LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :** BNA_P
**ClientSampleId :** CB9A6

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-BP032823.M
Title : SVOA CALIBRATION

Signal : TIC: BP014540.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 3.417 | 35 | 39 | 55 | rVB | 1400646 | 1764656 | 25.29% | 2.880% |
| 2 | 3.823 | 106 | 108 | 118 | rBV | 206420 | 337038 | 4.83% | 0.550% |
| 3 | 4.211 | 169 | 174 | 178 | rBV | 163315 | 201035 | 2.88% | 0.328% |
| 4 | 4.258 | 178 | 182 | 188 | rVB | 36368 | 52082 | 0.75% | 0.085% |
| 5 | 4.417 | 206 | 209 | 220 | rBV5 | 14489 | 39414 | 0.56% | 0.064% |
| 6 | 4.593 | 232 | 239 | 248 | rVB | 297284 | 448886 | 6.43% | 0.733% |
| 7 | 5.717 | 426 | 430 | 439 | rBV | 228158 | 345609 | 4.95% | 0.564% |
| 8 | 6.269 | 512 | 524 | 532 | rBV | 62663 | 114940 | 1.65% | 0.188% |
| 9 | 6.587 | 573 | 578 | 583 | rBV2 | 23409 | 40512 | 0.58% | 0.066% |
| 10 | 6.952 | 635 | 640 | 648 | rVB | 92348 | 150524 | 2.16% | 0.246% |
| 11 | 7.181 | 671 | 679 | 689 | rVV | 169665 | 365576 | 5.24% | 0.597% |
| 12 | 7.328 | 698 | 704 | 714 | rBV | 337030 | 538475 | 7.72% | 0.879% |
| 13 | 7.458 | 720 | 726 | 732 | rVV2 | 100768 | 180467 | 2.59% | 0.294% |
| 14 | 7.534 | 732 | 739 | 747 | rVV | 373775 | 658262 | 9.43% | 1.074% |
| 15 | 7.599 | 747 | 750 | 756 | rVV4 | 19299 | 40700 | 0.58% | 0.066% |
| 16 | 7.740 | 766 | 774 | 779 | rBV2 | 26097 | 43121 | 0.62% | 0.070% |
| 17 | 7.958 | 806 | 811 | 813 | rBV3 | 37153 | 50883 | 0.73% | 0.083% |
| 18 | 7.999 | 813 | 818 | 825 | rVV | 486297 | 774195 | 11.09% | 1.263% |
| 19 | 8.087 | 825 | 833 | 842 | rVB2 | 87302 | 190880 | 2.74% | 0.311% |
| 20 | 8.252 | 853 | 861 | 864 | rBV5 | 14974 | 36729 | 0.53% | 0.060% |
| 21 | 8.663 | 923 | 931 | 934 | rBV4 | 22066 | 43793 | 0.63% | 0.071% |
| 22 | 8.728 | 936 | 942 | 950 | rVV | 470567 | 854793 | 12.25% | 1.395% |
| 23 | 8.799 | 950 | 954 | 966 | rVB9 | 39828 | 97635 | 1.40% | 0.159% |
| 24 | 8.928 | 966 | 976 | 981 | rBV4 | 49305 | 109619 | 1.57% | 0.179% |
| 25 | 9.005 | 981 | 989 | 1002 | rBV | 4218330 | 6978101 | 100.00% | 11.387% |
| 26 | 9.116 | 1003 | 1008 | 1013 | rVV | 129440 | 224968 | 3.22% | 0.367% |
| 27 | 9.181 | 1013 | 1019 | 1027 | rVB | 449941 | 747016 | 10.71% | 1.219% |
| 28 | 9.381 | 1048 | 1053 | 1059 | rVB5 | 51223 | 95733 | 1.37% | 0.156% |
| 29 | 9.528 | 1071 | 1078 | 1084 | rBV4 | 94868 | 194868 | 2.79% | 0.318% |
| 30 | 9.652 | 1092 | 1099 | 1107 | rBV | 167191 | 302569 | 4.34% | 0.494% |
| 31 | 9.728 | 1107 | 1112 | 1115 | rBV2 | 49628 | 80507 | 1.15% | 0.131% |
| 32 | 9.822 | 1125 | 1128 | 1136 | rVB5 | 39663 | 66353 | 0.95% | 0.108% |
| 33 | 9.905 | 1136 | 1142 | 1154 | rBV2 | 359180 | 705806 | 10.11% | 1.152% |
| 34 | 10.028 | 1154 | 1163 | 1169 | rBV3 | 107170 | 289403 | 4.15% | 0.472% |
| 35 | 10.087 | 1169 | 1173 | 1184 | rVB | 161759 | 314740 | 4.51% | 0.514% |
| 36 | 10.234 | 1192 | 1198 | 1206 | rVB7 | 43896 | 94735 | 1.36% | 0.155% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
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Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

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-BP032823.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 10.363 | 1212 | 1220 | 1228 | rVB3 | 225006 | 481657 | 6.90% | 0.786% |
| 38 | 10.458 | 1229 | 1236 | 1254 | rBV | 532562 | 1258016 | 18.03% | 2.053% |
| 39 | 10.599 | 1255 | 1260 | 1269 | rBV5 | 77839 | 182632 | 2.62% | 0.298% |
| 40 | 10.769 | 1284 | 1289 | 1292 | rBV2 | 116117 | 187865 | 2.69% | 0.307% |
| 41 | 10.822 | 1292 | 1298 | 1304 | rVV | 704893 | 1185838 | 16.99% | 1.935% |
| 42 | 10.875 | 1304 | 1307 | 1312 | rVB2 | 77219 | 108918 | 1.56% | 0.178% |
| 43 | 10.975 | 1315 | 1324 | 1336 | rBV2 | 498612 | 1131327 | 16.21% | 1.846% |
| 44 | 11.087 | 1337 | 1343 | 1353 | rBV8 | 73702 | 194790 | 2.79% | 0.318% |
| 45 | 11.257 | 1366 | 1372 | 1375 | rBV4 | 65238 | 122827 | 1.76% | 0.200% |
| 46 | 11.346 | 1384 | 1387 | 1392 | rBV8 | 20176 | 40758 | 0.58% | 0.067% |
| 47 | 11.410 | 1392 | 1398 | 1402 | rBV9 | 46053 | 117492 | 1.68% | 0.192% |
| 48 | 11.493 | 1408 | 1412 | 1423 | rVB8 | 57301 | 159471 | 2.29% | 0.260% |
| 49 | 11.605 | 1426 | 1431 | 1433 | rBV3 | 46638 | 85731 | 1.23% | 0.140% |
| 50 | 11.652 | 1433 | 1439 | 1442 | rVV6 | 80507 | 207256 | 2.97% | 0.338% |
| 51 | 11.687 | 1442 | 1445 | 1452 | rVB5 | 102843 | 183816 | 2.63% | 0.300% |
| 52 | 11.757 | 1453 | 1457 | 1460 | rVV4 | 36994 | 61055 | 0.87% | 0.100% |
| 53 | 11.804 | 1461 | 1465 | 1472 | rVV6 | 97061 | 193440 | 2.77% | 0.316% |
| 54 | 11.875 | 1473 | 1477 | 1480 | rVV3 | 57319 | 80316 | 1.15% | 0.131% |
| 55 | 11.999 | 1494 | 1498 | 1500 | rVV5 | 49767 | 84478 | 1.21% | 0.138% |
| 56 | 12.028 | 1500 | 1503 | 1513 | rVB5 | 69375 | 143808 | 2.06% | 0.235% |
| 57 | 12.169 | 1519 | 1527 | 1533 | rBV | 168326 | 336028 | 4.82% | 0.548% |
| 58 | 12.275 | 1541 | 1545 | 1549 | rBV2 | 63783 | 86069 | 1.23% | 0.140% |
| 59 | 12.340 | 1551 | 1556 | 1560 | rBV2 | 88710 | 174919 | 2.51% | 0.285% |
| 60 | 12.387 | 1561 | 1564 | 1569 | rVB2 | 58306 | 103874 | 1.49% | 0.170% |
| 61 | 12.546 | 1587 | 1591 | 1600 | rVB7 | 62669 | 115113 | 1.65% | 0.188% |
| 62 | 12.722 | 1616 | 1621 | 1628 | rBV9 | 30124 | 83893 | 1.20% | 0.137% |
| 63 | 12.816 | 1632 | 1637 | 1643 | rVB | 185026 | 316109 | 4.53% | 0.516% |
| 64 | 12.916 | 1650 | 1654 | 1659 | rVB4 | 46983 | 75818 | 1.09% | 0.124% |
| 65 | 13.046 | 1672 | 1676 | 1681 | rVB4 | 94061 | 156230 | 2.24% | 0.255% |
| 66 | 13.451 | 1742 | 1745 | 1751 | rVB3 | 56091 | 79611 | 1.14% | 0.130% |
| 67 | 13.710 | 1784 | 1789 | 1795 | rVB3 | 169496 | 270955 | 3.88% | 0.442% |
| 68 | 13.775 | 1797 | 1800 | 1808 | rVV8 | 44159 | 91172 | 1.31% | 0.149% |
| 69 | 14.063 | 1844 | 1849 | 1859 | rVB | 1196554 | 1817533 | 26.05% | 2.966% |
| 70 | 14.346 | 1892 | 1897 | 1905 | rBV | 1188381 | 1867303 | 26.76% | 3.047% |
| 71 | 14.481 | 1914 | 1920 | 1924 | rBV3 | 209527 | 381287 | 5.46% | 0.622% |
| 72 | 14.581 | 1931 | 1937 | 1942 | rBV | 507808 | 753005 | 10.79% | 1.229% |
| 73 | 14.651 | 1944 | 1949 | 1958 | rVB | 1051569 | 1664100 | 23.85% | 2.716% |
| 74 | 14.746 | 1962 | 1965 | 1970 | rVV3 | 47939 | 69891 | 1.00% | 0.114% |
| 75 | 14.934 | 1993 | 1997 | 2001 | rBV4 | 94303 | 155236 | 2.22% | 0.253% |
| 76 | 15.263 | 2049 | 2053 | 2059 | rVB4 | 106551 | 188662 | 2.70% | 0.308% |

LSC Area Percent Report

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Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

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-BP032823.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 77 | 15.398 | 2072 | 2076 | 2078 | rBV | 116084 | 155028 | 2.22% | 0.253% |
| 78 | 15.651 | 2112 | 2119 | 2125 | rBV | 1563080 | 2328994 | 33.38% | 3.801% |
| 79 | 15.781 | 2138 | 2141 | 2145 | rBV | 145017 | 225149 | 3.23% | 0.367% |
| 80 | 15.910 | 2159 | 2163 | 2172 | rBV | 274311 | 476003 | 6.82% | 0.777% |
| 81 | 16.016 | 2178 | 2181 | 2191 | rVB | 240225 | 378996 | 5.43% | 0.618% |
| 82 | 16.116 | 2194 | 2198 | 2202 | rBV4 | 86677 | 128240 | 1.84% | 0.209% |
| 83 | 16.181 | 2203 | 2209 | 2213 | rBV | 1208632 | 1682216 | 24.11% | 2.745% |
| 84 | 16.328 | 2232 | 2234 | 2240 | rVB3 | 79508 | 110134 | 1.58% | 0.180% |
| 85 | 16.692 | 2291 | 2296 | 2302 | rBV | 692391 | 1005401 | 14.41% | 1.641% |
| 86 | 16.957 | 2338 | 2341 | 2344 | rVB | 86346 | 89288 | 1.28% | 0.146% |
| 87 | 17.416 | 2410 | 2419 | 2427 | rVB | 1519448 | 2299651 | 32.96% | 3.753% |
| 88 | 17.516 | 2431 | 2436 | 2442 | rVV | 1812725 | 2544683 | 36.47% | 4.153% |
| 89 | 18.287 | 2563 | 2567 | 2575 | rBV | 982526 | 1330762 | 19.07% | 2.172% |
| 90 | 18.498 | 2600 | 2603 | 2609 | rBV2 | 90592 | 139283 | 2.00% | 0.227% |
| 91 | 19.045 | 2693 | 2696 | 2702 | rVB | 455985 | 549770 | 7.88% | 0.897% |
| 92 | 19.510 | 2772 | 2775 | 2786 | rBV | 238801 | 328436 | 4.71% | 0.536% |
| 93 | 19.745 | 2810 | 2815 | 2820 | rBV | 2484527 | 3128123 | 44.83% | 5.105% |
| 94 | 19.792 | 2820 | 2823 | 2830 | rVB | 504385 | 662314 | 9.49% | 1.081% |
| 95 | 20.045 | 2863 | 2866 | 2874 | rVB | 216316 | 332172 | 4.76% | 0.542% |
| 96 | 21.110 | 3043 | 3047 | 3055 | rBV | 609802 | 1158131 | 16.60% | 1.890% |
| 97 | 21.175 | 3055 | 3058 | 3064 | rVB | 893201 | 1075313 | 15.41% | 1.755% |
| 98 | 21.504 | 3109 | 3114 | 3120 | rBV | 1895638 | 2428232 | 34.80% | 3.962% |
| 99 | 23.851 | 3507 | 3513 | 3528 | rVB2 | 1774464 | 3614082 | 51.79% | 5.898% |
| 100 | 24.016 | 3535 | 3541 | 3550 | rVB2 | 1231476 | 2537258 | 36.36% | 4.140% |

Sum of corrected areas: 61280581

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
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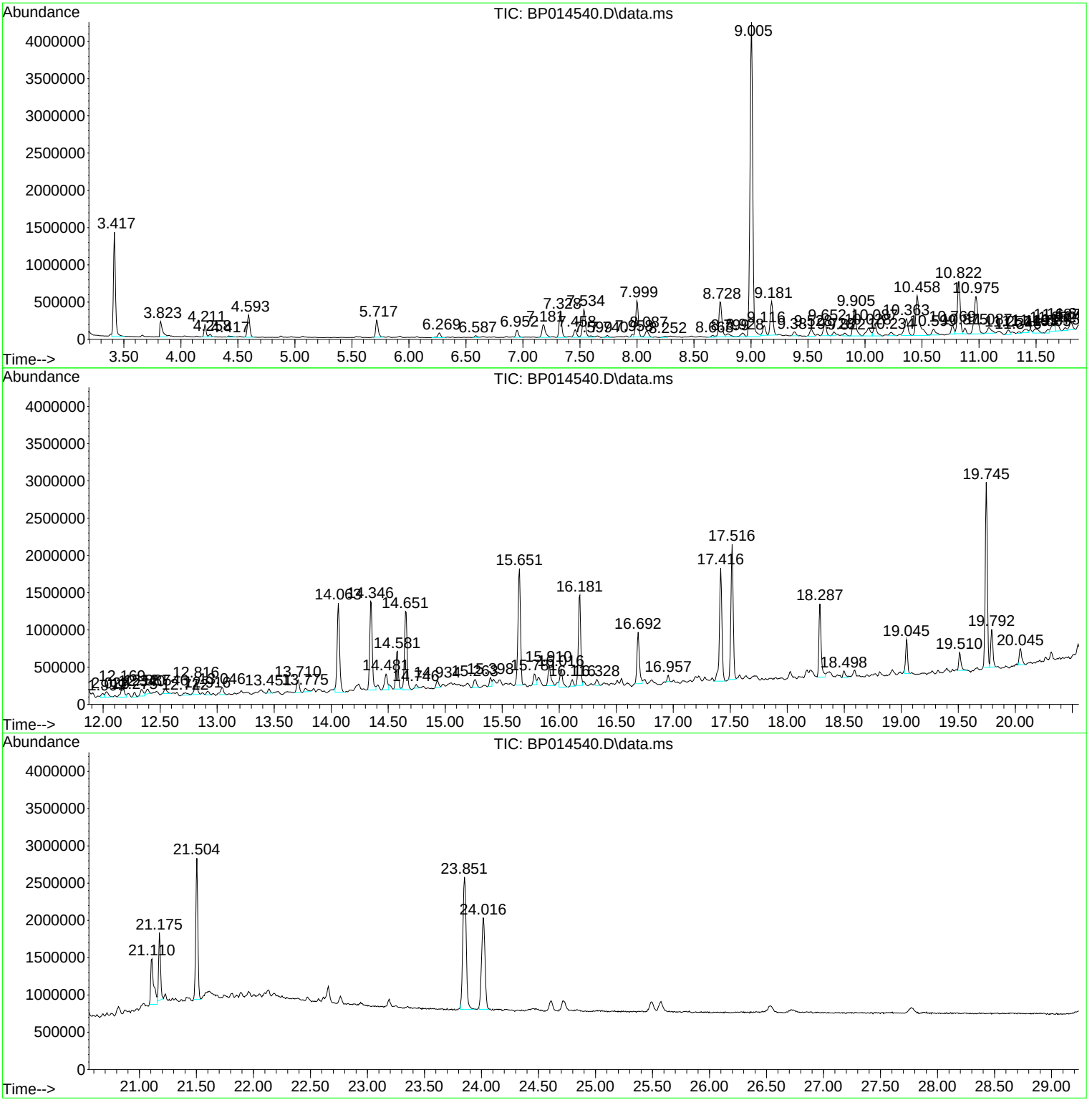
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

