LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017072.D
Acq On : 10 Sep 2023 03:15
Operator : MA/JU
Sample : 04227-06
Misc :
ALS Vial : 84 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BBKE5

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP017072.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 3.276 | 27 | 32 | 39 | rBV | 2692330 | 4173343 | 6.89% | 1.916% |
| 2 | 3.358 | 43 | 46 | 61 | rVB | 603135 | 838255 | 1.38% | 0.385% |
| 3 | 3.764 | 111 | 115 | 131 | rBV | 730517 | 1223107 | 2.02% | 0.562% |
| 4 | 3.958 | 143 | 148 | 155 | rVB | 78532 | 133555 | 0.22% | 0.061% |
| 5 | 4.593 | 250 | 256 | 260 | rBV | 30872004 | 60577641 | 100.00% | 27.811% |
| 6 | 6.652 | 601 | 606 | 614 | rBV | 242188 | 351834 | 0.58% | 0.162% |
| 7 | 7.122 | 680 | 686 | 697 | rVB | 472072 | 916546 | 1.51% | 0.421% |
| 8 | 7.311 | 707 | 718 | 728 | rBV | 2291701 | 3487877 | 5.76% | 1.601% |
| 9 | 7.487 | 741 | 748 | 757 | rBV | 2070751 | 3335445 | 5.51% | 1.531% |
| 10 | 7.569 | 757 | 762 | 770 | rVB3 | 29696 | 67617 | 0.11% | 0.031% |
| 11 | 7.969 | 822 | 830 | 841 | rBV | 1852598 | 2879642 | 4.75% | 1.322% |
| 12 | 8.699 | 943 | 954 | 961 | rBV | 1165645 | 2631672 | 4.34% | 1.208% |
| 13 | 8.758 | 961 | 964 | 970 | rVB | 228770 | 323484 | 0.53% | 0.149% |
| 14 | 9.169 | 1025 | 1034 | 1046 | rBV | 2617189 | 4355728 | 7.19% | 2.000% |
| 15 | 9.481 | 1079 | 1087 | 1100 | rVB | 10429285 | 17740393 | 29.29% | 8.144% |
| 16 | 9.881 | 1148 | 1155 | 1167 | rBV | 2172567 | 3481939 | 5.75% | 1.599% |
| 17 | 10.416 | 1235 | 1246 | 1263 | rBV | 3050733 | 5021237 | 8.29% | 2.305% |
| 18 | 10.793 | 1301 | 1310 | 1319 | rBV | 2589963 | 4263949 | 7.04% | 1.958% |
| 19 | 10.958 | 1330 | 1338 | 1351 | rBV | 2880445 | 4728182 | 7.81% | 2.171% |
| 20 | 12.375 | 1574 | 1579 | 1587 | rBV | 51587 | 87136 | 0.14% | 0.040% |
| 21 | 13.469 | 1759 | 1765 | 1771 | rVV | 108922 | 180399 | 0.30% | 0.083% |
| 22 | 13.604 | 1779 | 1788 | 1794 | rBV7 | 29078 | 65758 | 0.11% | 0.030% |
| 23 | 14.040 | 1855 | 1862 | 1875 | rBV | 5697525 | 7686954 | 12.69% | 3.529% |
| 24 | 14.328 | 1903 | 1911 | 1920 | rBV | 6331161 | 9235805 | 15.25% | 4.240% |
| 25 | 14.551 | 1944 | 1949 | 1955 | rBV | 107377 | 169934 | 0.28% | 0.078% |
| 26 | 14.628 | 1955 | 1962 | 1969 | rVB | 3778651 | 5479968 | 9.05% | 2.516% |
| 27 | 14.846 | 1989 | 1999 | 2014 | rBV | 403699 | 690387 | 1.14% | 0.317% |
| 28 | 15.169 | 2048 | 2054 | 2063 | rBV | 488303 | 685240 | 1.13% | 0.315% |
| 29 | 15.622 | 2123 | 2131 | 2139 | rBV | 8354293 | 11394931 | 18.81% | 5.231% |
| 30 | 15.746 | 2146 | 2152 | 2164 | rBV | 2286012 | 3011069 | 4.97% | 1.382% |
| 31 | 15.893 | 2173 | 2177 | 2186 | rVB | 232676 | 335284 | 0.55% | 0.154% |
| 32 | 16.251 | 2233 | 2238 | 2245 | rBV | 170177 | 243518 | 0.40% | 0.112% |
| 33 | 16.351 | 2251 | 2255 | 2260 | rVB2 | 70079 | 94923 | 0.16% | 0.044% |
| 34 | 16.998 | 2361 | 2365 | 2370 | rBV3 | 34008 | 64758 | 0.11% | 0.030% |
| 35 | 17.392 | 2425 | 2432 | 2441 | rBV | 4679677 | 6489388 | 10.71% | 2.979% |
| 36 | 17.492 | 2443 | 2449 | 2461 | rVV2 | 8825252 | 12179132 | 20.10% | 5.591% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017072.D
Acq On : 10 Sep 2023 03:15
Operator : MA/JU
Sample : 04227-06
Misc :
ALS Vial : 84 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BBKE5

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 18.234 | 2570 | 2575 | 2588 | rBV | 881368 | 1321691 | 2.18% | 0.607% |
| 38 | 19.298 | 2753 | 2756 | 2760 | rVB2 | 91623 | 109664 | 0.18% | 0.050% |
| 39 | 19.369 | 2764 | 2768 | 2777 | rVB | 122932 | 173769 | 0.29% | 0.080% |
| 40 | 19.469 | 2781 | 2785 | 2795 | rBV | 439689 | 643830 | 1.06% | 0.296% |
| 41 | 19.604 | 2805 | 2808 | 2812 | rVB | 78793 | 86686 | 0.14% | 0.040% |
| 42 | 19.728 | 2823 | 2829 | 2835 | rBV | 10005481 | 13437034 | 22.18% | 6.169% |
| 43 | 21.157 | 3068 | 3072 | 3079 | rBV | 434584 | 637552 | 1.05% | 0.293% |
| 44 | 21.486 | 3123 | 3128 | 3136 | rBV | 4409167 | 5799447 | 9.57% | 2.662% |
| 45 | 23.851 | 3523 | 3530 | 3545 | rVB2 | 5266034 | 10955413 | 18.08% | 5.030% |
| 46 | 24.022 | 3552 | 3559 | 3569 | rVB | 2492018 | 5285755 | 8.73% | 2.427% |
| 47 | 24.574 | 3649 | 3653 | 3663 | rVB | 371564 | 745623 | 1.23% | 0.342% |

