, the precise loading of the new linker on the solid phase is much easier to achieve when using the new Carbomoylation procedure, developed by Metkinen Chemistry (U.S. Patent Application Serial No 60/854,721 and International Patent Application No. PCT/ F12007/050575).

METKINEN *chemistry*



F-CS-supports: 2'-F-Nucleoside Bound Solid Supports for 2'-F- RNA synthesis

Catalogue number: 103-30

Description: Chemically modified Macroporous Aminomethyl Polystyrene. White to off-white powder.

Storage of dry compound: 1 year at +4°C

Loading: standard - 30 - 40 µmol/g; custom loading – from 5 to 80 µmol/g.

Oligosynthesis on CS supports: perform oligonucleotide assembly, using standard RNA protocols, recommended by your synthesizer manufacturer.

Cleavage: Cleave the oligo from the support using concentrated aqueous ammonium hydroxide at room temperature for 30 minutes. Alternatively, cleave the oligo from the support using concentrated aqueous ammonium hydroxide – 40 % aqueous methylamine (AMA) at room temperature for 15 minutes.



B*= bzAde, Thy, bzCyt, acCyt, ibuG; AMPS = macroporous aminomethyl polystyrene

Deprotection AFTER Cleavage: Following the cleavage step, proceed with oligonucleotide deprotection using the conditions appropriate for removal of the protecting groups on the nucleobases.



Metkinen Universal Solid Support

The universal support strategy offers the following clear advantages: eliminates the possibility of errors in parallel synthesis applications where up to 384 wells may contain different supports; eliminates the need for at least four supports for DNA synthesis and four for RNA synthesis; simplifies the preparation of oligonucleotides with modified or unusual nucleosides at the 3'-terminus.

Like our initial version of Universal Solid supports - USII, the new type of supports - Metkinen Universal Solid Support would be appropriate for the production of DNA oligos, long and short, as well as those requiring mild deprotection. It is compatible with the synthesis of RNA, (including siRNA) as well as virtually any oligonucleotide analogs (2'-F-RNA, 2'-OMe-RNA, LNA, oligonucleotide N3'•P5' phosphoramidates, etc.). The reagent used for the cleavage/dephosphorylation step is commercially available and the procedures described are fully compatible with high-throughput Synthesis.

The difference lies in the higher stability of Metkinen Universal Solid Support than that of USII upon prolong storage. In addition, the preparation of the new support appears to be more consistent and reliable.

Metkinen Universal Solid Support is subject to proprietary rights of Glen Research Corporation and is synthesized and sold under the following licensed patents: US Patent No.: 6,770,754 and European Patent No.: 1404695. The new carbomoylation chemistry, resulting in the stable urea fragment, bridging the Universal linker and aminoalkylated solid phase, is subject to proprietary rights of Metkinen Chemistry (U.S. Patent Application Serial No. 60/854,721; International Patent Application No. PCT/ F12007/050575).



Universal Solid Support

Metkinen Universal Solid Support for synthesis of DNA, RNA & any type of modified oligonucleotides.

Catalogue number: 103-00

Description: Chemically modified Macropourous Aminomethyl polystyrene. White to off-white powder.

Storage of dry compound: 1 year at +20°C

Loading: 20-80 µmol/g loadings is available. Please enquire for custom loading.

Oligo synthesis on USIII: Perform oligonucleotide assembly, using standard protocols, recommended by your synthesizer manufacturer. Upon the completion of synthesis wash the oligonucleotide bound support with pure acetonitrile. Do not perform any washing steps with solvents, containing basic reagents (diethylamine, triethylamine, dimethylamine, etc.) and water! **Cleavage:** Cleave the oligo from the support using 3.5N - 4.5N ammonia in methanol (dilute cold 7N ammonia in methanol, Aldrich Cat. No 499145-100ML, with cold anhydrous methanol) at room temperature for 30 minutes. Do not use aqueous ammonium hydroxide and/or mixtures of ammonium hydroxide and methanol for cleavage!

Deprotection AFTER Cleavage:

Standard: After Cleavage, add 1 volume of 3.5N - 4.5N ammonia in methanol, seal and deprotect for 8-15 hours at 60 °C for removal of the protecting groups on the nucleobases. or

Alternatively: add 1 volume of 30% ammonium hydroxide, seal and deprotect using the conditions appropriate for removal of the protecting groups on the nucleobases (e.g. at 55 °C for 5 hours).



Chemical Phosphorylation Reagent II

Chemical Phosphorylation Reagent (CPR II) contains a DMT group which can be left on the oligonucleotide and used for rapid purification of oligonucleotide 5'-phosphates by the popular DMTr-on technique, which employs disposable RP cartridges or "Trityl-on" RP HPLC purification. The DMTr group is removed with aqueous acid (e.g., 2%TFA in the case of Cartridge Purification) and the remaining linker is then eliminated after brief treatment with aqueous ammonium hydroxide (12 -15% ammonium hydroxide at room temperature for 15 minutes) to yield the 5'-phosphate. Catalogue number: 103-10

Description: amorphous colorless glass

Storage of dry compound: 1 year at -20°C

Coupling Conditions: 6 minute coupling time.

Omit the capping step after the addition of this reagent!