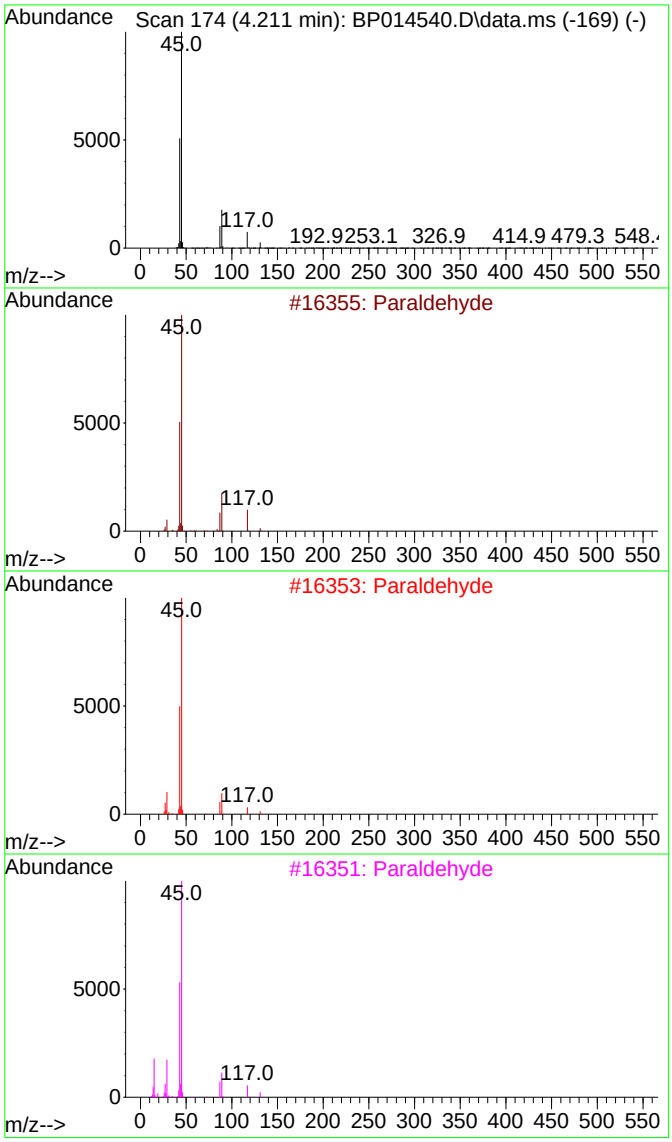
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

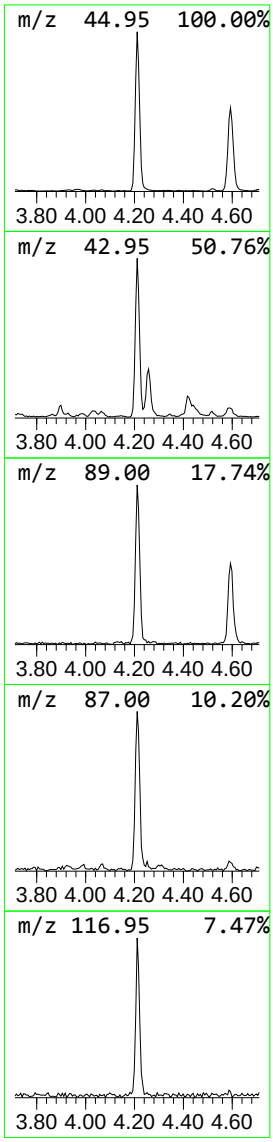
**********

Peak Number 1 Paraldehyde Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.211 | 5.19 ng/ul | 201035 | 1,4-Dichlorobenzene-d4 | 7.999 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Paraldehyde | 132 | C6H12O3 | 000123-63-7 | 86 |
| 2 | Paraldehyde | 132 | C6H12O3 | 000123-63-7 | 83 |
| 3 | Paraldehyde | 132 | C6H12O3 | 000123-63-7 | 83 |
| 4 | Paraldehyde | 132 | C6H12O3 | 000123-63-7 | 64 |
| 5 | Ethanol, 2-[2-(2-butoxyethoxy)et... | 206 | C10H22O4 | 000143-22-6 | 40 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

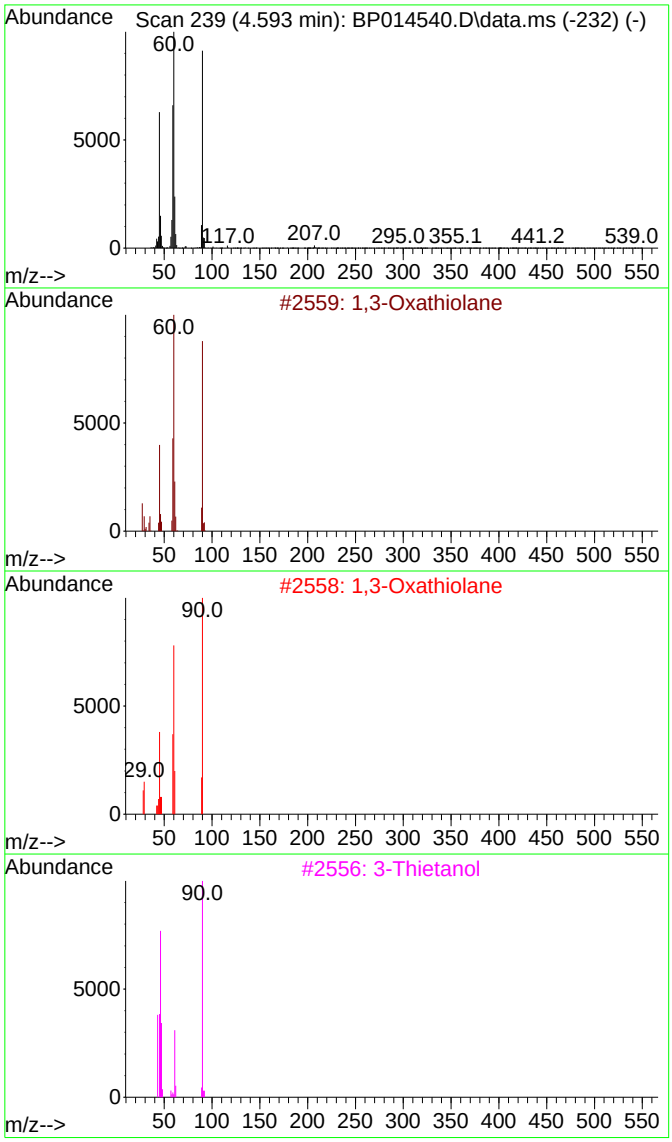
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

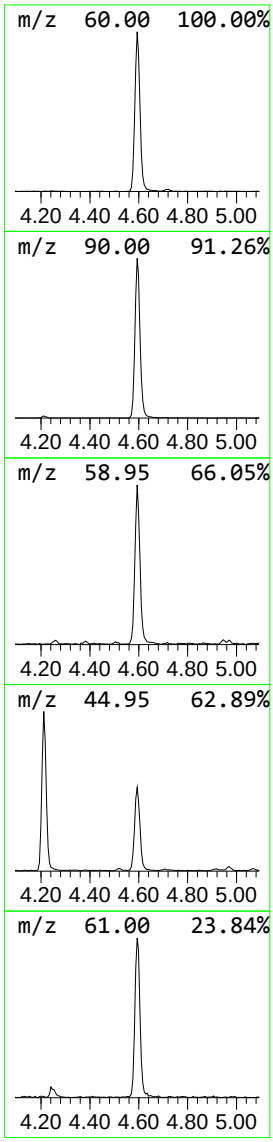
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 2 1,3-Oxathiolane Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.593 | 11.60 ng/ul | 448886 | 1,4-Dichlorobenzene-d4 | 7.999 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,3-Oxathiolane | 90 | C3H6OS | 002094-97-5 | 91 |
| 2 | 1,3-Oxathiolane | 90 | C3H6OS | 002094-97-5 | 64 |
| 3 | 3-Thietanol | 90 | C3H6OS | 010304-16-2 | 53 |
| 4 | Thiourea, methyl- | 90 | C2H6N2S | 000598-52-7 | 50 |
| 5 | Hydrazinecarboxylic acid, methyl... | 90 | C2H6N2O2 | 006294-89-9 | 45 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

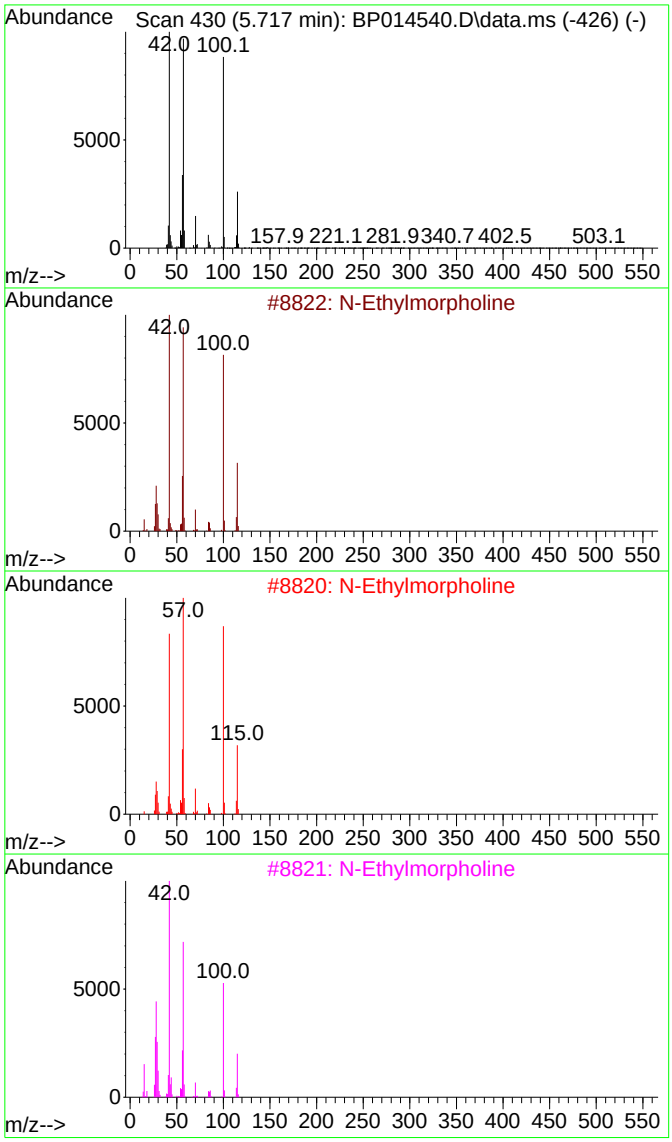
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

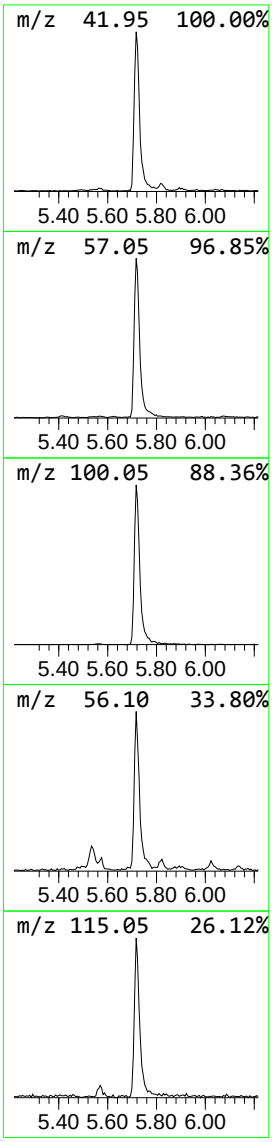
**********

Peak Number 3 N-Ethylmorpholine Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.717 | 8.93 ng/ul | 345609 | 1,4-Dichlorobenzene-d4 | 7.999 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | N-Ethylmorpholine | 115 | C6H13NO | 000100-74-3 | 94 |
| 2 | N-Ethylmorpholine | 115 | C6H13NO | 000100-74-3 | 94 |
| 3 | N-Ethylmorpholine | 115 | C6H13NO | 000100-74-3 | 90 |
| 4 | Sydnone, 3-methyl- | 100 | C3H4N2O2 | 006939-12-4 | 53 |
| 5 | 2H-1,4-Oxazine-4-acetic acid, te... | 145 | C6H11NO3 | 1000362-53-0 | 45 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

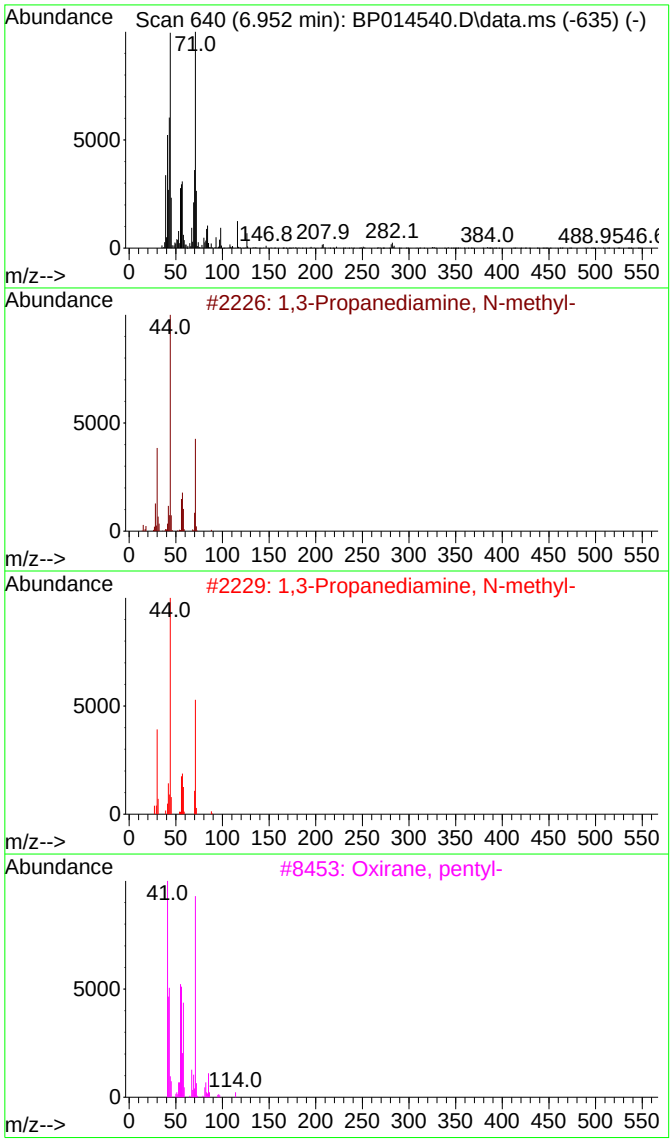
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

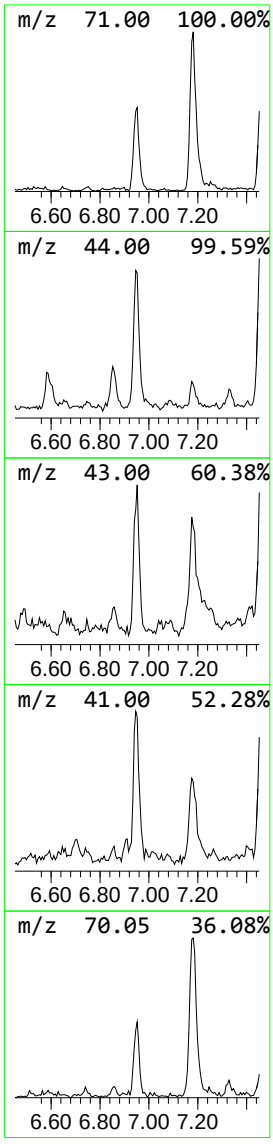
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 5 1,3-Propanediamine, N-methyl- Concentration Rank 23

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.952 | 3.89 ng/ul | 150524 | 1,4-Dichlorobenzene-d4 | 7.999 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,3-Propanediamine, N-methyl- | 88 | C4H12N2 | 006291-84-5 | 53 |
| 2 | 1,3-Propanediamine, N-methyl- | 88 | C4H12N2 | 006291-84-5 | 42 |
| 3 | Oxirane, pentyl- | 114 | C7H14O | 005063-65-0 | 27 |
| 4 | Sulfurous acid, nonyl 2-pentyl e... | 278 | C14H30O3S | 1010309-15-8 | 27 |
| 5 | 1-Butanol, 2,2-dimethyl- | 102 | C6H14O | 001185-33-7 | 25 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

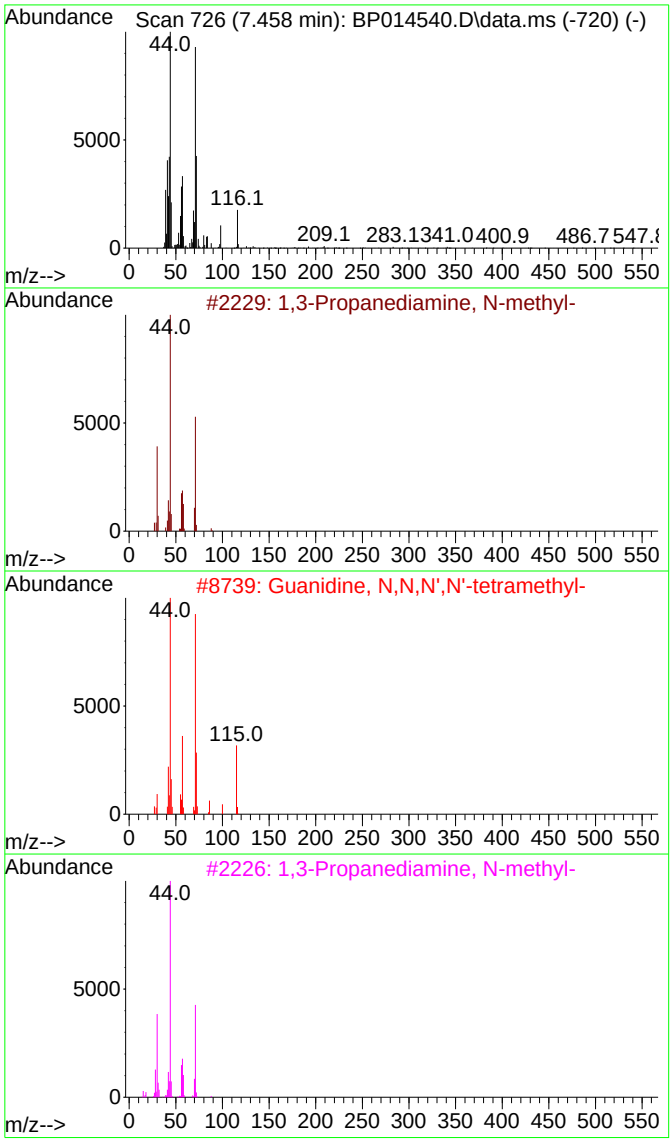
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

