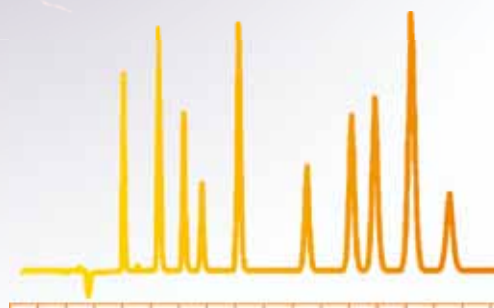


# Chromatography

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## Pesticides from water

MN Appl. No. 304250 / 304260

Gleis, C., GIU GmbH, Teningen

Matrix: water

**Columns:** CHROMABOND® HR-X, 3 mL, 200 mg  
REF 730931

**Sample pretreatment:** samples are spiked with 500 ng of each pesticide in 1000 mL water, adjusted to pH 2 with HCl or pH 7

**Conditioning:** 10 mL MeOH, 10 mL dist. water

**Sample application:** slowly pass 1000 mL spiked water sample through the column with the aid of a tubing adaptor (REF 730243).

**Elution:** after drying 5 mL MeOH / THF (1:1; v/v)

**Subsequent analysis:** HPLC

**Recovery rates [%]:**

compound	pH 2	compound	pH 7
metamitron	86	desisopropylatrazine	90
quinmerac	90	2,4-dichlorobenzamide	95
chloridazon	93	desethylatrazine	89
picloram	83	hexazinone	95
metribuzin	84	bromacil	103
cyanazine	83	simazine	91
metabenzthiazuron	94	desethylterbutylazine	89
chlortoluron	91	atrazine	88
isoproturon	89	metalaxyl	97
diuron	91	metazachlor	93
dimethenamid-P	89	propazine	88
linuron	94	terbutylazine	86
epoxyconazole	85	metolachlor	97
penconazole	90		
alachlor	93		
1-propiconazole	89		
flufenacet	91		
diflufenicam	58		
triallate	42		

## Extraction of paraquat and diquat from water

MN Appl. No. 305370

Matrix: water

**Column:** CHROMABOND® HR-XCW, 3 mL, 60 mg  
REF 730735

**Conditioning:** 1 mL MeOH, then 1 mL water (Do not let run the column dry!)

**Sample application:** 1 mL spiked water (0.5 µg/mL)

**Washing:** 1 mL MeOH

**Drying:** with nitrogen or air

**Elution:** 1 mL acetonitrile / water (1:1; v/v) + 2 % formic acid; Evaporation and reconstitution is suitable solvent.

**Subsequent analysis:** HPLC, e.g. NUCLEODUR® HILIC (see page 15)

**Substances:** paraquat; diquat

**Recovery rates:** paraquat: 103 %, diquat: 113 %

## Pesticides from water

MN Appl. No. 302060

Matrix: water

**Column:** CHROMABOND® Hydra, 6 mL, 2000 mg  
REF 730301

**Sample pretreatment:** adjust 1000 mL water to pH 7-8 with diluted NH<sub>3</sub> and add 100 µL of the internal standard (1 µg/L).

**Conditioning:** 2 x 5 mL MeOH, then 2 x 5 mL dist. water

**Sample application:** slowly force or aspirate the water sample through the column. Then dry for 2 h with 2 bar N<sub>2</sub>.

**Elution:** aspirate 10 mL MeOH slowly through the column. Evaporate the eluted solution to dryness in a tapered flask with a rotation evaporator at 30 °C and store it in a refrigerator for about 15 min. Dissolve the residue in 200 µL cold, fresh n-hexane and transfer the solution in a conical HPLC vial. Store the solution in a refrigerator until GC analysis.

**Subsequent analysis:** GC columns: OPTIMA® δ-3 or OPTIMA® δ-6; HPLC column: NUCLEODUR® PolarTec

**Substances:** desisopropylatrazine; desethylatrazine; desethylterbutylazine; simazine; atrazine; propazine; terbutylazine; sebuthylazine; desmetryn; ametryn; prometryne; terbutryn; cyanazine; metazachlor; napropamide; methoprotryn; hexazinone

**Recovery rates:** between 95 - 100 %

## Method development of an extraction of very polar pesticide metabolites - optimization of MeOH portion

MN Appl. No. 304400

Pantiru, M.E., Diss., Wissenschaftszentrum Weihenstephan für Ernährung, Landnutzung und Umwelt, TU München

**Column:** CHROMABOND® HR-P, 3 mL, 200 mg  
REF 730108

**Sample pretreatment:** samples of 100 mL phosphate buffer (pH 6) with 20, 10 or 5 % MeOH are spiked with 5 µg carbamate metabolites each

**Conditioning:** 3 mL MeOH, then 3 x 3 mL phosphate buffer (pH 6)

**Sample application:** with tubing adaptor (e.g. CHROMABOND® tubing adaptor, REF 730243) the prepared sample is passed through the column (vacuum: 4 - 5 mm Hg).

**Elution:** after drying in a nitrogen stream (15 min) it is eluted with 2 x 3 mL acetone

**Concentration:** the eluate is concentrated with nitrogen in a 5 mL test tube (30 °C). The residue is reconstituted with 500 µL MeOH and passed into a HPLC vial.

**Subsequent analysis:** HPLC

**Recovery rates [%]:**

compound	20 % MeOH	10 % MeOH	5 % MeOH
butocarboxim sulfoxid	5	82	103
aldicarb sulfon	98	83	102
ethiofencarb sulfon	102	88	101
methiocarb sulfoxid	97	91	102
carbofuran 3-OH	99	99	100
methiocarb sulfon	96	100	102
carbofuran 3-keto	99	99	101



## PAHs from drinking water (EPA 550)

MN Appl. No. 302170

Matrix: water

**Column:** CHROMABOND® C<sub>18</sub> ec, 6 mL, 1000 mg  
REF 730015

**Sample pretreatment:** add 100 mg Na<sub>2</sub>S<sub>2</sub>O<sub>3</sub> to 1000 mL water sample and adjust the pH value to pH 2 with 6 N HCl

**Conditioning:** 4 x 10 mL methylene chloride, 4 x 10 mL MeOH, then 4 x 10 mL ultra pure water

**Sample application:** suck or press the water sample through the column.

**Washing:** with 10 mL ultra pure water, then dry the column for 10 minutes with vacuum

**Elution:** suck 2 x 5 mL methylene chloride slowly through the column. Dry the combined fractions over Na<sub>2</sub>SO<sub>4</sub>. Filter the suspension and wash with 2 mL methylene chloride. Concentrate the sample to 1 mL under a nitrogen stream. For further analysis add 3 mL acetonitrile and concentrate to 0.5 mL with vacuum.

**Subsequent analysis:** HPLC, e.g. NUCLEODUR® C<sub>18</sub> PAH (see page 16)

## PAHs from water

MN Appl. No. 301250

Matrix: water

**Column:** CHROMABOND® C<sub>18</sub> PAH, 6 mL, 2000 mg  
REF 730166

**Conditioning:** 1 column volume MeOH, then 1 column volume dist. water

**Sample application:** aspirate 1000 mL water sample through the column (about 15 to 20 mL/min), then dry the column (stream of nitrogen or 24 h in a desiccator over P<sub>2</sub>O<sub>5</sub>).

**Elution:** elute with 4 mL acetonitrile / benzene\* (3:1; v/v) and evaporate or fill up to the volume required

\*: alternatively toluene can be used

**Subsequent analysis:** HPLC, e.g. NUCLEODUR® C<sub>18</sub> PAH (see page 16)

**Recovery rates [%]:**

compound (each 50 ng/L)	%
naphthalene	87
acenaphthylene	89
acenaphthene	90
fluorene	82
phenanthrene	85
anthracene	90
fluoranthene	89
pyrene	89
benz[a]anthracene	87
chrysene	95
benzo[b]fluoranthene	91
benzo[k]fluoranthene	89
benzo[a]pyrene	90
dibenz[ah]anthracene	97
benzo[ghi]perylene	91
indeno[1,2,3-cd]pyrene	96

## 16 EPA-PAHs from water with CHROMABOND® Easy

MN Appl. No. 302830

Matrix: water

**Column:** CHROMABOND® Easy, 6 mL, 200 mg  
REF 730755

**Sample pretreatment:** concentration of the standard: 1 µg/L

**Sample:** drinking water

**Conditioning:** 3 mL MeOH, 3 mL water

**Sample application:** 500 mL drinking water

**Washing:** 3 mL water / MeOH 5%

**Drying:** vacuum

**Elution:** 2 x 2 mL dichlormethane for GC analysis or for HPLC: concentrate dichlormethane in a nitrogen atmosphere and elute in ACN.

**Subsequent analysis:** HPLC, e.g. NUCLEODUR® C<sub>18</sub> PAH (see page 16)

## PAHs from water containing humic acids

MN Appl. No. 301260

Matrix: water

**Column:** CHROMABOND® NH<sub>2</sub> / C<sub>18</sub>, 6 mL,  
500/1000 mg (glass columns)  
REF 730620G

**Sample pretreatment:** add 25 mL 2-propanol to 500 mL water sample

**Conditioning:** 10 mL methylene chloride, 10 mL MeOH, then 10 mL dist. water / 2-propanol (9:1; v/v)

**Sample application:** 250 mL sample solution

**Washing:** 2 mL dist. water / 2-propanol (9:1; v/v)

**Drying:** about 20 min, vacuum

**Elution:** elute with 4x1 mL methylene chloride (let percolate first 1 mL into the column packing without vacuum, then apply light vacuum) and if necessary evaporate in a stream of nitrogen and fill up with a suitable solvent.

**Subsequent analysis:** HPLC, e.g. NUCLEODUR® C<sub>18</sub> PAH (see page 16)

**Recovery rates [%]:**

compound	%	compound	%
naphthalene	90	benzo[a]anthracene	88
acenaphthylene	89	chrysene	95
acenaphthene	86	benzo[b]fluoranthene	93
fluorene	87	benzo[k]fluoranthene	88
phenanthrene	87	benzo[a]pyrene	87
anthracene	89	dibenz[ah]anthracene	91
fluoranthene	90	benzo[ghi]perylene	90
pyrene	93	indeno[1,2,3-cd]pyrene	89



## 16 EPA-PAHs from soil with CHROMABOND® Easy

MN Appl. No. 302820

*Matrix:* sludge, soil

*Column:* CHROMABOND® Easy, 6 mL, 200 mg  
REF 730755

*Sample pretreatment:* homogenize 5 g soil sample in 30 mL 2-propanol, filter the solution and fill up to 250 mL with water

*Sample:* soil or sludge

*Conditioning:* 3 mL MeOH, 3 mL water

*Sample application:* 250 mL sample solution

*Washing:* 2 x 3 mL water / MeOH (95:5; v/v)

*Drying:* vacuum

*Elution:* 2 x 1 mL dichlormethane for GC analysis or for HPLC analysis: concentrate dichlormethane in a nitrogen atmosphere and dissolve in acetonitrile.

*Subsequent analysis:* HPLC, e.g. NUCLEODUR® C<sub>18</sub> PAH (see page 16)

## PCBs from soil, sludge or cement

MN Appl. No. 302040

Göldner, E., Inst. für Angewandte Chemie Gockel & Weischedel & Co.

*Matrix:* cement, sludge, soil

*Column:* CHROMABOND® NAN, 3 mL, 400/1400/400 mg  
REF 730109

*Sample pretreatment:* ca. 20 g of a dry soil, sludge, cement sample with 10 g sodium sulfate are extracted for 8 hours with 100 mL n-hexane in a soxhlet extractor. Then the extract is concentrated to a volume of 1-2 mL.

*Conditioning:* 2 mL n-hexane

*Sample application:* slowly force or aspirate the extract of the sample pretreatment through the column

*Elution:* aspirate 10 mL n-hexane slowly through the column. Then concentrate to 5 mL.

To gain a better peak shape in GC a subsequent acetonitrile extraction is performed: Extract 2 mL of the n-hexane-extract with 2 mL acetonitrile. After phase partition use ca. 1 mL of the acetonitrile phase for GC analysis.

*Subsequent analysis:* GC, e.g. OPTIMA® δ-6, as described in application number 250480 ([www.mn-net.com/apps](http://www.mn-net.com/apps))

*Recovery rates [%]:*

PCB	extraction rates n-hexane to acetonitril
28	30.8
52	36.9
101	26.8
153	19.1
138	23.4
180	15.9

## PAHs from soil

MN Appl. No. 301310

*Matrix:* soil

*Column:* CHROMABOND® CN/SiOH, 6 mL, 500/1000 mg  
REF 730135

*Sample pretreatment:* dry 30 g soil with sodium sulphate and reflux 4 hours with 250 mL petroleum ether in a soxhlet extractor. For low PAH contents (colorless or weakly colored extracts) concentrate extract to 1/10 of its volume in a rotation evaporator.

*Conditioning:* 4 mL petroleum ether

*Sample application:* force or aspirate 20 mL of the extract through the column

*Elution:* 2 x 2 mL acetonitrile / benzene (3:1; v/v), then evaporate or fill up to the volume required (alternative 2 x 2 mL acetonitrile / toluene (3:1; v/v))

*Subsequent analysis:* HPLC, e.g. NUCLEODUR® C<sub>18</sub> PAH (see page 16)

*Recovery rates [%]:*

compound	%	compound	%
naphthalene	85	benzo[a]anthracene	84
acenaphthylene	92	chrysene	96
acenaphthene	89	benzo[b]fluoranthene	95
fluorene	87	benzo[k]fluoranthene	90
phenanthrene	83	benzo[a]pyrene	90
anthracene	88	dibenz[ah]anthracene	96
fluoranthene	87	benzo[ghi]perylene	87
pyrene	90	indeno[1,2,3-cd]pyrene	97

## Phenols in aqueous samples (EPA 528)

MN Appl. No. 302370

Ciube, R.; Spangenberg, J.; Wild, G.; Meyer, Th., GIT Fachz. Lab 8/96

*Matrix:* water

*Column:* CHROMABOND® HR-P (e.g. 6 mL, 200 mg  
REF 730119 or sorbent REF 730615)

*Sample pretreatment:* water samples with two different concentration ranges (0.1 and 0.003 mg/kg) were prepared without pH adjustment or desalting with MeOH as auxiliary solvent.

*Conditioning:* 10 mL MeOH and then 4 mL water. The column must not run dry.

*Sample application:* 50 mL (0.1 mg/kg) or 500 mL (0.003 mg/kg) of the prepared samples are sucked through the column. Dry the column by sucking air for 1 minute.

*Elution:* 4 mL MeOH

*Concentration:* the eluate is concentrated to 1 or 0.5 mL by heating to 70 °C.

*Subsequent analysis:* GC

*Recovery rates [%]:*

compound	0.1 mg/L	0.003 mg/L
phenol	99.0 +/-1.0	98.8 +/-0.9
o-cresol	99.1 +/-0.7	97.9 +/-0.8
m- + p-cresol	99.2 +/-0.7	97.5 +/-0.8
2,6-xylenol	97.8 +/-1.2	99.2 +/-0.7
2,4- + 2,5-xylenol	99.7 +/-0.7	98.6 +/-0.8
3,4-xylenol	98.9 +/-0.6	97.4 +/-0.7
2,3-xylenol	99.0 +/-0.4	98.9 +/-0.6
3,5-xylenol	99.4 +/-0.3	98.5 +/-0.7



## Enrichment of explosive compounds from water samples with SPE

MN Appl. No. 302910

Bausinger, T., Geographisches Inst. Universität Mainz

Matrix: water

**Column:** CHROMABOND® HR-P, 6 mL, 200 mg  
REF 730119

**Sample pretreatment:** test mixture contains 5 µg/L of each compound

**Conditioning:** 5 mL MeOH and then 3 mL acetonitrile and 10 mL water. The column must not run dry.

**Sample application:** suck 400 mL of the sample with 10-15 mL/min through the column, wash with water and suck the column to dryness for 40-60 min

**Elution:** apply 1 mL MeOH / acetonitrile (50:50; v/v) onto the column. The sorbent should become wet but do not suck the solvent through the column. Wait for 5 min and elute with 1.5 mL of the same solvent.

**Subsequent analysis:** fill the eluate up to 4 mL with water and then HPLC, e.g. NUCLEODUR® PolarTec (see page 17)

**Recovery rates [%]:**

compound	%
octogene	114 +/- 6
2,4-diamino-6-nitrotoluene	63 +/- 25
hexogene	99 +/-14
1,3,5-trinitrobenzene	105 +/-2
2-amino-6-nitrotoluene	102 +/-1
2-amino-4-nitrotoluene	101 +/-1
1,3-dinitrobenzene	106 +/-1
2,4,6 trinitrotoluene	103 +/-1
4-amino-2,6-dinitrotoluene	107 +/-1
2-amino-4,6-dinitrotoluene	107 +/-2
2,6-dinitrotoluene	98 +/-2
2,4-dinitrotoluene	97 +/-1
2-nitrotoluene	103 +/-2
4-nitrotoluene	87 +/-1
3-nitrotoluene	92 +/-6



## AOX determination in water with high content of salt or organic compounds according to DIN 38409 - H 22

MN Appl. No. 302080

Matrix: water

**Column:** CHROMABOND® HR-P-AOX, 6 mL, 500 mg  
REF 730111.AOX

**Conditioning:** 5 mL MeOH, then 10 mL dist. Water, the column should not run dry.

**Sample application:** force or aspirate 100 mL of the original or diluted water sample (pH 1) through the column (3-5 mL/min). The column should not run dry. Discard the eluted solution. Wash with 50 mL of a nitrate washing solution (dissolve 17 g NaNO<sub>3</sub> in 100 mL dist. water, add 1.4 mL HNO<sub>3</sub> 10 M), fill up with dist. water to 1000 mL take off 50 mL and fill up with dist. water to 1000 mL). Discard the eluted solution.

**Elution:** slowly force or aspirate 1 x 1 mL, then 1 x 4 mL MeOH and 10 mL dist. water through the column. Collect eluted solutions in 100 mL measuring flask and fill up with dist. water to 100 mL.

**Note:** we recommend to test the recovery with a standard solution. Is the recovery to low, an increase of elution solvent is recommended (2 x 5 mL MeOH or 2 x 5 mL MeOH/ ethyl acetate 1:1; v/v). If the sample has only a small contamination (conc. DOC < 500 mg/L) the adsorbent weight in the cartridge can be reduced for a better recovery, i.e. with CHROMABOND® HR-P-AOX, 6 mL, 200 mg, REF 730119.AOX

## Hydrocarbons in water according to ISO DIS 9377-4 / DIN H-53

MN Appl. No. 302090

Matrix: water

**Column:** CHROMABOND® Na<sub>2</sub>SO<sub>4</sub> / Florisil®,  
6 mL, 2000/2000 mg, glass column  
REF 730249G

**Internal standard solution:** solve 20 mg n-tetracontane (C<sub>40</sub>H<sub>82</sub>) in cyclohexane, add 20 mL n-nonane (C<sub>9</sub>H<sub>20</sub>) and fill up to 1 L with cyclohexane. To prepare the extraction solution dilute the standard solution with cyclohexane 1:10.