[3-(4,4'-Dimethoxytrityloxy)-2,2dicarboxyethyl]propyl-(2-cyanoethyl)-(N,N-diisopropyl)-phosphoramidite

Trinucleotide (TRIMER CODON) Phosphoramidites

Catalogue number: 103-20

Description: white to off-white powder

Storage of dry compound: 1 year at -20°C

Protein mutagenesis can be used to fine tune a variety of properties, such as improved stability to high temperatures, denaturants, or nonaqueous solvents; higher affinity binding to a target molecule; increased rates of enzymatic reactions; or changes of specificities. However, generating and finding these improved proteins can be a difficult task. One of the most popular methods is to make pools of degenerate oligonucleotides, which can be incorporated into the genes as cassettes or by PCR by using the degenerate oligo as a primer.1 Degenerate oligonucleotides are synthesized as a mixture of A/C/G/T phosphoramidites (N) at the site of the codons to be mutated. Problems arise, though, from using an equimolar solution of each base. First there is a coding bias. Out of the 64 possible codon combinations of A. C. G and T. 18 code for leucine, arginine or serine, but only 2 for tryptophan or methionine. As a result, only 3% of the mutagenic oligonucleotides will contain methionine or tryptophan, and over 28% will contain either leucine, arginine or serine. In addition, the three nonsense codons will lead to chain termination in 4.7% of the sequences. There are ways to improve this situation. For instance, using two degenerate mixes of bases,



N and G/C, on the DNA synthesizer to insert NNG/C into the sequence will halve the number of the most degenerate codons, but still code for all 20 amino acids. However, still 59% of the clones will code for just eight amino acids and 3% will have a stop codon inserted.



Trinucleotide (TRIMER CODON) Phosphoramidites

The generation of redundant sequences and stop codons makes searching a clonal library inefficient. However, it is possible to improve the efficiency of this process by using a mixture of trinucleotide (trimer) phosphoramidites.^{2–5} By synthesizing a set of trimers that cover all 20 amino acids, the mutation of a gene can be carried out at the codon level rather

than at individual bases. Therefore, unlike other methods of mutagenesis, trimer phosphoramidites lead to no codon bias, no frame-shift mutations, and no production of stop codons, making them one of the most efficient tools to explore sequence space in protein regions that are important for function 6 – even in nonsaturating conditions.^{7,8}

References:

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New 2'-Deoxynucleoside bound Solid Supports (CS-supports)

The CS Solid Supports incorporate a new linker that allows 3–4 times faster cleavage of target oligonucleotide off the solid phase under the the regular cleavage/deprotection conditions. This feature make them indispensable, when Ultra-FAST Deprotection ULtraMILD Deprotection is necessary in order to insure the convenient isolation of modified and dye labelled oligonucleotides that appear not stable to deprotection with ammonium hydroxide or AMA at regular conditions (e.g. oligonucleotides, labled with Cy5 or Cy 5.5 dyes) The novel CS supports appear absolutely similar, if not better than the conventional nucleoside bound supports (incorporating the traditional succinate linker) in terms of performance as solid phases for Oligonucleotide synthesis. Noteworthy, the precise loading of the new linker on the solid phase is much easier to achieve when using the new Carbomoylation procedure, developed by Metkinen Chemistry (U.S. Patent Application Serial No 60/854,721 and International Patent Application No. PCT/FI2007/050575).



New 2'-Deoxynucleoside bound Solid Supports (CS-supports)

DMT rO

CS –**Supports:** 2'-Deoxynucleoside Bound Solid Supports for DNA synthesis

Catalogue number: 103-31

Description: Chemically modified Macroporous Aminomethyl Polystyrene. White to off-white powder.

Storage of dry compound: 1 year at +4°C

Loading: standard - 30 - 40 µmol/g; custom loading – from 5 to 80 µmol/g.

Oligosynthesis on CS supports: perform oligonucleotide assembly, using standard DNA protocols, recommended by your synthesizer manufacturer.

Cleavage: Cleave the oligo from the support using concentrated aqueous ammonium hydroxide at room temperature for 30 minutes. Alternatively, cleave the oligo from the support using



B*= bzAde, Thy, bzCyt, acCyt, ibGua; AMPS = macroporous aminomethyl polystyrene

concentrated aqueous ammonium hydroxide - 40 % aqueous methylamine (AMA) at room temperature for 15 minutes.

Deprotection AFTER Cleavage: Following the cleavage step, proceed with oligonucleotide deprotection using the conditions appropriate for removal of the protecting groups on the nucleobases.



2'-Fluoroaraninooligonucleotides (FANA) can be used as antisense oligonucleotides,¹ siRNA² and aptamers.³Today Metkinen Chemistry offers synthetically synthesized pyrimidine 2'-fluoroarabinonucleosides as well as purine 2'-fluoroarabinonucleosides, prepared by bioconversion of synthetic 2'-F-arabinothymidine, employing evolved recombinant E.coli strains. We guarantee the complete absence of ANY BSE/TSE risk, associated with our 2'-F-fluoroarabinonucleosides. Starting from October 2009 Metkinen Chemistry will launch production of a full set of 2'-F-arabinonucleoside CE-phosphoramidites for the chemical synthesis and a full set of 2'-Farabinonucleoside 5'-triphosphates for enzymatic preparation of 2'-FANA.

References:

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METKINEN chemistry

















Biotin-II-UTP

5-[3-(6-(Biotinylamido)hexanoylamido)propenyl]-uridine-5'-triphosphate, triethylammonium or tetralithium (optional) salt



Catalogue number: 303-31

Diluent: dd-H₂O

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: 1 month as 1 mM solution in water or neutral buffers at +4°C, 6 months as 1 mM solution in water or neutral buffers at -20°C.