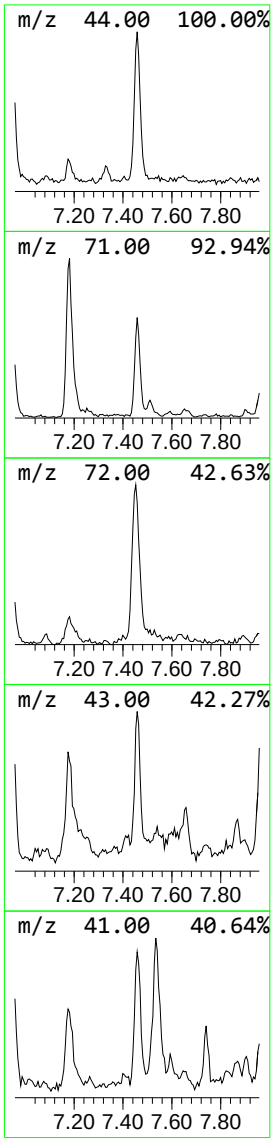
**********

Peak Number 6 unknown-01 Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.458 | 4.66 ng/ul | 180467 | 1,4-Dichlorobenzene-d4 | 7.999 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,3-Propanediamine, N-methyl- | 88 | C4H12N2 | 006291-84-5 | 40 |
| 2 | Guanidine, N,N,N',N'-tetramethyl- | 115 | C5H13N3 | 000080-70-6 | 38 |
| 3 | 1,3-Propanediamine, N-methyl- | 88 | C4H12N2 | 006291-84-5 | 35 |
| 4 | 1,3-Propanediamine, N-methyl- | 88 | C4H12N2 | 006291-84-5 | 33 |
| 5 | Tris(2-aminoethyl)amine | 146 | C6H18N4 | 004097-89-6 | 25 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

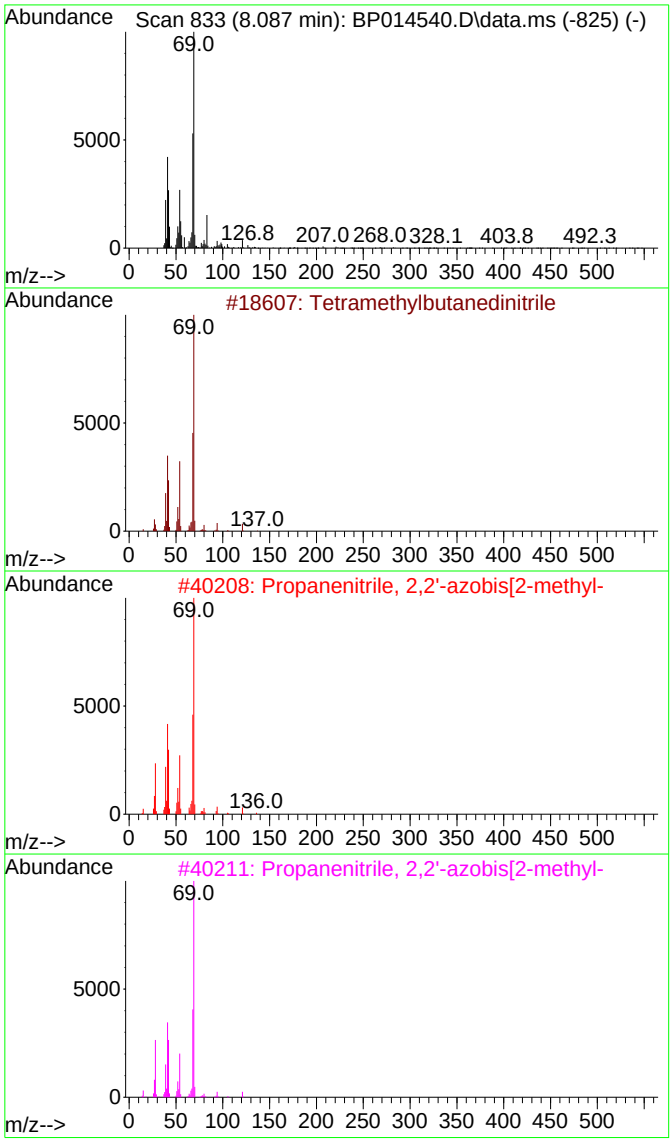
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

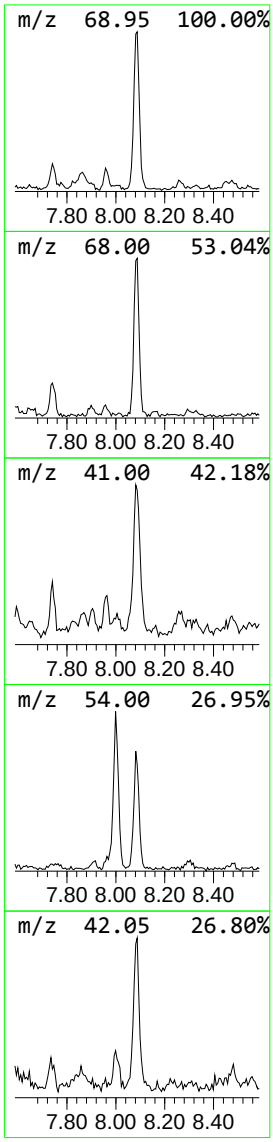
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 7 Tetramethylbutanedinitrile Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.087 | 4.93 ng/ul | 190880 | 1,4-Dichlorobenzene-d4 | 7.999 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Tetramethylbutanedinitrile | 136 | C8H12N2 | 003333-52-6 | 86 |
| 2 | Propanenitrile, 2,2'-azobis[2-me... | 164 | C8H12N4 | 000078-67-1 | 86 |
| 3 | Propanenitrile, 2,2'-azobis[2-me... | 164 | C8H12N4 | 000078-67-1 | 86 |
| 4 | Tetramethylbutanedinitrile | 136 | C8H12N2 | 003333-52-6 | 72 |
| 5 | Propanenitrile, 2,2'-azobis[2-me... | 164 | C8H12N4 | 000078-67-1 | 72 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

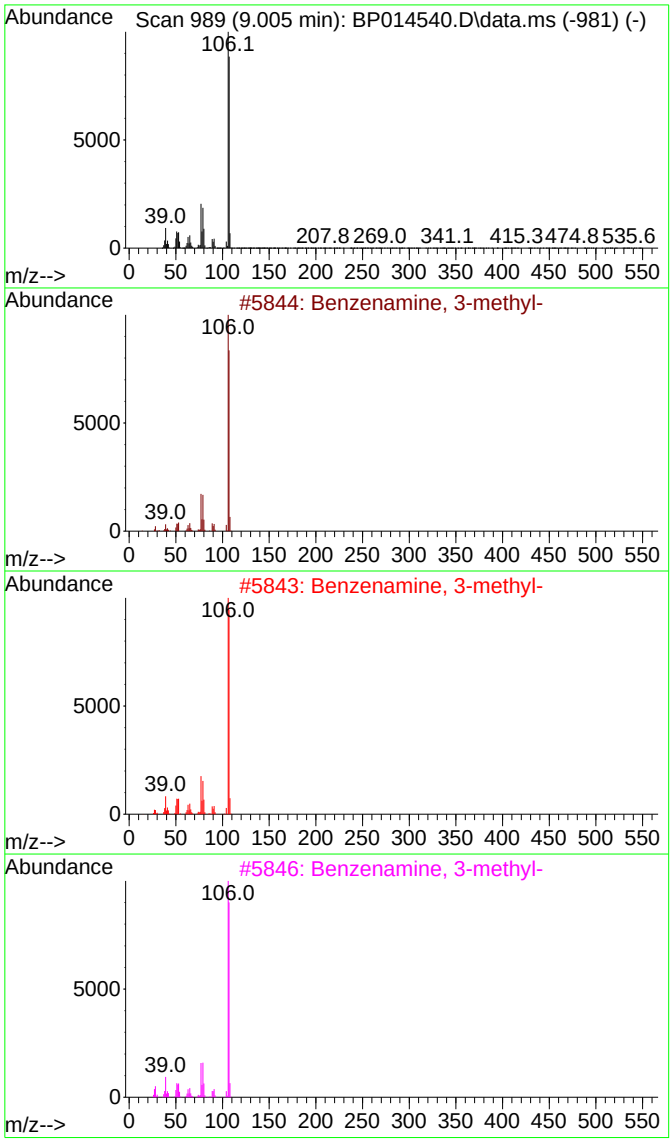
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

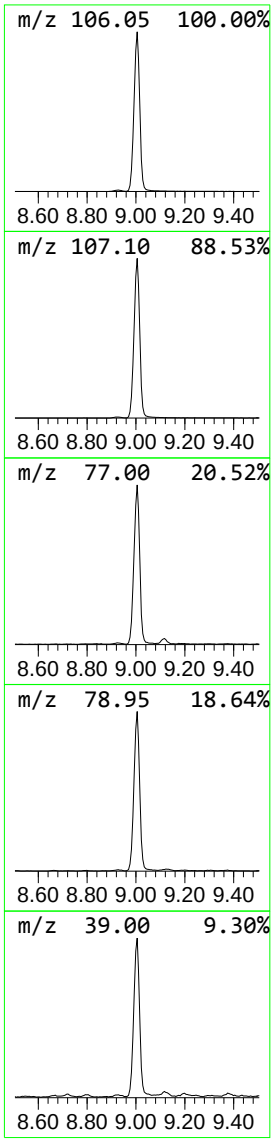
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 10 Benzenamine, 3-methyl- Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.005 | 180.27 ng/ul | 6978100 | 1,4-Dichlorobenzene-d4 | 7.999 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzenamine, 3-methyl- | 107 | C7H9N | 000108-44-1 | 97 |
| 2 | Benzenamine, 3-methyl- | 107 | C7H9N | 000108-44-1 | 97 |
| 3 | Benzenamine, 3-methyl- | 107 | C7H9N | 000108-44-1 | 97 |
| 4 | Benzenamine, 3-methyl- | 107 | C7H9N | 000108-44-1 | 96 |
| 5 | o-Toluidine | 107 | C7H9N | 000095-53-4 | 94 |



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Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

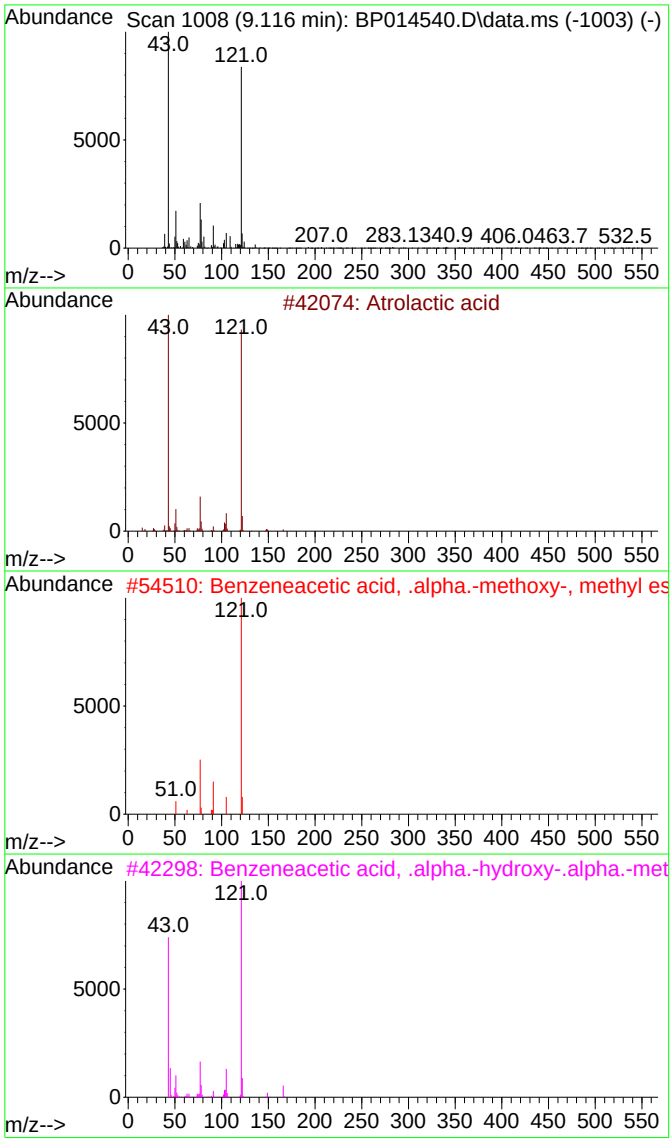
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

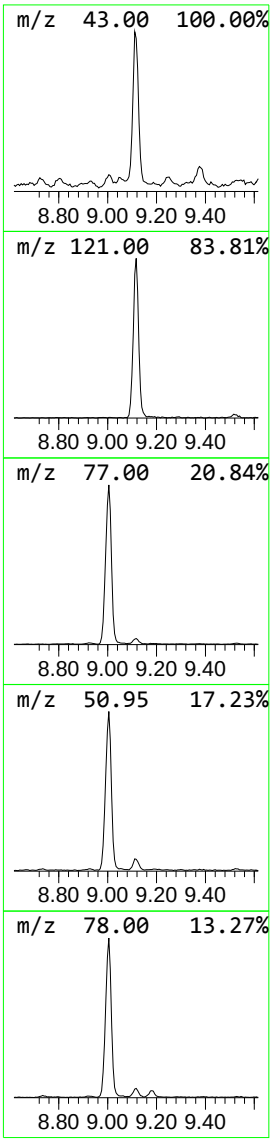
**********

Peak Number 11 Atrolactic acid Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.116 | 5.81 ng/ul | 224968 | 1,4-Dichlorobenzene-d4 | 7.999 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Atrolactic acid | 166 | C9H10O3 | 000515-30-0 | 72 |
| 2 | Benzeneacetic acid, .alpha.-meth... | 180 | C10H12O3 | 056143-21-6 | 72 |
| 3 | Benzeneacetic acid, .alpha.-hydr... | 166 | C9H10O3 | 013113-71-8 | 72 |
| 4 | .alpha.-Ethyl-.alpha.-methylbenz... | 150 | C10H14O | 001565-75-9 | 64 |
| 5 | 3-Hydroxy-3-phenylbutan-2-one | 164 | C10H12O2 | 003155-01-9 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

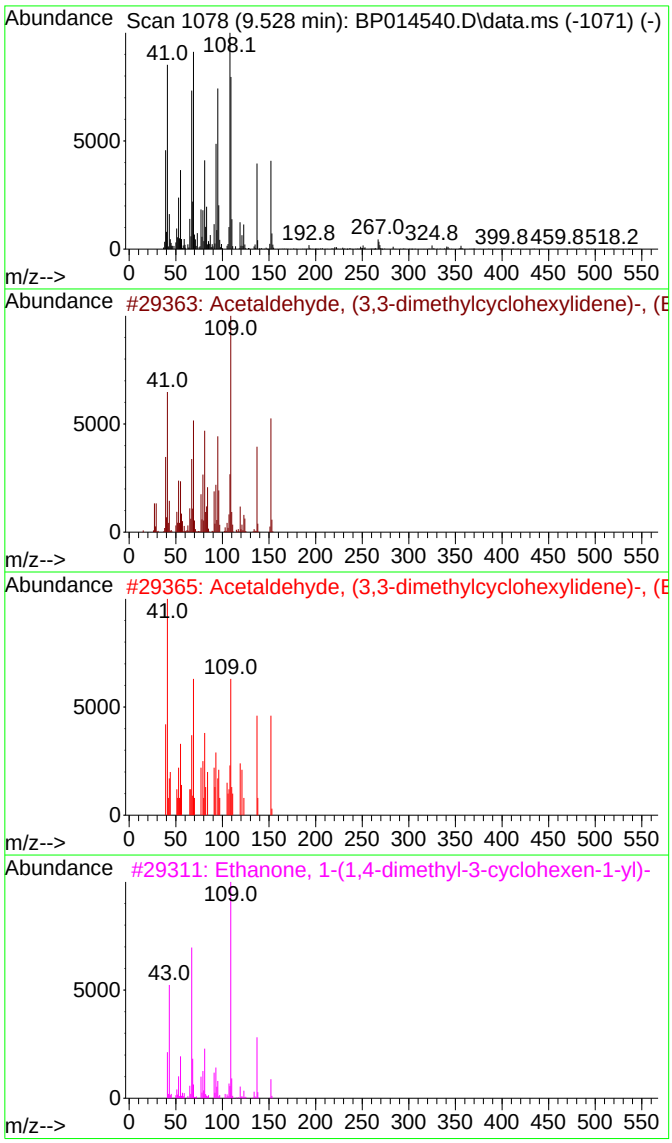
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