**Sample pretreatment:** adjust 900 mL water (10 °C) with HCl (12 mol/L) to pH 2 and add 80 g MgSO<sub>4</sub>. Add 50 mL of the extraction solution, close the bottle and stir the suspension intensely for 30 min. Add enough distilled water to separate the organic from the aqueous phase.

**Conditioning:** 5 mL cyclohexane

**Sample application:** slowly force or aspirate the organic solution through the column.

**Elution:** wash with 10 mL cyclohexane. Evaporate the combined organic solutions carefully to 1 mL or less. If necessary, fill up to 1 mL exactly. (Evaporation to 1 mL can be unnecessary, if the hydrocarbon content is high.)

**Recovery rates:** must be > 80 % for n-tetracontane.

**Subsequent analysis:** GC, according to application number 210600 on OPTIMA® 1 ([www.mn-net.com/apps](http://www.mn-net.com/apps))



## Plasticizers (phthalates and adipates) from drinking water (EPA 506)

**MN Appl. No. 302160**

*Matrix:* water

**Column:** CHROMABOND® C18 ec, 3 mL, 500 mg (glass column) REF 730013G

*Sample pretreatment:* add 5 mL MeOH to 1000 mL water sample.

*Conditioning:* 2 x 10 mL methylene chloride, 2 x 10 mL MeOH, then 10 mL ultra pure water.

*Sample application:* suck or press up to 1000 mL water sample through the column.

*Washing:* with 10 mL ultra pure water

*Elution:* suck 10 mL methylene chloride slowly through the column. Concentrate the sample under nitrogen stream to about 0.5 mL and dry over Na<sub>2</sub>SO<sub>4</sub>.

*Subsequent analysis:* GC according to EPA 606 on OPTIMA® 1 or OPTIMA® 5 column, see GC application numbers 201210 and 201220 ([www.mn-net.com/apps](http://www.mn-net.com/apps))

## Extraction of perfluorinated surfactants from water

**MN Appl. No. 305140**

*Matrix:* water

**Column:** CHROMABOND® HR-XAW, 3 mL, 60 mg REF 730747

*Conditioning:* 2 mL MeOH + 5% ammonia, then 2 mL MeOH and at last 2 mL water. Do not let run the column dry.

*Sample application:* 500 mL water sample, spiked with 1 mL standard solution (concentration: 20 µg/L of each compound)

*Washing:* 2 mL water, then 2 mL acetone / acetonitrile / formic acid, (50:50:1; v/v/v) and at last 2 mL MeOH

*Elution:* 2 mL MeOH + 5 % ammonia

*Evaporation:* nitrogen stream with slight heating to dryness

*Reconstitution:* HPLC solvent

*Subsequent analysis:* LC-MS, e.g. NUCLEODUR® Sphinx RP (see page 18)

*Recovery rates [%]:*

compound	%
perfluoropropionic acid (PFPrA)	103
perfluoropentanoic acid (PFPeA)	94
perfluorohexanoic acid	94
perfluorooctanoic acid	95
perfluorooctane sulfonate potassium salt (PFOS)	81
perfluorododecanoic acid (PFDoDA)	82

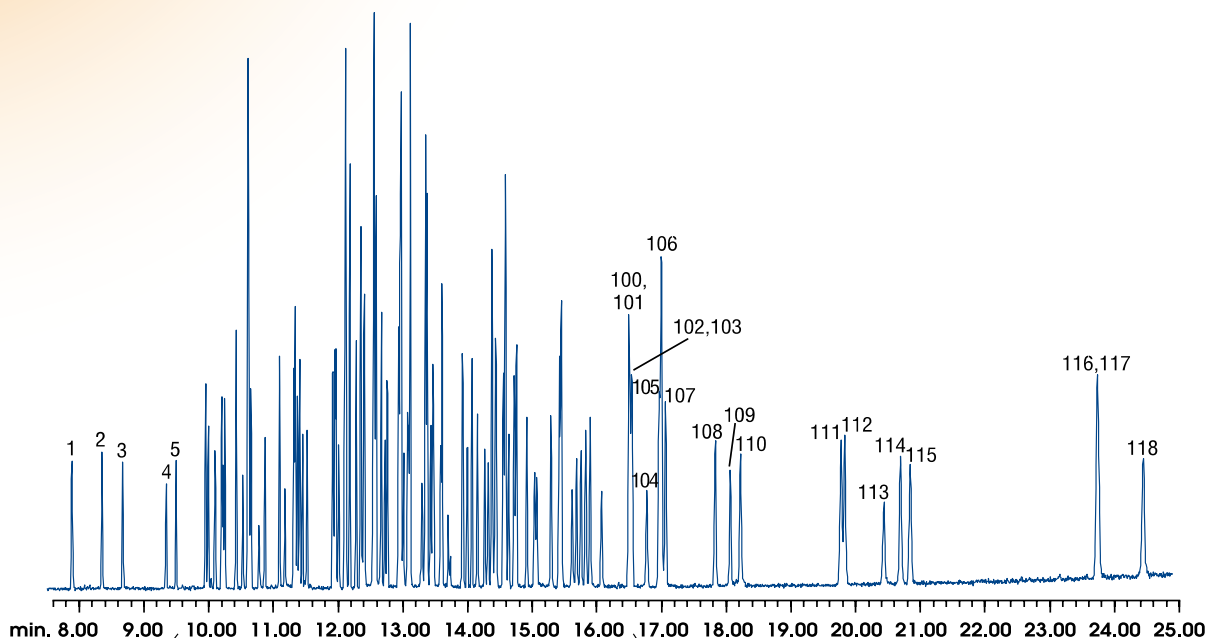




# Pesticide analysis

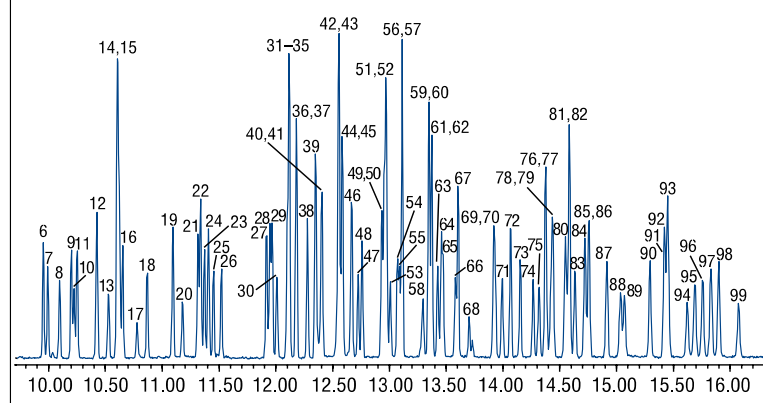
## Separation of PAHs, pesticides, phthalates (EPA 525)

MN Appl. No. 212810



### Peaks:

1. isophorone, 2. 2-nitro-m-xylene, 3. dichlorvos,
4. hexachlorocyclopentadiene, 5. EPTC, 6. butylate,
7. mevinphos, 8. vernolate, 9. pebulate, 10. etridiazole (terrazole),
11. dimethylphthalate, 12. acenaphthene,
13. 2,6-dinitrotoluene, 14. acenaphthene-d10,
15. 2-chlorobiphenyl, 16. chloroneb, 17. tebuthiuron,
18. molinate, 19. diethyl phthalate, 20. 2,4-dinitrotoluene,
21. propachlor, 22. fluorene, 23. ethoprop, 24. cycloate,
25. trifluralin, 26. chlorpropham, 27. 2,3-dichlorobiphenyl,
28. atraton, 29. prometon, 30. a-BHC, 31. hexachlorobenzene,
32. propazine, 33. simazine, 34. atrazine, 35. metribuzin,
36. diazinon, 37. terbufos, 38. pronamide, 39. pentachlorophenol,
40. b-BHC, 41. disulfoton, 42. terbacil, 43. phenanthrene-d10,
44. methyl parathion, 45. phenanthrene, 46. anthracene,
47. g-BHC (lindane), 48. 2,4,5-trichlorobiphenyl, 49. alachlor,
50. prometryne, 51. ametryn, 52. simetryn, 53. d-BHC,
54. heptachlor, 55. chlorothalonil, 56. di-n-butylphthalate,
57. terbutryn, 58. bromacil, 59. chlorpyrifos, 60. metolachlor,
61. DCPA methyl ester (Dacthal®), 62. 2,2',4,4'-tetrachlorobiphenyl,
63. aldrin, 64. triadimefon, 65. cyanazine (Bladex), 66. MGK-264,
67. diphenamid, 68. merphos, 69. 2,2',3',4,6-penta-chlorobiphenyl,
70. heptachloro epoxide (isomer B), 71. heptachloro epoxide (isomer A), 72. butachlor, 73. stirofos (tetrachlorvinphos),
74. fenamiphos, 75. a-chlordane, 76. napropamide,
77. g-chlordane, 78. endosulfan I, 79. trans-nonachlor,
80. pyrene-d10, 81. pyrene, 82. 4,4'-DDE
83. 2,2',4,4',5,6'-hexachloro-biphenyl, 84. p-terphenyl-d14,
85. Dieldrin, 86. carboxin, 87. chlorbenzilate, 88. tricyclazole,
89. endrin, 90. 4,4'-DDD, 91. bis(2-ethylhexyl)adipate,
92. butyl benzyl phthalate, 93. endosulfan II, 94. endrin aldehyde,
95. norflurazon, 96. 4,4'-DDT, 97. triphenylphosphate,
98. hexazinone, 99. endosulfan sulfate,
100. bis(2-ethylhexyl)phthalate, 101. methoxychlor,
102. 2,2',3,3',4,5',6,6'-octachloro-biphenyl,
103. 2,2',3,3',4,4',6-heptachloro-biphenyl, 104. endrin ketone,
105. benzo(a)anthracene, 106. chrysene-d12, 107. chrysene,
108. fenarimol, 109. cis-permethrin, 110. trans-permethrin,
111. benzo(b)fluoranthene, 112. benzo(k)fluoranthene,
113. fluridone (Sonar®), 114. benzo(a)pyrene, 115. perylene-d12,
116. dibenzo(a,h)anthracene, 117. indeno(1,2,3-cd)pyrene,
118. benzo(ghi)perylene



### OPTIMA® XLB, 30 m, 0.25mm ID, 0.25 µm

REF 725850.30

**Sample:** US EPA method 525 standards, 1 µL, 5 ng / compound

**Injection:** pressure pulsed (0.4 min 30 psi), splitless (for 0.4 min)

**Inj. temperature:** 300°C

**Carrier gas:** helium, 1.0 mL/min

**Temperature:** 35 °C (for 2 min) → 260 °C at 20 °C/min  
260 °C → 330 °C at 6 °C/min, 330 °C (for 5 min)

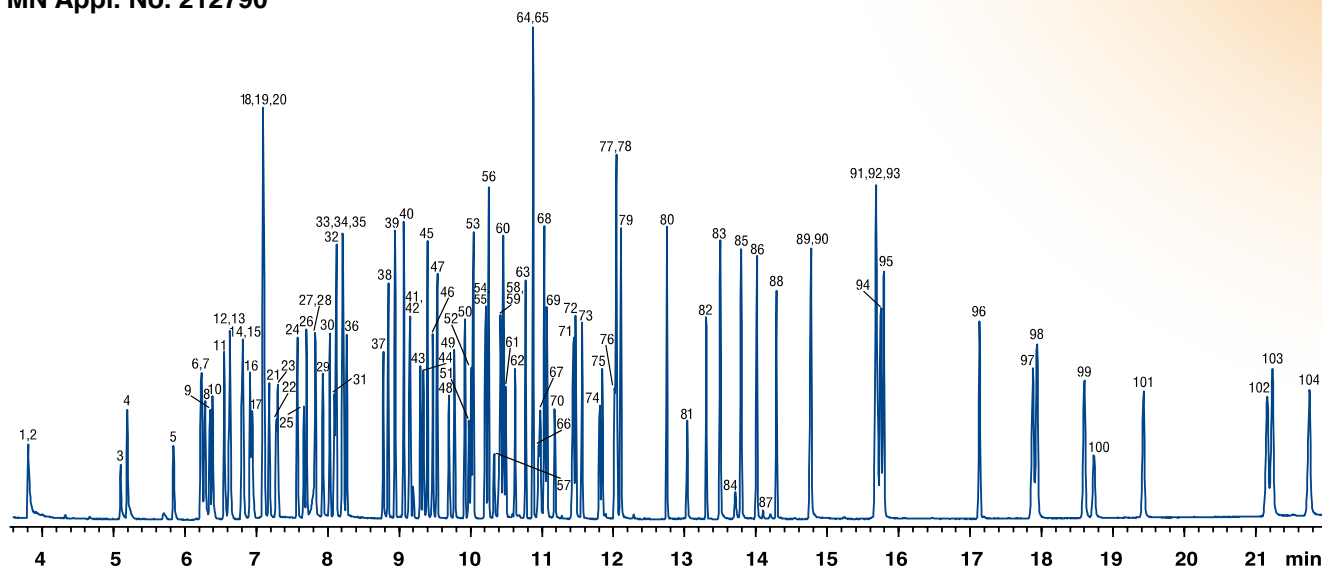
**Detection:** GC / MS, 280 °C, scan range: 45 - 550 amu





## Separation of pesticides (EPA 8270)

MN Appl. No. 212790



**OPTIMA® 5 MS Accent, 30 m, 0.25 mm ID, 0.25 µm**

REF 725820.30

**Sample:** 16 µg/mL in methylene chloride

**Injection:** 1 µL splitless (for 0.4 min)

**Inj. temperature:** 300 °C

**Carrier gas:** helium, 1.0 mL/min

**Temperature:** 35 °C (for 2 min) → 260 °C at 20 °C/min  
→ 330 °C at 6 °C/min (for 1 min)

**Detection:** GC/MS, 280 °C, scan range: 35-550 amu

### Peaks:

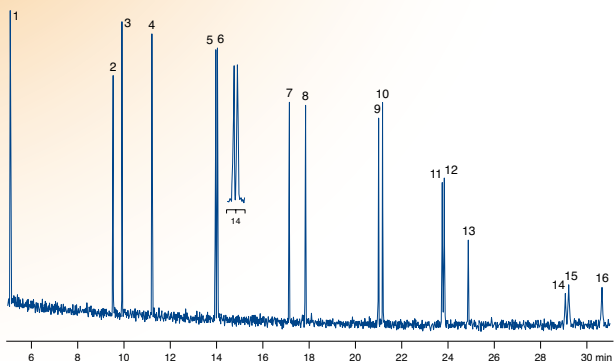
1. N-nitrosodimethylamine, 2. pyridine, 3. methyl methanesulfonate, 4. 2-fluorophenol, 5. ethyl methanesulfonate, 6. phenol-d6, 7. phenol, 8. aniline, 9. bis(2-chloroethyl)ether, 10. 2-chlorophenol, 11. 1,3-dichlorobenzene, 12. 1,4-dichlorobenzene-d4, 13. 1,4-dichlorobenzene, 14. 1,2-dichlorobenzene, 15. benzyl alcohol, 16. 2-methylphenol, 17. bis(2-chloroisopropyl)ether, 18. acetophenone, 19a. 4-methylphenol, 19b. 3-methylphenol, 20. N-nitroso-di-n-propylamine, 21. hexachloroethane, 22. nitrobenzene-d5, 23. nitrobenzene, 24. isophorone, 25. 2-nitrophenol, 26. 2,4-dimethylphenol, 27. bis(2-chloro-ethoxy) methane, 28. benzoic acid, 29. 2,4-dichlorophenol, 30. 1,2,4-trichlorobenzene, 31. naphthalene-d8, 32. naphthalene, 33. 2,6-dichlorophenol, 34. 4-chloroaniline, 35. hexachloropropene, 36. hexachlorobutadiene, 37. 4-chloro-3-methylphenol, 38. Isosafrole, 39. 2-methyl-naphthalene, 40. 1-methylnaphthalene, 41. hexachlorocyclopentadiene, 42. 1,2,4,5-tetrachlorobenzene, 43. 2,4,6-trichlorophenol, 44. 2,4,5-trichlorophenol, 45. 2-fluorobiphenyl, 46. safrole, 47. 2-chloronaphthalene, 48. 2-nitroaniline, 49. 1,4-naphthoquinone, 50. dimethylphthalate, 51. 1,3-dinitrobenzene, 52. 2,6-dinitrotoluene, 53. acenaphthylene, 54. acenaphthene-d10, 55. 3-nitroaniline, 56. acenaphthene, 57. 2,4-dinitrophenol, 58. pentachlorobenzene, 59. 4-nitrophenol, 60. dibenzofurane, 61. 2,4-dinitrotoluene, 62. 2,3,4,6-tetrachlorophenol, 63. diethyl phthalate, 64. fluorene, 65. 4-chlorophenyl phenyl ether, 66. 4-nitro-aniline, 67. 4,6-dinitro-2-methylphenol, 68. di-phenylamine, 69. azobenzene, 70. 2,4,6-tribromophenol, 71. phenacetin, 72. 4-bromo-phenyl phenyl ether, 73. hexachlorobenzene, 74. pentachlorophenol, 75. pentachloronitrobenzene, 76. phenanthrene-d10, 77. dinoseb, 78. phenanthrene, 79. anthracene, 80. di-n-butylphthalate, 81. 4-nitro-quinolin-1-oxide, 82. isodrin, 83. fluoranthene, 84. benzidine, 85. pyrene, 86. p-terphenyl-d14, 87. aramite, 88. chlorbenzilate, 89. kepone, 90. butyl benzyl phthalate, 91. benzo(a)anthracene, 92. 3,3'-dichloro-benzidine, 93. chrysene-d12, 94. chrysene, 95. bis(2-ethylhexyl)phthalate, 96. di-n-octylphthalate, 97. benzo(b)fluor-anthene, 98. benzo(k)fluoranthene, 99. benzo(a)pyrene, 100. perylene-d12, 101. 3-methylcholanthrene, 102. indeno(1,2,3-cd)pyrene, 103. dibenzo(a,h)anthracene, 104. benzo(ghi)perylene





## PAHs acc. to EPA 610

MN Appl. No. 213190



**OPTIMA® 35 MS, 30 m, 0.25mm ID, 0.25 µm**  
REF 726154.30

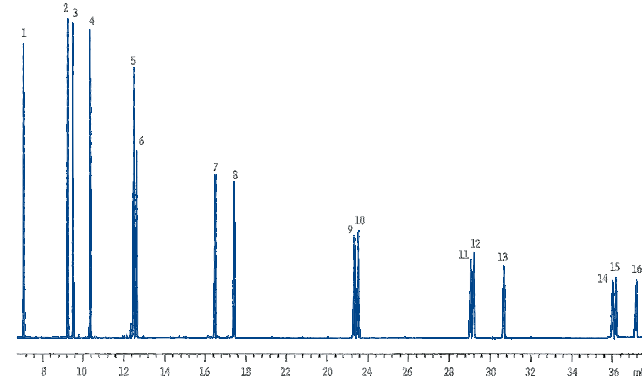
**Injection:** 1.0 µL H<sub>2</sub>, split 1:10  
**Carrier gas:** helium, 0.6 bar  
**Temperature:** 100 °C (for 3 min) → 300 °C at 6 °C/min,  
300 °C (for 10 min)  
**Detector:** MSD