Catalogue number: 303-11

Diluent: dd-H,O

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: 1 month as 1 mM solution in water or neutral buffers at +4°C, 6 months as 1 mM solution in water or neutral buffers at -20°C.









Diluent: dd-H₂O

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: 1 month as 1 mM solution in water or neutral buffers at +4°C, 6 months as 1 mM solution in water or neutral buffers at -20°C.



Biotin-16-dUTP

5-[3-{3-(6-(Biotinylamido)hexanoylamido)bytyrylamido}propenyl]-2'-deoxyuridine-5'-triphosphate, triethylammonium or tetralithium (optional) salt



Catalogue number: 303-14

Diluent: dd-H₂O

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: 1 month as 1 mM solution in water or neutral buffers at +4°C, 6 months as 1 mM solution in water or neutral buffers at -20°C.







6-FAM-II-dUTP

5-[3-(6-(fluoresceinyl-6-carboxamido)hexanoylamido)-propenyl]-2'-deoxyuridine-5'triphosphate, triethylammonium or tetralithium (optional) salt



Catalogue number: 303-12

Diluent: dd-H₂O

Storage: Freezer storage in dark, -10 to -30°C, dry.

Stability in Solution: 1 month as 1 mM solution in water or neutral buffers at +4°C, 6 months as 1 mM solution in water or neutral buffers at -20°C in dark.



6-FAM-11-dCTP 5-[3-(6-(fluoresceinyl-6-carboxamido)hexanoylamido)-propenyl]-2'-deoxycytidine-5'triphosphate, triethylammonium or tetralithium (optional) salt HC ΟН o N O ö

NH2

οн

Catalogue number: 303-22

Diluent: dd-H₂O

Storage: Freezer storage in dark, -10 to -30°C, dry.

Stability in Solution: 1 month as 1 mM solution in water or neutral buffers at +4°C in dark, 6 months as 1 mM solution in water or neutral buffers at -20°C in dark.



6-FAM-II-UTP

5-[3-(6-(fluoresceinyl-6-carboxamido)hexanoylamido)-propenyl]- uridine-5'-triphosphate, triethylammonium or tetralithium (optional) salt



Catalogue number: 303-32

Diluent: dd-H₂O

Storage: Freezer storage in dark, -10 to -30°C, dry.

Stability in Solution: 1 month as 1 mM solution in water or neutral buffers at +4°C in dark, 6 months as 1 mM solution in water or neutral buffers at -20°C in dark.






SIMA phosphoramidite

[(3',6'-dipivaloyl-2',7'-diphenyl-4,7-dichlorofluoresceinyl)-6-carboxamidohexyl]-I-O-(2-cyanoethyl)-(N,N-diisopropyl)-phosphoramidite

Catalog number: 303-55

SIMA is a full spectral analog of HEX in all fluorescent applications, but has significantly improved stability. It has a higher quantum yield and molecular extinction than HEX. SIMA is also more versatile in oligo synthesis especially due to improved stability during deprotection and simple purification of labeled oligos



Formula: C₅₈H₆₄Cl₂N₃O₁₀P M.W.: 1065.0

Diluent: Anhydrous Acetonitrile

Coupling: No changes needed from standard method recommended by synthesizer manufacturer.

Deprotection: No changes needed from standard method recommended by synthesizer manufacturer. SIMA labeled oligos are stable at 55°C in ammonium hydroxide (up to 6-8 hours) and can be deprotected with AMA.

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: 2-3 days, <90% efficient after 4 days

Store labeled oligo in the dark, either dry or in a neutral aqueous media at -20°C. Do not store crude fluorescently labeled oligonucleotides in ammonia.





Deprotection: No changes needed from standard method recommended by synthesizer manufacturer. (55 C in ammonium hydroxide up to 6-8 hours) and can be deprotected with AMA

Storage: Freezer storage, -10 to -30°C, dry. 1 year of storage guarantied





Catalogue Number: 303-57

Diluent: Anhydrous Acetonitrile

Coupling: No changes needed from standard method recommended by synthesizer manufacturer.

Deprotection: No changes needed from standard method recommended by synthesizer manufacturer. SIMA labeled oligos are stable at 55°C in ammonium hydroxide (up to 6-8 hours) and can be deprotected with AMA.

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: 2-3 days, <90% efficient after 4 days





Catalogue Number: 303-58

Deprotection: No changes needed from standard method recommended by synthesizer manufacturer. (55 C in ammonium hydroxide up to 6-8 hours)) and can be deprotected with AMA



5'-Fluorescein phosphoramidite (6-FAM)



Catalogue number: 303-41

Diluent: Anhydrous Acetonitrile

Coupling: 3 minute coupling time recommended

Deprotection: No changes needed from standard method recommended by synthesizer manufacturer.

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: 2-3 days, <90% efficient after 4 days

Store labeled oligo in the dark, either dry or in a neutral aqueous media at -20°C. Do not store crude fluorescently labeled oligonucleotides in ammonia.



5'-Fluorescein phosphoramidite (5-FAM)

[(3',6'-dipivaloylfluoresceinyl)-5-carboxamidohexyl]-I-O-(2-cyanoethyl)-(N,N-diisopropyl)-phosphoramidite



Catalogue number: 303-42

Diluent: Anhydrous Acetonitrile

Coupling: 3 minute coupling time recommended

Deprotection: No changes needed from standard method recommended by synthesizer manufacturer.