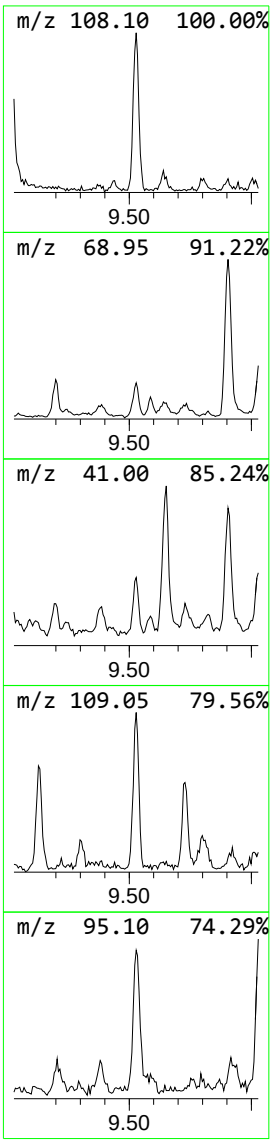
**********

Peak Number 13 unknown-02 Concentration Rank 26

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.528 | 3.29 ng/ul | 194868 | Naphthalene-d8 | 10.822 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Acetaldehyde, (3,3-dimethylcyclo... | 152 | C10H16O | 026532-25-2 | 42 |
| 2 | Acetaldehyde, (3,3-dimethylcyclo... | 152 | C10H16O | 026532-25-2 | 38 |
| 3 | Ethanone, 1-(1,4-dimethyl-3-cycl... | 152 | C10H16O | 043219-68-7 | 27 |
| 4 | 2H-Inden-2-one, octahydro-3a-met... | 152 | C10H16O | 013351-29-6 | 25 |
| 5 | 1(2H)-Naphthalenone, octahydro-,... | 152 | C10H16O | 021370-71-8 | 25 |



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Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

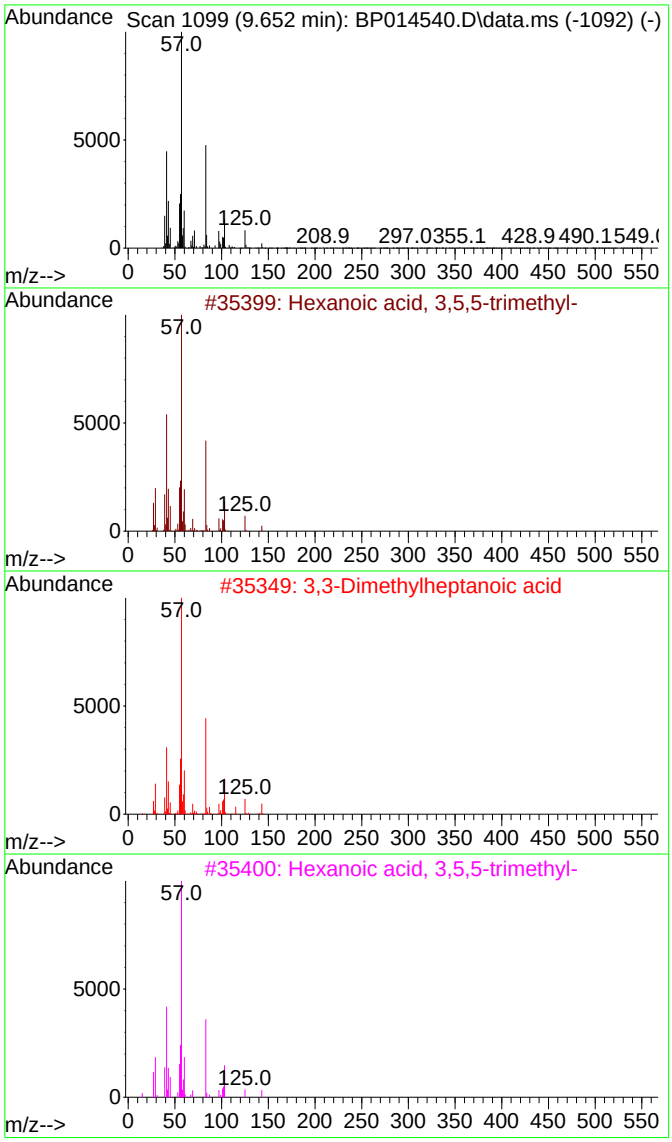
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

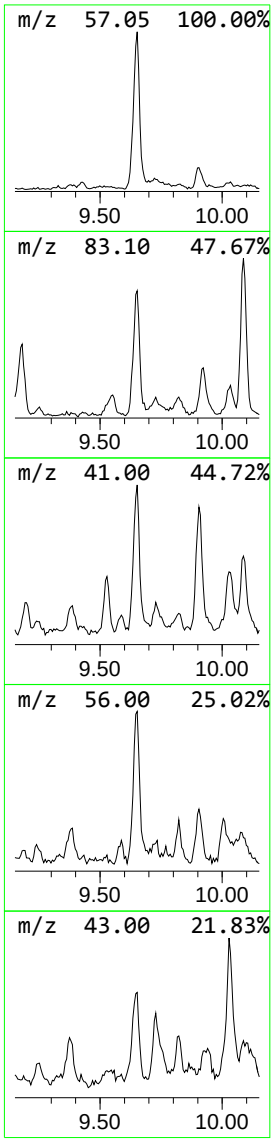
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 14 Hexanoic acid, 3,5,5-trimet... Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.652 | 5.10 ng/ul | 302569 | Naphthalene-d8 | 10.822 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexanoic acid, 3,5,5-trimethyl- | 158 | C9H18O2 | 003302-10-1 | 90 |
| 2 | 3,3-Dimethylheptanoic acid | 158 | C9H18O2 | 067061-30-7 | 83 |
| 3 | Hexanoic acid, 3,5,5-trimethyl- | 158 | C9H18O2 | 003302-10-1 | 83 |
| 4 | Hexanoic acid, 3,5,5-trimethyl- | 158 | C9H18O2 | 003302-10-1 | 83 |
| 5 | Hexanoic acid, 3,5,5-trimethyl-,... | 512 | C30H56O6 | 056554-53-1 | 83 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

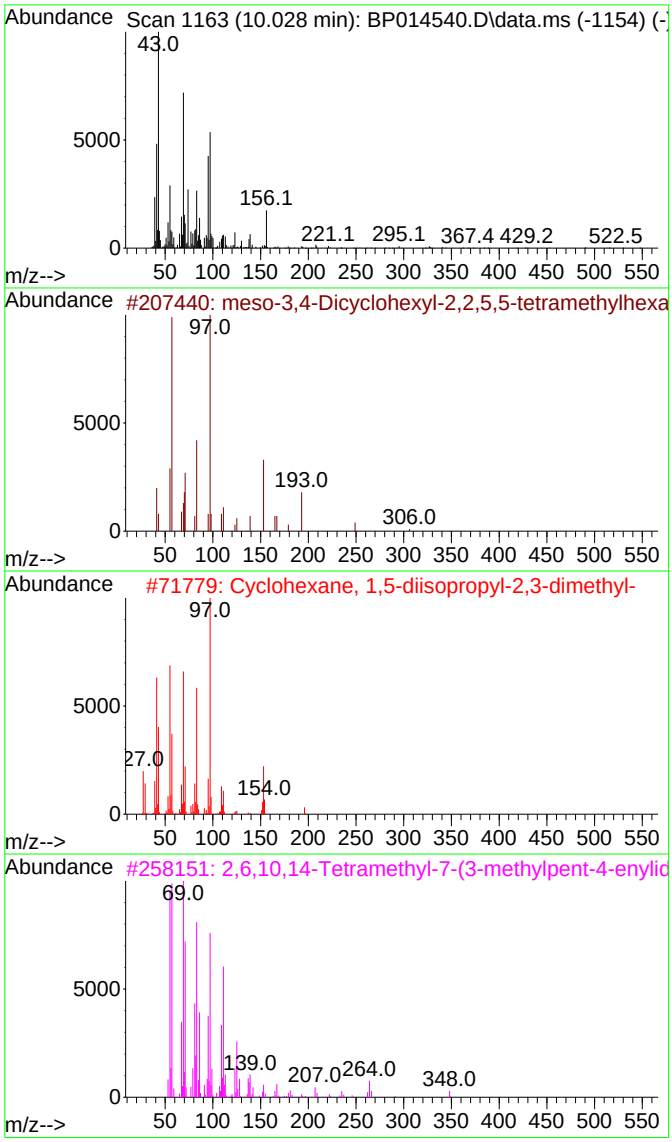
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

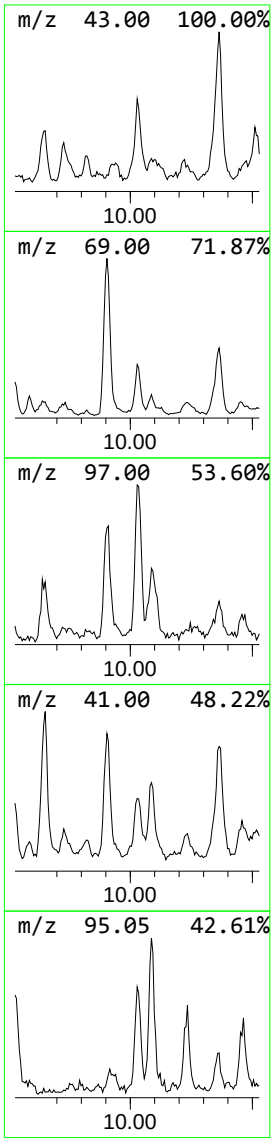
**********

Peak Number 15 unknown-03 Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.028 | 4.88 ng/ul | 289403 | Naphthalene-d8 | 10.822 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | meso-3,4-Dicyclohexyl-2,2,5,5-te... | 306 | C22H42 | 065149-86-2 | 27 |
| 2 | Cyclohexane, 1,5-diisopropyl-2,3... | 196 | C14H28 | 1000149-58-8 | 27 |
| 3 | 2,6,10,14-Tetramethyl-7-(3-methy... | 348 | C25H48 | 1000370-41-6 | 27 |
| 4 | 3-Thiopheneacetic acid, methyl e... | 156 | C7H8O2S | 058414-52-1 | 18 |
| 5 | Oxalic acid, cyclohexylmethyl no... | 312 | C18H32O4 | 1000309-68-6 | 14 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

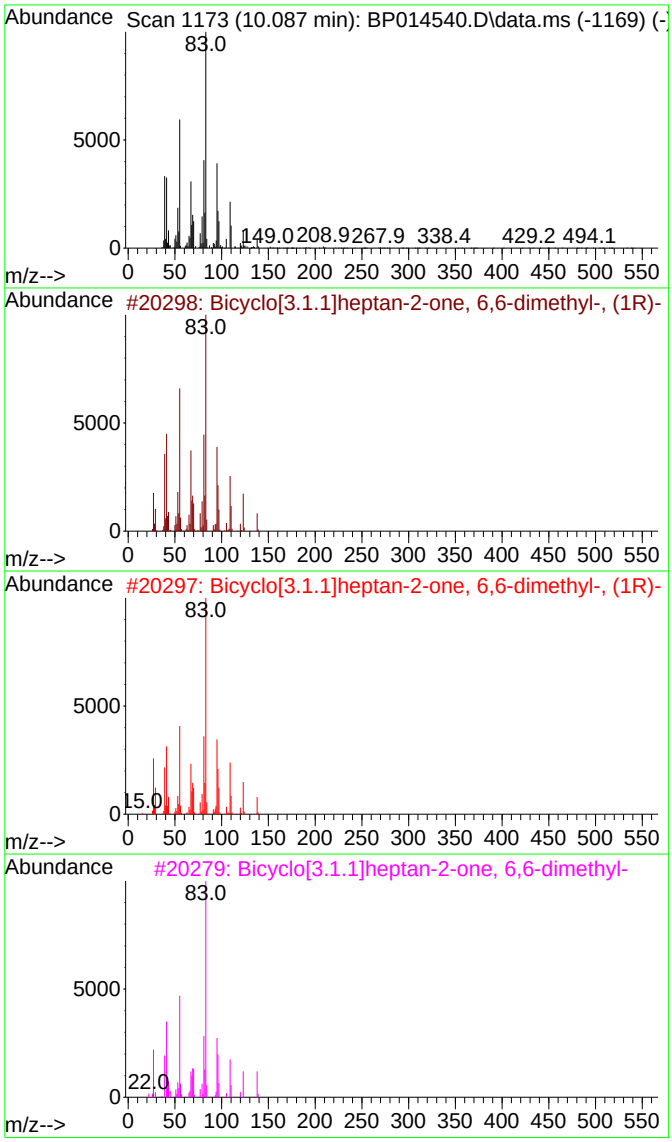
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

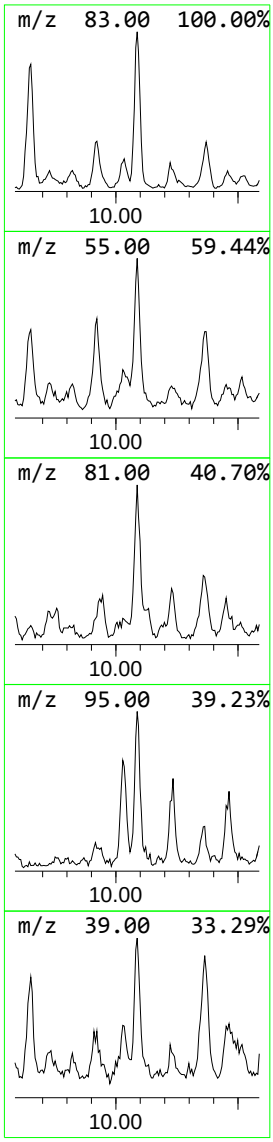
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 16 Bicyclo[3.1.1]heptan-2-one,... Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.087 | 5.31 ng/ul | 314740 | Naphthalene-d8 | 10.822 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Bicyclo[3.1.1]heptan-2-one, 6,6-... | 138 | C9H14O | 038651-65-9 | 91 |
| 2 | Bicyclo[3.1.1]heptan-2-one, 6,6-... | 138 | C9H14O | 038651-65-9 | 78 |
| 3 | Bicyclo[3.1.1]heptan-2-one, 6,6-... | 138 | C9H14O | 024903-95-5 | 53 |
| 4 | Spiro[2.3]hexane-5-carboxylic ac... | 278 | C18H30O2 | 1000152-25-7 | 45 |
| 5 | 1-Cyclohexylethanol, trifluoroac... | 224 | C10H15F3O2 | 1000365-19-2 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

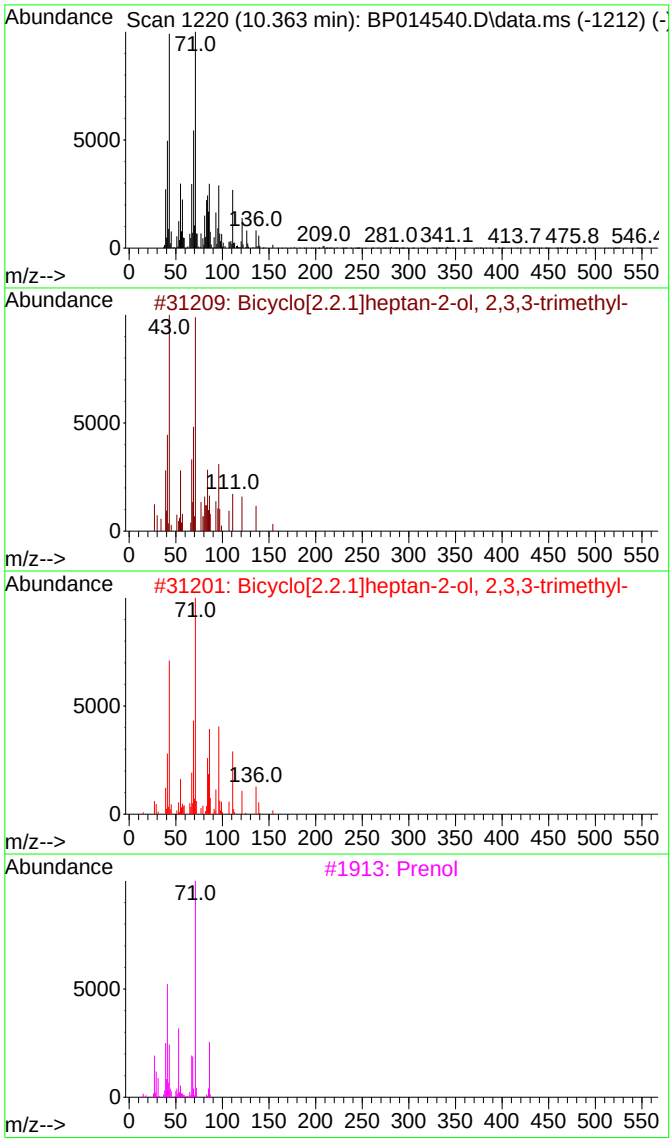
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

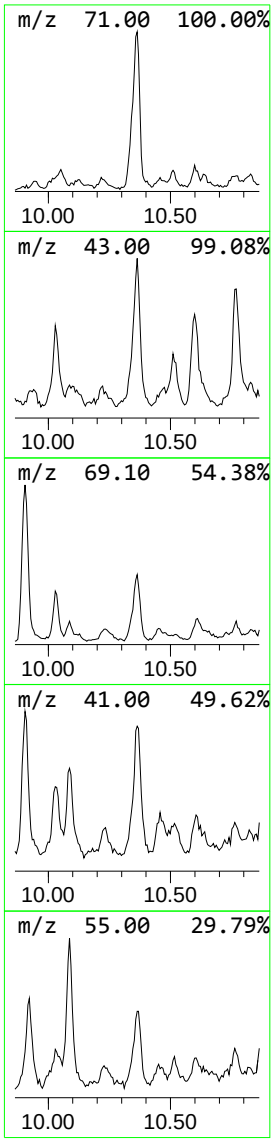
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 17 Bicyclo[2.2.1]heptan-2-ol, ... Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.363 | 8.12 ng/ul | 481657 | Naphthalene-d8 | 10.822 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Bicyclo[2.2.1]heptan-2-ol, 2,3,3... | 154 | C10H18O | 000465-31-6 | 81 |
| 2 | Bicyclo[2.2.1]heptan-2-ol, 2,3,3... | 154 | C10H18O | 000465-31-6 | 46 |
| 3 | Prenol | 86 | C5H10O | 000556-82-1 | 43 |
| 4 | 1-Propene, 3-methoxy-2-methyl- | 86 | C5H10O | 022418-49-1 | 35 |
| 5 | 1,3-Dimethylbutyl butyrate | 172 | C10H20O2 | 005332-88-7 | 35 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