### Peaks:

1. naphthalene, 2. acenaphthylene, 3. acenaphthene, 4. fluorene, 5. phenanthrene, 6. anthracene, 7. fluoranthene, 8. pyrene, 9. benzo[a]anthracene, 10. chrysene, 11. benzo[b]fluoranthene, 12. benzo[k]fluoranthene, 13. benzo[a]pyrene, 14. indeno[1,2,3-cd]pyrene, 15. dibenz[ah]anthracene, 16. benzo[ghi]perylene

## Separation of PAHs

MN Appl. No. 212800



**OPTIMA® 5 MS Accent, 30 m, 0.25 mm ID, 0.25 µm**  
REF 725820.30

**Sample:** 1 µL of 20 ng/µL, PAH Mix  
**Injection:** splitless (for 1 min)  
**Inj. temperature:** 300 °C  
**Carrier gas:** hydrogen, 40 cm/sec.  
**Temperature:** 40 °C (for 1 min) → 200 °C at 20 °C/min,  
200 °C → 310 °C at 4 °C/min,  
310 °C (for 5 min)  
**Detection:** FID, 310 °C

### Peaks:

1. naphthalene, 2. acenaphthylene, 3. acenaphthene, 4. fluorene, 5. phenanthrene, 6. anthracene, 7. fluoranthene, 8. pyrene, 9. benzo(a)anthracene, 10. chrysene, 11. benzo(b)fluoranthene, 12. benzo(k)fluoranthene, 13. benzo(a)pyrene, 14. indeno(1,2,3-cd)pyrene, 15. dibenzo(a,h)anthracene, 16. benzo(ghi)perylene

## Determination of PCBs and PAHs with PCB 28 / PCB 31 separation in less than 10 min

MN Appl. No. 212920

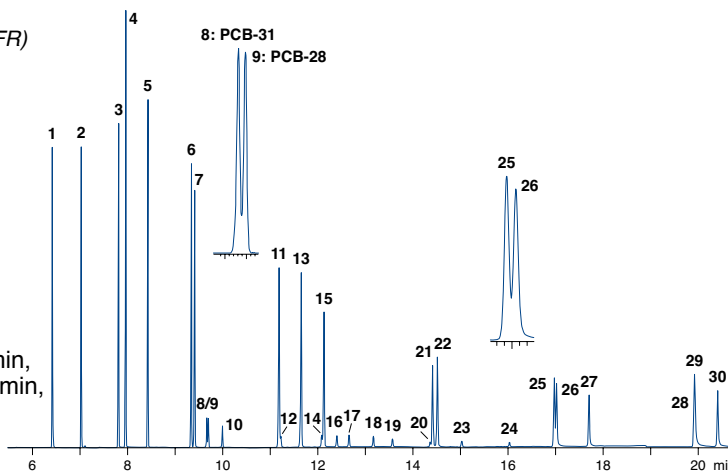
Centre d'Analyses de Recherche, Lab. D'Hydrologie, Illkirch (FR)

**OPTIMA® XLB, 30 m, 0.25mm ID, 0.25 µm**  
REF 725850.30

**Inj. volume:** 1 µL, standard 0.005 ng/µL  
**Inj. temperature:** 250 °C  
**Injection:** pulsed, splitless  
**Inj. pulse:** 1.38 bar in 1 min  
**Purge flow:** 60 mL/min  
**Carrier gas:** helium  
**Temperature:** 40 °C (for 2 min) → 240 °C at 30 °C/min,  
240 °C (for 2 min) → 340 °C at 10 °C/min,  
340 °C (for 5 min)  
**Detection:** MS source: 230 °C  
Interface: 280 °C  
Quadrupol: 150 °C

### Peaks:

1. naphthalene, 2. 2-methylnaphthalene, 3. acenaphthylene, 4. acenaphthene, 5. fluorene, 6. phenanthrene, 7. anthracene, 8. PCB 31, 9. PCB 28, 10. PCB 52, 11. Fluoranthene, 12. PCB 101, 13. Pyrene, 14. PCB 77, 15. 2-methylfluoranthene, 16. PCB 118, 17. PCB 153, 18. PCB 138, 19. PCB 126, 20. PCB 180, 21. Benzo(a)anthracene, 22. Chrysene, 23. PCB 169, 24. PCB 194, 25. Benzo(b)fluoranthene, 26. Benzo(k)fluoranthene, 27. Benzo(a)pyrene, 28. Dibenzo(ah)anthracene, 29. Indeno(123cd)pyrene, 30. Benzo(ghi)perylene



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## Analysis of phenols in accordance with EPA 604 on medium-polar ultra low bleed column

MN Appl. No. 213600

OPTIMA® 17 MS, 30 m, 0.25mm ID, 0.25 µm

REF 726162.30

Sample: phenol-mix 604

Inj. volume: 1 µL

Inj. temperature: 230 °C

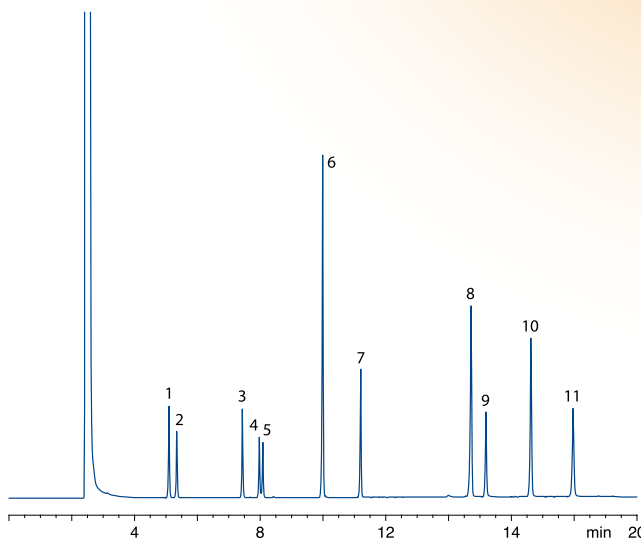
Carrier gas: helium, 0.8 bar, split 1:30

Temperature: 100 °C → 250 °C at 10 °C/min

Detection: FID, 280 °C

### Peaks:

1. phenol, 2. 2-chlorophenol, 3. 2,4-dimethylphenol,  
4. 2-nitrophenol, 5. 2,4-dichlorophenol, 6. 4-chloro-3-methylphenol,  
7. 2,4,6-trichlorophenol, 8. 4-nitrophenol, 9. 2,4-dinitrophenol,  
10. 2-methyl-4,6-dinitrophenol, 11. pentachlorophenol



## Analysis of isomeric phenols

MN Appl. No. 250060

OPTIMA® 5-3, 60 m, 0.25 mm ID, 0.25 µm

REF 726420.60

Injection: 1.0 µL, split 1:80

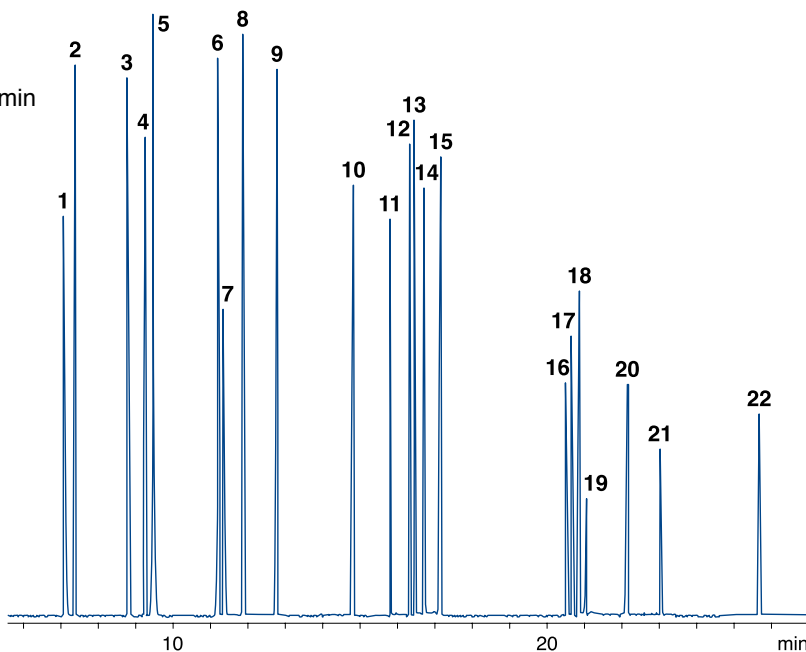
Carrier gas: helium, 1.3 bar

Temperature: 60 °C (3 min) → 320 °C, 6 °C/min

Detector: MSD HP 5971

### Peaks:

1. phenol, 2. 2-chlorophenol, 3. 2-methylphenol,  
4. 4-methylphenol, 5. 3-methylphenol,  
6. 2,4-dimethylphenol, 7. 2-nitrophenol,  
8. 2,4-dichlorophenol, 9. 2,6-dichlorophenol,  
10. 4-chloro-3-methylphenol, 11. 2,3,5-trichlorophenol,  
12. 2,4,6-trichlorophenol, 13. 2,4,5-trichlorophenol,  
14. 2,3,4-trichlorophenol, 15. 2,3,6-trichlorophenol,  
16. 2,3,5,6-tetrachlorophenol,  
17. 2,3,4,5-tetrachlorophenol,  
18. 2,3,4,6-tetrachlorophenol,  
19. 2,4-dinitrophenol, 20. 3,4,5-trichlorophenol,  
21. 2-methyl-4,6-dinitrophenol,  
22. 2-isopropyl-4,6-dinitrophenol





# Volatiles analysis

## EPA 502 / EPA 524 volatile organics calibration mix (VOCs)

MN Appl. No. 211280

OPTIMA® 624, 50 m, 0.25 mm, 1.40 µm

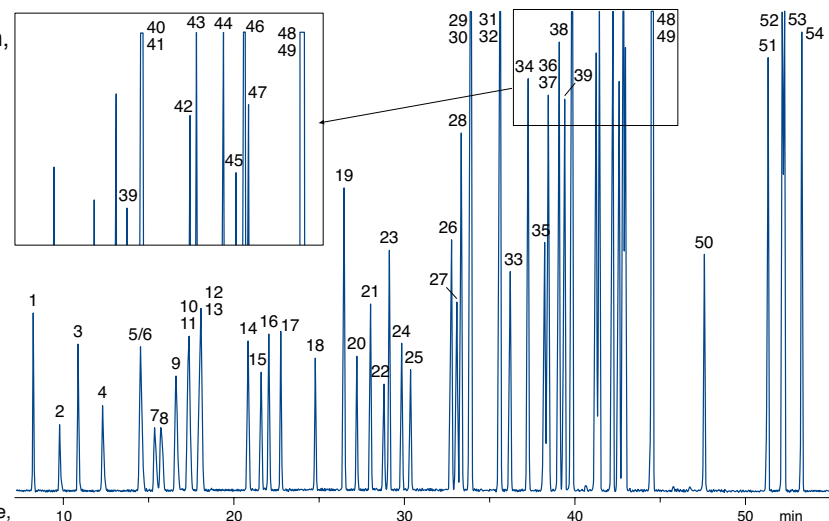
REF 726785.50

**Injection:** 1 µL, 280 °C, split 1: 50

**Carrier gas:** helium, 1.5 bar

**Temperature:** 40 °C (for 5 min) → 70 °C at 2.5 °C/min,  
70 °C → 100 °C at 3.0 °C/min,  
100 °C → 220 °C at 4.0 °C/min,  
220 °C (for 5 min)

**Detector:** MSD 5971 [HP]  
Scanmode, thres: 300,  
EMV: 2400



### Peaks:

1. 1,1-dichloroethene, 2. dichloromethane (methylene chloride), 3. trans-1,2-dichloroethene,
4. 1,1-dichloroethane, 5. 2,2-dichloropropane,
6. is-1,2-dichloroethene, 7. bromochloromethane,
8. trichloromethane (chloroform), 9. 1,1,1-trichloroethane,
10. 1,1-dichloropropene, 11. tetrachloromethane (carbon tetrachloride), 12. 1,2-dichloroethane, 13. benzene,
14. trichloroethene, 15. 1,2-dichloropropane, 16. bromodichloromethane, 17. dibromomethane, 18. cis-1,3-dichloropropene, 19. toluene,
20. trans-1,3-dichloropropene, 21. 1,1,2-trichloroethane, 22. tetrachlorethene, 23. 1,3-dichloropropane, 24. dibromochloromethane, 25. 1,2-dibromoethane,
26. chlorobenzene, 27. ethylbenzene, 28. 1,1,1,2-tetrachloroethane, 29. m-xylene, 30. p-xylene, 31. o-xylene, 32. styrene, 33. tribromomethane (bromoform),
34. isopropylbenzene, 35. 1,1,2,2-tetrachloroethane, 36. 1,2,3-trichloropropane, 37. bromobenzene, 38. n-propylbenzene, 39. 2-chlorotoluene,
40. 1,3,5-trimethylbenzene, 41. 4-chlorotoluene, 42. tert-butylbenzene, 43. 1,2,4-trimethylbenzene, 44. sec-butylbenzene, 45. p-isopropyltoluene,
46. 1,3-dichlorobenzene, 47. 1,4-dichlorobenzene, 48. n-butylbenzene, 49. 1,2-dichlorobenzene 50. 1,2-dibromo-3-chloropropane, 51. 1,2,4-trichlorobenzene,
52. hexachlorobutadiene, 53. naphthalene, 54. 1,2,3-trichlorobenzene

## EPA 502 / EPA 524 volatile organics calibration mix (VOCs)

MN Appl. No. 211300

OPTIMA® δ-3, 50 m, 0.20 mm, 0.20 µm

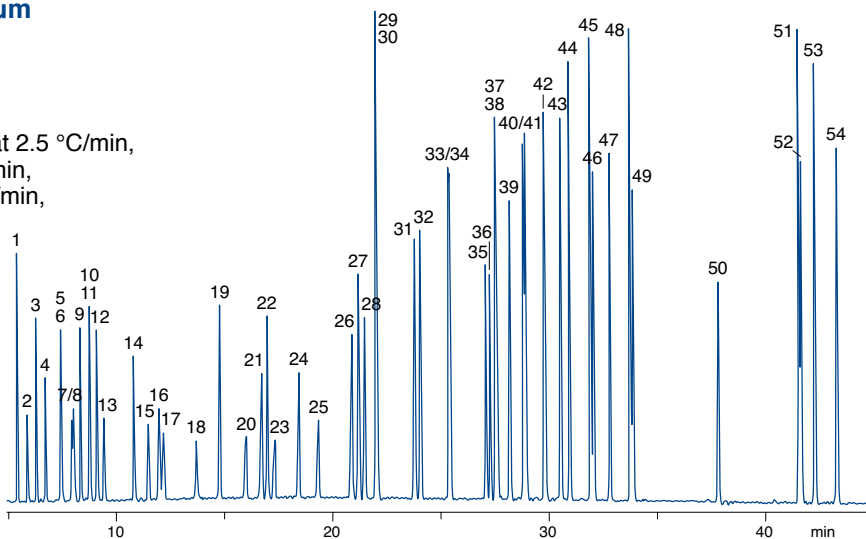
REF 726400.50

**Injection:** 1 µL, 280 °C, split 1: 40

**Carrier gas:** helium, 1.8 bar

**Temperature:** 40 °C (for 5 min) → 70 °C at 2.5 °C/min,  
70 °C → 100 °C at 3.0 °C/min,  
100 °C → 160 °C at 4.0 °C/min,  
160 °C (for 5 min)

**Detector:** MSD 5971 [HP]  
Scanmode, thres: 400,  
EMV: 1976



### Peaks:

1. 1,1-dichloroethene, 2. dichloromethane (methylene chloride), 3. trans-1,2-dichloroethene,
4. 1,1-dichloroethane, 5. 2,2-dichloropropane,
6. is-1,2-dichloroethene, 7. bromochloromethane,
8. trichloromethane (chloroform), 9. 1,1,1-trichloroethane, 10. 1,1-dichloropropene,
11. tetrachloromethane (carbon tetrachloride), 12. 1,2-dichloroethane, 13. benzene,
14. trichloroethene, 15. 1,2-dichloropropane, 16. bromodichloromethane, 17. dibromomethane, 18. cis-1,3-dichloropropene, 19. toluene, 20. trans-1,3-dichloropropene,
21. 1,1,2-trichloroethane, 22. tetrachlorethene, 23. 1,3-dichloropropane, 24. dibromochloromethane, 25. 1,2-dibromoethane, 26. chlorobenzene,
27. ethylbenzene, 28. 1,1,1,2-tetrachloroethane, 29. m-xylene, 30. p-xylene, 31. o-xylene, 32. styrene, 33. tribromomethane (bromoform), 34. isopropylbenzene,
35. 1,1,2,2-tetrachloroethane, 36. 1,2,3-trichloropropane, 37. bromobenzene, 38. n-propylbenzene, 39. 2-chlorotoluene, 40. 1,3,5-trimethylbenzene
41. 4-chlorotoluene, 42. tert-butylbenzene, 43. 1,2,4-trimethylbenzene, 44. sec-butylbenzene, 45. p-isopropyltoluene, 46. 1,3-dichlorobenzene,
47. 1,4-dichlorobenzene, 48. n-butylbenzene, 49. 1,2-dichlorobenzene 50. 1,2-dibromo-3-chloropropane, 51. 1,2,4-trichlorobenzene, 52. hexachlorobutadiene,
53. naphthalene, 54. 1,2,3-trichlorobenzene



## Analysis of hydrocarbons

MN Appl. No. 210780

Inst. für Meereskunde, Kiel

**OPTIMA® δ-6, 30 m, 0.25 mm ID, 0.25 µm**

REF 726470.30

**Sample:** 10 ng/µL

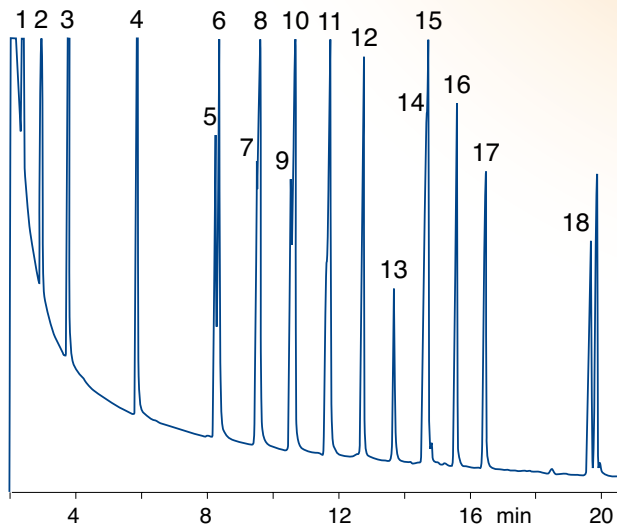
**Flow:** 36 cm/sec

**Carrier gas:** hydrogen

**Temperature:** 80 °C (for 1 min) → 320 °C at 6 °C/min,  
320 °C (for 10 min)

### Peaks:

1. C-10, 2. C-11, 3. C-12, 4. C-14, 5. C-16:1, 6. C-16, 7. C-17:1,  
8. C-17, 9. C-18:1, 10. C-18, 11. C-19, 12. C-20, 13. C-21, 14. C-22:1,  
15. C-22, 16. C-23, 17. C-24, 18. C-28