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: 2-3 days, <90% efficient after 4 days

Store labeled oligo in the dark, either dry or in a neutral aqueous media at -20°C. Do not store crude fluorescently labeled oligonucleotides in ammonia.



Fluorescein phosphoramidite (I)

(2S,4R)-N-(6-(3',6'-dipivaloylfluoresceinyl-5-carboxamido)hexanoylamido)-4-O-[(2-cyanoethyl)-(N,N-diisopropyl)-phosphoramidite]-2-(dimethoxytrityloxymethyl)pyrrolidine

Catalogue number: 303-43

Diluent: Anhydrous Acetonitrile

Coupling: 10 minute coupling time recommended

Deprotection: No changes needed from standard method recommended by synthesizer manufacturer.

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: 7-10 days, <90% efficient after 14 days



Store labeled oligo in the dark, either dry or in a neutral aqueous media at -20°C. Do not store crude fluorescently labeled oligonucleotides in ammonia.



Fluorescein phosphoramidite (II)

(2S,4R)-N-(6-(3',6'-dipivaloylfluoresceinyl-6-carboxamido)hexanoylamido)-4-O-[(2-cyanoethyl)-(N,N-diisopropyl)-phosphoramidite]-2-(dimethoxytrityloxymethyl)pyrrolidine



Store labeled oligo in the dark, either dry or in a neutral aqueous media at -20°C. Do not store crude fluorescently labeled oligonucleotides in ammonia.



6-Fluorescein dT phosphoramidite

5'-Dimethoxytrityl-5-[3-(6-(fluoresceinyl-6-carboxamido)hexanoylamido)-propenyl] uridine, 3'- (2-cyanoethyl)-(N,N-diisopropyl)-phosphoramidite



Diluent: Anhydrous Acetonitrile

Coupling: 10 minute coupling time recommended

Deprotection: No changes needed from standard method recommended by synthesizer manufacturer.

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: 7-10 days, <90% efficient after 14 days

Store labeled oligo in the dark, either dry or in a neutral aqueous media at -20°C. Do not store crude fluorescently labeled oligonucleotides in ammonia.



5-Fluorescein dT phosphoramidite

5'-Dimethoxytrityl-5-[3-(6-(fluoresceinyl-5-carboxamido)hexanoylamido)-propenyl] uridine-3'-(2-cyanoethyl)-(N,N-diisopropyl)-phosphoramidite



Store labeled oligo in the dark, either dry or in a neutral aqueous media at -20°C. Do not store crude fluorescently labeled oligonucleotides in ammonia.



HN

CPG

FAM Phosphoramidites and CPG

3'-(6-Fluorescein)-CPG

(2S,4R)-N-(6-(3',6'-dipivaloylfluoresceinyl-6-carboxamido)hexanoylamido)-4-O-(diglycoyl-long chain alkylamino-CPG)-2-(dimethoxytrityloxymethyl)pyrrolidine

Catalogue number: 303-47

Diluent: Not Applicable

Coupling: This support should be used in a manner identical to normal protected nucleoside support since it contains the DMT group.

Deprotection: No changes needed from standard method recommended by synthesizer manufacturer.

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: Not Applicable

Store labeled oligo in the dark, either dry or in a neutral aqueous media at -20°C. Do not store crude fluorescently labeled oligonucleotides in ammonia.

DMTrO



3'-(5-Fluorescein)-CPG

(2S,4R)-N-(6-(3',6'-dipivaloylfluoresceinyl-5-carboxamido)hexanoylamido)-4-O-(diglycoyl-long chain alkylamino-CPG)-2-(dimethoxytrityloxymethyl)pyrrolidine

I-Dimethoxytrityloxy-3-aza-3-[I-oxo-3-{(di-O-pivaloyl-fluorescein)-5-carboxamido} propyl]heptyl-6-Odiglycoyl- long chain alkylamino-CPG

Catalogue number: 303-48

Diluent: Not Applicable

Coupling: This support should be used in a manner identical to normal protected nucleoside support since it contains the DMT group.

Deprotection: No changes needed from standard method recommended by synthesizer manufacturer.

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: Not Applicable

Store labeled oligo in the dark, either dry or in a neutral aqueous media at -20°C. Do not store crude fluorescently labeled oligonucleotides in ammonia.

DMTrO



HEX Phosphoramidites

HEX phosphoramidite

[(4,7,2',4',5',7'-hexachloro-3',6'-dipivaloylfluoresceinyl)-6-carboxamidohexyl]-I-O-(2-cyanoethyl)-(N,N-diisopropyl)-phosphoramidite



Catalogue number: 303-51

Diluent: Anhydrous Acetonitrile

Coupling: 3 minute coupling time recommended

Deprotection: Ammonium Hydroxide for 24 hrs at room temperature.

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: 2-3 days, <90% efficient after 4 days

Store labeled oligo in the dark, either dry or in a neutral aqueous media at -20°C. Do not store crude fluorescently labeled oligonucleotides in ammonia.



HEX Phosphoramidites





Catalogue number: 303-59

Diluent: Anhydrous Acetonitrile

Coupling: 3 minute coupling time recommended

Deprotection: Ammonium Hydroxide for 24 hrs at room temperature.