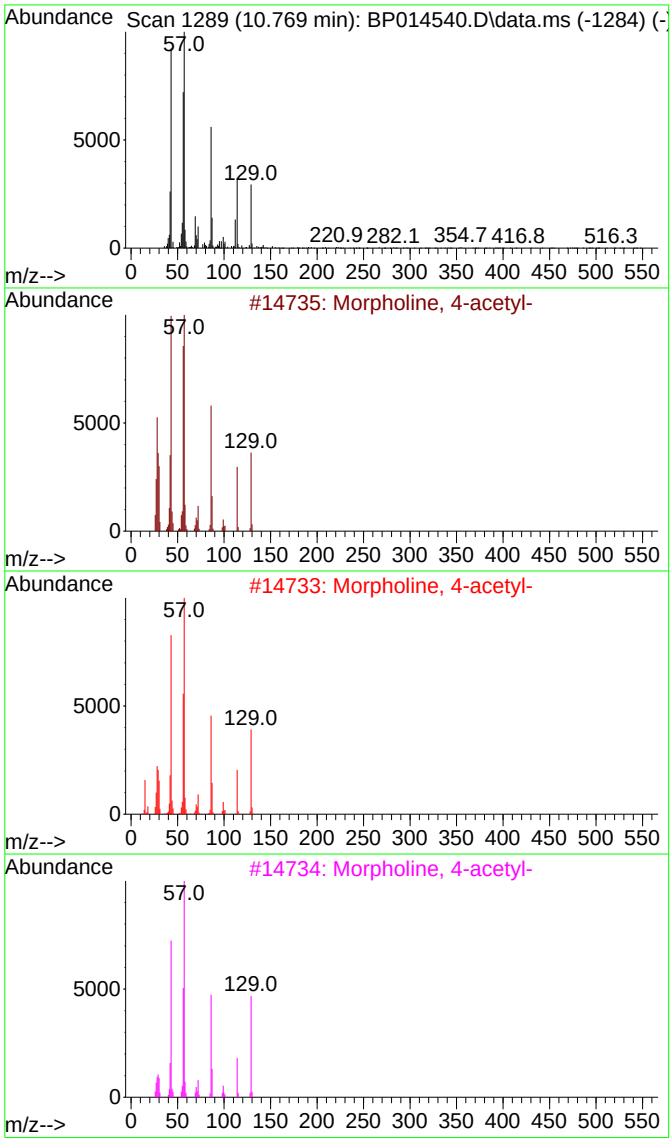
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

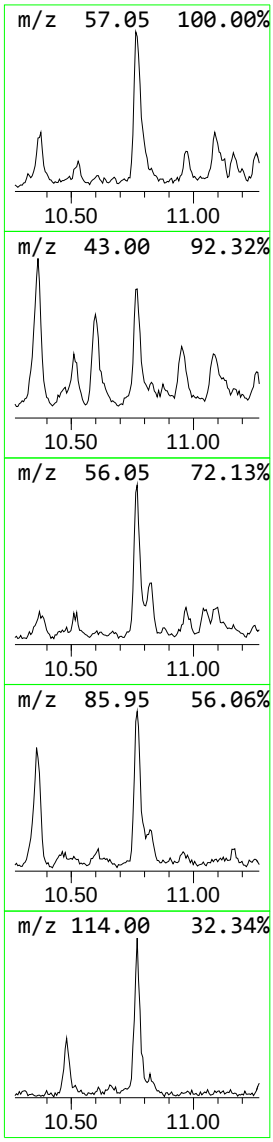
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 19 Morpholine, 4-acetyl- Concentration Rank 30

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.769 | 3.17 ng/ul | 187865 | Naphthalene-d8 | 10.822 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Morpholine, 4-acetyl- | 129 | C6H11NO2 | 001696-20-4 | 94 |
| 2 | Morpholine, 4-acetyl- | 129 | C6H11NO2 | 001696-20-4 | 83 |
| 3 | Morpholine, 4-acetyl- | 129 | C6H11NO2 | 001696-20-4 | 43 |
| 4 | 2-Methylpropionic acid, morpholide | 157 | C8H15NO2 | 1000307-31-7 | 42 |
| 5 | Cyclohexanemethanol, 2-amino-, t... | 129 | C7H15NO | 005691-21-4 | 37 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

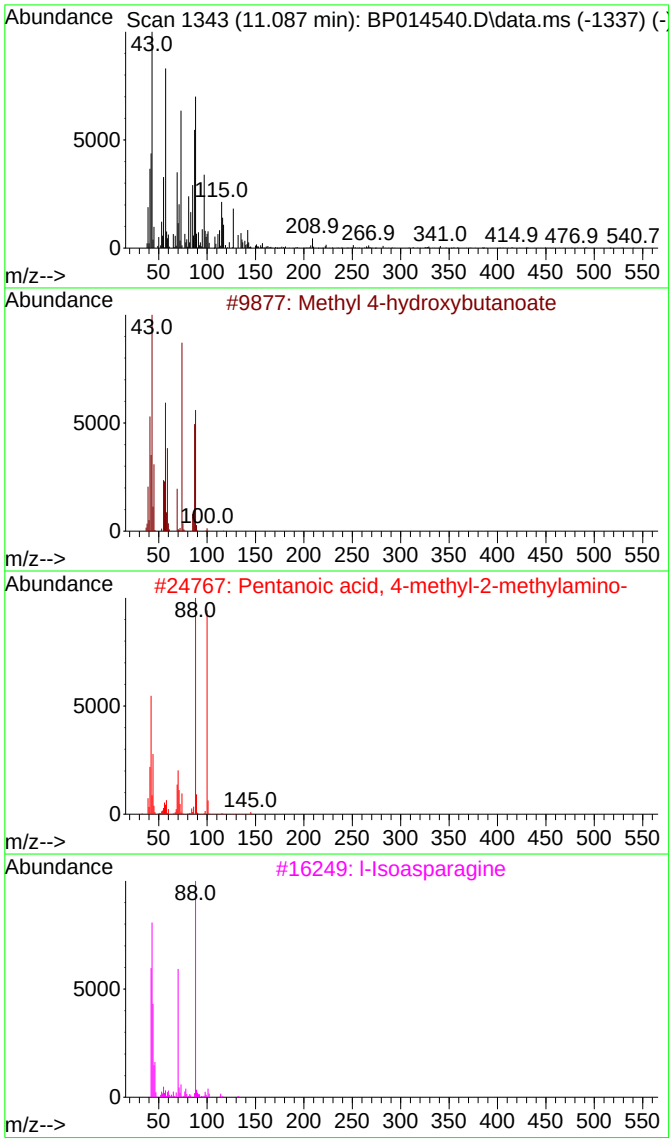
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

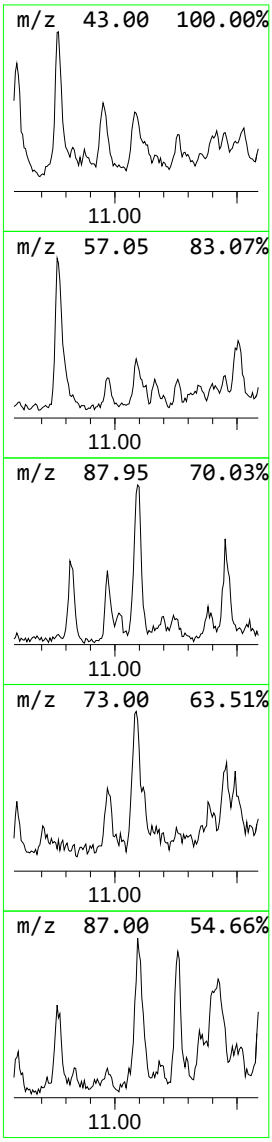
**********

Peak Number 20 unknown-04 Concentration Rank 27

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.087 | 3.29 ng/ul | 194790 | Naphthalene-d8 | 10.822 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Methyl 4-hydroxybutanoate | 118 | C5H10O3 | 000925-57-5 | 37 |
| 2 | Pentanoic acid, 4-methyl-2-methy... | 145 | C7H15NO2 | 004125-95-5 | 25 |
| 3 | l-Isoasparagine | 132 | C4H8N2O3 | 1000132-80-6 | 22 |
| 4 | Methyldiethanolamine | 119 | C5H13NO2 | 000105-59-9 | 22 |
| 5 | Acetic acid, diethyl- | 116 | C6H12O2 | 000088-09-5 | 14 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

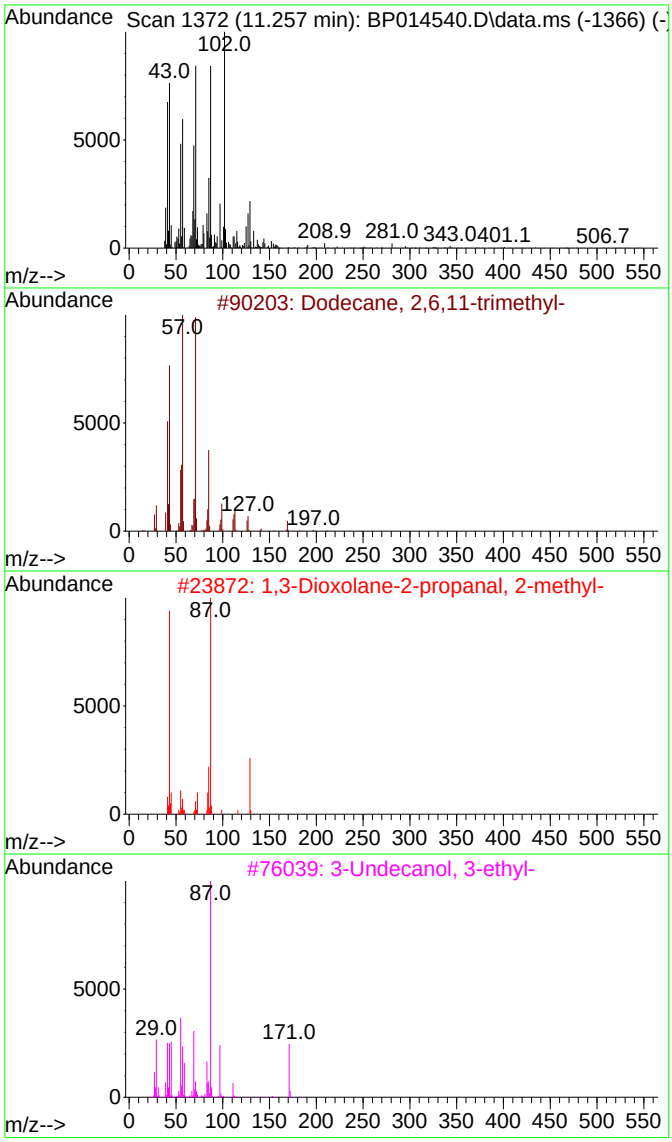
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

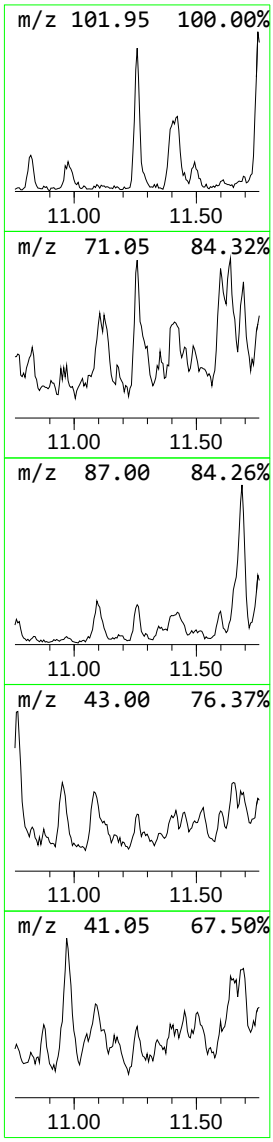
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 43

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.258 | 2.07 ng/ul | 122827 | Naphthalene-d8 | 10.822 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Dodecane, 2,6,11-trimethyl- | 212 | C15H32 | 031295-56-4 | 35 |
| 2 | 1,3-Dioxolane-2-propanal, 2-methyl- | 144 | C7H12O3 | 024108-29-0 | 22 |
| 3 | 3-Undecanol, 3-ethyl- | 200 | C13H28O | 062101-31-9 | 14 |
| 4 | Ethyl 3-methyl-2-oxobutyrate | 144 | C7H12O3 | 020201-24-5 | 11 |
| 5 | Propanoic acid, 2-methyl-, 1-met... | 158 | C9H18O2 | 054340-93-1 | 10 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

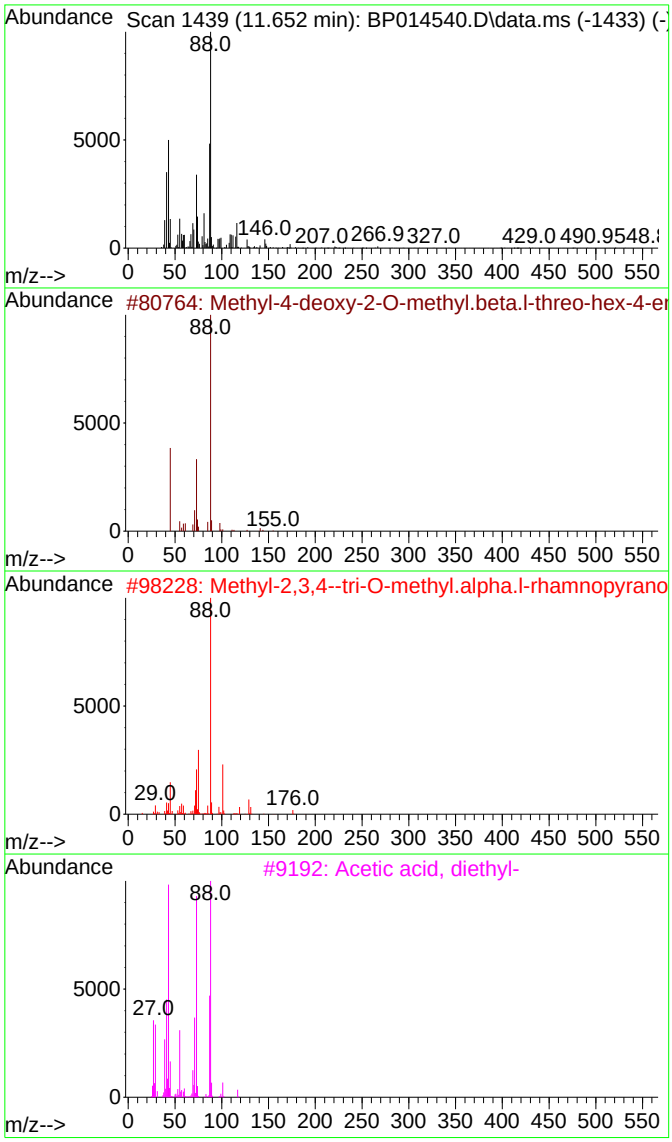
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

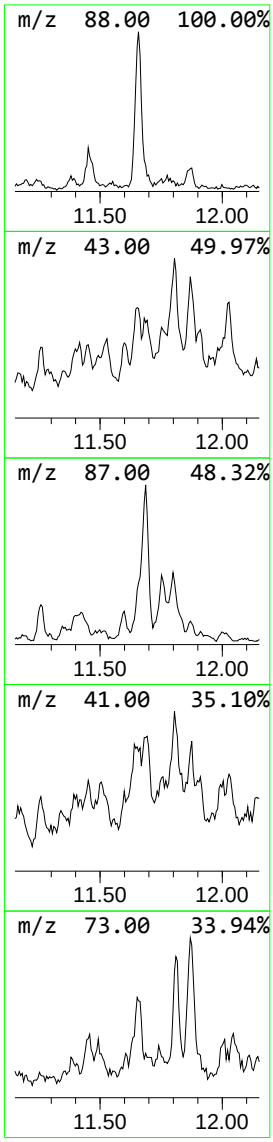
**********

Peak Number 23 unknown-05 Concentration Rank 25

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.652 | 3.50 ng/ul | 207256 | Naphthalene-d8 | 10.822 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Methyl-4-deoxy-2-O-methyl.beta.l... | 204 | C8H12O6 | 1000312-10-0 | 43 |
| 2 | Methyl-2,3,4--tri-O-methyl.alpha... | 220 | C10H20O5 | 003253-23-4 | 43 |
| 3 | Acetic acid, diethyl- | 116 | C6H12O2 | 000088-09-5 | 40 |
| 4 | Methyl(methyl-4-deoxy-2-O-methyl... | 218 | C9H14O6 | 052545-23-0 | 38 |
| 5 | Glycine, N-methoxycarbonyl-, tri... | 315 | C17H33NO4 | 1000313-49-3 | 38 |



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Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