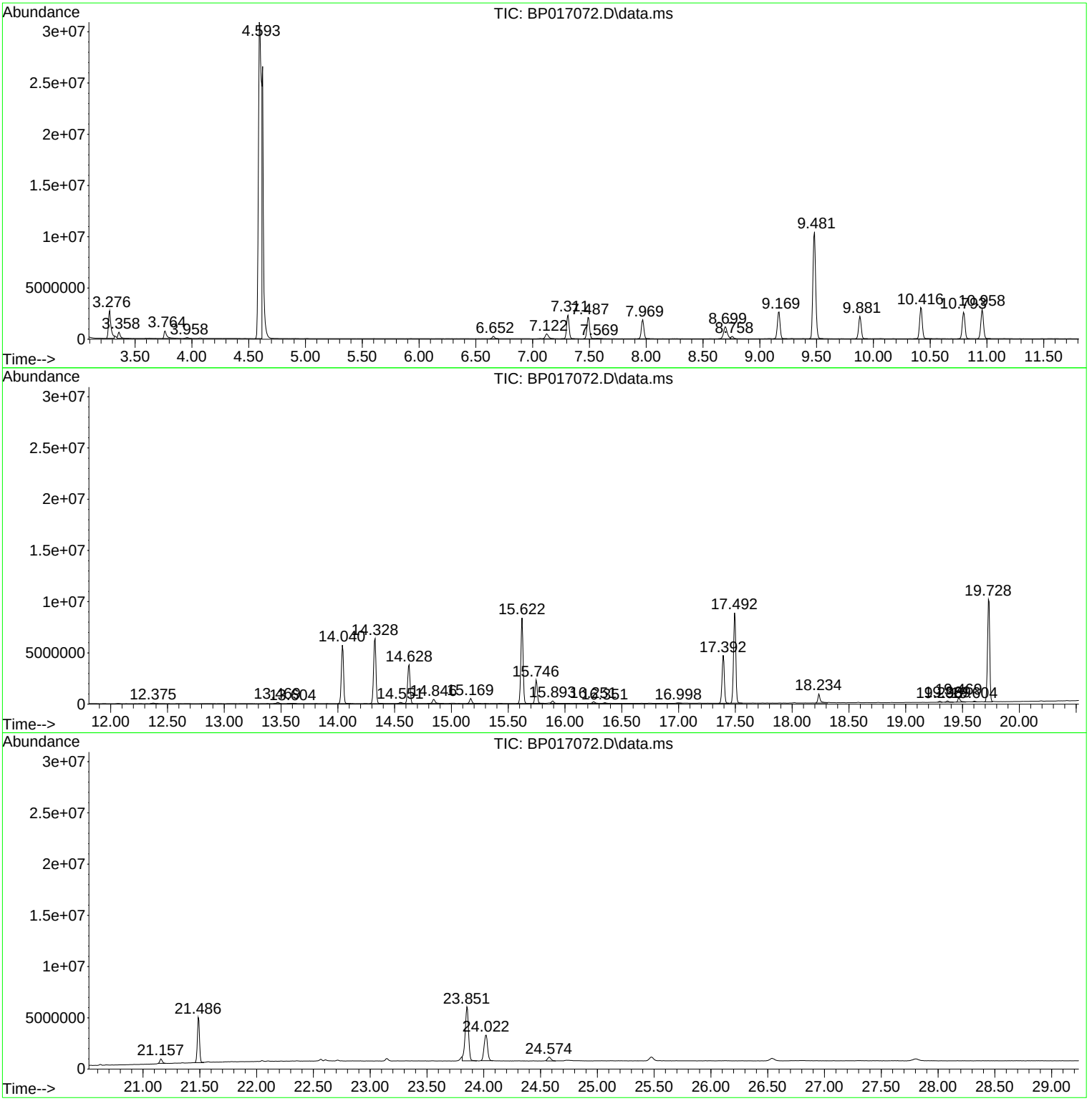
Sum of corrected areas: 217822494

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017072.D
Acq On : 10 Sep 2023 03:15
Operator : MA/JU
Sample : 04227-06
Misc :
ALS Vial : 84 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BBKE5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017072.D
Acq On : 10 Sep 2023 03:15
Operator : MA/JU
Sample : 04227-06
Misc :
ALS Vial : 84 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BBKE5

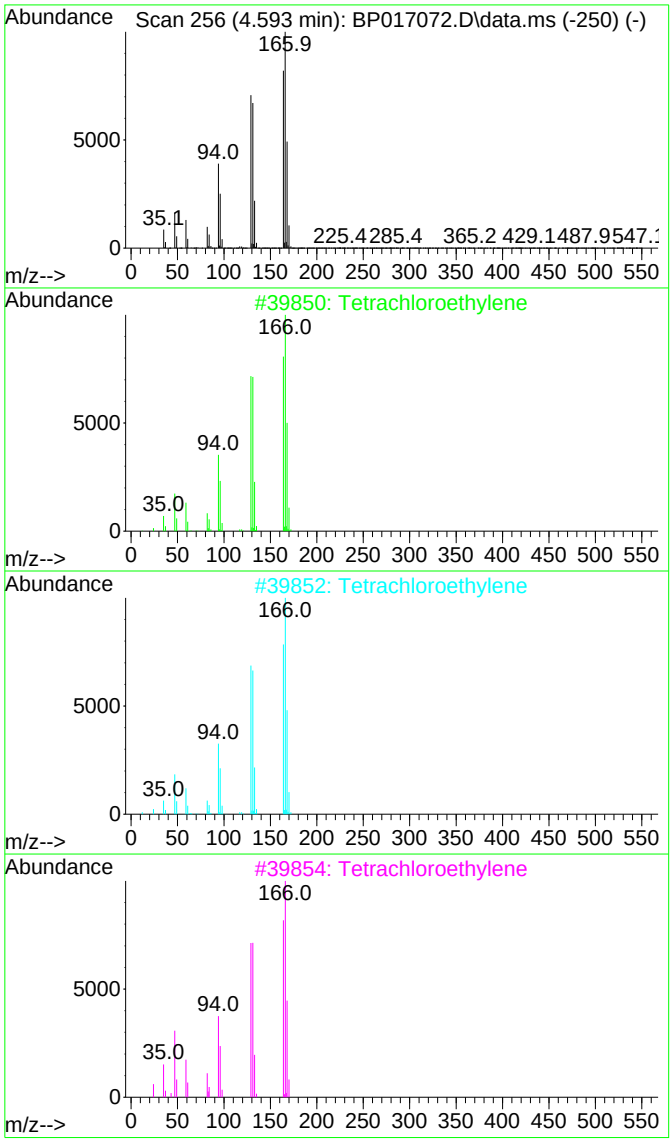
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