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: 2-3 days, <90% efficient after 4 days

Store labeled oligo in the dark, either dry or in a neutral aqueous media at -20°C. Do not store crude fluorescently labeled oligonucleotides in ammonia.



TET Phosphoramidites

TET phosphoramidite

[(4,7,2',7'-tetrachloro-3',6'-dipivaloylfluoresceinyl)-6-carboxamidohexyl]-I-O-(2-cyanoethyl)-(N,N-diisopropyl)-phosphoramidite



Catalogue number: 303-61

Diluent: Anhydrous Acetonitrile

Coupling: 3 minute coupling time recommended

Deprotection: Ammonium Hydroxide for 24 hrs at room temperature.

Storage: Freezer storage, -10 to -30°C, dry.

Stability in Solution: 2-3 days, <90% efficient after 4 days

Store labeled oligo in the dark, either dry or in a neutral aqueous media at -20°C. Do not store crude fluorescently labeled oligonucleotides in ammonia.



TET Phosphoramidites



Store labeled oligo in the dark, either dry or in a neutral aqueous media at -20°C. Do not store crude fluorescently labeled oligonucleotides in ammonia



Biotin Phosphoramidites





Purine 3'-Amino -2', 3'-Dideoxynucleosides & Derivatives





Purine 3'-Amino -2', 3'-Dideoxynucleosides & Derivatives





Purine 3'-Amino -2', 3'-Dideoxynucleosides & Derivatives



Catalogue number: 203-30

IUPAC name: 3'- Amino-2',3'-dideoxy-2-fluoroadenosine

Purity: > 99%

Description: white to off-white crystals

Storage of dry compound: 3 years at +4°C

Resistant to enzymatic deamination

METKINEN chemistry

Purine 3'-Amino -2', 3'-Dideoxynucleosides & Derivatives

3'-Amino-2',3'-Dideoxy-2,6 Diaminopurineriboside

9-(3-Amino-2,3-dideoxy-ß-D-ribofuranosyl)- 2,6-diaminopurine



Catalogue number: 203-12

IUPAC name: 3'-Amino-2',3'-dideoxy-2,6-diaminopurineriboside

Purity: > 99%

Description: white to off-white crystals

Storage of dry compound: 3 years at +4°C

For bulk quantities please enquire















2'-Amino-2'-deoxyguanosine 9-(2-amino-2-deoxy-B-D-ribofuranosyl)-guanine NH₂ HO Catalogue number: 203-29 **CAS number:** 60966-26-9 IUPAC name: 2'-Amino-2'-deoxyguanosine **Purity:** > 99% **Description:** white to off-white crystals Storage of dry compound: 3 years at +4°C For bulk quantities please enquire





Catalogue number: 203-41

IUPAC name: 2'-Amino-2'-deoxyinosine

Purity: > 99%

Description: white to off-white crystals

Storage of dry compound: 3 years at +4°C

For bulk quantities please enquire







${\bf Arabino fur an osyl-6-B} enzylam in opurine$

9-(B-D-arabinoribofuranosyl)-6-Benzylaminopurine



Catalogue number: 203-20

IUPAC name: Arabinofuranosyl-6-Benzylaminopurine

Purity: > 99%

Description: white to off-white crystals

Storage of dry compound: 3 years at +4°C

For bulk quantities please enquire









Catalogue number: 203-32

CAS number: 21679-14-1

IUPAC name: Arabinofuranosyl-2-fluoroadenine

Purity: > 99%

Description: white to off-white crystals

Storage of dry compound: 3 years at +4°C

Resistant to enzymatic deamination





Purity: > 99%

Description: white to off-white crystals

Storage of dry compound: 3 years at +4°C

For bulk quantities please enquire





Purity: > 98%

Description: white to off-white crystals

Storage of dry compound: 3 years at +4°C

For bulk quantities please enquire







2-Chloro-2'-deoxyadenosine (2-CdA, cladribine) 9-(2-deoxy-β-D-ribofuranosyl)-2-chloro-adenine

CAS number: 4291-63-8

IUPAC name: 2-Chloro-2'-deoxyadenosine

Purity: > 99%

Description: white to off-white crystals

Storage of dry compound: 3 years at +4°C

For bulk quantities please enquire









Catalogue number: 203-36

IUPAC name: 2'-deoxy-2'-fluoro-2-fluoro-adenosine

Purity: > 99%

Description: white to off-white crystals

Storage of dry compound: 3 years at +4°C

Resistant to enzymatic deamination.


2'-Fluoro-2'-Deoxy-2,6-Diaminopurineriboside 9-(2-Deoxy-2-fluoro-B-D-ribofuranosyl)-2,6-diaminopurine NH_2 Ν ≥N HO NH_2 н Catalogue number: 203-18 IUPAC name: 2'-Deoxy-2'-fluoro-2,6diaminopurineriboside **Purity:** > 99% Description: white to off-white crystals Storage of dry compound: 3 years at +4°C For bulk quantities please enquire













For bulk quantities please enquire











Modifying nucleosides

Pyrene Derivative

3'-O-(diisopropylamino-2-cyanethoxyphosphinyl)-5'-O-(4,4'-dimethoxytrityl)-5-(pyren-I-ylethynyl)-2'-deoxyuridine



Diluent: Mixture CH3CN - CH2Cl2 (1:1), v/v

Coupling: 5 minute coupling time recommended

Storage: Freezer storage, -10 to -30°C, dry.