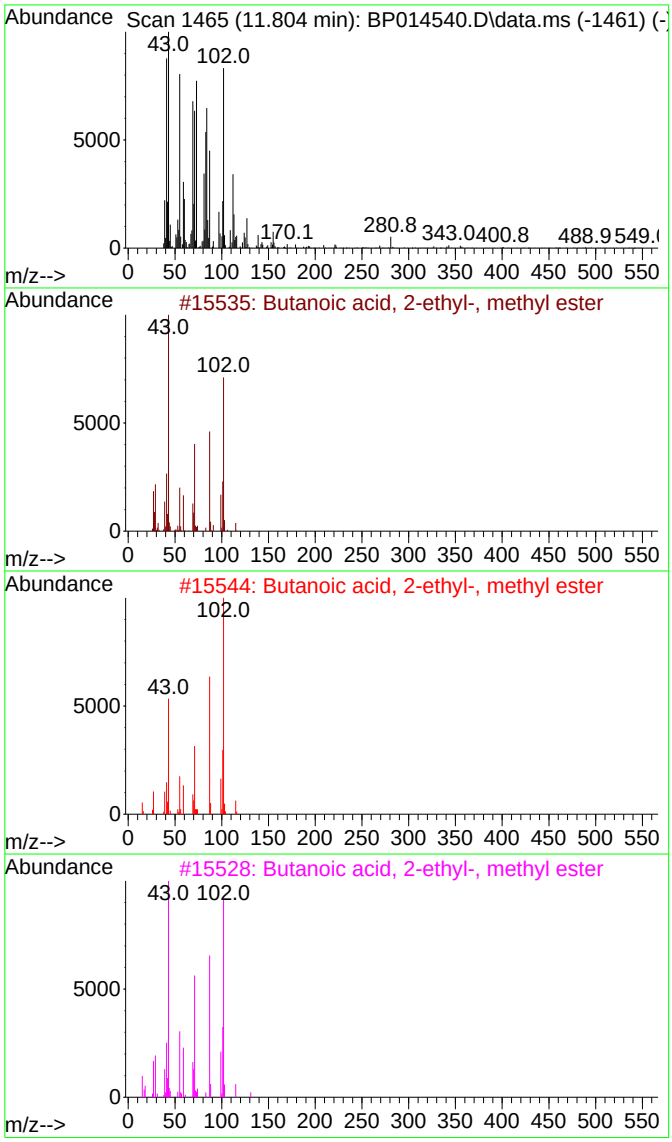
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

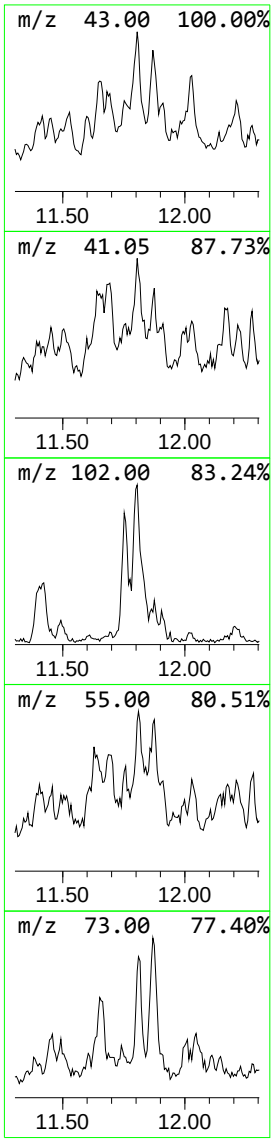
**********

Peak Number 25 unknown-06 Concentration Rank 28

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.805 | 3.26 ng/ul | 193440 | Naphthalene-d8 | 10.822 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butanoic acid, 2-ethyl-, methyl ... | 130 | C7H14O2 | 000816-11-5 | 43 |
| 2 | Butanoic acid, 2-ethyl-, methyl ... | 130 | C7H14O2 | 000816-11-5 | 38 |
| 3 | Butanoic acid, 2-ethyl-, methyl ... | 130 | C7H14O2 | 000816-11-5 | 38 |
| 4 | l-Norleucenine ethyl ester | 226 | C10H14N2O4 | 180995-53-3 | 22 |
| 5 | Butanoic acid, 4-(2-oxocyclopent... | 170 | C9H14O3 | 003296-46-6 | 15 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

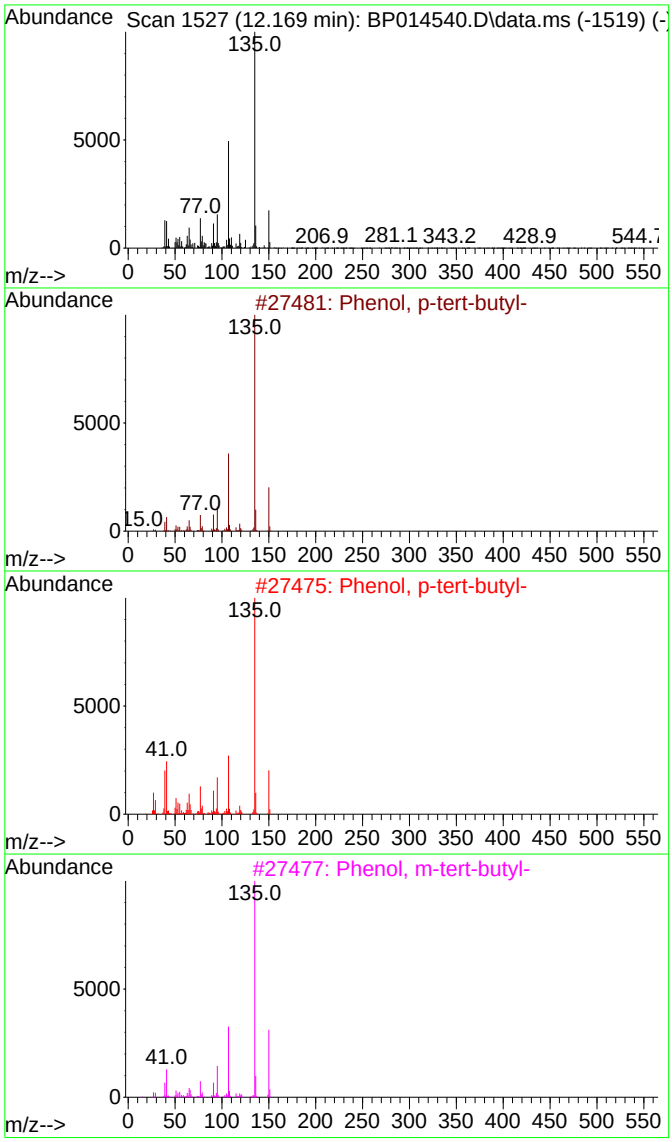
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

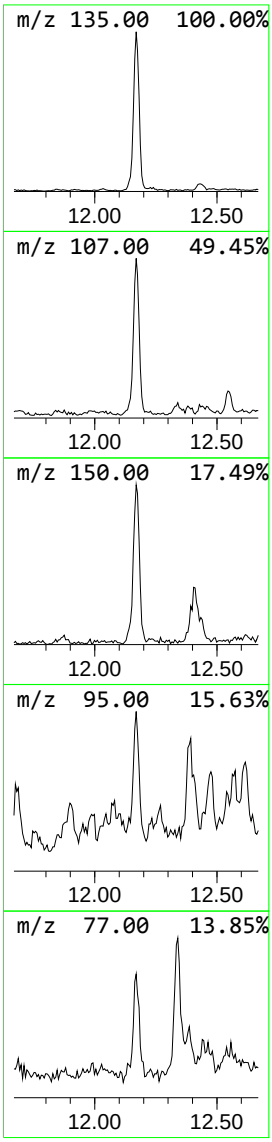
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 27 Phenol, p-tert-butyl- Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.169 | 5.67 ng/ul | 336028 | Naphthalene-d8 | 10.822 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 95 |
| 2 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 93 |
| 3 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 90 |
| 4 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 90 |
| 5 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 90 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

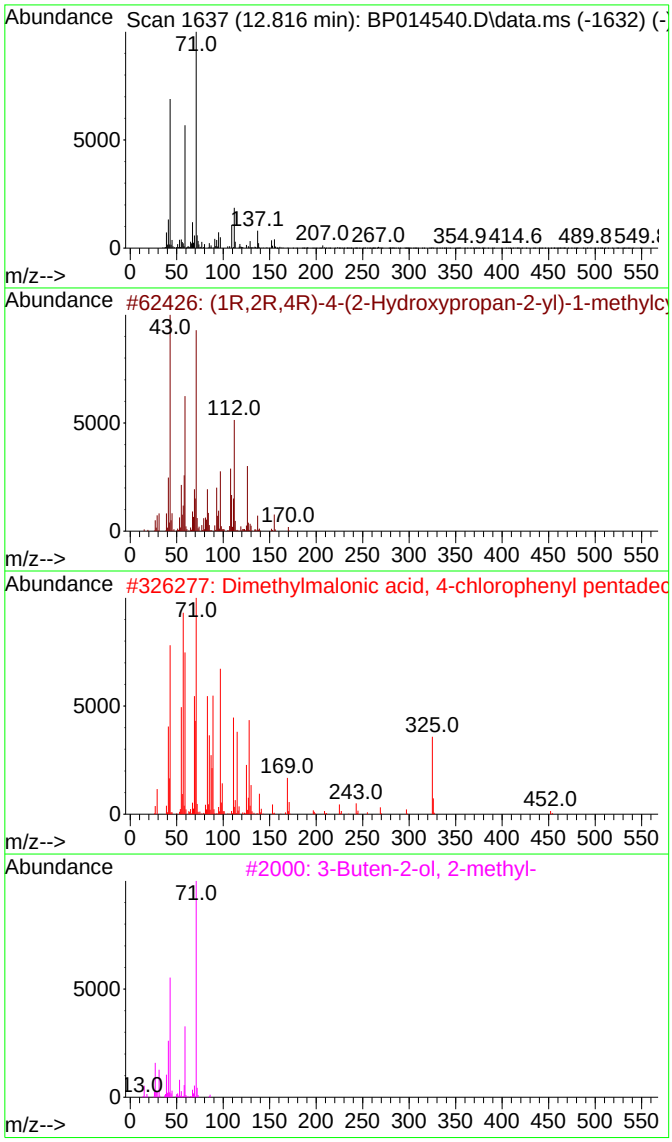
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

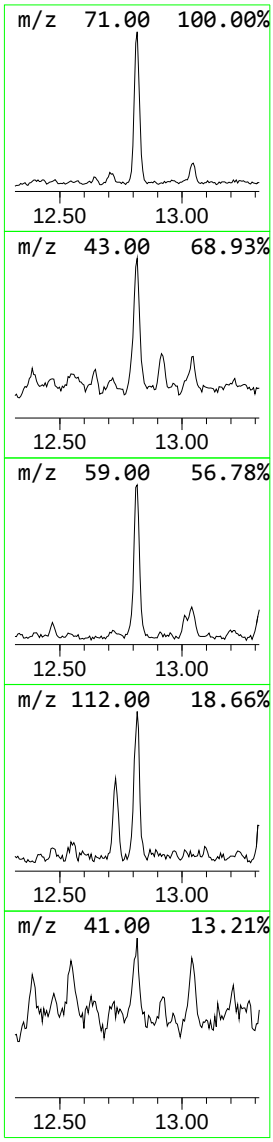
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 29 (1R,2R,4R)-4-(2-Hydroxyprop... Concentration Rank 24

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.816 | 3.80 ng/ul | 316109 | Acenaphthene-d10 | 14.651 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | (1R,2R,4R)-4-(2-Hydroxypropan-2-... | 188 | C10H20O3 | 093861-31-5 | 56 |
| 2 | Dimethylmalonic acid, 4-chloroph... | 452 | C26H41ClO4 | 1000361-98-3 | 53 |
| 3 | 3-Buten-2-ol, 2-methyl- | 86 | C5H10O | 000115-18-4 | 42 |
| 4 | 2-METHYLHEPT-1-EN-3-OL | 128 | C8H16O | 013019-19-7 | 38 |
| 5 | Ether, 1-hexadecenyl methyl, (Z)- | 254 | C17H34O | 020246-43-9 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

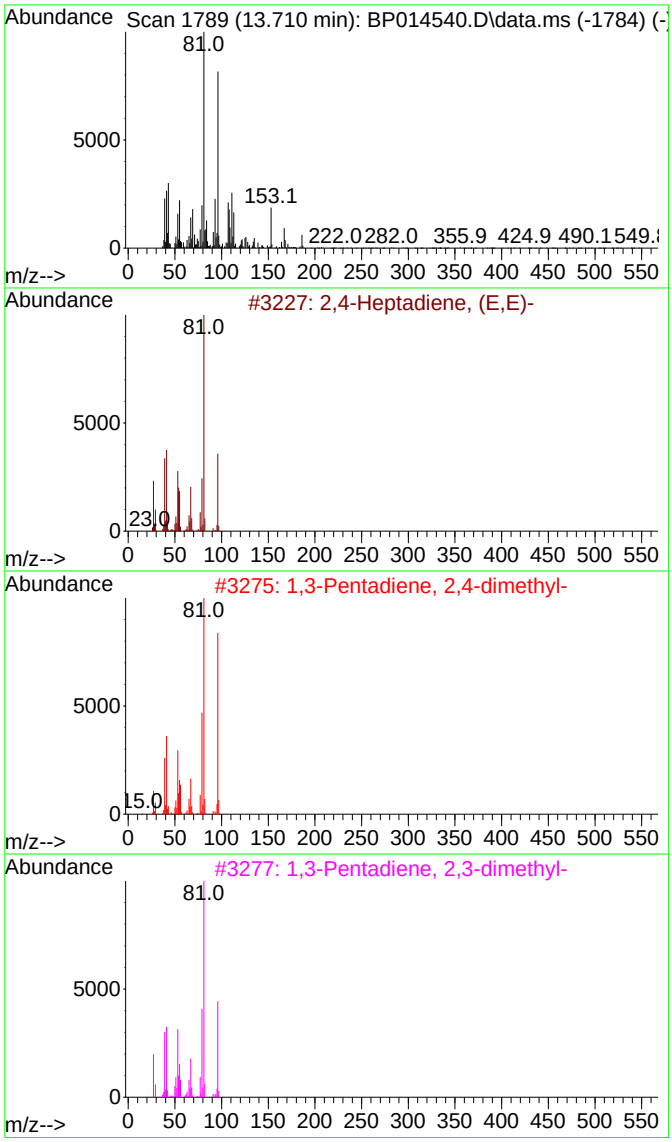
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

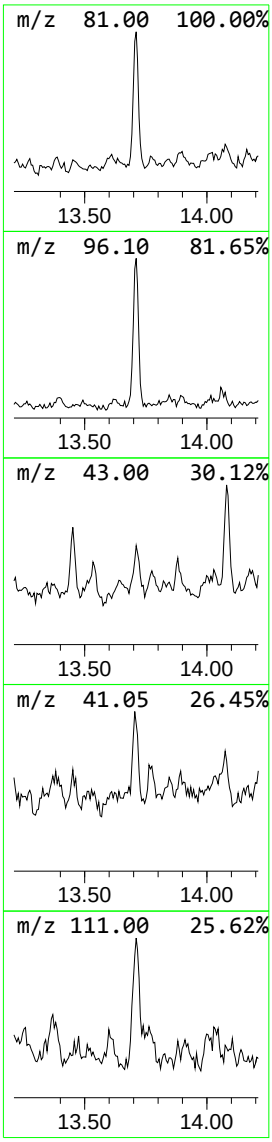
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 30 2,4-Heptadiene, (E,E)- Concentration Rank 29

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.710 | 3.26 ng/ul | 270955 | Acenaphthene-d10 | 14.651 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,4-Heptadiene, (E,E)- | 96 | C7H12 | 002384-94-3 | 53 |
| 2 | 1,3-Pentadiene, 2,4-dimethyl- | 96 | C7H12 | 001000-86-8 | 49 |
| 3 | 1,3-Pentadiene, 2,3-dimethyl- | 96 | C7H12 | 001113-56-0 | 47 |
| 4 | Cyclohexene, 1-methyl- | 96 | C7H12 | 000591-49-1 | 47 |
| 5 | (S)-2,5-Dimethyl-3-vinylhex-4-en... | 154 | C10H18O | 035671-15-9 | 47 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

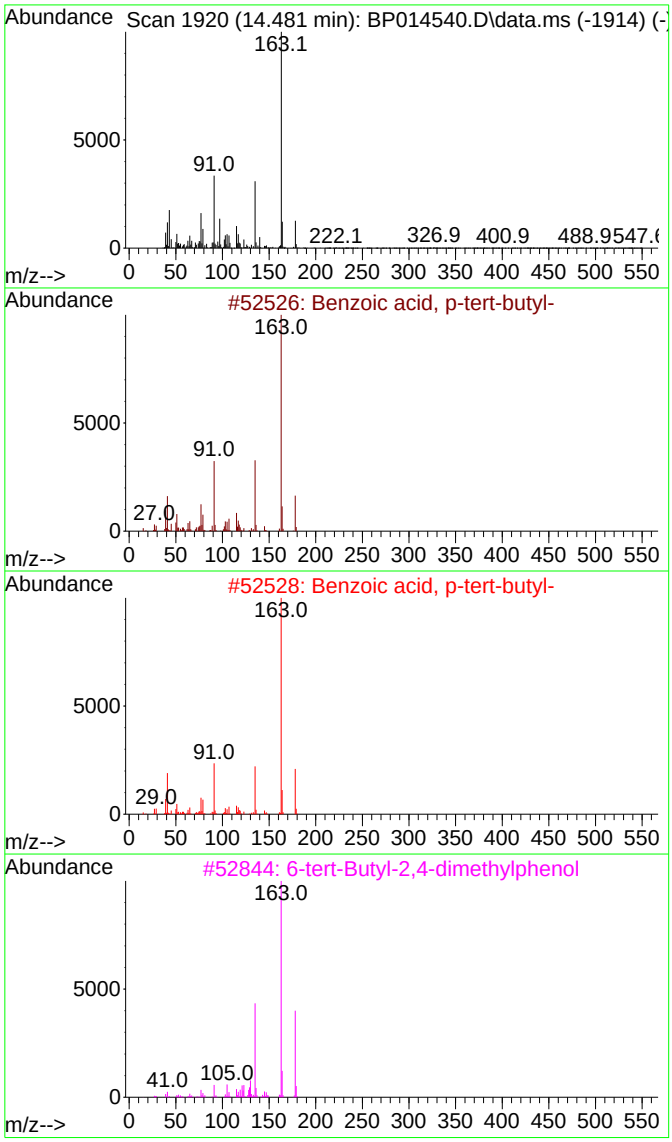
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