## Analysis of phthalates in accordance with EPA 8060 on medium-polar ultra low bleed column

MN Appl. No. 213610

**OPTIMA® 17 MS, 30 m, 0.25 mm ID, 0.25 µm**

REF 726162.30

**Injection:** 1.0 mL, 280 °C, 0.5 min splitless, 25 mL/min

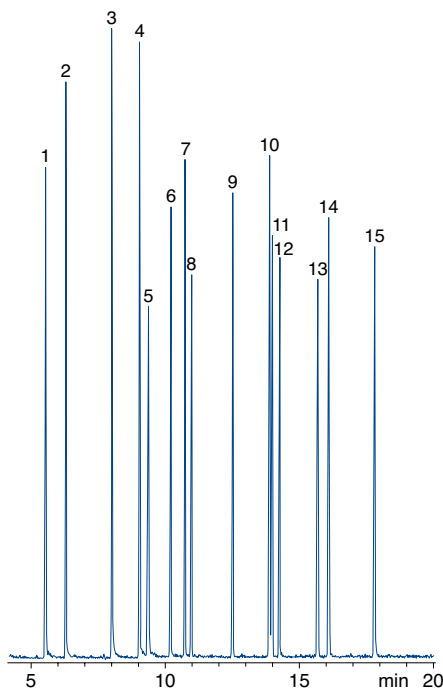
**Carrier gas:** helium, 0.6 bar

**Temperature:** 120 °C → 220 °C at 25 °C/min,  
220 °C → 330 °C at 8 °C/min, 330 °C (for 10 min)

**Detector:** MSD

### Peaks:

1. Dimethyl phthalate, 2. Diethyl phthalate,  
3. Di-isobutyl phthalate, 4. Di-n-butyl phthalate,  
5. Bis (4-methyl-2-pentyl) phthalate, 6. Bis (2-methoxyethyl) phthalate  
7. Di-n-pentyl phthalate, 8. Bis (2-ethoxyethyl) phthalate, 9. Di-n-hexyl phthalate,  
10. Bis (2-ethylhexyl) phthalate, 11. Benzyl-butyl phthalate,  
12. Bis-(2-butoxyethyl) phthalate, 13. Di-cyclohexyl phthalate,  
14. Di-n-octyl phthalate, 15. Di-n-nonyl phthalate





# Phthalates analysis

## Separation of phthalates (EPA 606)

MN Appl. No. 213160

OPTIMA®  $\delta$ -3, 30 m, 0.25 mm ID, 0.25  $\mu$ m

REF 726420.30

*Inj. volume:* 1  $\mu$ L, split 1:10

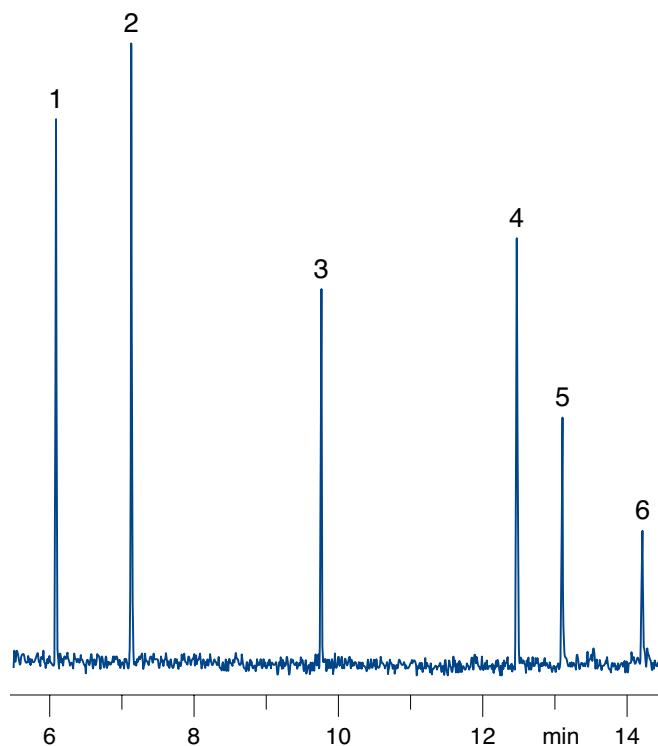
*Carrier gas:* hydrogen, 1.4 mL/min

*Temperature:* 100 °C  $\rightarrow$  320 °C at 15 °C/min

*Detection:* MSD

### Peaks:

1. dimethyl phthalate,
2. diethyl phthalate,
3. dibutyl phthalate,
4. benzyl butyl phthalate,
5. bis(2-ethylhexyl) phthalate,
6. di-n-octyl phthalate





## Pesticides

MN Appl. No. 124480

**Column:** EC 250/3 NUCLEODUR® PolarTec, 3 µm  
REF 760479.30

**Conditions**

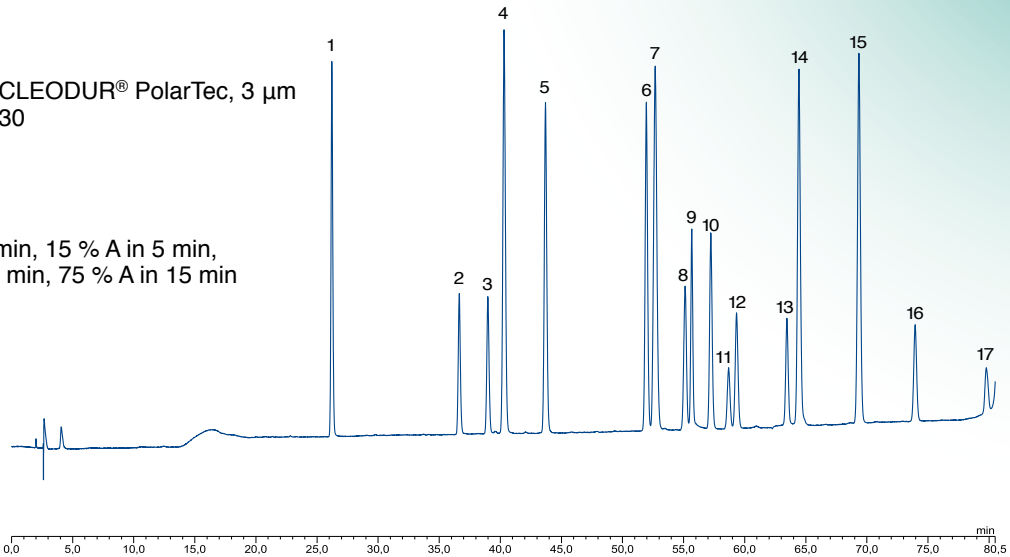
**Eluent:** A) MeOH  
B) water

**Gradient:** 2 % A for 10 min, 15 % A in 5 min,  
43 % A for 60 min, 75 % A in 15 min

**Flow rate:** 0.56 mL/min

**Temperature:** 35 °C

**Detection:** 230 nm



**Peaks:**

1. desethylatrazine, 2. metoxuran, 3. hexazinone, 4. simazine, 5. cyanazine, 6. methabenzthiazuron, 7. atrazine, 8. chlorotoluron, 9. monolinuron, 10. isoproturon, 11. metazachlor, 12. diuron, 13. metobromuron, 14. sebuthylazine, 15. terbuthylazine, 16. linuron, 17. metolachlor

## Separation of paraquat and diquat on NUCLEODUR® HILIC with LC-MS detection

MN Appl. No. 123060

**Column:** EC 125/2 NUCLEODUR® HILIC, 3 µm  
REF 760531.20

**Substances:** paraquat; diquat

**Concentration:** 500 µg/mL

**Conditions**

**Eluent:** acetonitrile / 50 mM ammonium formate,  
pH 3.2 (80:20; v/v)

**Flow rate:** 0.3 mL/min

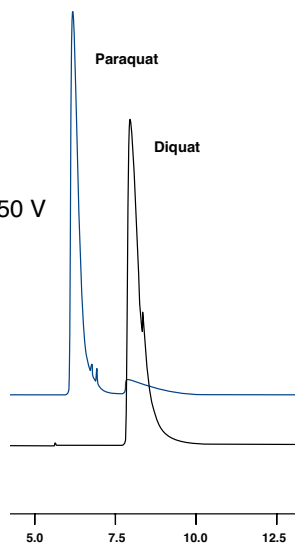
**Inj. volume:** 1.0 µL

**Detection:** LC-MS

**Mode:** ESI+

**Sample temp.:** 450 °C

**Needle:** 3.0 kV  
Cone voltage: 50 V



## Separation of carbamate pesticides (EPA 531.1)

MN Appl. No. 124220

**Column:** EC 150/4.6 NUCLEODUR® C<sub>18</sub> HTec, 3 µm  
REF 760325.46

**Conditions**

**Eluent:** A) water  
B) acetonitrile

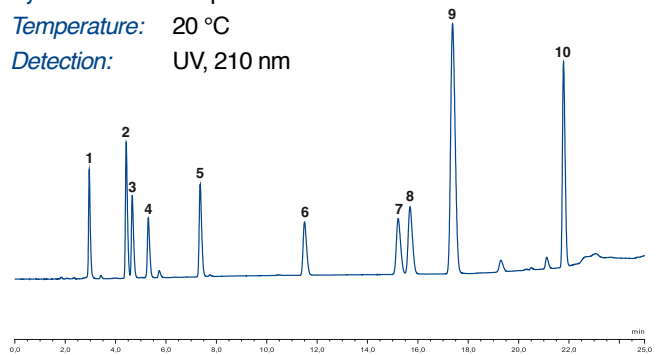
**Gradient:** 17.5 - 32.5 % B in 6 min, 32.5 - 45.0 % B  
in 10 min, 45.0 - 90.0 % B in 9 min

**Flow rate:** 0.8 mL/min

**Inj. volume:** 2.5 µL

**Temperature:** 20 °C

**Detection:** UV, 210 nm



**Peaks:**

1. aldicarb sulfoxide, 2. aldicarb sulfone, 3. oxamyl, 4. methomyl, 5. 3-hydroxycarbofuran, 6. aldicarb, 7. propoxur, 8. carbofuran, 9. carbaryl, 10. methiocarb

A separation with UV detection is shown by Appl. No.: 123050  
([www.mn-net.com/apps](http://www.mn-net.com/apps))



## Fast separation of 16 EPA PAHs on NUCLEODUR® C<sub>18</sub> PAH

MN Appl. No. 123820

**Column:** EC 100/4 NUCLEODUR® C<sub>18</sub> PAH, 3 µm  
REF 760783.40

**Conditions**

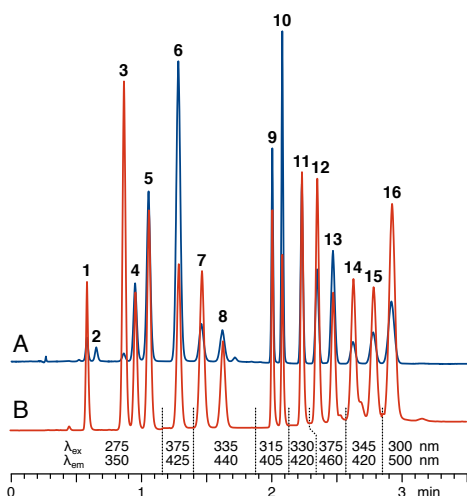
**Eluent:** A) MeOH/water (80:20; v/v)  
B) acetonitrile

**Gradient:** 2 - 20 % B in 1.2 min, 20 - 100 % B in 0.5 min, 100 % B for 2.5 min, 100 - 2 % B in 0.4 min, 2 % B for 5 min (equilibration)

**Flow rate:** 2.5 mL/min

**Temperature:** 35 °C

**Detection:** A: UV, 254 nm  
B: Fluorescence, see chromatogram



**Peaks:**  
see below

## Separation of 16 EPA PAHs with MeOH/water gradient on NUCLEODUR® C<sub>18</sub> PAH

MN Appl. No. 123830

**Column:** EC 125/4 NUCLEODUR® C<sub>18</sub> PAH, 3 µm  
REF 760784.40

**Conditions**

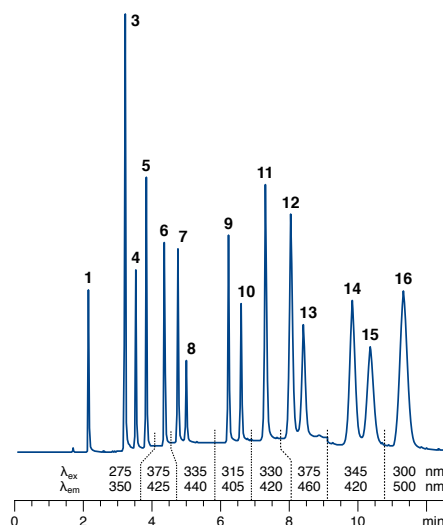
**Eluent:** A) water  
B) MeOH

**Gradient:** 65 - 97 % B in 6 min, 97 % B for 5 min, 97 - 65 % B in 0.5 min

**Flow rate:** 2.0 mL/min

**Temperature:** 35 °C

**Detection:** Fluorescence



**Peaks:**  
see below

## Separation of 18 PAHs on NUCLEODUR® C<sub>18</sub> PAH

MN Appl. No. 123840

**Column:** EC 125/4 NUCLEODUR® C<sub>18</sub> PAH, 3 µm  
REF 760784.40

**Conditions**

**Eluent:** A) MeOH/water (70:30; v/v)  
B) acetonitrile

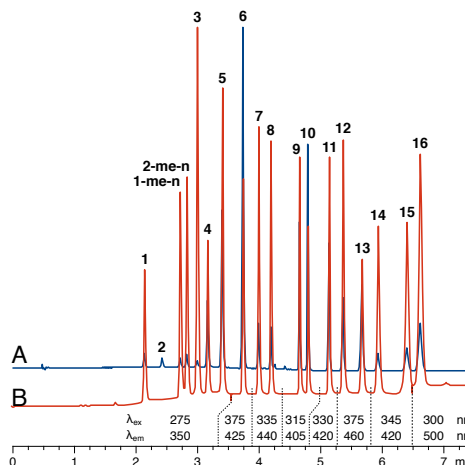
**Gradient:** 0 - 20 % B in 1.5 min, 20 - 50 % B in 1.5 min, 50 - 100 % B in 1.0 min, 100 % B for 3.0 min, 100 - 0 % B in 0.5 min

**Flow rate:** 1.5 mL/min

**Temperature:** 35 °C

**Inj. volumes:** UV, 1 µL (10 ng per compound)  
Fluorescence, 0.5 µL (5 ng per compound)

**Detection:** A: UV, 254 nm,  
B: Fluorescence, see chromatogram



**Peaks:**

1. naphthalene, 2. acenaphthylene (not detectable with fluorescence), 3. acenaphthene, 4. fluorene, 5. phenanthrene, 6. anthracene, 7. fluoranthene, 8. pyrene, 9. benz[a]anthracene, 10. chrysene, 11. benzo[b]fluoranthene, 12. benzo[k]fluoranthene, 13. benzo[a]pyrene, 14. dibenz[ah]anthracene, 15. benzo[ghi]perylene, 16. indeno[1,2,3-cd]pyrene, 1-me-n: 1-methylnaphthalene, 2-me-n: 2-methylnaphthalene





## Separation of nitroaromatics (EPA 8330 Mix A/B)

MN Appl. No. 124490 and 124500

**Column:** EC 150/3 NUCLEODUR® PolarTec, 5 µm  
REF 760488.30

**Conditions**

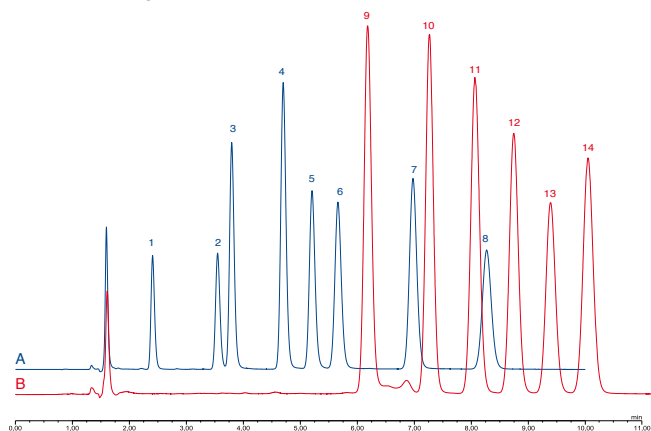
**Eluent:** Mix A: MeOH/water (50:50; v/v)  
Mix B: water + 0.1 % CHOOH/  
MeOH (55:45; v/v)

**Flow rate:** 0.46 mL/min

**Inj. temperature:** Mix A: 50 °C; Mix B: 60 °C

**Note:**

Please also view application number 124510 showing Mix A and Mix B in a single run ([www.mn-net.com/apps](http://www.mn-net.com/apps))



**Peaks:**

**Mix A:** 1. octogen (HMX), 2. hexogen (RDX), 3. 1,3,5-trinitrobenzene, 4. 1,3-dinitrobenzene, 5. nitrobenzene, 6. 2,4,6-trinitrotoluene, 7. 2-amino-4,6-dinitrotoluene, 8. 2,4-dinitrotoluene,

**Mix B:** 9. n-methyl-n-2,4,6,-tetranitroaniline (tetryl), 10. 4-amino-2,6-dinitrotoluene, 11. 2,6-dinitrotoluene 12. 2-nitrotoluene, 13. 4-nitrotoluene, 14. 3-nitrotoluene

## Separation of carbonyl DNPH compounds (EPA TO-11 A, EPA 8315)

MN Appl. No. 123620

**Column:** EC 250/4 NUCLEODUR® C<sub>18</sub> HTec, 5 µm  
REF 760316.40

**Sample:**

commercial standard test mixture

**Conditions**

**Eluent:** A) acetonitrile  
B) water

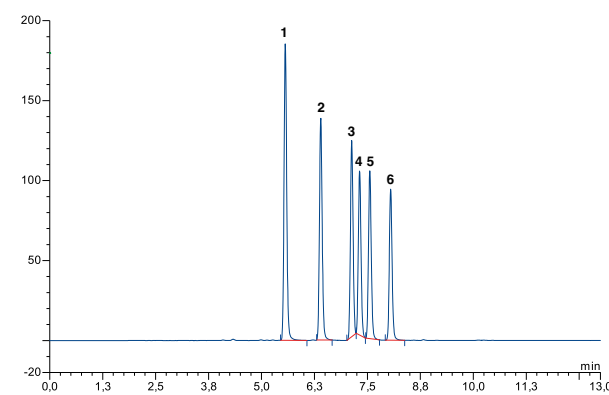
**Gradient:** 44 - 12 % B in 7.5 min,  
12 % B for 7.5 min,  
12 - 44 % B in 5 min

**Flow rate:** 0.65 mL/min

**Temperature:** 60 °C

**Inj. volume:** 5 µL

**Detection:** UV, 360 nm



**Peaks:**

1. formaldehyde-2,4-DNPH, 2. acetaldehyde-2,4-DNPH, 3. acrolein-2,4-DNPH, 4. acetone-2,4-DNPH, 5. propionaldehyde-2,4-DNPH, 6. crotonaldehyde-2,4-DNPH