Stability: 3-4 days, <90% efficient after 5 days

Store labeled oligo in the dark.

5-(Pyren-1-ylethynyl)-2'-deoxyuridine contains pyrene chromophore conjugated with nucleobase through the triple bond. Its fluorescent properties are different from those of non-conjugated pyrene thus making this nucleoside suitable for hybridization studies including FRET, mismatch detection, electron injection studies, and other applications.



Modifying nucleosides

Pyrelene Derivative

3'-O-(N,N-diisopropylamino-2-cyanethoxyphosphinyl)-5'-O-(4,4'-dimethoxytrityl)-5-(perylen-3-ylethynyl)-2'-deoxyuridine



Appearance: Orange amorphous solid

Diluent: Mixture CH3CN - CH2Cl2 (1:1), v/v

Coupling: 5 minute coupling time recommended

Storage: Freezer storage, -10 to -30°C, dry.

Stability: 3-4 days, <90% efficient after 5 days

Store labeled oligo in the dark.

5-(Perylen-3-ylethynyl)-2'-deoxyuridine contains perylene conjugated to uracil. Emission wavelength maximum is 490 nm. In DNA probes this compound shows increase in fluorescence intensity.















C6 MMTr-Aminolinker Phosphoramidite $MMTrNH \longrightarrow O - P < \frac{N(iPr)_2}{OCNEt}$ Catalogue number: 501-04 Purity: > 90% Formula: $C_{35}H_{48}N_3O_3P$ Dilution (for 0.1M): 0.25g/4.2 ml 1.0g/17 ml

Description: Colorless to light yellow oil

Storage: Refrigerated, 2-8°C



Spacer Phosphoramidites





Acetylenic Amidites



A hydroxyprolinol-derived reagent (enantiomerically pure). Suitable for 5' and internal labeling. Stable to all ONS steps and ammonolysis. Oligonucleotides synthesized undergo smooth conjugation to azides in the presence of copper(I) species.



Acetylenic Amidites



Catalogue number: 501-07

Purity: > 95%

A 2,4-dihydroxybutyramide-derived reagent, enantiomerically pure. Suitable for 5' and internal labeling. Stable to all ONS steps. When on 5', stable to ammonolysis with DMT-on. Oligonucleotides synthesized undergo smooth conjugation to azides in the presence of copper(I) species. When the label is 5'-terminal, oligonucleotide should be subjected to ammonolysis only in DMT-on mode.

References:

- Dioubankova, N.N.; Malakhov, A.D.; Stetsenko, D.A.; Gait, M.J.; Korshun, V.A. Phosphoramidites and solid supports based on N-substituted 2,4-dihydroxybutyramides: universal reagents for synthesis of modified oligonucleotides. Tetrahedron, 2006, 62, 6762–6773.
- Ustinov, A.V.; Dubnyakova, V.V.; Korshun, V.A. Perylene Diimide-Oligonucleotide Conjugates Constructed by Click Chemistry. Nucleosides Nucleotides & Nucleic Acids, 2007, 26, 751–754.

 Ustinov, A.V.; Dubnyakova, V.V.; Korshun, V.A. A convenient "click chemistry" approach to perylene diimide–oligonucleotide conjugates. Tetrahedron, 2008, in press.



Acetylenic Amidites



Catalogue number: 501-08

Purity: > 95%

A 3,3-dimethyl-2,4-dihydroxybutyramide-derived reagent, enantiomerically pure. Suitable for 5' and internal labeling. Stable to all ONS steps. When on 5', stable to ammonolysis with DMTon. When the label is 5'-terminal, oligonucleotide should be subjected to ammonolysis only in DMT-on mode.

References:

 Dioubankova, N.N.; Malakhov, A.D.; Stetsenko, D.A.; Gait, M.J.; Korshun, V.A. Phosphoramidites and solid supports based on N-substituted 2,4-dihydroxybutyramides: universal reagents for synthesis of modified oligonucleotides. Tetrahedron, 2006, 62, 6762–6773.

ince	3						
Cat.#	Name	Page	Unit	Price, EUR	Price, USD*		
2'-Fluoro-2'-deoxynucleosides & Derivatives							
203-15	2'-Fluoro-2'-Deoxyguanosine	3	g 0 g 00 g kg	225 1500 4670 request	315 2100 6538 request		
203-17	2'-Fluoro-2'-Deoxyadenosine	4	lg 10g 100g lkg	210 1400 4000 request	295 1960 5600 request		
203-50	2'-Fluoro-2'-Deoxyguanosine-						
203-51	3'-CE-phosphoramidite	5	lg 10g 100g Ikg	225 2000 14000 request	315 2800 18900 request		
203-31	2'-Fluoro-2'-Deoxyadenosine- 3'-CE-phosphoramidite	6	lg 10g 100g lkg	220 2000 13500 request	308 2800 18300 request		
203-52	2'-Fluoro-2'-Deoxyuridine-		0				
203-53	2'-Fluoro-2'-Deoxycytidine-	7	l g 10 g 100 g 1 kg	80 650 4500 request	0 9 0 6300 request		
	3'-CE-phosphoramidite	8	l g 10 g 100 g 1 kg	80 650 4500 request	0 9 0 6300 request		
104-01	2'-Fluoro-2'-deoxyadenosine-5'-Triphosphate	9	10 µmol	75	105		
104-01	2'-Fluoro-2'-deoxycytidine-5'-Triphosphate	9	10 μmol	75	105		
104-02	2'-Fluoro-2'-deoxyguanosine-5'-Triphosphate	9 9	10 μmol	75 75	105		
104-03	2'-Fluoro-2'-deoxygualosille-5'-Triphosphate	9	10 μmol	75	105		
104-04	2 -Fluoro-2 -deoxyuridine-5 - Triphosphate 2'-Fluoro-2'-deoxynucleoside-5'-Triphosphate Kit, containing 5 µmol each	9 9	10 μmol 20 μmol, (4x5 μmol)	75 145	205		
103-30	2'-F-Nucleoside bound Solid Supports		(וטווון באד)	J	205		
103-30	(F-CS-supports)	П	lg I0g	40 35	55 50		