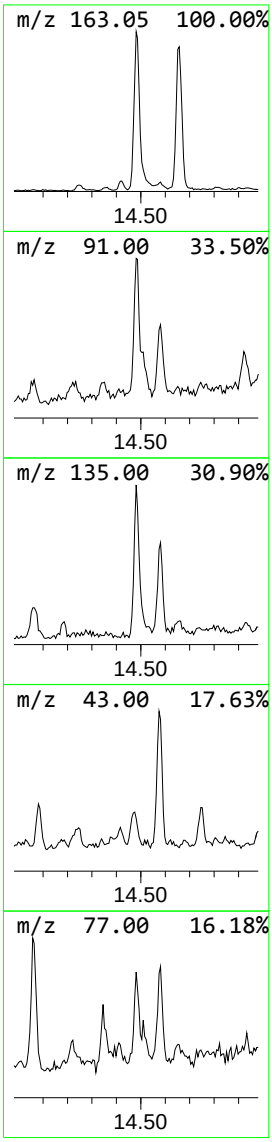
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 31 Benzoic acid, p-tert-butyl- Concentration Rank 21

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.481 | 4.58 ng/ul | 381287 | Acenaphthene-d10 | 14.651 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzoic acid, p-tert-butyl- | 178 | C11H14O2 | 000098-73-7 | 93 |
| 2 | Benzoic acid, p-tert-butyl- | 178 | C11H14O2 | 000098-73-7 | 90 |
| 3 | 6-tert-Butyl-2,4-dimethylphenol | 178 | C12H18O | 001879-09-0 | 64 |
| 4 | 4-tert-Butyl-2,6-dimethylphenol | 178 | C12H18O | 000879-97-0 | 59 |
| 5 | 2-Propenal, 3-(2,6,6-trimethyl-1... | 178 | C12H18O | 004951-40-0 | 53 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

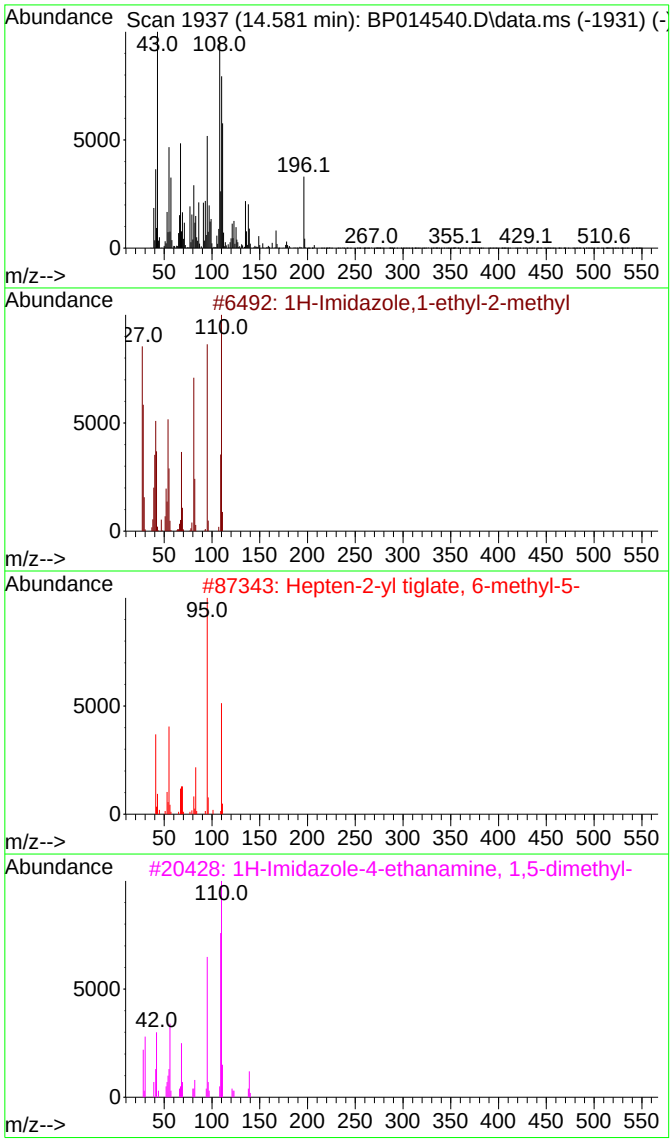
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

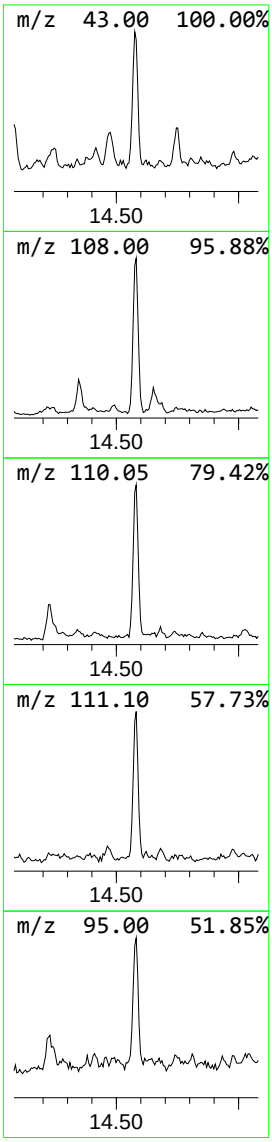
**********

Peak Number 32 unknown-07 Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.581 | 9.05 ng/ul | 753005 | Acenaphthene-d10 | 14.651 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1H-Imidazole,1-ethyl-2-methyl | 110 | C6H10N2 | 021202-52-8 | 27 |
| 2 | Hepten-2-yl tiglate, 6-methyl-5- | 210 | C13H22O2 | 1000383-64-5 | 25 |
| 3 | 1H-Imidazole-4-ethanamine, 1,5-d... | 139 | C7H13N3 | 053966-45-3 | 22 |
| 4 | 1H-Imidazole, 5-octanoyl- | 194 | C11H18N2O | 061985-27-1 | 22 |
| 5 | 2,4-Hexadiene, 2,3-dimethyl- | 110 | C8H14 | 005678-98-8 | 22 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

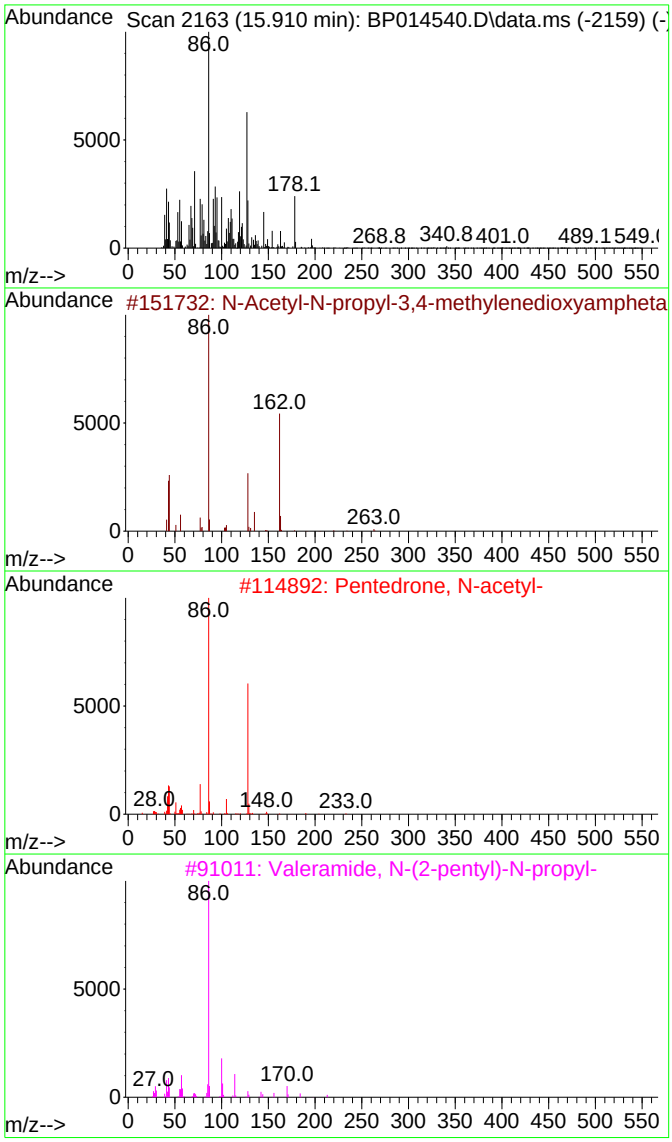
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

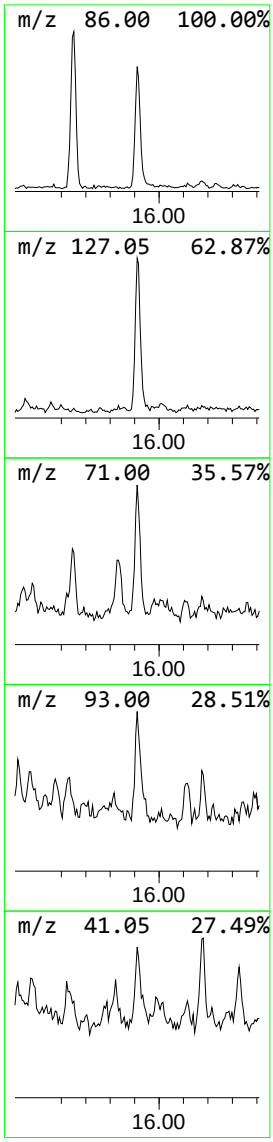
**********

Peak Number 34 unknown-08 Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.910 | 5.72 ng/ul | 476003 | Acenaphthene-d10 | 14.651 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | N-Acetyl-N-propyl-3,4-methylened... | 263 | C15H21NO3 | 1000123-03-1 | 27 |
| 2 | Pentedrone, N-acetyl- | 233 | C14H19NO2 | 1000448-38-9 | 27 |
| 3 | Valeramide, N-(2-pentyl)-N-propyl- | 213 | C13H27NO | 1000489-76-7 | 27 |
| 4 | N-Acetyl-DL-leucine, 2-methylpro... | 229 | C12H23NO3 | 1000453-80-3 | 27 |
| 5 | 4-Amino-3-chloro-N-[2-(diethylam... | 311 | C15H22ClN3O2 | 227014-59-7 | 22 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

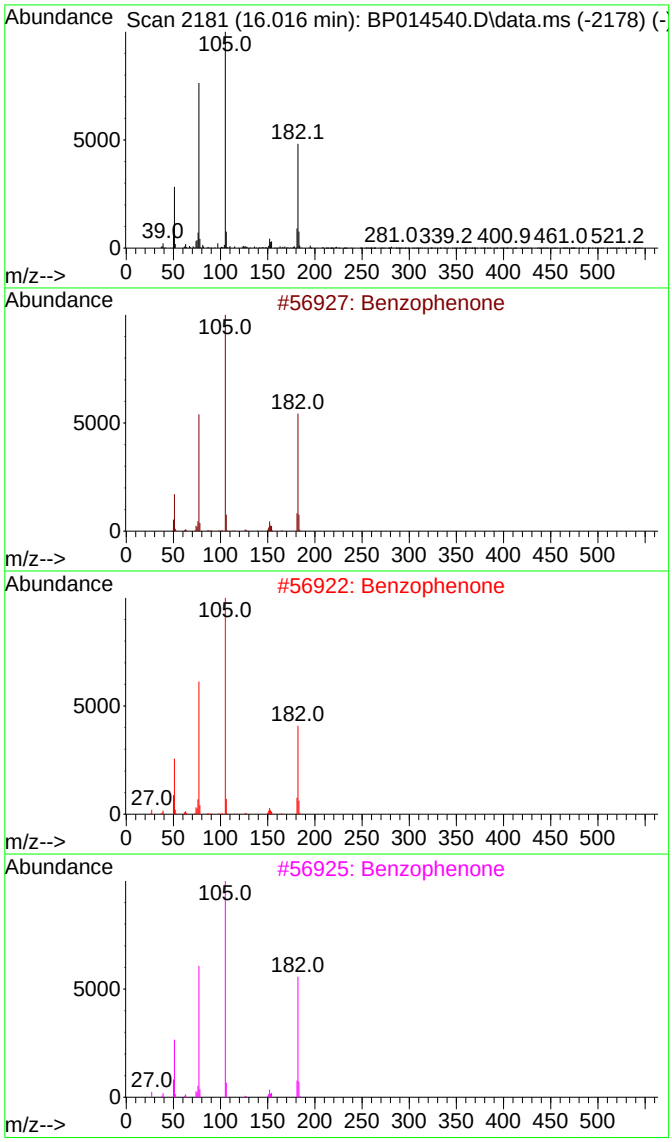
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

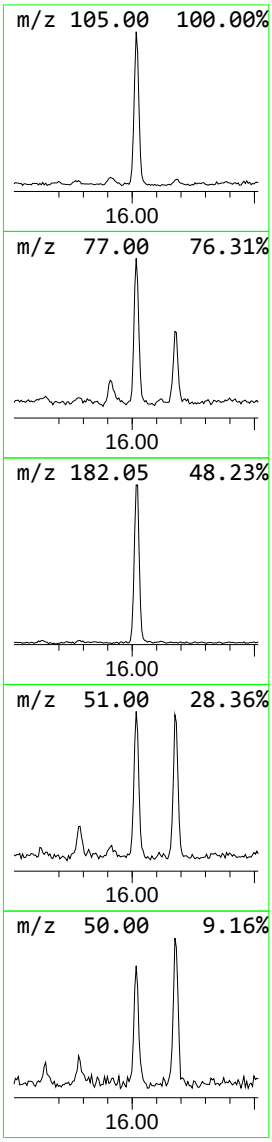
**********

Peak Number 35 Benzophenone Concentration Rank 22

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.016 | 4.55 ng/ul | 378996 | Acenaphthene-d10 | 14.651 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzophenone | 182 | C13H10O | 000119-61-9 | 95 |
| 2 | Benzophenone | 182 | C13H10O | 000119-61-9 | 95 |
| 3 | Benzophenone | 182 | C13H10O | 000119-61-9 | 93 |
| 4 | Benzophenone | 182 | C13H10O | 000119-61-9 | 91 |
| 5 | 1,2,4,5-Tetroxane, 3,3,6,6-tetra... | 396 | C26H20O4 | 016204-36-7 | 78 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

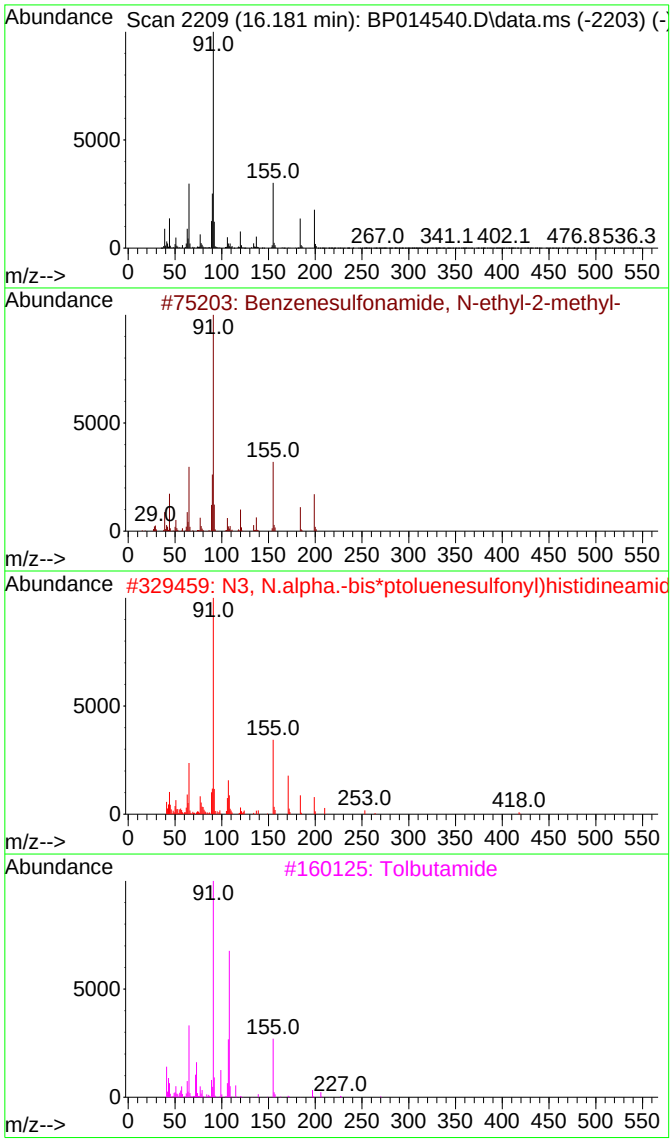
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