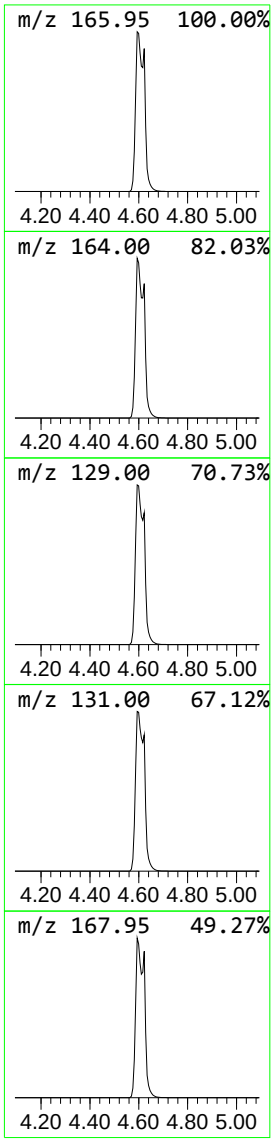
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 1 Tetrachloroethylene Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.593 | 420.73 ng/ul | 60577600 | 1,4-Dichlorobenzene-d4 | 7.969 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Tetrachloroethylene | 164 | C2Cl4 | 000127-18-4 | 99 |
| 2 | Tetrachloroethylene | 164 | C2Cl4 | 000127-18-4 | 99 |
| 3 | Tetrachloroethylene | 164 | C2Cl4 | 000127-18-4 | 97 |
| 4 | Tetrachloroethylene | 164 | C2Cl4 | 000127-18-4 | 97 |
| 5 | 2,4-Dichloro-6-fluoropyrimidine | 166 | C4HCl2FN2 | 003833-57-6 | 35 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017072.D
Acq On : 10 Sep 2023 03:15
Operator : MA/JU
Sample : 04227-06
Misc :
ALS Vial : 84 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BBKE5

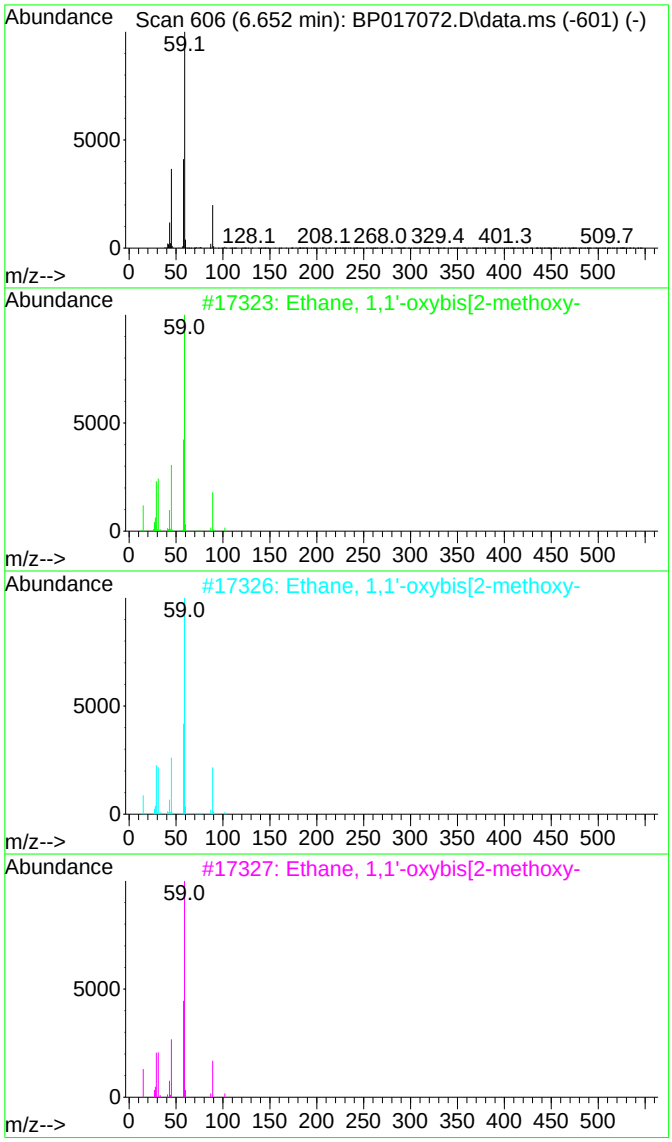
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

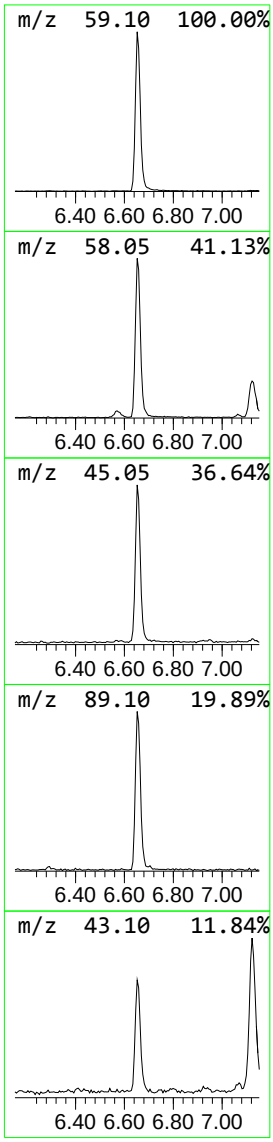
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 2 Ethane, 1,1'-oxybis[2-methoxy- Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.652 | 2.44 ng/ul | 351834 | 1,4-Dichlorobenzene-d4 | 7.969 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Ethane, 1,1'-oxybis[2-methoxy- | 134 | C6H14O3 | 000111-96-6 | 83 |
| 2 | Ethane, 1,1'-oxybis[2-methoxy- | 134 | C6H14O3 | 000111-96-6 | 83 |
| 3 | Ethane, 1,1'-oxybis[2-methoxy- | 134 | C6H14O3 | 000111-96-6 | 74 |
| 4 | Ethane, 1,1'-oxybis[2-methoxy- | 134 | C6H14O3 | 000111-96-6 | 74 |
| 5 | 2,5,8,11-Tetraoxadodecane | 178 | C8H18O4 | 000112-49-2 | 72 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017072.D
Acq On : 10 Sep 2023 03:15
Operator : MA/JU
Sample : 04227-06
Misc :
ALS Vial : 84 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BBKE5