Prices

Cat.#	Name	Page	Unit	Price, EUR	Price, USD*		
Reagents for RNA and DNA synthesis							
03-00 03-10 03-20 03-21 03-3	Universal Solid Support III Chemical Phosphorylation Reagent II Trimer (Codon) Phosphoramidites Dimer Phosphoramidites 2'-Deoxynucleoside bound Solid Supports (CS-supports)	13 14 15 17 19		request request request request request	request request request request request		
2'-Flu	oro-2'-deoxyarabinonucleosides (FA	NA) 8	& Derivati	ves			
105-01	2'-FANA-A	21	0,5 g I g I0 g	300 550 3500	420 770 4900		
105-02	2'-FANA-DAP	22	0,5 g I g I0 g	310 570 3700	435 790 5180		
105-03	2'-FANA-G	23	0,5 g I g I0 g	330 600 3800	460 840 5320		
105-04	2'-FANA-I	24	0,5 g I g I0 g	320 590 3690	450 825 5165		
105-05	2'-FANA-T	25	0,5 g I g I0 g	0 90 90	155 165 1665		
105-06	2'-FANA-C	26	0,5 g I g I0 g	150 280 1750	210 390 2450		

Price	s						
Cat.#	Name	Page	Unit	Price, EUR	Price, USD*		
Biotin Labeled Triphosphates							
303-31	Biotin-11-UTP	27	I μMol I0 μMol I00 μMol	360 2755 11020	505 3860 15430		
303-11	Biotin-11-dUTP	28	I μMol I0 μMol I00 μMol	230 2175 8700	320 3045 12180		
303-21	Biotin-11-dCTP	29	I μMol I0 μMol I00 μMol	230 2175 8700	320 3045 12180		
303-71	Biotin-11-dATP	30	0,25 μMol Ι μMol Ι0 μMol	380 1450 8265	530 2030 11570		
303-14	Biotin-16-dUTP	31	I μMol I0 μMol I00 μMol	230 2175 8700	320 3045 12180		
303-72	Biotin-11-dGTP	32	0,25 μMol Ι μMol Ι0 μMol	380 1450 8265	530 2030 11570		
Fluors	cein Labeled Triphosphates						
303-12	6-FAM-11-dUTP	33	0,5 μMol Ι μMol Ι0 μMol Ι00 μMol	300 550 2755 8265	420 770 3860 15430		
303-22	6-FAM-11-dCTP	34	0,5 μMol Ι μMol Ι0 μMol Ι00 μMol	300 550 2755 11020	420 770 3860 15430		

Price	s						
Cat.#	Name	Page	Unit	Price, EUR	Price, USD*		
303-32	6-FAM-11-UTP	35	0,5 μMol Ι μMol Ι0 μMol Ι00 μMol	300 550 2755 11020	420 770 3860 15430		
303-80	6-FAM-11-dATP	36	0,25 μMol Ι μMol Ι0 μMol	425 1650 8265	595 2310 11570		
SIMA Phosphoramidites and CPG							
303-55	SIMA phosphoramidite	37	0,5 g I g I0 g	230 435 4250	320 610 5950		
303-56	6-SIMA-dT	38	250 mg I g	360 1160	505 1625		
303-57	5-SIMA phosphoramidite	39	0.5 g I g I0 g	230 435 4250	320 610 5950		
303-58	6-SIMA-CPG	40	l g	725	1015		
FAM F	Phosphoramidites and CPG						
303-41	5'-Fluorescein phosphoramidite (6-FAM)	41	l g 10 g	290 2900	405 4060		
303-42	5'-Fluorescein phosphoramidite (5-FAM)	42	l g 10 g	290 2900	405 4060		
303-43	Fluorescein phosphoramidite (I)	43	0,5 g I g I0 g	230 435 3265	320 610 5075		
303-44	Fluorescein phosphoramidite (II)	44	0,5 g I g I0 g	230 435 3625	320 610 5075		

Prices								
Cat.#	Name	Page	Unit	Price, EUR	Price, USD*			
303-45	6-Fluorescein dT phosphoramidite	45	0,5 g I g I0 g	360 680 5075	505 950 7105			
303-46	5-Fluorescein dT phosphoramidite	46	0,5 g I g I0 g	360 680 5075	505 950 7105			
303-47	3'-(6-Fluorescein)-CPG	47	l g 10 g	220 1450	310 2030			
303-48	3'-(5-Fluorescein)-CPG	48	l g 10 g	220 1450	310 2030			
HEX F	HEX Phosphoramidites							
303-51	HEX phosphoramidite	49	0,5 g I g I0 g	230 435 4350	320 610 6090			
303-59	5-HEX phosphoramidite	50	l g 10 g	290 2900	405 4060			
TET P	hosphoramidites							
303-61	TET phosphoramidite	51	0,5 g I g I0 g	230 435 4350	320 610 6090			
303-65	5-TET phosphoramidite	52	l g 10 g	290 2900	410 4060			
Biotin	Phosphoramidites							
303-70	5'-Biotin Phosphoramidite	53	250 mg I g	110 290	155 405			