## Determination of perfluorinated surfactants by LC-MS

MN Appl. No. 123340

**Matrix:** water

**Column:** EC 125/2 NUCLEODUR® Sphinx RP, 3 µm  
REF 760807.20

**Conditions**

**Eluent:** A) 10 mM ammonium acetate in water/MeOH (75:25; v:v)  
B) 10 mM ammonium acetate in acetonitrile/MeOH (75:25; v:v)

**Gradient:** 10 % - 30 % B in 3 min, 30 % - 55 % B in 8 min, 55 % - 10 % B in 4 min

**Flow rate:** 0.30 mL/min

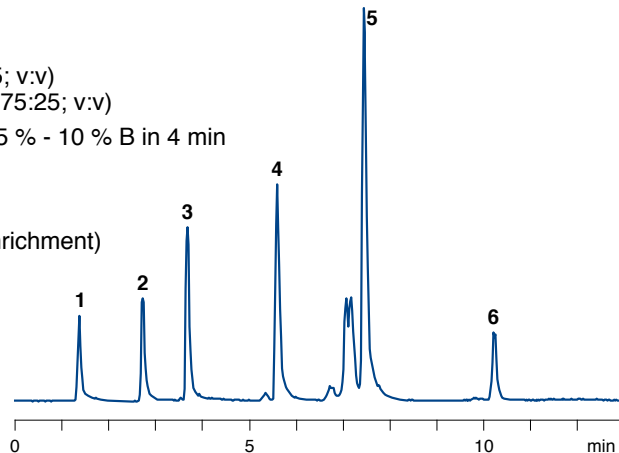
**Temperature:** 50 °C

**Inj. volume:** 2.5 µL (each concentration = 5 mg/L, after SPE enrichment)

**Detection:** MS, ESI negative

**Needle voltage:** -3.00 kV

**Sample temperature:** 450 °C



**Peaks:**

1. perfluoropropionic acid (PFPrA, pentafluoropropionic acid), 2. perfluoropentanoic acid (PFPeA), 3. perfluorohexanoic acid (PFHxA, undecafluorohexanoic acid), 4. perfluorooctanoic acid (PFOA), 5. perfluorooctane sulfonate potassium salts (PFOS, heptafluorooctanesulfonic acid), 6. perfluorododecanoic acid (PFDoDA)

## Separation of Phthalates on NUCLEODUR® C<sub>18</sub> HTec, 3 µm

MN Appl. No. 124300

**Column:** EC 125/2 NUCLEODUR® C<sub>18</sub> HTec, 3 µm  
REF 760324.20

**Sample concentration:** 0.5 mg/mL of each compound

**Conditions**

**Eluent:** A) acetonitrile  
B) water

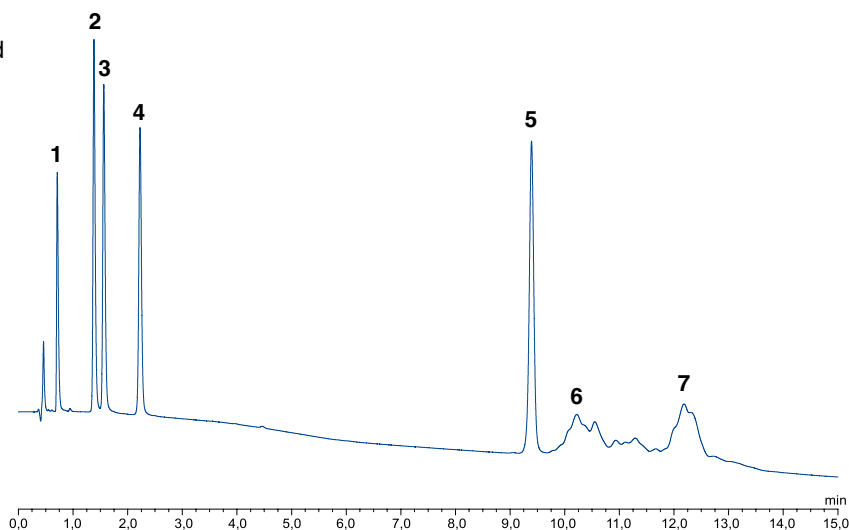
**Gradient:** 27.5 - 22.5 % B in 2.5 min,  
22.5 - 15 % B in 2.0 min,  
15 - 0 % B in 10.5 min

**Flow rate:** 0.5 mL/min

**Inj. volume:** 1 µL

**Temperature:** 45 °C

**Detection:** UV, 225 nm



**Peaks:**

1. dimethyl phthalate, 2. dipropyl phthalate,  
3. diphenyl phthalate, 4. dibutyl phthalate,  
5. diisooctyl phthalate, 6. diisononyl phthalate,  
7. diisodecyl phthalate



## CHROMABOND® HR-Xpert...

the innovative concept of five polymer-based RP- and mixed-mode ion exchange phases for SPE:

- ⇒ **CHROMABOND® HR-X** hydrophobic PS/DVB copolymer
- ⇒ **CHROMABOND® HR-XC** strong mixed-mode cation exchanger
- ⇒ **CHROMABOND® HR-XA** strong mixed-mode anion exchanger
- ⇒ **CHROMABOND® HR-XCW** weak mixed-mode cation exchanger
- ⇒ **CHROMABOND® HR-XAW** weak mixed-mode anion exchanger

All **CHROMABOND® HR-Xpert** phases are based on pure and spherical polymeric resin and provide:

- ⇒ Excellent recovery rates and highest reproducibility
- ⇒ Reliable and cost-efficient analyses
- ⇒ Robust retention mechanism even for aggressive washing procedures
- ⇒ Low limits of detection also for critical matrices

## HR-X spherical, hydrophobic polystyrene-divinylbenzene adsorbent resin

- ⇒ hydrophobic polystyrene-divinylbenzene copolymer  
pH stability 1 – 14  
high-purity material with highest reproducibility and lowest blank values due to a novel manufacturing process  
spherical particles, size 85 µm; pore size 55 – 60 Å  
very high surface 1000 m<sup>2</sup>/g  
capacity 390 mg/g (caffeine in water)  
excellent recovery rates especially for the enrichment of pharmaceuticals/active ingredients due to the spherical structure of the particles, very homogeneous surface, and optimized pore structure

- ⇒ recommended application:  
pharmaceuticals/active ingredients from tablets, creams and water/waste water  
drugs and pharmaceuticals from urine, blood, serum and plasma  
trace analysis of pesticides, herbicides, phenols, PAHs and PCBs from water

Volume	Adsorbent weight						Pack of
<b>CHROMABOND® HR-X polypropylene columns</b>							
	30 mg	60 mg	100 mg	200 mg	500 mg	1 g	
1 mL	<b>730934</b>		<b>730935</b>				30
3 mL		<b>730936</b>		<b>730931</b>	<b>730937</b>		30
6 mL				<b>730938</b>	<b>730939</b>		30
15 mL					<b>730940</b>	<b>730941</b>	20

## HR-XC

- ⇒ strong acidic benzenesulphonic acid cation exchanger  
exchange capacity 1.0 meq/g, pKa < 1  
base material PS/DVB copolymer (HR-X)  
pH stability 1 – 14  
high purity material, highest reproducibility and lowest blank values due to an optimized production process  
spherical particles, size 85 µm; pore size 65 – 75 Å  
very large specific surface 800 m<sup>2</sup>/g; pore volume 1.4 cm<sup>3</sup>/g  
RP capacity 300 mg/g (caffeine in water)  
outstanding recovery rates especially for the enrichment of basic analytes

## strong cation exchanger

- ⇒ recommended application:  
basic active ingredients from heavily matrix-contaminated samples like e.g. urine, plasma, serum  
fungicides from food, melamine from milk  
basic analytes like e.g. amines  
bases with **pKa 2 – 10**

Volume	Adsorbent weight						Pack of
<b>CHROMABOND® HR-XC polypropylene columns</b>							
	30 mg	60 mg	100 mg	150 mg	200 mg	500 mg	
1 mL	<b>730969</b>		<b>730049</b>				30
3 mL		<b>730956</b>			<b>730952</b>	<b>730953</b>	30
6 mL				<b>730957</b>		<b>730955</b>	30

Glass columns, LV columns, CHROMAFIX® cartridges, MULTI 96 and adsorbent on request!




## HR-XA

- ◆ strong basic quaternary ammonium anion exchanger  
 exchange capacity 0.25 meq/g, pKa ~ 18  
 base material PS/DVB copolymer (HR-X)  
 pH stability 1 – 14  
 high purity material with highest reproducibility and  
 lowest blank values due to an optimized production process  
 spherical particles, size 85 µm; pore size 55 – 65 Å  
 very large specific surface 850 m<sup>2</sup>/g; pore volume 1.4 cm<sup>3</sup>/g  
 RP capacity 350 mg/g (caffeine in water)  
 outstanding recovery rates especially for the enrichment of acidic analytes

## strong anion exchanger

- ◆ recommended application:  
 acidic active ingredients from  
 heavily matrix-contaminated  
 samples like e. g. urine,  
 plasma, serum  
 phenolic acids  
 acidic herbicides  
 weak/medium-strength acids  
 with **pKa 2 – 8**


	Volume	Adsorbent weight					Pack of	
	<b>CHROMABOND® HR-XA polypropylene columns</b>							
		30 mg	60 mg	100 mg	150 mg	200 mg	500 mg	
	1 mL	<b>730968</b>		<b>730727</b>				30
	3 mL		<b>730950</b>			<b>730951</b>	<b>730954</b>	30
	6 mL				<b>730958</b>	<b>730966</b>		30

## HR-XCW

- ◆ weak acidic carboxylic acid cation exchanger  
 exchange capacity > 0.7 meq/g, pKa ~ 5  
 base material spherical PS/DVB copolymer (HR-X)  
 pH stability 1 – 14  
 high purity material, highest reproducibility and  
 lowest blank values due to an optimized production process  
 spherical particles, size 85 µm; pore size 50 – 60 Å  
 very large specific surface 850 m<sup>2</sup>/g; pore volume 1.2 – 1.4 cm<sup>3</sup>/g  
 RP capacity 350 mg/g (caffeine in water)  
 outstanding recovery rates especially for enrichment of strongly basic analytes

## weak cation exchanger

- ◆ recommended application:  
 basic compounds like quater-  
 nary amines  
 active ingredients from heavily  
 matrix-contaminated samples  
 like e.g. urine, plasma, serum  
 strong bases with **pKa > 10**

	Volume	Adsorbent weight					Pack of	
	<b>CHROMABOND® HR-XCW polypropylene columns</b>							
		30 mg	60 mg	100 mg	150 mg	200 mg	500 mg	
	1 mL	<b>730731</b>		<b>730733</b>				30
	3 mL		<b>730735</b>			<b>730739</b>	<b>730741</b>	30
	6 mL				<b>730737</b>	<b>730743</b>		30

Glass columns, LV columns, CHROMAFIX® cartridges, MULTI 96 and adsorbent on request!

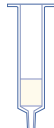


## HR-XAW

- weak basic secondary and tertiary ammonium anion exchanger, exchange capacity > 0.5 meq/g, pKa ~ 6
- base material spherical PS/DVB copolymer (HR-X)
- pH stability 1 – 14
- high purity material with highest reproducibility and lowest blank values due to an optimized production process
- spherical particles, size 85 µm; pore size 55 – 65 Å
- very large specific surface 850 m<sup>2</sup>/g; pore volume 1.2 – 1.4 cm<sup>3</sup>/g
- RP capacity 350 mg/g (caffeine in water)
- outstanding recovery rates especially for enrichment of acidic analytes

## weak anion exchanger

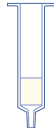
- recommended application:
  - perfluorinated surfactants
  - acidic compounds like sulfonates
  - active ingredients from heavily matrix-contaminated samples like e.g. urine, plasma, serum
  - strong acids with pKa < 1

	Volume	Adsorbent weight					Pack of	
	<b>CHROMABOND® HR-XAW polypropylene columns</b>							
		30 mg	60 mg	100 mg	150 mg	200 mg	500 mg	
	1 mL	<b>730728</b>		<b>730729</b>				30
	3 mL		<b>730747</b>			<b>730748</b>	<b>730744</b>	30
	6 mL			<b>730749</b>		<b>730745</b>		30

## Easy polar, bifunctionally modified polystyrene-divinylbenzene copolymer

- polar modified polystyrene-divinylbenzene copolymer with a weak anion exchanger
- pH stability 1 – 14, particle size 80 µm, pore size 50 Å, specific surface 650 – 700 m<sup>2</sup>/g,
- The Easy effect:**
  - without preconditioning
  - due to bifunctional modification much more hydrophilic than conventional polystyrene-divinylbenzene polymers
  - easily wettable with water

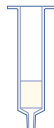
- recommended application:
  - polar herbicides / pesticides from water (acidic, neutral, basic)
  - polar phenols from water
  - polyaromatic compounds
  - polychlorinated biphenyls
  - drug analysis from urine, blood, serum, plasma, pharmaceuticals / active ingredients from tablets, creams

	Volume	Adsorbent weight					Pack of	
	<b>CHROMABOND® Easy polypropylene columns</b>							
		30 mg	60 mg	100 mg	200 mg	500 mg	1 g	
	1 mL	<b>730751</b>		<b>730752</b>				30
	3 mL		<b>730753</b>		<b>730754</b>	<b>730759</b>		30
	6 mL			<b>730755</b>	<b>730756</b>			30

## HR-P polystyrene-divinylbenzene adsorbent resin

- highly porous polystyrene-divinylbenzene copolymer
- specific surface 1200 m<sup>2</sup>/g
- particle size 50 – 100 µm
- very high binding capacity, up to 30% of adsorbent weight (for comparison: silica adsorbents about 3%)

- recommended application:
  - aromatic compounds
  - phenols from water
  - nitroaromatics from water
  - pesticides from water
  - PAHs from oil

	Volume	Adsorbent weight				Pack of	
	<b>CHROMABOND® HR-P polypropylene columns</b>						
			100 mg	200 mg	500 mg	1 g	
	1 mL		<b>730280</b>				30
	3 mL			<b>730108</b>	<b>730117</b>		30
	6 mL			<b>730119</b>	<b>730111</b>	<b>730118</b>	30

Glass columns, LV columns, CHROMAFIX® cartridges, MULTI 96 and adsorbent on request!



## HR-P-AOX

### AOX from waters with high salt loads (DIN 38409 – H22)

special PS/DVB phase

recommended application:  
extraction of AOX (adsorbable organically bonded halogens) from waters containing high salt loads / organic pollutants in accordance with DIN 38409 – H22

	Volume	Adsorbent weight			Pack of
	<b>CHROMABOND® HR-P-AOX polypropylene columns</b>				
	6 mL	200 mg 730119.AOX	500 mg 730111.AOX		30

## C<sub>18</sub> ec

### octadecyl silica, endcapped

base material silica, pore size 60 Å, particle size 45 µm, specific surface 500 m<sup>2</sup>/g, pH stability 2 – 8  
octadecyl phases, endcapped, carbon content 14 %  
very nonpolar, hydrophobic interactions with a wide variety of organic compounds  
advantageous for clean-up of samples with large structural variations (polarity differences)

recommended application:  
nonpolar compounds  
aflatoxins, amphetamines, antibiotics, antiepileptics, barbiturates, caffeine, drugs, preservatives, fatty acids, nicotine, PAHs, pesticides, PCBs, heavy metals, vitamins  
very well suited for desalting of samples

	Volume	Adsorbent weight					Pack of
	<b>CHROMABOND® C<sub>18</sub> ec polypropylene columns</b>						
		100 mg	200 mg	500 mg	1 g	2 g	
	1 mL	730011					100
	3 mL		730012	730013			50
	6 mL			730014	730015	730141	30
	<b>CHROMABOND® C<sub>18</sub> ec glass column</b>						
	3 mL			730013G			50

## C<sub>18</sub> Hydra

### octadecyl silica for polar analytes

base material silica, pore size 60 Å, particle size 45 µm, specific surface 500 m<sup>2</sup>/g, pH stability 2 – 8  
special octadecyl phase for polar analytes, not endcapped, carbon content 15 %

recommended application:  
more polar compounds like pesticides and their polar degradation products, phenols, phenoxy-carboxylic acids, nitroaromatics, pharmaceuticals

	Volume	Adsorbent weight					Pack of
	<b>CHROMABOND® C<sub>18</sub> Hydra polypropylene columns</b>						
		50 mg	100 mg	200 mg	500 mg	1 g	2 g
	1 mL	730294	730295				
	3 mL			730296	730297	730298	
	6 mL			730299	730300	730301	
							100
							50
							30

Glass columns, LV columns, CHROMAFIX® cartridges, MULTI 96 and adsorbent on request!

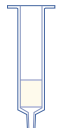


## C<sub>18</sub> PAH

- base material silica, pore size 60 Å, particle size 45 µm, specific surface 500 m<sup>2</sup>/g, pH stability 2 – 8
- special octadecyl modification for enrichment of PAH, not endcapped, carbon content 14 %

## octadecyl silica for PAH analysis

- recommended application: PAHs from water

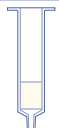
Volume	Adsorbent weight			Pack of
	<b>CHROMABOND® C<sub>18</sub> PAH polypropylene columns</b>			
	6 mL	2 g 730166		30

## NH<sub>2</sub>/C<sub>18</sub>

- special combination phase: aminopropyl phase for removal of interfering humic acids
- octadecyl phase for enrichment of PAH

## combination phase for PAH analysis

- recommended application: PAHs from water containing humic acids

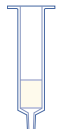
Volume	Adsorbent weight			Pack of
	<b>CHROMABOND® NH<sub>2</sub>/C<sub>18</sub> polypropylene columns · glass column</b>			
		500/500 mg	500 mg/1 g	
	6 mL PP	730618	730620	30
6 mL glass		730620G	30	

## NH<sub>2</sub>

- base material silica, pore size 60 Å, particle size 45 µm, specific surface 500 m<sup>2</sup>/g, pH stability 2 – 8
- aminopropyl phase, carbon content 3.5 %
- polar, weak anion exchanger

## aminopropyl silica

- recommended application: trace elements, lipids

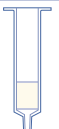
Volume	Adsorbent weight				Pack of
	<b>CHROMABOND® NH<sub>2</sub> polypropylene columns</b>				
		100 mg	200 mg	500 mg	1 g
	1 mL	730031			
	3 mL		730413	730033	
	6 mL			730180	730626
					100
					50
					30

## SiOH

- unmodified, weakly acidic silica, pore size 60 Å, particle size 45 µm, specific surface 500 m<sup>2</sup>/g, pH stability 2 – 8
- very polar
- adsorbs humidity from air, for this reason it should be kept well closed and if necessary dried before use
- due to its high affinity for polar compounds it should not be conditioned with polar (e. g. MeOH) or water-containing solvents

## unmodified silica

- recommended application: aflatoxins, chloramphenicol, pesticides, steroids, vitamins

Volume	Adsorbent weight					Pack of
	<b>CHROMABOND® SiOH polypropylene columns</b>					
		100 mg	200 mg	500 mg	1 g	2 g
	1 mL	730071				
	3 mL		730214	730073		
	6 mL			730070	730075	730107
	15 mL					730217
						100
						50
						30
						20

Glass columns, LV columns, CHROMAFIX® cartridges, MULTI 96 and adsorbent on request!