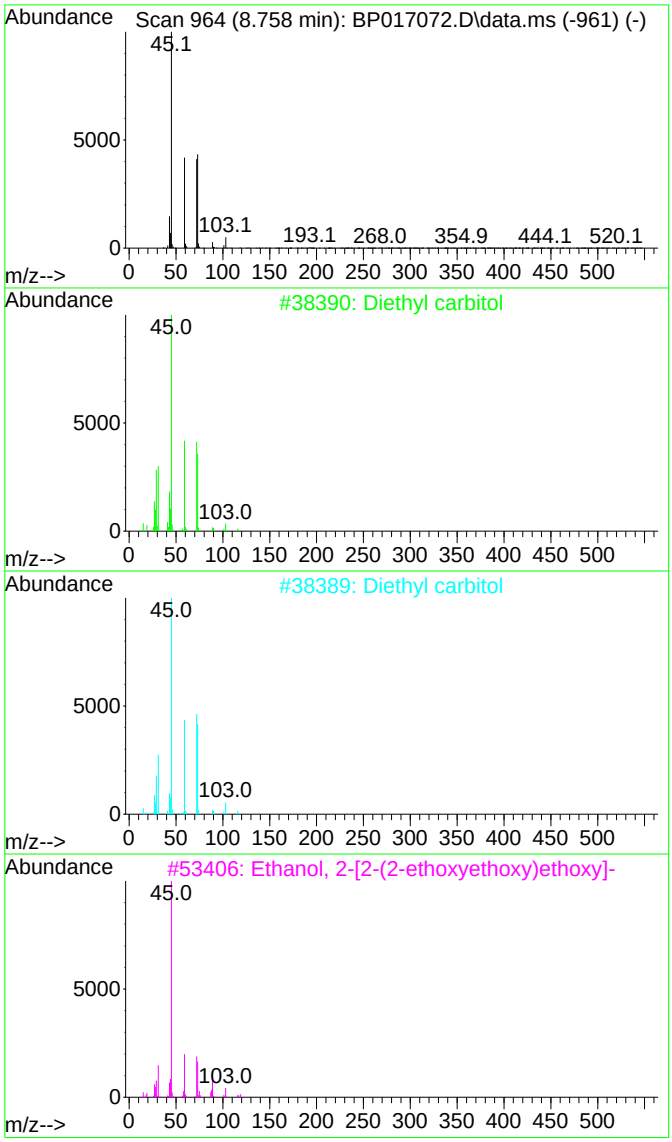
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

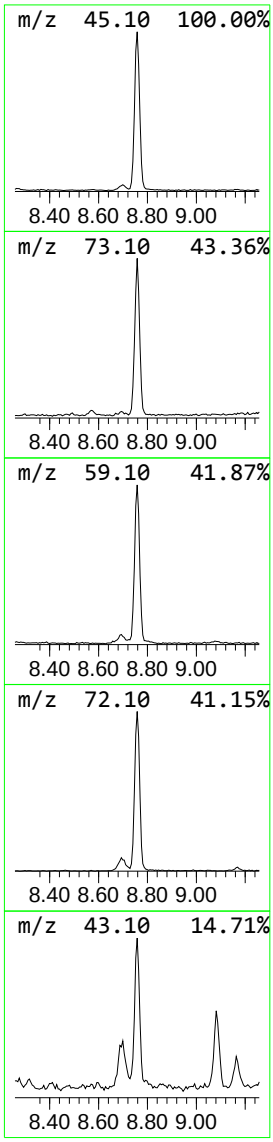
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 3 Diethyl carbitol Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.758 | 2.25 ng/ul | 323484 | 1,4-Dichlorobenzene-d4 | 7.969 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Diethyl carbitol | 162 | C8H18O3 | 000112-36-7 | 83 |
| 2 | Diethyl carbitol | 162 | C8H18O3 | 000112-36-7 | 83 |
| 3 | Ethanol, 2-[2-(2-ethoxyethoxy)et... | 178 | C8H18O4 | 000112-50-5 | 78 |
| 4 | Ethanol, 2-(2-ethoxyethoxy)- | 134 | C6H14O3 | 000111-90-0 | 64 |
| 5 | Ethanol, 2-(2-ethoxyethoxy)- | 134 | C6H14O3 | 000111-90-0 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017072.D
Acq On : 10 Sep 2023 03:15
Operator : MA/JU
Sample : 04227-06
Misc :
ALS Vial : 84 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BBKE5