Price	s							
Cat.#	Name	Page	Unit	Price, EUR	Price, USD*			
Purine 3'-Amino-2',3'-Dideoxynucleosides & Derivatives								
203-11	3'- Amino-2',3'-dideoxyadenosine	54	250 mg I g I0 g	40 90 600	55 125 840			
203-10	3'- Amino-2',3'-dideoxyguanosine	55	250 mg I g I0 g	50 120 800	70 170 1120			
203-30	3'- Amino-2',3'-dideoxy-2-fluoroadenosine	56	250 mg I g I0 g	115 270 1800	160 380 2520			
203-12	3'-Amino-2',3'-dideoxy- 2,6-diaminopurineriboside	57	250 mg I g 10 g	45 105 700	60 150 980			
Purine	e Nucleosides							
203-27	2'-Amino-2'-deoxyadenosine	58	250 mg I g I0 g	150 345 2300	210 485 3220			
203-31	2'-Amino-2'-deoxy-2-fluoroadenosine	59	250 mg I g I0 g	220 525 3500	310 735 4900			
203-28	2'-Amino-2'-deoxy-2,6-diaminopurineriboside	60	250 mg I g I0 g	130 300 2000	180 420 2800			
203-29	2'-Amino-2'-deoxyguanosine	61	250 mg I g I0 g	160 375 2500	225 525 3500			
203-41	2'-Amino-2'-deoxyinosine	62	250 mg I g I0 g	130 300 2000	180 420 2800			

Prices

Cat.#	Name	Page	Unit	Price, EUR	Price, USD*
203-05	Arabinofuranosyl-adenine	63	250 mg I g I0 g	20 45 300	30 65 420
203-20	Arabinofuranosyl-6-Benzylaminopurine	64	250 mg I g I0 g	85 195 1300	120 275 1820
203-06	Arabinofuranosyl-2,6- diaminopurine	65	250 mg I g I0 g	20 45 300	30 65 420
203-32	Arabinofuranosyl-2-fluoroadenine	66	250 mg I g I0 g	100 225 1500	140 315 2100
203-03	Arabinofuranosyl-guanine	67	250 mg I g I0 g	35 75 500	50 105 700
203-37	Arabinofuranosyl-isoguanine	68	250 mg I g I0 g	290 675 4500	405 945 6300
203-21	6-Benzylaminopurine 2'-deoxyriboside	69	250 mg I g I0 g	85 195 1300	120 275 1820
203-25	2-Chloro-2'-deoxyadenosine	70	250 mg I g I0 g	80 220 1500	115 310 2100
203-35	2'-deoxy-2-fluoroadenosine	71	250 mg I g I0 g	85 195 1300	120 275 1820
203-36	2'-deoxy-2'-fluoro-2-fluoroadenosine	72	250 mg I g I0 g	290 675 4500	410 945 6300

Prices

Cat.#	Name	Page	Unit	Price, EUR	Price, USD*
203-18	2'-Deoxy-2'-fluoro-2,6-diaminopurineriboside	73	250 mg I g I0 g	100 240 1600	140 340 2240
203-14	2'-Deoxy-2'-fluoroinosine	74	250 mg I g I0 g	85 195 1300	120 275 1820
203-38	2'-Deoxyisoguanosine	75	250 mg I g I0 g	130 300 2000	180 420 2800
203-40	2'-Deoxy-2'-fluoroisoguanosine	76	250 mg I g I0 g	290 675 4500	410 945 6300
203-23	2,6-Diaminopurine 2'-deoxyriboside	77	250 mg I g I0 g	35 75 500	50 105 700
203-39	Isoguanosine	78	250 mg I g I0 g	170 390 2600	240 545 3640
Modify	ying Nucleosides				
403-03	Pyrene Derivative	79	250 mg I g I0 g	200 670 6000	280 940 8400
403-06	Pyrelene Derivative	80	250 mg I g I0 g	280 960 8700	395 1345 12180

Prices

Cat.#	Name	Page	Unit	Price, EUR	Price, USD*			
Amino	Aminolinker Phosphoramidites							
501-01 501-02 501-03 501-04	C6 TFA Aminolinker Phosphoramidite DMT C6-Aminolinker Phosphoramidite C3 MMTr-Aminolinker Phosphoramidite C6 MMTr-Aminolinker Phosphoramidite	81 82 83 84	250 mg 250 mg 250 mg 250 mg	25 40 55 40	35 55 80 55			
Space	r Phosphoramidites							
501-05	Spacer 18 Phosphoramidite	85	250 mg	85	120			
Acety	lenic Amidites							
501-06	Reagent I-amidite	86	250 mg I g I0 g	50 540 4800	210 755 6720			
501-07	Reagent 2-amidite	87	250 mg I g 10 g	130 430 3900	180 600 5460			
501-08	Reagent 3-amidite	88	250 mg I g 10 g	100 360 2900	140 505 4060			



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