# SPE · Solid Phase Extraction

## Florisol®

matrix magnesium silicate (MgO - SiOH 15:85), high purity, particle size 150 – 250 µm

## magnesium silicate

recommended application:  
organic tin compounds,  
aliphatic carboxylic acids,  
PCBs, PAHs


Volume	Adsorbent weight				Pack of	
	<b>CHROMABOND® Florisol® polypropylene columns</b>					
		200 mg	500 mg	1 g	2 g	
	3 mL	730457	730081			50
	6 mL		730238	730082	730239	30
<b>CHROMABOND® Florisol® glass columns</b>						
6 mL			1 g	2 g		
			730082G	730239G	30	

## Na<sub>2</sub>SO<sub>4</sub>/Florisol®

## hydrocarbons from water acc. to DIN H-53/ISO DIS 9377-4

special combination phase of sodium sulphate and Florisol®

recommended application:  
hydrocarbons from drinking, surface  
and waste waters


Volume	Adsorbent weight		Pack of
	<b>CHROMABOND® Na<sub>2</sub>SO<sub>4</sub> / Florisol® glass columns</b>		
	6 mL	2 g/2 g	730249G

## CN/SiOH

## combination phase for PAH analysis

special combination phase  
cyanopropyl phase for selective adsorption of polycyclic aromatics  
via π-π interactions  
unmodified silica phase for removal of polar compounds

recommended application:  
extraction of the 16 PAHs according  
to EPA from soil samples

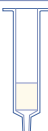
Volume	Adsorbent weight		Pack of	
	<b>CHROMABOND® CN/SiOH polypropylene columns</b>			
		500 mg/1 g		
	3 mL	730112		50
6 mL	730135		30	

## NAN

## special phase for PCB analysis

special combination phase:  
N: sodium sulphate for removal of trace water;  
A: SiOH/AgNO<sub>3</sub> phase for removal of sulphur, sulphur-containing and  
polar compounds

recommended application  
extraction of PCB from sludge

Volume	Adsorbent weight		Pack of	
	<b>CHROMABOND® NAN polypropylene columns</b>			
		400/1400/400 mg	700/2000/700 mg	
	3 mL	730109		50
6 mL		730149	30	

Glass columns, LV columns, CHROMAFIX® cartridges, MULTI 96 and adsorbent on request!



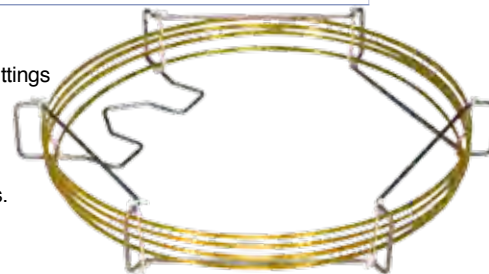


Phase	Composition	max. Temp. <sup>1</sup>	USP	Similar phases <sup>2</sup>
<b>OPTIMA® 1</b>	100 % dimethylpolysiloxane	340/360 °C	G1 G2 G38	PERMABOND® SE-30, OV-1, DB-1, SE-30, HP-1, SPB™-1, CP-Sil 5 CB, Rtx®-1, 007-1, BP1, MDN-1, AT™-1, ZB-1, OV-101
<b>OPTIMA® 1 MS</b> <b>OPTIMA® 1 MS Accent</b>	100 % dimethylpolysiloxane	340/360 °C	G1 G2 G38	Ultra-1, DB-1MS, HP-1MS, Rxi®-1MS, Rtx®-1MS, Equity™-1, AT™-1MS, VF-1MS, CP-Sil 5 CB MS
<b>OPTIMA® 5</b>	5 % phenyl – 95 % methylpolysiloxane	340/360 °C	G27 G36	PERMABOND® SE-52, SE-54, SE-52, HP-5, SPB™-5, CP-Sil 8, Rtx®-5, 007-5, BP5, MDN-5, AT™-5, ZB-5
<b>OPTIMA® 5 MS</b>	5 % diphenyl – 95 % dimethylpolysiloxane	340/360 °C	G27 G36	DB-5, DB-5MS, HP-5MS, Ultra-2, Equity™-5, CP-Sil 8CB low bleed/MS, Rxi®-5MS, Rtx®-5SIL-MS, Rtx®-5MS, 007-5MS, BPX™5, MDN-5S, AT™-5MS, VF-5MS
<b>OPTIMA® 5 MS Accent</b>	silarylene phase with selectivity similar to 5 % diphenyl – 95 % dimethylpolysiloxane	340/360 °C	G27 G36	
<b>OPTIMA® XLB</b>	silarylene phase as above, optimized silarylene content	340/360 °C	–	DB-XLB, Rxi®-XLB, Rtx®-XLB, MDN-12, VF-XMS
<b>OPTIMA® δ-3</b>	phase with autoselectivity	340/360 °C	G49	no similar phases
<b>OPTIMA® δ-6</b>	phase with autoselectivity	340/360 °C	–	no similar phases
<b>OPTIMA® 1301</b>	6 % cyanopropylphenyl – 94 % dimethylpolysiloxane	300/320 °C	G43	HP-1301, DB-1301, SPB™-1301, Rtx®-1301, CP-1301, 007-1301
<b>OPTIMA® 624</b>	6 % cyanopropylphenyl – 94 % dimethylpolysiloxane	280/300 °C	G43	HP-624, HP-VOC, DB-624, DB-VRX, SPB™-624, CP-624, Rtx®-624, Rtx®-Volatiles, 007-624, BP624, VOCOL
<b>OPTIMA® 624 LB</b>	as above, low bleed phase	280/300 °C	G43	
<b>OPTIMA® 1701</b>	14 % cyanopropylphenyl – 86 % dimethylpolysiloxane	300/320 °C	G46	OV-1701, DB-1701, CP-Sil 19 CB, HP-1701, Rtx®-1701, SPB™-1701, 007-1701, BP10, ZB-1701
<b>OPTIMA® 35 MS</b>	silarylene phase with selectivity similar to a 35 % diphenyl – 65 % dimethylpolysiloxane phase	360/370 °C	G42	DB-35 MS, HP-35, SPB™-35, Rxi®-35SIL MS, Rtx-35, 007-35, BPX™-35, MDN-35, AT™-35 MS, ZB-35, OV-11, VF-35 MS
<b>OPTIMA® 17</b>	phenylmethylpolysiloxane, 50 % phenyl	320/340 °C	G3	OV-17, DB-17, HP-50+, HP-17, SPB™-50, SP-2250, Rxi®-17, Rtx®-50, CP-Sil 24 CB, 007-17, ZB-50
<b>OPTIMA® 17 MS</b>	silarylene phase with selectivity similar to 50 % phenyl, 50 % methylpolysiloxane	340/360 °C	G3	OV-17, AT™-50, BPX™-50, DB-17, DB-18ms, HP-50+, HP-17, SPB™-50, SPB™-17, SP-2250, Rtx®-50, CP-Sil 24 CB, 007-17, VF-17ms, ZB-50
<b>OPTIMA® 210</b>	trifluoropropylmethylpolysiloxane (50 % trifluoropropyl)	260/280 °C	G6	OV-210, DB-210, Rtx®-200, 007-210
<b>OPTIMA® 225</b>	50 % cyanopropylmethyl – 50 % phenylmethylpolysiloxane	260/280 °C	G7 G19	DB-225, HP-225, OV-225, Rtx®-225, CP-Sil 43, 007-225, BP225
<b>OPTIMA® 240</b>	33 % cyanopropylmethyl – 67 % dimethylpolysiloxane	260/280 °C	–	no similar phases
<b>OPTIMA® WAX</b>	polyethylene glycol 20000 daltons	250/260 °C	G16	PERMABOND® CW 20 M, DB-Wax, Supelcowax™, HP-Wax, HP-INNOWax, Rtx®-Wax, CP-Wax 52 CB, Stabilwax, 007-CW, BP20, AT™-Wax, ZB-Wax
<b>OPTIMA® FFAP</b>	polyethylene glycol 2-nitro-terephthalate	250/260 °C	G25 G35	PERMABOND® FFAP, DB-FFAP, HP-FFAP, CP-SIL 58 CB, 007-FFAP, CP-FFAP CB, Nukol

<sup>1</sup> **first temperature for isothermal operation, second value for short isotherms in a temperature program**  
Please note, that for columns with 0.53 mm ID and for columns with thicker films temperature limits are generally lower. For details refer to the description of individual phases.

<sup>2</sup> **phases which provide a similar selectivity based on chemical and physical properties**

Each Column is individually tested and supplied with test certificate and test chromatogram, but without fittings or ferrules. Column ends are melted or closed with septa, and thus protected from atmospheric oxygen. Additionally, we supply the corresponding test mixture with each column. On request, all columns can be supplied on a **5 inch (13 mm) cage** for the Agilent GC 6850: For ordering, please add an E at the end of the REF number (e.g. 726470.30E). For a considerably longer lifetime, even for contaminated or matrix-containing samples; MN offers the option of **integrated precolumns**. Please contact us for details.



**MACHEREY-NAGEL** Trademarks: CHROMABOND, CHROMAFIL, CHROMAFIX, ChromCart, NUCLEODUR, OPTIMA, PERMABOND  
Trademarks of the other companies: DB / SE, Agilent Technologies Inc. / J&W Scientific Inc. (USA); CP / VF, Agilent / Varian Inc. (USA); HP, Agilent Technologies Inc. (USA); AT, Alltech Associates Inc. (USA); OV, Ohio Valley Specialty Company (USA); ZB, Phenomenex Inc. (USA); 007, Quadrex Corp. (USA); Rtx / Rxi, Restek Corp. (USA); BPX, SGE Analytical Science Pty Ltd. (Australia); Equity / SPB / Nukol / MDN, Sigma-Aldrich Biotechnology LP / Sigma-Aldrich Co. / Supelco (USA); Florisil / U.S. Silica Co. (USA).



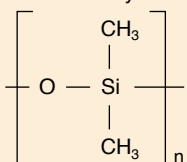
The following selection of GC column dimensions is only an excerpt of our product range. For the full range of GC columns ask for our chromatography catalog or visit:

[www.mn-net.com](http://www.mn-net.com)

Custom made dimensions can be obtained on request.

## OPTIMA® 1 MS Accent

selectivity identical to OPTIMA® 1



increased sensitivity due to an unmatched low background level

100 % dimethylpolysiloxane



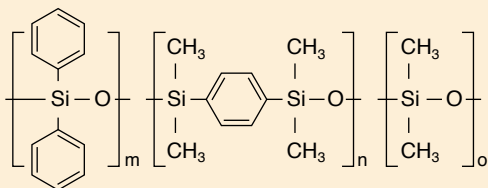
max. temperature for isothermal operation 340 °C,  
max. temperature for short isotherms in a temperature program 360 °C

- lowest column bleed, nonpolar phase, ideal for ion trap and quadrupol MS detectors
- perfect inertness for basic compounds
- solvent rinsing for removal of impurities applicable
- application areas: all-round phase for environmental analyses, trace analyses, EPA methods, pesticides, PCB, food and drug analyses
- similar phases: Ultra-1, DB-1 MS, HP-1 MS, Rxi-1 MS, Rtx-1 MS, Equity-1, AT-1 MS, VF-1 MS, CP-Sil 5 CB MS
- USP G1 / G2 / G38

Length →	15 m	25 m	30 m	50 m	60 m
<b>0.2 mm ID (0.4 mm OD)</b>					
0.20 µm film		725801.25		725801.50	
<b>0.25 mm ID (0.4 mm OD)</b>					
0.25 µm film	725805.15		725805.30		725805.60
0.50 µm film			725806.30		725806.60
<b>0.32 mm ID (0.5 mm OD)</b>					
0.25 µm film			725802.30		725802.60

## OPTIMA® 5 MS Accent

chemically bonded, cross-linked silarylene phase with polarity similar to a 5% diphenyl - 95% dimethylpolysiloxane phase



increased sensitivity due to an unmatched low background level

silarylene phase



max. temperature for isothermal operation 340 °C, max. temperature for short isotherms in a temperature program 360 °C, for columns with films > 0.5 µm max. temperatures are 320 and 340 °C, respectively

- lowest column bleed, nonpolar phase, ideal for ion trap and quadrupol MS detectors

solvent rinsing for removal of impurities applicable

application areas: all-round phase for environmental analyses, trace analyses, EPA methods, pesticides, PCB, food and drug analyses

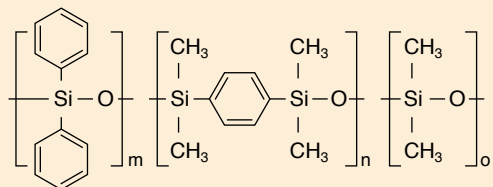
similar phases: DB-5 MS, HP-5 MS, Ultra-2, Equity-5, CP-Sil 8 CB low bleed/MS, Rxi-5 MS, Rtx-5SIL-MS, Rtx-5 MS, 007-5 MS, BPX5, MDN-5S, AT-5 MS, VF-5 MS

- USP G27 / G36

Length →	12 m	15 m	25 m	30 m	50 m	60 m
<b>0.2 mm ID (0.4 mm OD)</b>						
0.20 µm film			725810.25		725810.50	
0.35 µm film	725815.12				725815.50	
<b>0.25 mm ID (0.4 mm OD)</b>						
0.25 µm film		725820.15		725820.30		725820.60
0.50 µm film				725825.30		725825.60
1.00 µm film				725826.30		725826.60
<b>0.32 mm ID (0.5 mm OD)</b>						
0.25 µm film				725811.30		725811.60
0.50 µm film				725813.30		
1.00 µm film			725812.25			725812.60

**OPTIMA® XLB**

chemically bonded, cross-linked silarylene phase,  
optimized silarylene content for lowest column bleed



similar phases: DB-XLB, Rxi-XLB, Rtx-XLB, MDN-12,  
VF-XMS

**silarylene phase**

max. temperature for isothermal operation 340 °C,  
max. temperature for short isotherms in a temperature program 360 °C,



lowest column bleed, nonpolar phase, ideal for ion trap and quadrupole MS detectors  
perfect inertness for basic compounds  
solvent rinsing for removal of impurities applicable  
application areas: ultra low bleed phase, highly selective for environmental and trace analyses, pesticides  
recommended phase for PCB separations

Length →	30 m	60 m
<b>0.25 mm ID (0.4 mm OD)</b>		
0.25 µm film	<b>725850.30</b>	<b>725850.60</b>

**OPTIMA® δ-3**

- medium polar without CN groups  
analyses determine the polarity of the phase  
unique from MN, no similar phase  
ideal for MSD and PND detectors
- USP G49



max. temperature for isothermal operation 340 °C,  
max. temperature for short isotherms in a temperature program 360 °C  
for 0.53 mm ID columns the max. temperatures are 320 and 340 °C, resp.



autoselectivity resulting in a wide range of polarities from approximately the non-polar OPTIMA® 5 to the midpolar OPTIMA® 1701

**polysiloxane phase with autoselectivity**

Length →	25 m	30 m	50 m	60 m
<b>0.2 mm ID (0.4 mm OD)</b>				
0.20 µm film	<b>726400.25</b>		<b>726400.50</b>	
<b>0.25 mm ID (0.4 mm OD)</b>				
0.25 µm film		<b>726420.30</b>		<b>726420.60</b>
0.50 µm film		<b>726421.30</b>		
<b>0.32 mm ID (0.5 mm OD)</b>				
0.25 µm film		<b>726440.30</b>		<b>726440.60</b>
0.35 µm film		<b>726441.30</b>		<b>726441.60</b>
1.00 µm film		<b>726442.30</b>		<b>726442.60</b>

**OPTIMA® δ-6**

- medium polar without CN groups  
analyses determine the polarity of the phase  
unique from MN, no similar phase  
ideal for MSD and PND detectors



max. temperature for isothermal operation 340 °C,  
max. temperature for short isotherms in a temperature program 360 °C  
for 0.53 mm ID columns the max. temperatures are 320 and 340 °C, resp.



autoselectivity resulting in a wide range of polarities from approximately the mid-polar OPTIMA® 17 to the polar OPTIMA® 210

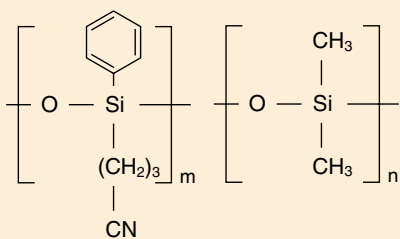
**polysiloxane phase with autoselectivity**

Length →	25 m	30 m	50 m	60 m
<b>0.2 mm ID (0.4 mm OD)</b>				
0.20 µm film	<b>726465.25</b>		<b>726465.50</b>	
<b>0.25 mm ID (0.4 mm OD)</b>				
0.25 µm film		<b>726470.30</b>		<b>726470.60</b>
<b>0.32 mm ID (0.5 mm OD)</b>				
0.25 µm film		<b>726480.30</b>		<b>726480.60</b>
0.35 µm film		<b>726481.30</b>		<b>726481.60</b>
1.00 µm film		<b>726482.30</b>		<b>726482.60</b>



## OPTIMA® 624

medium polar phase



similar phases: HP-624, HP-VOC, DB-624, DB-VRX, SPB-624, CP-624, Rtx-624, Rtx-Volatiles, 007-624, BP624, VOCOL

## 6 % cyanopropyl-phenyl – 94 % dimethylpolysiloxane

- max. temperature for isothermal operation 280 °C, max. temperature for short isotherms in a temperature program 300 °C
- recommended for environmental analyses for corresponding columns with lower film thickness see OPTIMA® 1301
- USP G43

## OPTIMA® 624 LB

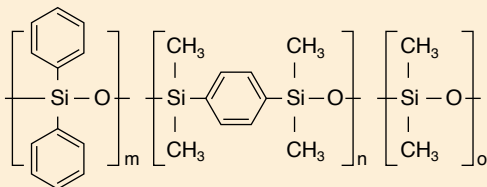
excellent Low Bleed columns for halogenated hydrocarbons, volatiles, aromatic compounds, solvents etc.