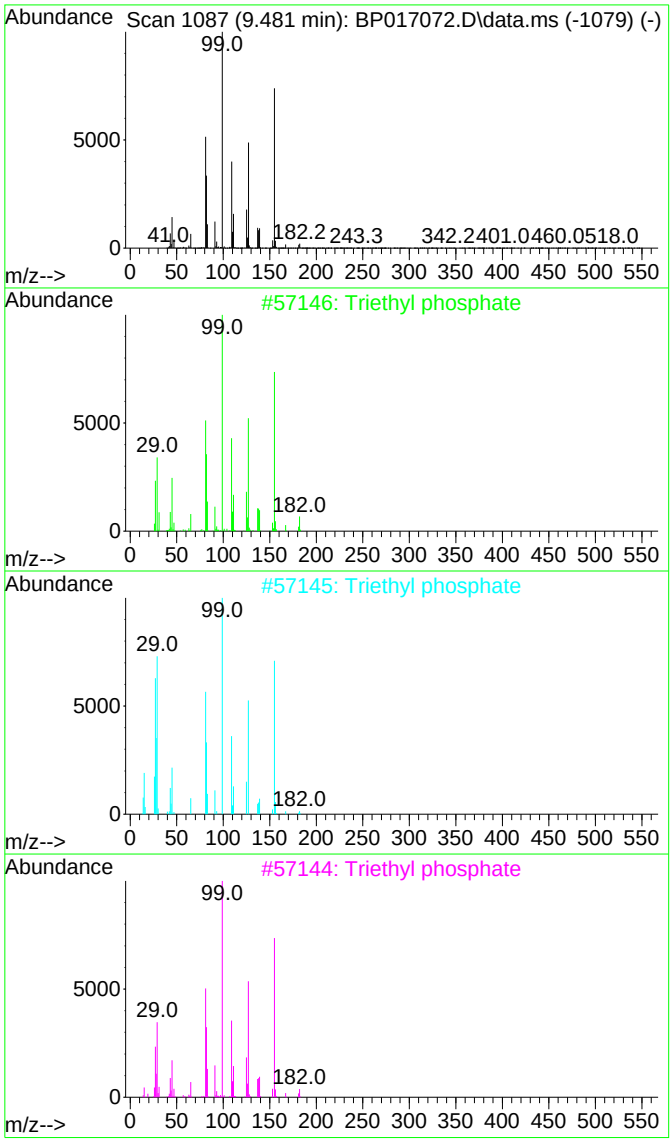
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

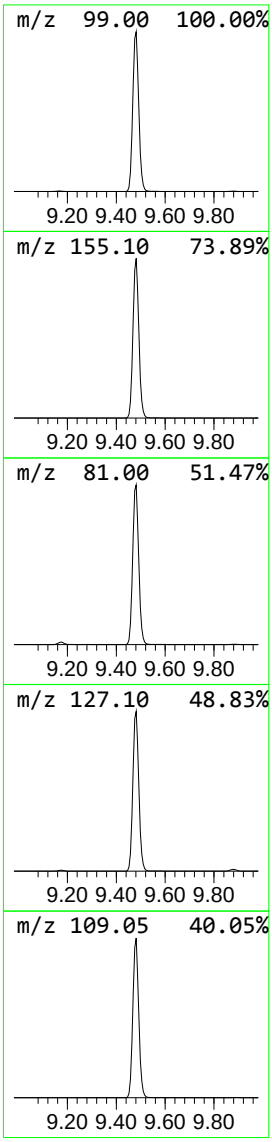
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 4 Triethyl phosphate Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.481 | 83.21 ng/ul | 17740400 | Naphthalene-d8 | 10.793 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Triethyl phosphate | 182 | C6H15O4P | 000078-40-0 | 94 |
| 2 | Triethyl phosphate | 182 | C6H15O4P | 000078-40-0 | 94 |
| 3 | Triethyl phosphate | 182 | C6H15O4P | 000078-40-0 | 93 |
| 4 | Triethyl phosphate | 182 | C6H15O4P | 000078-40-0 | 83 |
| 5 | Butyl diethyl phosphate | 210 | C8H19O4P | 1000298-58-0 | 42 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017072.D
Acq On : 10 Sep 2023 03:15
Operator : MA/JU
Sample : 04227-06
Misc :
ALS Vial : 84 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BBKE5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

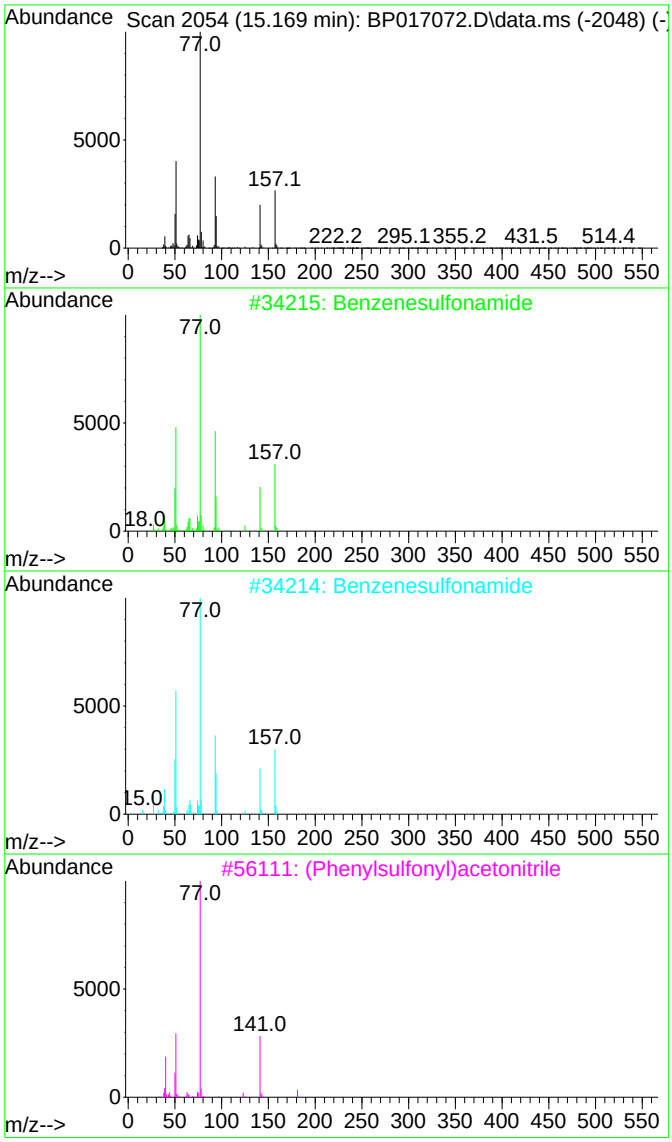
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

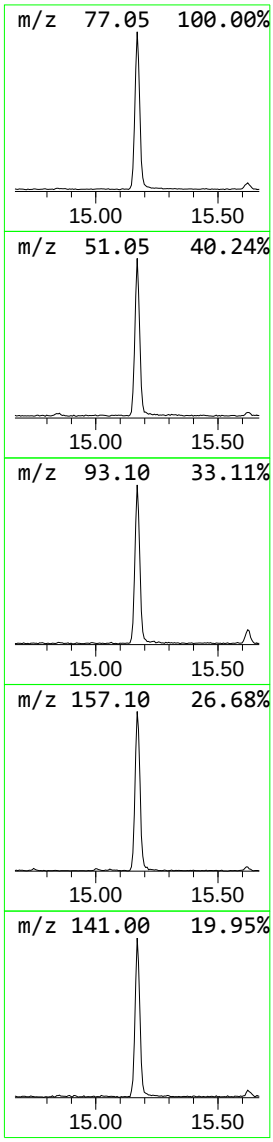
**********

Peak Number 5 Benzenesulfonamide Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.169 | 2.50 ng/ul | 685240 | Acenaphthene-d10 | 14.628 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzenesulfonamide | 157 | C6H7NO2S | 000098-10-2 | 97 |
| 2 | Benzenesulfonamide | 157 | C6H7NO2S | 000098-10-2 | 70 |
| 3 | (Phenylsulfonyl)acetonitrile | 181 | C8H7NO2S | 007605-28-9 | 40 |
| 4 | Ethanone, 2-amino-1-phenyl- | 135 | C8H9NO | 000613-89-8 | 38 |
| 5 | Benzenesulfonyl chloride | 176 | C6H5ClO2S | 000098-09-9 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017072.D
Acq On : 10 Sep 2023 03:15
Operator : MA/JU
Sample : 04227-06
Misc :
ALS Vial : 84 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BBKE5

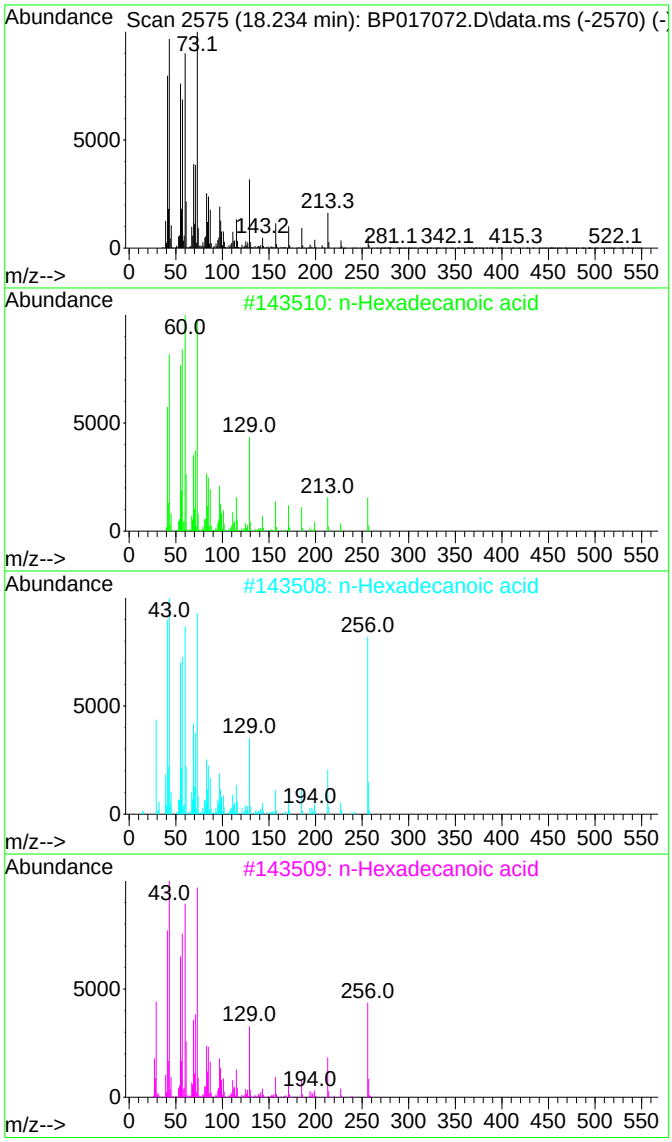
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

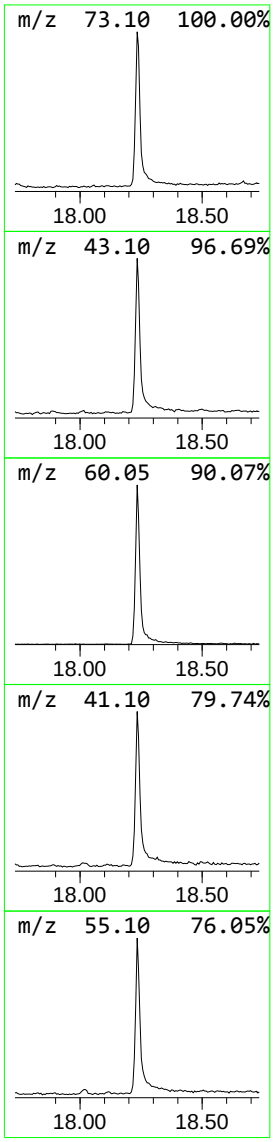
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 6 n-Hexadecanoic acid Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.234 | 4.07 ng/ul | 1321690 | Phenanthrene-d10 | 17.392 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 99 |
| 2 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98 |
| 3 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98 |
| 4 | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 95 |
| 5 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 93 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017072.D
Acq On : 10 Sep 2023 03:15
Operator : MA/JU
Sample : 04227-06
Misc :
ALS Vial : 84 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BBKE5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