## 6 % cyanopropyl-phenyl – 94 % dimethylpolysiloxane

Length →		25 m	30 m	50 m	60 m
<b>OPTIMA® 624</b>	<b>0.2 mm ID (0.4 mm OD)</b>				
	1.10 µm film	<b>726784.25</b>			
	<b>0.25 mm ID (0.4 mm OD)</b>				
	1.40 µm film	<b>726785.25</b>	<b>726785.30</b>	<b>726785.50</b>	<b>726785.60</b>
	<b>0.32 mm ID (0.5 mm OD)</b>				
	1.80 µm film	<b>726787.25</b>	<b>726787.30</b>	<b>726787.50</b>	<b>726787.60</b>
	<b>0.53 mm ID (0.8 mm OD)</b>				
	3.00 µm film	<b>726789.25</b>	<b>726789.30</b>		
<b>OPTIMA® 624 LB</b>	<b>0.25 mm ID (0.5 mm OD)</b>				
	1.40 µm film		<b>726791.30</b>		<b>726791.60</b>
	<b>0.32 mm ID (0.5 mm OD)</b>				
	1.80 µm film		<b>726786.30</b>	<b>726786.50</b>	

## OPTIMA® 35 MS

chemically bonded cross-linked silarylene phase with selectivity similar to 35 % phenyl / 65 % methyl polysiloxane



similar phases: DB-35 MS, HP-35, SPB-35, Rxi-35SIL MS, Rtx-35, 007-35, BPX-35, MDN-35, AT-35 MS, ZB-35, OV-11, VF-35 MS

## silarylene phase

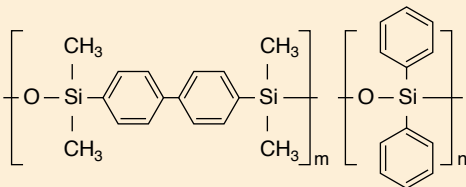
- max. temperature for isothermal operation 360 °C, max. temperature for short isotherms in a temperature program 370 °C
- very low column bleeding, midpolar phase, recommended for ion-trap detectors
- optimum column for confirmation of analytical results in combination with a 1 MS or 5 MS
- polymer without CN groups
- recommended application: allround phase for environmental analyses, ultra trace analyses, EPA methods, pesticides, PCBs, food and drug analyses
- USP G42

Length →		30 m	60 m
<b>0.25 mm ID (0.4 mm OD)</b>	0.25 µm film	<b>726154.30</b>	<b>726154.60</b>
	<b>0.32 mm ID (0.5 mm OD)</b>		
	0.25 µm film	<b>726157.30</b>	<b>726157.60</b>



## OPTIMA® 17 MS

- medium polar silarylene phase with selectivity analogue to 50% phenyl – 50% methylpolysiloxane



similar phases: OV-17, AT-50, BPX-50, DB-17, DB-17ms, HP-50+, HP-17, SPB-50, SPB-17, SP-2250, Rtx-50, CP-Sil 24 CB, 007-17, VF-17ms, ZB-50

## silarylene phase

- max. temperature for isothermal operation 340 °C, max. temperature for short isotherms in a temperature program 360 °C
- ideal for ion trap detectors
- optimum reference column in combination with a 1 MS or 5 MS
- no CN groups in the polymer
- recommended applications: all-round phase for environmental analyses, ultra-trace analyses, EPA methods, pesticides, PCBs, food and drug analyses
- USP G3

	Length →	30 m	60 m
<b>0.25 mm ID (0.4 mm OD)</b>			
0.25 µm film		<b>726162.30</b>	<b>726162.60</b>
<b>0.32 mm ID (0.5 mm OD)</b>			
0.25 µm film		<b>726165.30</b>	<b>726165.60</b>

## GC Application Guide

- explaining basics and principles of GC: phase selection by column properties, important GC parameters, helpful hints for troubleshooting
- 280 selected applications** from the fields
  - ✓ environmental pollutants
  - ✓ solvents · chemicals
  - ✓ fragrances · food and cosmetic components
  - ✓ drugs · pharmaceutical ingredients
  - ✓ petrochemical products
  - ✓ chiral separations
- latest and more applications at [www.mn-net.com/apps](http://www.mn-net.com/apps)





## NUCLEODUR® C<sub>18</sub> HTec

base-deactivated octadecyl phase






◆ **key features:**

- reliable and durable standard RP phase, suited for LC/MS
- outstanding base deactivation
- excellent stability and high loadability

◆ **technical characteristics:**

high density octadecyl modification (C<sub>18</sub>)  
 pore size 110 Å; particle sizes 1.8 µm, 3 µm and 5 µm for analytical separations  
 carbon content 18%, pH stability 1 – 11, USP L1

eluent in column acetonitrile / water

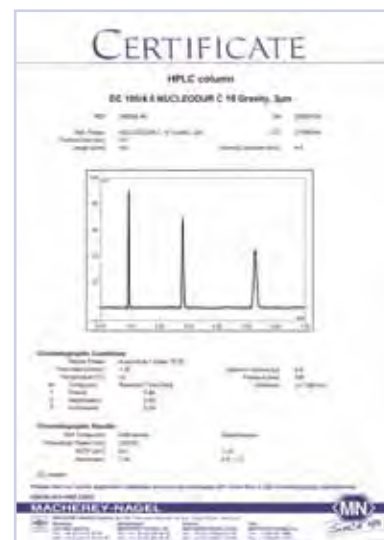
Length →	30 mm	50 mm	75 mm	100 mm	125 mm	150 mm	250 mm	Guard columns
<b>NUCLEODUR® C<sub>18</sub> HTec, 1.8 µm</b>								particle size 1.8 µm
<b>EC columns</b>								
	2 mm ID	760301.20	760305.20	760304.20	760306.20		760308.20	
	3 mm ID	760301.30	760305.30					
	4 mm ID	760301.40	760305.40					
	4.6 mm ID	760301.46	760305.46					
<b>NUCLEODUR® C<sub>18</sub> HTec, 3 µm</b>								particle size 3 µm
	2 mm ID		760321.20		760324.20	760325.20	760326.20	761120.30
	3 mm ID		760321.30		760324.30	760325.30	760326.30	761120.30
	4 mm ID		760321.40		760324.40	760325.40	760326.40	761120.40
	4.6 mm ID		760321.46	760322.46	760323.46	760324.46	760325.46	760326.46
								761120.40
<b>NUCLEODUR® C<sub>18</sub> HTec, 5 µm</b>								particle size 5 µm
	2 mm ID		760311.20		760314.20	760315.20	760316.20	761110.30
	3 mm ID		760311.30		760314.30	760315.30	760316.30	761110.30
	4 mm ID		760311.40		760314.40	760315.40	760316.40	761110.40
	4.6 mm ID		760311.46	760312.46	760313.46	760314.46	760315.46	760316.46
								761110.40

## Our HPLC QC policy

- ◆ **highest production standard**  
our facilities are EN ISO 9001:2008 certified
- ◆ **strict quality specifications** for outstanding reliability
- ◆ **perfect reproducibility** within each batch and from lot to lot
- ◆ Each column is individually tested and supplied with test chromatogram and test conditions

### Test mixture for reversed phase columns

Designation	Pack of	REF
Test mixture for reversed phase columns in acetonitrile	1 mL	722394





## NUCLEODUR® Sphinx RP

bifunctional RP phase






key features:

- distinct selectivity based on well-balanced bifunctional surface coverage
- widens the scope for method development based on additional  $\pi$ - $\pi$  interactions
- suitable for LC/MS due to low bleeding characteristics

technical characteristics:

octadecyl and propylphenyl modified silica; pore size 110 Å; particle sizes 1.8  $\mu$ m, 3  $\mu$ m and 5  $\mu$ m; carbon content 15 %; pH stability 1 – 10; high reproducibility and consistent quality, USP L1/L11



eluent in column acetonitrile / water

Length →	30 mm	50 mm	75 mm	100 mm	125 mm	150 mm	250 mm	Guard columns
<b>NUCLEODUR® Sphinx RP, 1.8 <math>\mu</math>m</b>								particle size 1.8 $\mu$ m
<b>EC columns</b>								
	2 mm ID	760821.20	760822.20	760825.20	760823.20		760824.20	
	3 mm ID	760821.30	760822.30					
	4 mm ID	760821.40	760822.40					
	4.6 mm ID	760821.46	760822.46					
<b>NUCLEODUR® Sphinx RP, 3 <math>\mu</math>m</b>								particle size 3 $\mu$ m
	2 mm ID		760806.20		760807.20	760805.20	760808.20	761557.30
	3 mm ID		760806.30		760807.30	760805.30	760808.30	761557.30
	4 mm ID		760806.40		760807.40	760805.40	760808.40	761557.40
	4.6 mm ID		760806.46	760813.46	760812.46	760807.46	760808.46	761557.40
<b>NUCLEODUR® Sphinx RP, 5 <math>\mu</math>m</b>								particle size 5 $\mu$ m
	2 mm ID		760800.20		760801.20	760802.20	760803.20	761550.30
	3 mm ID		760800.30		760801.30	760802.30	760803.30	761550.30
	4 mm ID		760800.40		760801.40	760802.40	760803.40	761550.40
	4.6 mm ID		760800.46	760815.46	760809.46	760801.46	760802.46	760803.46

## NUCLEODUR® C<sub>18</sub> PAH

special octadecyl phase for PAH analyses

- base material NUCLEODUR® silica, particle size 3  $\mu$ m, pore size 110 Å; polymeric coating · USP L1
- eluent in column acetonitrile / water 70:30
- allows efficient gradient separation of the 16 PAH according to EPA
- detection of the separated PAH by UV (250 to 280 nm), with diode array or with fluorescence detection at different wavelengths for excitation and emission (acenaphthylene cannot be analysed with fluorescence detection)

Length →	50 mm	100 mm	125 mm	150 mm	250 mm	Guard columns
<b>NUCLEODUR® C<sub>18</sub> PAH, 1,8 <math>\mu</math>m</b>						particle size 1.8 $\mu$ m
<b>EC columns</b>						
	2 mm ID	760771.20	760773.20			
	3 mm ID	760771.30	760773.30			
	4 mm ID	760771.40	760773.40			
<b>NUCLEODUR® C<sub>18</sub> PAH, 3 <math>\mu</math>m</b>						particle size 3 $\mu$ m
	3 mm ID		760783.30	760784.30	760785.30	760786.30
	4 mm ID		760783.40	760784.30	760785.30	760786.30
<b>PAH standard according to EPA for HPLC</b>						
<b>PAH standard for HPLC</b>	16 PAH according to EPA method 610 in acetonitrile (1 ml)					722393

As guard columns for EC columns use ChromCart® guard column cartridges with guard column adaptor EC (REF 721359).  
8 mm ChromCart® guard column cartridges in packs of 3, analytical main columns in packs of 1.  
Microbore and VarioPrep columns on request.



## NUCLEODUR® PolarTec

RP phase with embedded polar group




🔸 **key features:**

- hydrophobic phase with pronounced hydrophilic properties
- separation mechanism based on hydrophobic (van der Waals) and polar interactions
- suitable for LC/MS due to low bleeding characteristics

🔸 **technical characteristics:**

embedded polar group, endcapped · 15.5 % C · USP L1/L60  
 suited for 100 % aqueous eluents  
 pore size 110 Å, available particle sizes 3 µm and 5 µm  
 pH stability 1 – 10

eluent in column acetonitrile / water

Length →	50 mm	75 mm	100 mm	125 mm	150 mm	250 mm	Guard columns
<b>NUCLEODUR® PolarTec, 3 µm</b>							particle size 3 µm
<b>EC columns</b>							
	2 mm ID	760473.20		760477.20	760478.20	760479.20	761160.30
	3 mm ID	760473.30		760477.30	760478.30	760479.30	761160.30
	4 mm ID	760473.40		760477.40	760478.40	760479.40	761160.40
	4.6 mm ID	760473.46	760475.46	760476.46	760477.46	760478.46	761160.40
<b>NUCLEODUR® PolarTec, 5 µm</b>							particle size 5 µm
	2 mm ID	760483.20		760487.20	760488.20	760489.20	761161.30
	3 mm ID	760483.30		760487.30	760488.30	760489.30	761161.30
	4 mm ID	760483.40		760487.40	760488.40	760489.40	761161.40
	4.6 mm ID	760483.46	760485.46	760486.46	760487.46	760488.46	761161.40

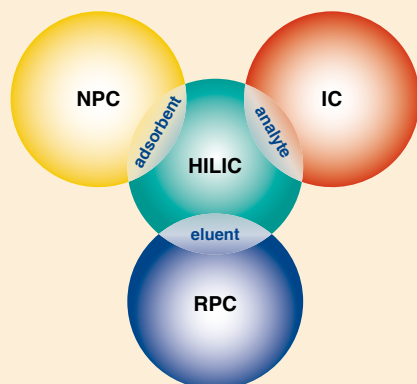
As guard columns for EC columns use ChromCart® guard column cartridges with guard column adaptor EC (REF 721359).  
 8 mm ChromCart® guard column cartridges in packs of 3, analytical main columns in packs of 1.  
 Microbore and VarioPrep columns on request.





## NUCLEODUR® HILIC

zwitterionic phase






### key features:

- ideal for reproducible and stable chromatography of highly polar analytes
- suitable for analytical and preparative applications as well as LC-MS
- very short column conditioning period

### technical characteristics:

ammonium - sulfonic acid modified silica; pore size 110 Å; particle sizes 1.8 µm, 3 µm and 5 µm; carbon content 7%; pH stability 2 – 8.5

eluent in column acetonitrile – water (80:20; v/v)

Length →	30 mm	50 mm	75 mm	100 mm	125 mm	150 mm	250 mm	Guard columns
<b>NUCLEODUR® HILIC, 1.8 µm</b>								particle size 1.8 µm
<b>EC columns</b>								
	2 mm ID	760521.20	760523.20	760525.20	760526.20		760528.20	
	3 mm ID	760521.30	760523.30					
	4 mm ID	760521.40	760523.40					
	4.6 mm ID	760521.46	760523.46					
<b>NUCLEODUR® HILIC, 3 µm</b>								particle size 3 µm
	2 mm ID		760532.20		760531.20	760533.20	760530.20	761580.30
	3 mm ID		760532.30		760531.30	760533.30	760530.30	761580.30
	4 mm ID		760532.40		760531.40	760533.40	760530.40	761580.40
	4.6 mm ID		760532.46		760534.46	760531.46	760533.46	760530.46
<b>NUCLEODUR® HILIC, 5 µm</b>								particle size 5 µm
	2 mm ID		760552.20		760551.20	760553.20	760550.20	761590.30
	3 mm ID		760552.30		760551.30	760553.30	760550.30	761590.30
	4 mm ID		760552.40		760551.40	760553.40	760550.40	761590.40
	4.6 mm ID		760552.46		760554.46	760551.46	760553.46	760550.46

As guard columns for EC columns use ChromCart® guard column cartridges with guard column adaptor EC (REF 721359).  
8 mm ChromCart® guard column cartridges in packs of 3, analytical main columns in packs of 1.  
Microbore and VarioPrep columns on request.



## Derivatization reagents

- for improved volatility, better thermal stability or a lower limit of detection in gas chromatography  
prerequisite: quantitative, rapid and reproducible formation of only one derivative  
halogen atoms introduced by derivatization (e.g. trifluoroacetates) allow specific detection (ECD) with the advantage of high sensitivity  
elution orders and fragmentation patterns in MS can be influenced by a specific derivatization
- reagents for **silylation**, **acylation** and **alkylation** (methylation) available



For the full range of derivatization reagents and method development kits please ask for our chromatography catalog "Columns and Supply" or visit [www.mn-net.com](http://www.mn-net.com)

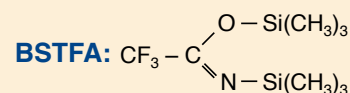
## Silylation reagents · BSTFA · SILYL-991

### N,O-bis-trimethylsilyl-trifluoroacetamide

m.w. 257.4, Bp 40 °C (12 mm Hg), density  $d_{20}^{40} = 0.961$

BSTFA is nonpolar (less polar than MSTFA), and can be mixed with acetonitrile for improved solubility.

For compounds, which are difficult to silylate (like secondary alcohols and amines), we recommend BSTFA + 1 % trimethylchlorosilane (TMCS), available under the designation SILYL-991.



## Silylation with BSA, BSTFA or SILYL-991 (BSTFA + 1 % TMCS)

### Procedure:

add 0.5 mL of the silylation reagent to 1 – 10 mg sample; if necessary, add some solvent (normally pyridine or DMF [dimethylformamide] are used). Heat to 60 – 80 °C for 20 min to increase the reaction rate. 1 – 2 drops of TMCS (trimethylchlorosilane) or TSIM will also speed up the reaction.

BSA MN Appl. No. 213091 · BSTFA MN Appl. No. 213092  
SILYL-991 MN Appl. No. 213093

	20 x 1 mL	1 x 10 mL	Packing unit 5 x 10 mL	1 x 50 mL	1 x 100 mL
BSTFA	<b>701220.201</b>	<b>701220.110</b>	<b>701220.510</b>		
SILYL-991 (BSTFA – TMCS (99:1))	<b>701490.201</b>			<b>701490.150</b>	<b>701490.1100</b>

Due to their purpose, derivatization reagents are very reactive substances. For this reason they should be stored cool and protected from moisture. Our derivatization reagents are supplied in vials with crimp caps for easy access with a syringe. Vials with pierced sealing disks have limited stability and should be used soon.



## MSTFA · MSHFBA · MBDSTFA

### ⬢ N-methyl-N-trimethylsilyl-trifluoroacetamide

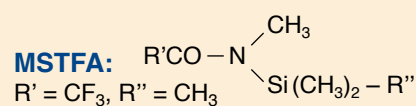
m.w. 199.1, Bp 70 °C (75 mm Hg), density  $d_{20}^{20}/4^{\circ} = 1.11$

### ⬢ N-methyl-N-trimethylsilyl-heptafluorobutyramide

m.w. 299.1, Bp 148 °C (760 mm Hg)

### ⬢ N-methyl-N-tert-butyl dimethylsilyl-trifluoroacetamide

m.w. 241.3, Bp 168 – 170 °C (760 mm Hg), density  $d_{20}^{20}/4^{\circ} = 1.121$



**MSHFBA:**  $\text{R}' = \text{C}_3\text{F}_7, \text{R}'' = \text{CH}_3$

### **MBDSTFA (MTB-TFA):**

$\text{R}' = \text{CF}_3, \text{R}'' = \text{C}_4\text{H}_9$

## Silylation with MSTFA, MSHFBA or MBDSTFA

### Procedure:

Dissolve 10 – 15 mg sample in 0.8 mL solvent, then add 0.2 mL of the silylation reagent. The reaction mixture can be heated to 60 – 70 °C for up to 1 hour and can be analysed directly. If TFA is used as a solvent, proceed as follows [M. Donike, J. Chromatogr. 85 (1973) 1 – 7]: dissolve 1 – 2 mg sample in 100  $\mu\text{L}$  TFA. Dropwise add 0.9 mL of the silylating reagent. After cooling the sample can be chromatographed directly.

MSTFA MN Appl. No. 213111 · MSHFBA MN Appl. No. 213112 · MBDSTFA MN Appl. No. 213113

	Packing unit							
	10 x 1 mL	20 x 1 mL	1 x 10 mL	5 x 10 mL	1 x 100 mL	6 x 50 mL	6 x 100 mL	12 x 100 mL
MSHFBA		<b>701260.201</b>	<b>701260.110</b>	<b>701260.510</b>	<b>701260.1100</b>		<b>701260.6100</b>	
MSTFA		<b>701270.201</b>	<b>701270.110</b>	<b>701270.510</b>	<b>701270.1100</b>	<b>701270.650</b>	<b>701270.6100</b>	<b>701270.12100</b>
MBDSTFA	<b>701440.101</b>	<b>701440.201</b>						

## Acylation reagents

### ⬢ Anhydrides

#### Heptafluorobutyric acid anhydride

**HFBA:**  $\text{C}_3\text{F}_7 - \text{CO} - \text{O} - \text{CO} - \text{C}_3\text{F}_7$

m.w. 410.06, Bp 106 – 107 °C (760 mm Hg),  
density  $d_{20}^{20}/4^{\circ} = 1.665$

### ⬢ Bisacylamides

#### N-methyl-bis(trifluoroacetamide)

**MBTFA:**  $\text{CF}_3 - \text{CO} - \text{N}(\text{CH}_3) - \text{CO} - \text{CF}_3$

m.w. 223.08, Bp 123 – 124 °C (760 mm Hg),  
density  $d_{20}^{20}/4^{\circ} = 1.55$

## Methods for acylation

### Acylation with fluorinated acid anhydrides:

Acylation with HFBA can be used for alcohols, phenols, carboxylic acids, amines, amino acids and steroids forming volatile, stable derivatives suited for FID as well as for ECD detection.