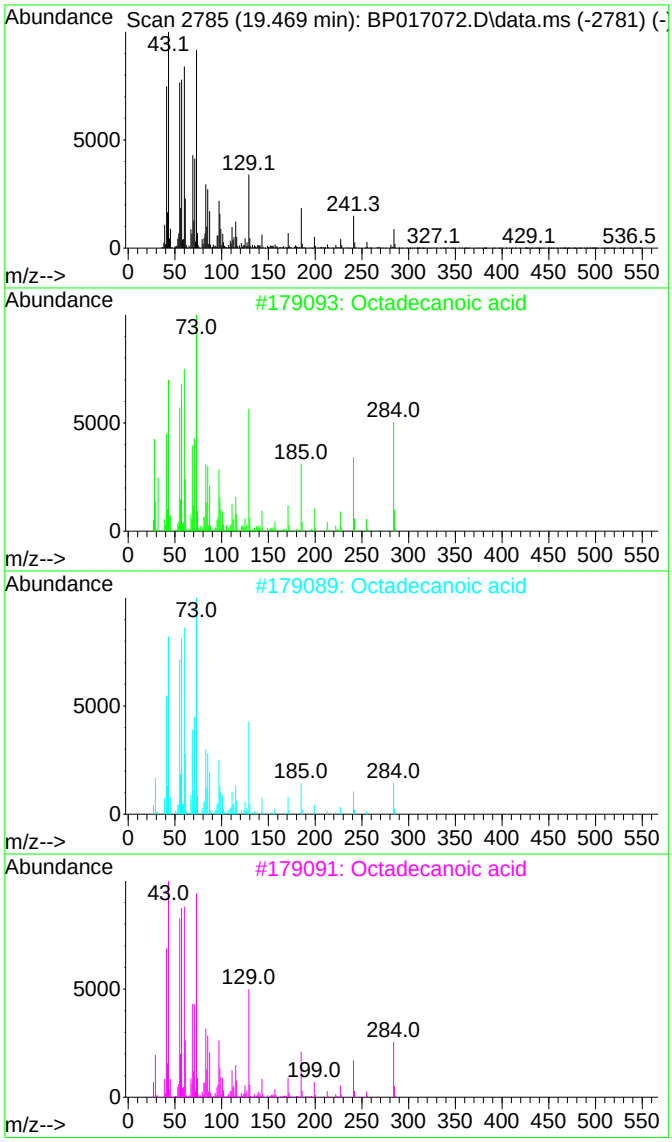
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

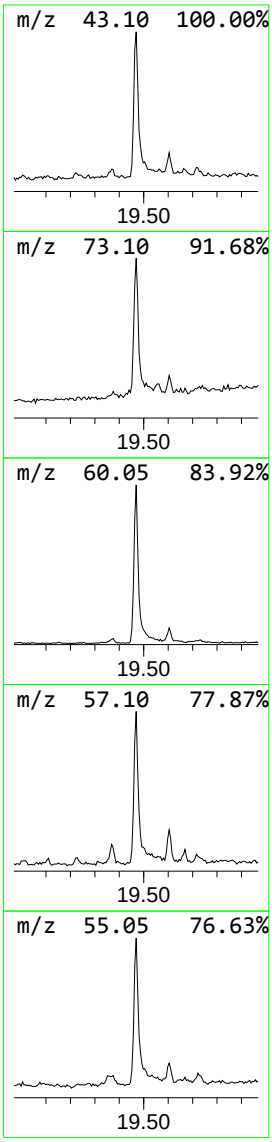
**********

Peak Number 7 Octadecanoic acid Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.469 | 2.22 ng/ul | 643830 | Chrysene-d12 | 21.486 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 99 |
| 2 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 98 |
| 3 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 97 |
| 4 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 95 |
| 5 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 93 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017072.D
Acq On : 10 Sep 2023 03:15
Operator : MA/JU
Sample : 04227-06
Misc :
ALS Vial : 84 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BBKE5

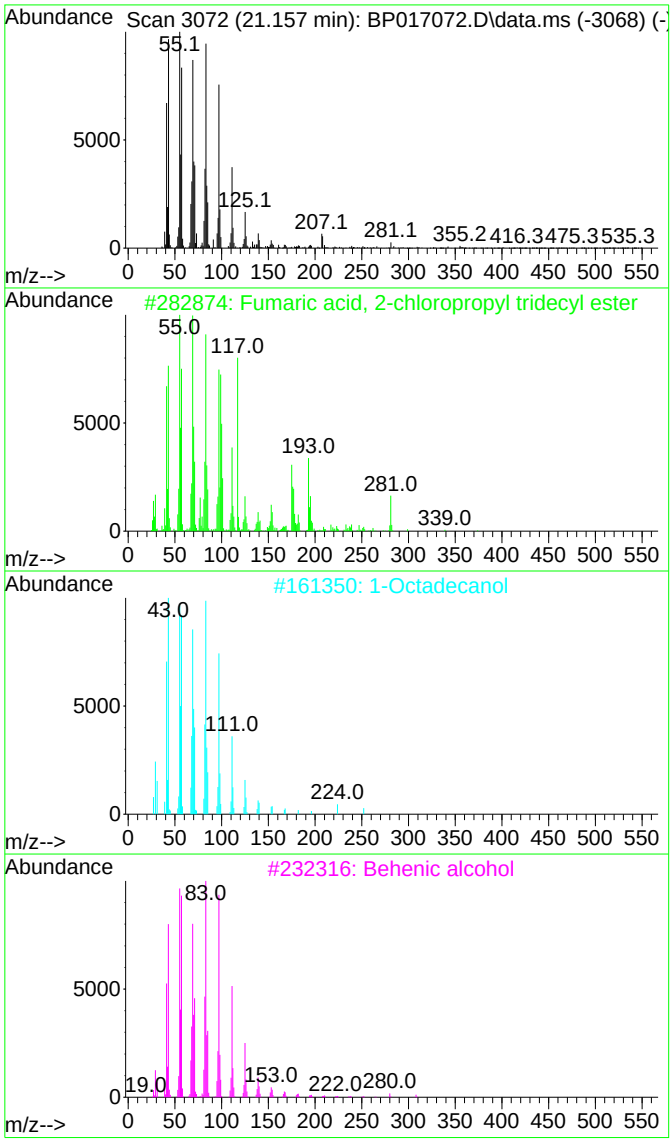
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

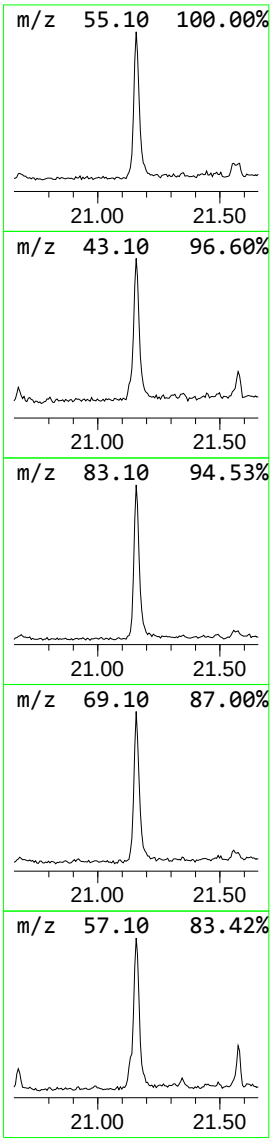
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 8 Fumaric acid, 2-chloropropy... Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.157 | 2.20 ng/ul | 637552 | Chrysene-d12 | 21.486 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Fumaric acid, 2-chloropropyl tri... | 374 | C20H35ClO4 | 1010348-57-1 | 98 |
| 2 | 1-Octadecanol | 270 | C18H38O | 000112-92-5 | 95 |
| 3 | Behenic alcohol | 326 | C22H46O | 000661-19-8 | 94 |
| 4 | n-Heptadecanol-1 | 256 | C17H36O | 001454-85-9 | 94 |
| 5 | n-Nonadecanol-1 | 284 | C19H40O | 001454-84-8 | 94 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017072.D
Acq On : 10 Sep 2023 03:15
Operator : MA/JU
Sample : 04227-06
Misc :
ALS Vial : 84 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BBKE5