### Procedure:

Dissolve 0.1 to 1 mg of the sample in 0.1 mL solvent, add 0.1 mL of the respective anhydride and heat to 60 – 70 °C for 1 – 2 hours. If the sample need not be concentrated prior to the analysis and if there is no danger of catalytically induced side reactions, pyridine is used as solvent. The reaction solution can be injected directly into the gas chromatograph. Otherwise use a volatile solvent and evaporate solvent, excess reagent and acid in a stream of nitrogen. Dissolve the residue in 50  $\mu\text{L}$  hexane, chloroform etc. and inject aliquot portions.

TFAA MN Appl. No. 213041 · HFBA MN Appl. No. 213042

### Acylation with fluorinated acid amides:

This method is recommended for alcohols, primary and secondary amines as well as for thiols under mild, neutral conditions.

### Procedure:

Add 0.5 mL MBTFA to about 2 mg sample. If there is no reaction at ambient temperature, heat the reaction mixture to 120 °C. Compounds which are difficult to dissolve, can be trifluoroacetylated in suitable solvent mixtures. It is recommended to use a ratio of solvent to MBTFA of 4 : 1. The reaction mixture can be chromatographed directly.

MBTFA MN Appl. No. 213051 · MBHFBA MN Appl. No. 213052

	Packing unit		
	20 x 1 mL	1 x 10 mL	5 x 10 mL
HFBA	<b>701110.201</b>	<b>701110.110</b>	<b>701110.510</b>
MBTFA	<b>701410.201</b>	<b>701410.110</b>	<b>701410.510</b>

Due to their purpose, derivatization reagents are very reactive substances. For this reason they should be stored cool and protected from moisture. Our derivatization reagents are supplied in vials with crimp caps for easy access with a syringe. Vials with pierced sealing disks have limited stability and should be used soon.

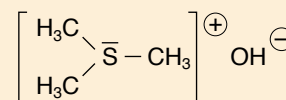


## Alkylation reagents

In GC generally methylation is the main type of alkylation used.

### Trimethylsulphonium hydroxide

**TMSH** (0.2 M in MeOH) · m.w. 94.06



## Methods for methylation

### Methylation with TMSH

Methylation with TMSH is recommended for free acids, chlorophenoxyacetic acids, their salts and derivatives as well as for phenols and chlorophenols.

#### Procedure:

Dissolve 100 mg sample (e. g. butter) in 5 mL of a suitable solvent (e. g. *tert.*-butyl methyl ether). Add 50 µl reagent to 100 µl of this solution. The mixture is injected directly. The temperature of the injector must be at least 250 °C.

MN Appl. No. 213060

	Packing unit			
	10 x 1 mL	20 x 1 mL	1 x 10 mL	5 x 10 mL
TMSH	<b>701520.101</b>	<b>701520.201</b>	<b>701520.110</b>	<b>701520.510</b>

## Useful consumables for your GC

### Ferrules for GC

- Graphite ferrules provide the highest temperature stability (up to 450 °C). They are reusable when handled with care. We also offer 1/16" graphite ferrules specially designed for Carlo Erba / Fisons or for Agilent gas chromatographs.
- Vespel ferrules come in three types: pure Vespel, Vespel with 15 % graphite and Vespel with 40 % graphite. All versions are stable up to 400 °C and reusable.
- PTFE ferrules can only be used up to 250 °C. They are not reusable and not recommended for temperature programming. However, they show the best chemical inertness of all ferrules.



(packing unit 10 ferrules)

Bore (= column OD)	Graphite		Vespel		PTFE
	max. temp. →	plain	+ 15 % graphite	+ 40 % graphite	
	450 °C	400 °C	400 °C	400 °C	250 °C
<b>1/16" ferrules</b>					
no bore	<b>708336</b>	<b>706187</b>	<b>706167</b>		<b>706177</b>
0.4 mm	<b>708309</b>			<b>706246</b>	
0.5 mm	<b>708308</b>			<b>706247</b>	
0.8 mm	<b>708301</b>			<b>706248</b>	
1.0 mm	<b>708302</b>				
1.2 mm	<b>708303</b>				
1/16"	<b>706155</b>	<b>706180</b>	<b>706160</b>	<b>706190</b>	<b>706170</b>



## Septa for GC

Designation	Material	Thickness	Hardness	max. Temp.
Standard septa (ST)	beige silicone rubber	4 mm	60 shore	
High temperature septa (HT)	red, specially pretreated, non-bleeding silicone rubber	3 mm	60 shore	320 °C *
Silicone septa, soft	transparent silicone rubber	3 mm	45 shore	250 °C
Silicone septa PTFE	white silicone rubber, one side coated with grey PTFE	3 mm		200 °C

\* When used at considerably higher temperatures – and working without septum purge – interfering peaks can occur due to thermal decomposition of the material.

Septum grade (packs of 50 septa)	Outer diameter					
	9 mm N 9	10 mm N 10	11 mm N 11	12 mm N 12	13 mm N 13	17 mm N 17
Standard septa (ST)	702609	702610	702611	702612	702613	
High temperature septa (HT)	702619	702620	702621	702622	702623	702632
Silicone septa, soft	702602		702604	702605	702606	
Silicone septa PTFE		702625	702626	702627	702628	
Septum remover (tool for removing septa which have become baked into the injection port of the gas chromatograph)						706141

## Tools and general accessories for GC

- Diamond file:**  
 a useful tool for cutting capillaries and smoothing ends of capillaries. Square capillary ends without protruding particles are especially important for butt connections (e.g. in Valco unions).
- Magnifying lens:**  
 a very versatile tool for any laboratory. In capillary GC it is often important to inspect column integrity or check cut ends of capillaries. When closing a column by melting the magnifying lens can be used to check whether the column is really closed or whether an open channel has been formed in the sealed end. Our lens provides 8fold magnification and is supplied with a scale as pictured in the figure below. The space between lines corresponds to 1/10 mm.



Lens with scale



Diamond file

Description	Specification	Pack of	REF
<b>Tools for capillary GC</b>			
Diamond file	for cutting capillaries and straightening capillary ends	1	708300
Magnifying lens with scale	magnification 8x	1	706296

Description	Pack of	REF	Specification
<b>Universal capillary glass connectors</b>			
linear	5 connectors	707971	
linear	10 connectors	707972	
Y splitter	1 connector	707973	
PTFE shrink- ing tube, thin-walled	1 m	708305	for connecting capillaries, min ID expanded 1.17 mm, max. ID shrunk 0.40 mm



## Polyamide (PA) = Nylon

- ◈ hydrophilic membrane
- ◈ for aqueous and organic/aqueous medium polar liquids

	Type	Pore size [µm]	Membrane diameter [mm]	Color code		Standard pack		BIG-BOX	
				top	bottom	filters/pack	REF	filters/pack	REF
	Xtra PA-20/25	0.20	25	labelled	–	100	<b>729212</b>	400	<b>729212.400</b>
	Xtra PA-45/25	0.45	25	labelled	–	100	<b>729213</b>	400	<b>729213.400</b>
	AO-20/15 MS*	0.20	15	yellow	green	100	<b>729048</b>	800	<b>729048.800</b>
	AO-45/15 MS*	0.45	25	colorless	green	100	<b>729049</b>	800	<b>729049.800</b>
	AO-20/3	0.20	3	colorless	colorless	100	<b>729010</b>		
	AO-45/3	0.45	3	colorless	colorless	100	<b>729011</b>		

\* MS = minispikes on filter exit



## Polytetrafluoroethylene (PTFE)

- ◈ hydrophobic membrane
- ◈ for nonpolar liquids and gases
- ◈ very resistant towards all kinds of solvents as well as acids and bases  
flushing with alcohol, followed by water, makes the originally hydrophobic membrane more hydrophilic

	Type	Pore size [µm]	Membrane diameter [mm]	Color code		Standard pack		BIG-BOX	
				top	bottom	filters/pack	REF	filters/pack	REF
	Xtra PTFE-20/25	0.20	25	labelled		100	<b>729207</b>	400	<b>729207.400</b>
	Xtra PTFE-45/25	0.45	25	labelled		100	<b>729205</b>	400	<b>729205.400</b>
	Xtra PTFE-100/25	0.1	25	labelled	colorless	100	<b>729247</b>	400	<b>729247.400</b>
	O-20/15 MS*	0.20	15	yellow	colorless	100	<b>729008</b>	800	<b>729008.800</b>
	O-45/15 MS*	0.45	15	colorless	colorless	100	<b>729009</b>	800	<b>729009.800</b>
	O-20/3	0.20	3	colorless	colorless	100	<b>729014</b>		
	O-45/3	0.45	3	colorless	colorless	100	<b>729015</b>		

\* MS = minispikes on filter exit



## Regenerated Cellulose (RC)

- hydrophilic membrane with very low adsorption
- for aqueous and organic/aqueous liquids, i. e. polar and medium polar sample solutions
- binding capacity for proteins 84 µg/filter of 25 mm diameter
- RC filter with integrated glass fibre prefilter (**GF/RC**): recommended for solutions with a high load of particulate matter or for highly viscous solutions

	Type	Pore size [µm]	Membrane diameter [mm]	Color code		Standard pack		BIG-BOX	
				top	bottom	filters/pack	REF	filters/pack	REF
	Xtra RC-20/25	0.20	25	labelled		100	<b>729230</b>	400	<b>729230.400</b>
	Xtra RC-45/25	0.45	25	labelled		100	<b>729231</b>	400	<b>729231.400</b>
	RC-20/15 MS*	0.20	15	yellow	blue	100	<b>729036</b>	800	<b>729036.800</b>
	RC-45/15 MS*	0.45	15	colorless	blue	100	<b>729037</b>	800	<b>729037.800</b>
<b>Combi Filters</b>									
	GF/RC-20/25	1.0/0.20	25	blue	blue	100	<b>729050</b>	400	<b>729050.400</b>
	GF/RC-45/25	1.0/0.45	25	black	blue	100	<b>729051</b>	400	<b>729051.400</b>

\* MS = minispikes on filter exit



## Polyester (PET)

- hydrophilic multipurpose membrane
- for polar as well as nonpolar solvents  
**the HPLC filter**, especially suited for mixtures of water and organic solvents  
 for TOC/DOC determination, not cytotoxic, does not inhibit the growth of microorganisms and higher cells
- polyester filter with integrated glass fibre prefilter (**GF/PET**): recommended for solutions with a high load of particulate matter or for highly viscous solutions





	Type	Pore size [µm]	Membrane diameter [mm]	Color code		Standard pack		BIG-BOX	
				top	bottom	filters/pack	REF	filters/pack	REF
	Xtra PET-20/25	0.20	25	labelled		100	<b>729221</b>	400	<b>729221.400</b>
	Xtra PET-45/25	0.45	25	labelled		100	<b>729220</b>	400	<b>729220.400</b>
	Xtra PET-120/25	1.2	25	labelled		100	<b>729229</b>	400	<b>729229.400</b>
	PET-20/15 MS*	0.20	15	yellow	orange	100	<b>729022</b>	800	<b>729022.800</b>
	PET-45/15 MS*	0.45	15	colorless	orange	100	<b>729023</b>	800	<b>729023.800</b>
<b>Combi Filters</b>									
	GF/PET-20/25	1.0/0.20	25	blue	orange	100	<b>729032</b>	400	<b>729032.400</b>
	GF/PET-45/25	1.0/0.45	25	black	orange	100	<b>729033</b>	400	<b>729033.400</b>

\* MS = minispikes on filter exit








# Vials and Caps



## Screw neck vials and inserts N 9

702283	702284	702813	702818
			
1.5 mL	1.5 mL	0.2 mL	0.1 mL
11.6 x 32 mm	11.6 x 32 mm	6 x 31 mm	5.7 x 29 mm
clear, label + scale	amber, label + scale	clear	clear
flat bottom	flat bottom	conical, 15 mm tip	with plastic spring
100 / PP box	100 / PP box	100 / PE bag	100 / PE bag

## Ready assembled screw closures N 9 and single septa N 9

REF	(Scale 1:1.6)	Cap description	Septum description	Hardness	Thickness	Pack of
702732		N 9 PP screw cap, color as indicated, center hole	Red Rubber / FEP colorless	40° shore A	1.0 mm	100 / PE bag
702033		as above, but with closed top				
702287.1		N 9 PP screw cap, color as indicated, center hole	Silicone white / PTFE red	40° shore A	1.0 mm	100 / PE bag
702288.1		N 9 PP screw cap, color as indicated, center hole	Silicone white / PTFE blue, slit	40° shore A	1.0 mm	100 / PE bag
702035		N 9 PP screw cap, color as indicated, center hole	PTFE red / Silicone white / PTFE red	40° shore A	1.0 mm	100 / PE bag

## Bonded screw closures N 9 (septa firmly connected with the cap; cannot be separated)

REF		Cap description	Septum description	Hardness	Thickness	Pack of
702026		N 9 PP bonded screw cap, blue, center hole	Silicone beige / PTFE white	45° shore A	1.3 mm	100 / PE bag
702027		N 9 PP bonded screw cap, blue, center hole	Silicone beige / PTFE white, slit	45° shore A	1.3 mm	100 / PE bag





## Crimp neck vials and inserts N 11

702885	702892	702813	702818
1.5 mL	1.5 mL	0.2 mL	0.1 mL
11.6 x 32 mm	11.6 x 32 mm	6 x 31 mm	5.7 x 29 mm
clear, label + scale	amber, label + scale	clear	clear
flat bottom	flat bottom	conical, 15 mm tip	with plastic spring
100 / PP box	100 / PP box	100 / PE bag	100 / PE bag

## Ready assembled crimp closures N 11, plain crimp caps N 11 and single septa N 11

REF	Cap description	Septum description	Hardness	Thickness	Pack of
70256	N 11 aluminium crimp cap, silver, center hole	Natural rubber / Butyl red-orange / TEF colorless	45° shore A	1.0 mm	100 / PE bag
70288	N 11 aluminium crimp cap, silver, center hole	Silicone white / PTFE red	40° shore A	1.3 mm	100 / PE bag
702823	N 11 aluminium crimp cap, silver, center hole	Silicone white / PTFE blue, cross-slit	40° shore A	1.5 mm	100 / PE bag
702995	N 11 aluminium crimp cap, silver, center hole	PTFE red / Silicone white / PTFE red	40° shore A	1.0 mm	100 / PE bag
702879	N 11 magnetic crimp cap, gold, center hole	Silicone white / PTFE red	55° shore A	1.0 mm	100 / PE bag

## Crimping tools N 11

REF	Type of crimping tool	Pack of
735111	Manual crimper, height adjustable, for 11 mm aluminium crimp caps	1 / box
735911	Manual decapper for 11 mm aluminium crimp caps	1 / box

## Ready assembled magnetic screw closures N 18

REF	Cap description	Septum description	Hardness	Thickness	Pack of
702055	N 18 magnetic screw cap, silver, center hole	Silicone white / PTFE blue	55° shore A	1.5 mm	100 / PE bag



# Vials and Caps

## Screw neck Headspace vials N 18










702826

20 mL
22.5 x 75.5 mm
clear
rounded bottom
100 / PP box

## Crimp neck vials N 20: 20 and 50 mL

REF 70254	REF 702261	REF 702263
		
PerkinElmer	DANI, Agilent	CTC, Varian
20 mL	20 mL	20 mL
23 x 75.5 mm	22.5 x 75.5 mm	22.5 x 75.5 mm
clear	clear	clear
rounded bottom	flat bottom	rounded bottom
bevelled HS crimp neck	flat DIN crimp neck	flat DIN crimp neck
100 / PP box	100 / PP box	100 / PP box

## Ready assembled crimp closures, plain crimp caps and single septa N 20

REF	Cap description	Septum description	Hardness	Thickness	Pack of
 <b>Center hole caps</b>					
702775	 N 20 aluminium crimp cap, silver, center hole	Butyl light grey / PTFE dark grey	50° shore A	3 mm	100 / PE bag
70234	 N 20 aluminium crimp cap, silver, center hole	Butyl dark grey / PTFE grey (only centrally laminated, typically called Pharma-Fix)	50° shore A	3 mm	100 / PE bag
702817	 N 20 aluminium crimp cap, silver, center hole	Silicone blue / PTFE colorless	40° shore A	3 mm	100 / PE bag
 <b>Bi-metal crimp caps</b>					
702838	 N 20 bi-metal crimp cap, blue/silver, center hole	Butyl light grey / PTFE dark grey	50° shore A	3 mm	100 / PE bag
702834	 N 20 bi-metal crimp cap, blue/silver, center hole	Silicone blue / PTFE colorless	40° shore A	3 mm	100 / PE bag
 <b>Magnetic crimp caps</b>					
702929	 N 20 magnetic crimp cap, silver, center hole 8 mm	Silicone blue / PTFE colorless	40° shore A	3 mm	100 / PE bag



## Screw neck vials N 24 (EPA)

702021	702022	702023	702024
20 mL	20 mL	40 mL	40 mL
27.5 x 57 mm	27.5 x 57 mm	27.5 x 95 mm	27.5 x 95 mm
clear, flat bottom	amber, flat bottom	clear, flat bottom	amber, flat bottom
100 / PP box	100 / PP box	100 / PP box	100 / PP box

## Bonded screw closures N 24, plain screw caps N 24 and single septa N 22

(images scale 1:2)

702058	702059	702060	702061	702062
<b>Bonded screw closures N 24</b> (septa firmly connected with the cap; cannot be separated)		<b>Plain screw caps N 24</b>		<b>Septum N 22</b>
N 24 PP bonded screw cap, white, center hole	N 24 PP bonded screw cap, white, closed top	N 24 PP screw cap, white, center hole	N 24 PP screw cap, white, closed top	Silicone natural / PTFE beige, 45° shore A, 3.2 mm
Silicone white / PTFE beige, 45° shore A, 3.2 mm	Silicone white / PTFE beige, 45° shore A, 3.2 mm	no liner	no liner	100 / PE bag
100 / PE bag	100 / PE bag	100 / PE bag	100 / PE bag	100 / PE bag

## Crimping tools N 20

REF	Type of crimping tool	Pack of
735120	Manual crimper, height adjustable, for 20 mm aluminium crimp caps	1 / box
735920	Manual decapper for 20 mm aluminium crimp caps	1 / box





HPLC



GC



TLC



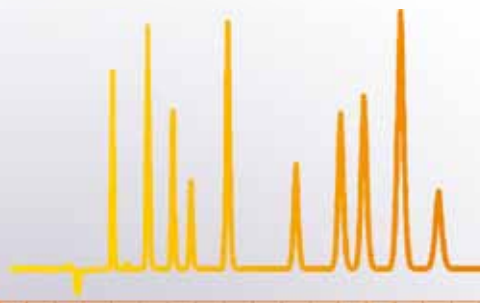
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KÄTEN20087  
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