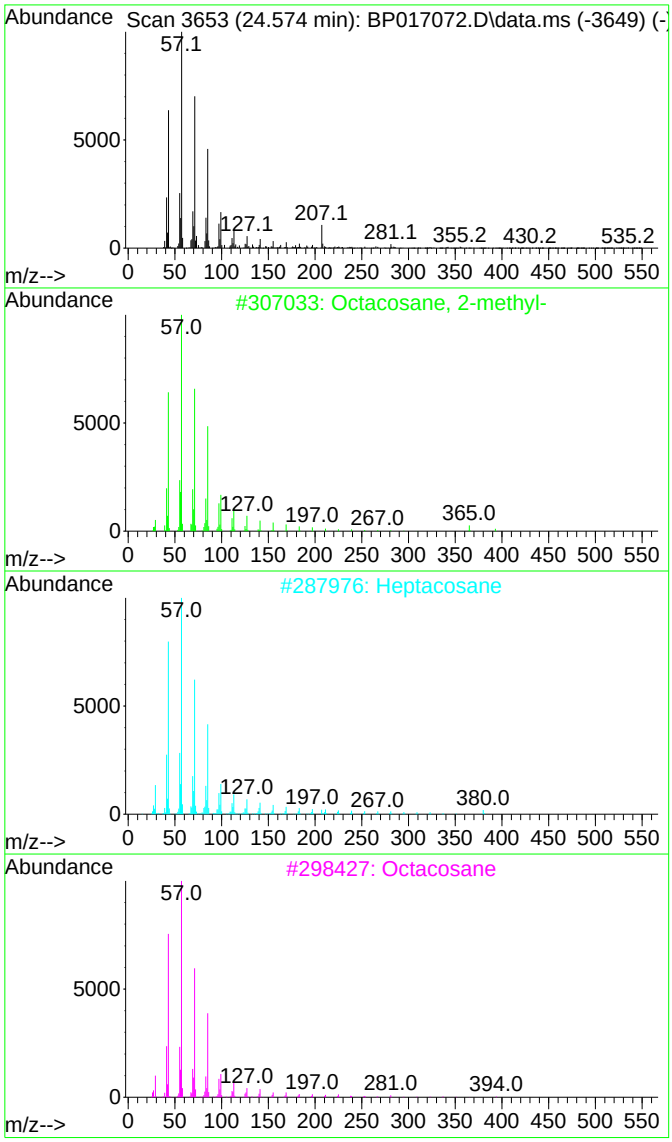
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

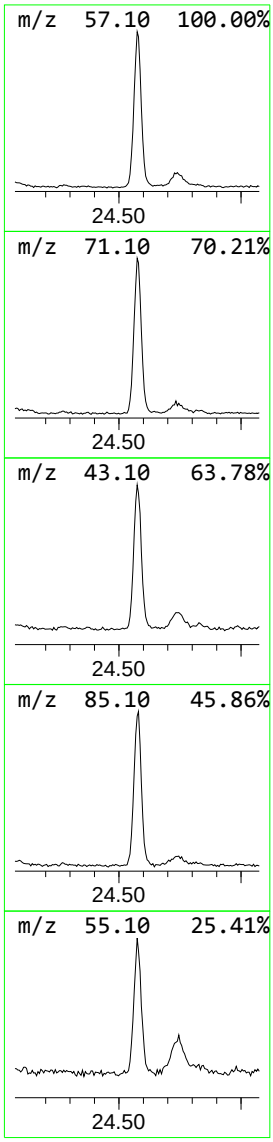
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 24.575 | 2.82 ng/ul | 745623 | Perylene-d12 | 24.022 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octacosane, 2-methyl- | 408 | C29H60 | 001560-98-1 | 90 |
| 2 | Heptacosane | 380 | C27H56 | 000593-49-7 | 90 |
| 3 | Octacosane | 394 | C28H58 | 000630-02-4 | 90 |
| 4 | Tetratetracontane | 619 | C44H90 | 007098-22-8 | 90 |
| 5 | Triacontane | 422 | C30H62 | 000638-68-6 | 90 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017072.D
Acq On : 10 Sep 2023 03:15
Operator : MA/JU
Sample : 04227-06
Misc :
ALS Vial : 84 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
BBKE5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

| | | | | | --Internal Standard--- | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| Tetrachloroethy... | 4.593 | 420.7 | ng/ul | 60577600 | 1 | 7.969 | 2879640 | 20.0 |
| Ethane, 1,1'-ox... | 6.652 | 2.4 | ng/ul | 351834 | 1 | 7.969 | 2879640 | 20.0 |
| Diethyl carbitol | 8.758 | 2.3 | ng/ul | 323484 | 1 | 7.969 | 2879640 | 20.0 |
| Triethyl phosphate | 9.481 | 83.2 | ng/ul | 17740400 | 2 | 10.793 | 4263950 | 20.0 |
| Benzenesulfonamide | 15.169 | 2.5 | ng/ul | 685240 | 3 | 14.628 | 5479970 | 20.0 |
| n-Hexadecanoic ... | 18.234 | 4.1 | ng/ul | 1321690 | 4 | 17.392 | 6489390 | 20.0 |
| Octadecanoic acid | 19.469 | 2.2 | ng/ul | 643830 | 5 | 21.486 | 5799450 | 20.0 |
| Fumaric acid, 2... | 21.157 | 2.2 | ng/ul | 637552 | 5 | 21.486 | 5799450 | 20.0 |
| (DEL) Alkane: S... | 24.575 | 2.8 | ng/ul | 745623 | 6 | 24.022 | 5285760 | 20.0 |