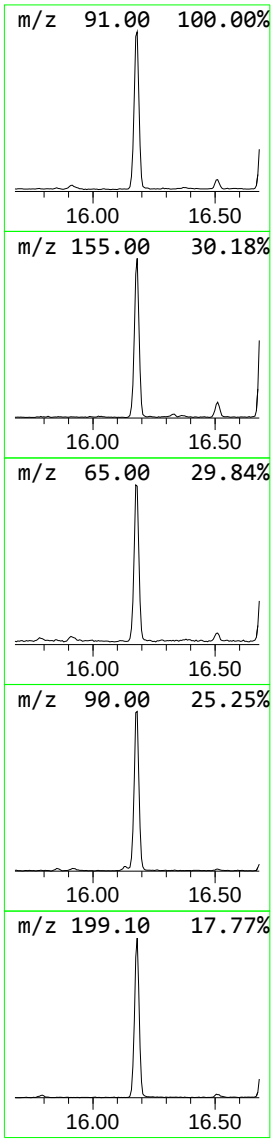
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 36 Benzenesulfonamide, N-ethyl... Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.181 | 14.63 ng/ul | 1682220 | Phenanthrene-d10 | 17.416 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzenesulfonamide, N-ethyl-2-me... | 199 | C9H13NO2S | 001077-56-1 | 98 |
| 2 | N3, N.alpha.-bis*ptoluenesulfony... | 462 | C20H22N4O5S2 | 1000130-21-9 | 50 |
| 3 | Tolbutamide | 270 | C12H18N2O3S | 000064-77-7 | 50 |
| 4 | Benzenesulfonamide, N-ethyl-2-me... | 199 | C9H13NO2S | 001077-56-1 | 49 |
| 5 | p-Toluenesulfonylacetonitrile | 195 | C9H9NO2S | 005697-44-9 | 47 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

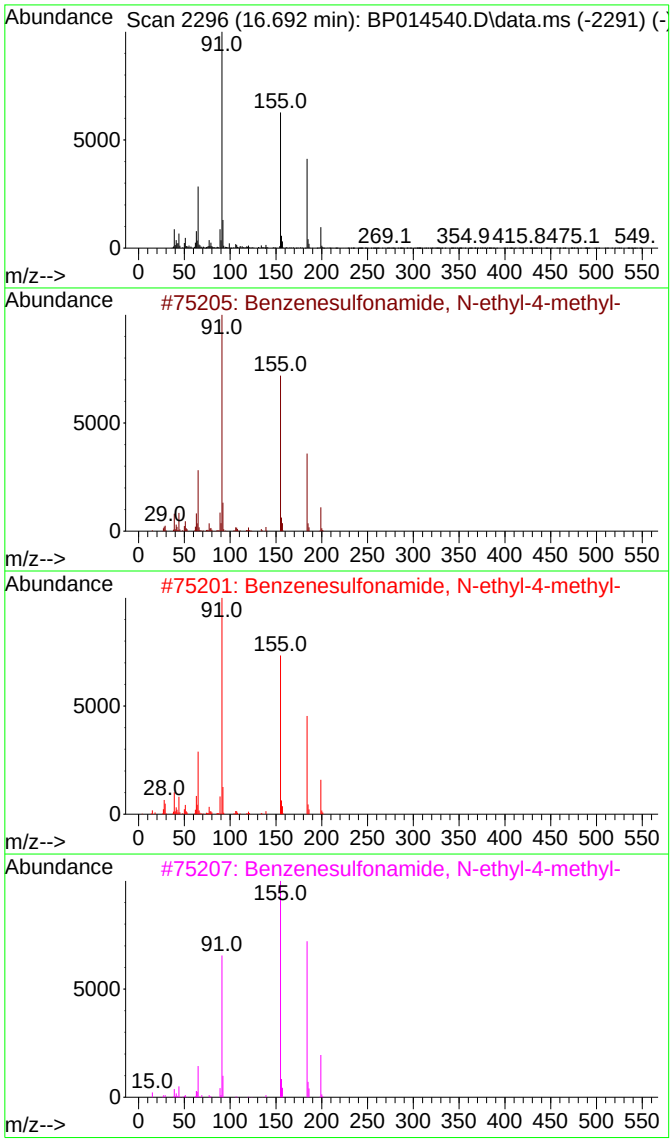
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

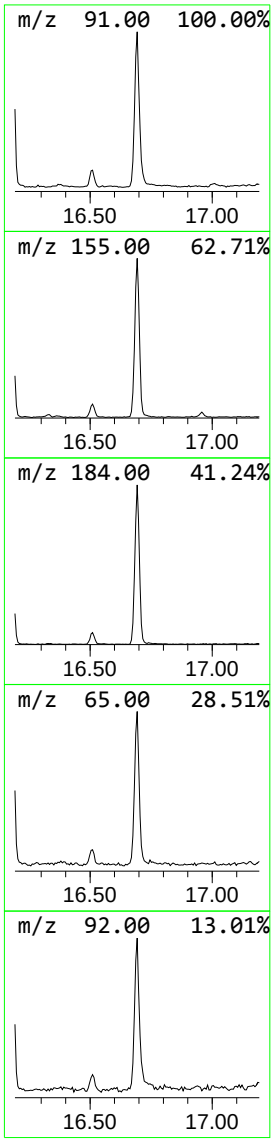
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 37 Benzenesulfonamide, N-ethyl... Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.692 | 8.74 ng/ul | 1005400 | Phenanthrene-d10 | 17.416 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzenesulfonamide, N-ethyl-4-me... | 199 | C9H13NO2S | 000080-39-7 | 97 |
| 2 | Benzenesulfonamide, N-ethyl-4-me... | 199 | C9H13NO2S | 000080-39-7 | 96 |
| 3 | Benzenesulfonamide, N-ethyl-4-me... | 199 | C9H13NO2S | 000080-39-7 | 91 |
| 4 | Benzenesulfonamide, N-ethyl-4-me... | 199 | C9H13NO2S | 000080-39-7 | 91 |
| 5 | N-isobutyl-4-methylbenzenesulfon... | 227 | C11H17NO2S | 023705-38-6 | 86 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

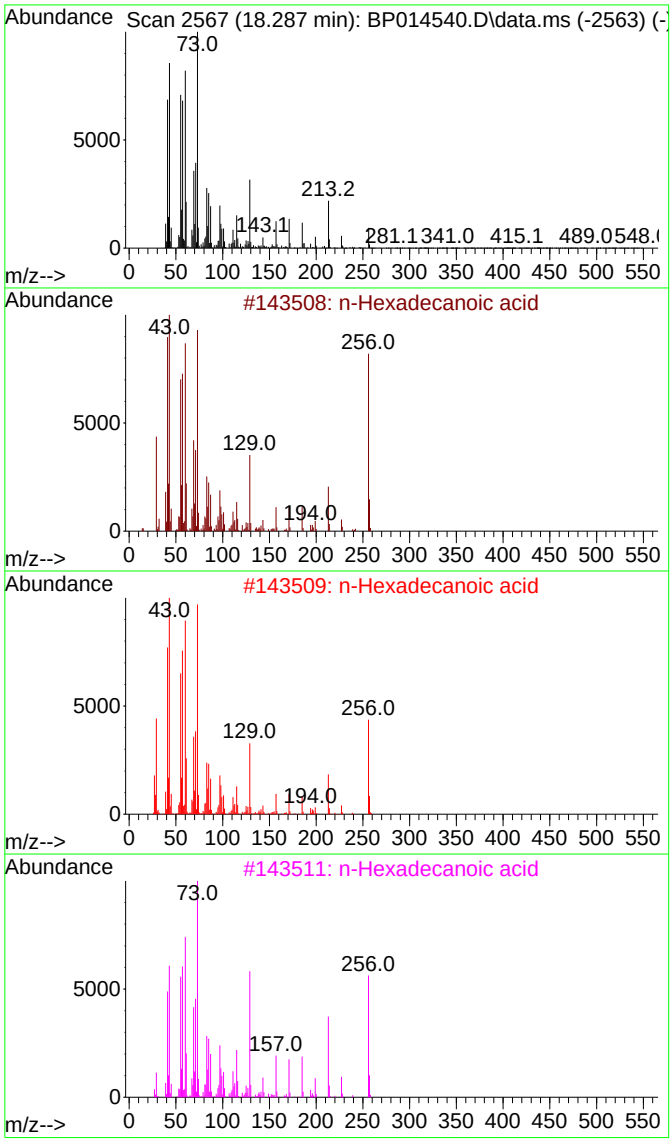
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

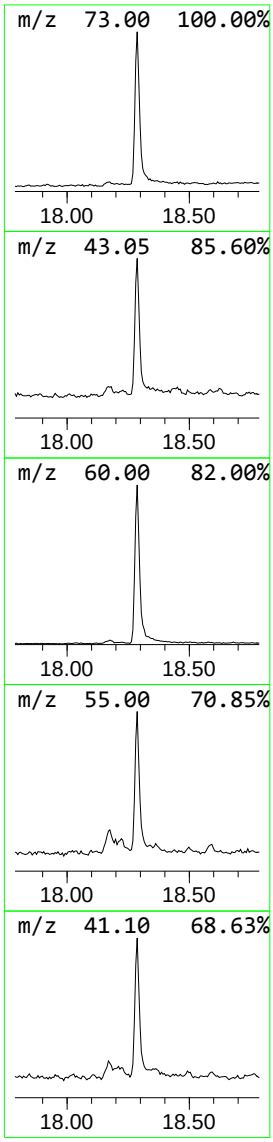
**********

Peak Number 38 n-Hexadecanoic acid Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.287 | 11.57 ng/ul | 1330760 | Phenanthrene-d10 | 17.416 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98 |
| 2 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98 |
| 3 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 97 |
| 4 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 97 |
| 5 | Tridecanoic acid | 214 | C13H26O2 | 000638-53-9 | 95 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

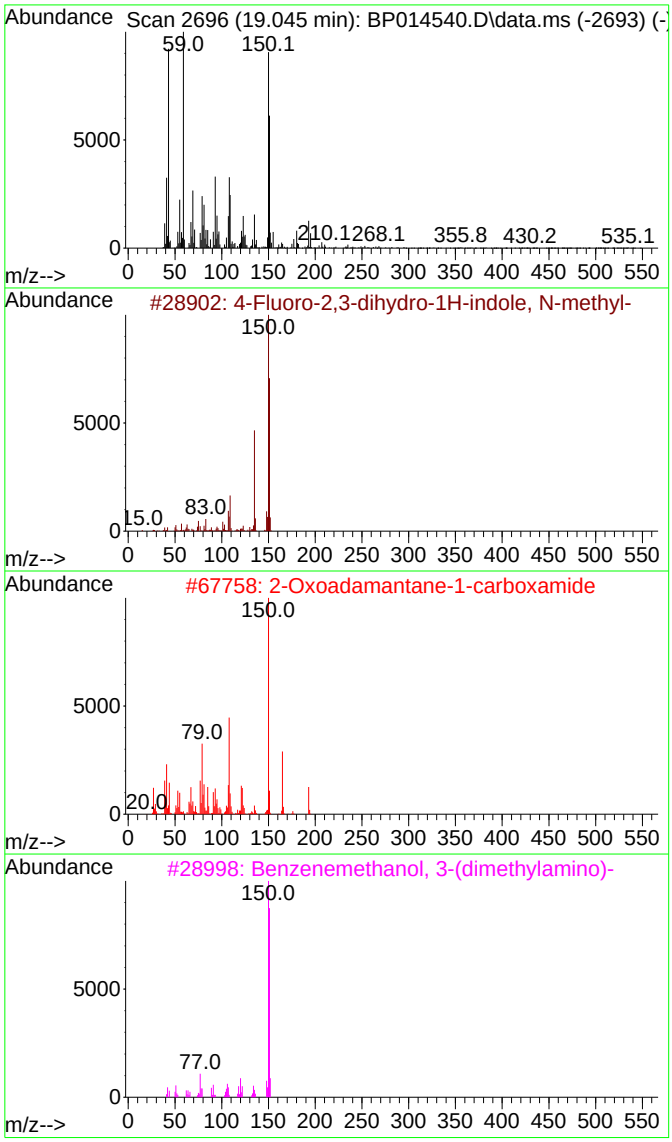
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

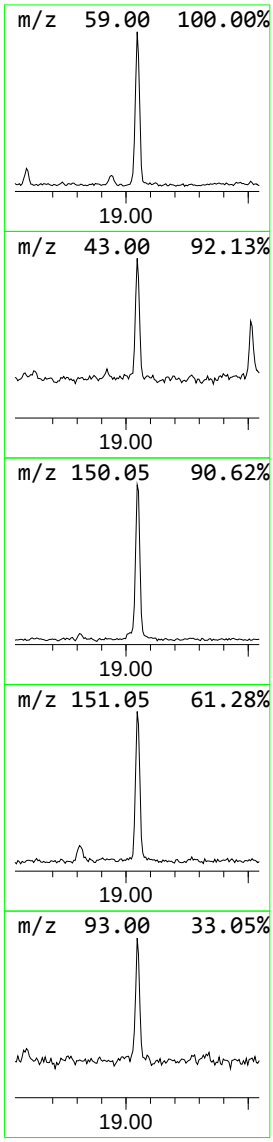
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 39 4-Fluoro-2,3-dihydro-1H-ind... Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.045 | 4.78 ng/ul | 549770 | Phenanthrene-d10 | 17.416 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4-Fluoro-2,3-dihydro-1H-indole, ... | 151 | C9H10FN | 1851835-07-8 | 60 |
| 2 | 2-Oxoadamantane-1-carboxamide | 193 | C11H15NO2 | 041171-77-1 | 35 |
| 3 | Benzenemethanol, 3-(dimethylamino)- | 151 | C9H13NO | 023501-93-1 | 30 |
| 4 | 6-Benzothiazolamine | 150 | C7H6N2S | 000533-30-2 | 25 |
| 5 | 4,6-Diaminopyrazolo[3,4-d]pyrimi... | 150 | C5H6N6 | 005413-80-9 | 22 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

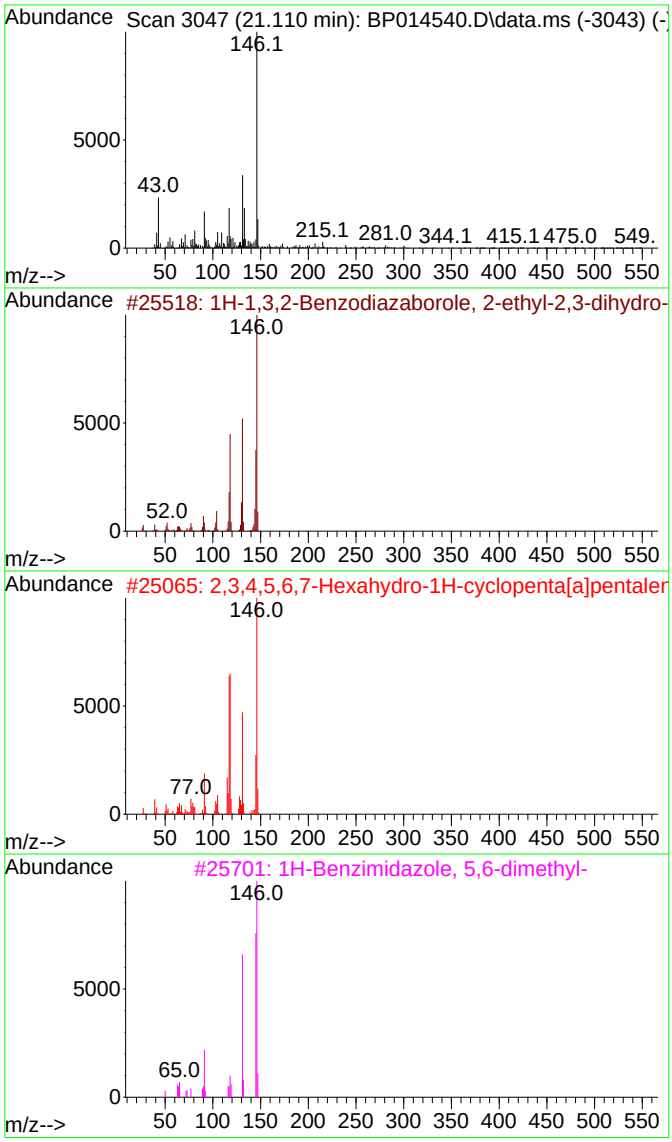
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

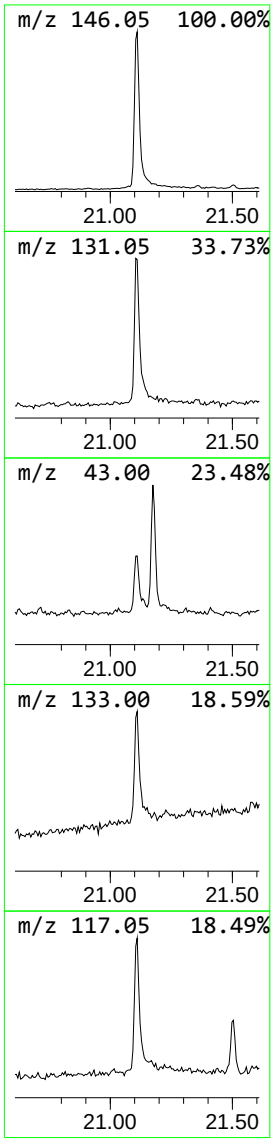
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 42 1H-1,3,2-Benzodiazaborole, ... Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.110 | 9.54 ng/ul | 1158130 | Chrysene-d12 | 21.504 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1H-1,3,2-Benzodiazaborole, 2-eth... | 146 | C8H11BN2 | 074663-81-3 | 64 |
| 2 | 2,3,4,5,6,7-Hexahydro-1H-cyclope... | 146 | C11H14 | 1010189-31-0 | 59 |
| 3 | 1H-Benzimidazole, 5,6-dimethyl- | 146 | C9H10N2 | 000582-60-5 | 59 |
| 4 | 2-Isoquinolineacetamide, N-cyclo... | 258 | C16H22N2O | 1000338-10-0 | 59 |
| 5 | 3-(Nitromethyl)phthalide | 193 | C9H7NO4 | 003598-68-3 | 53 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

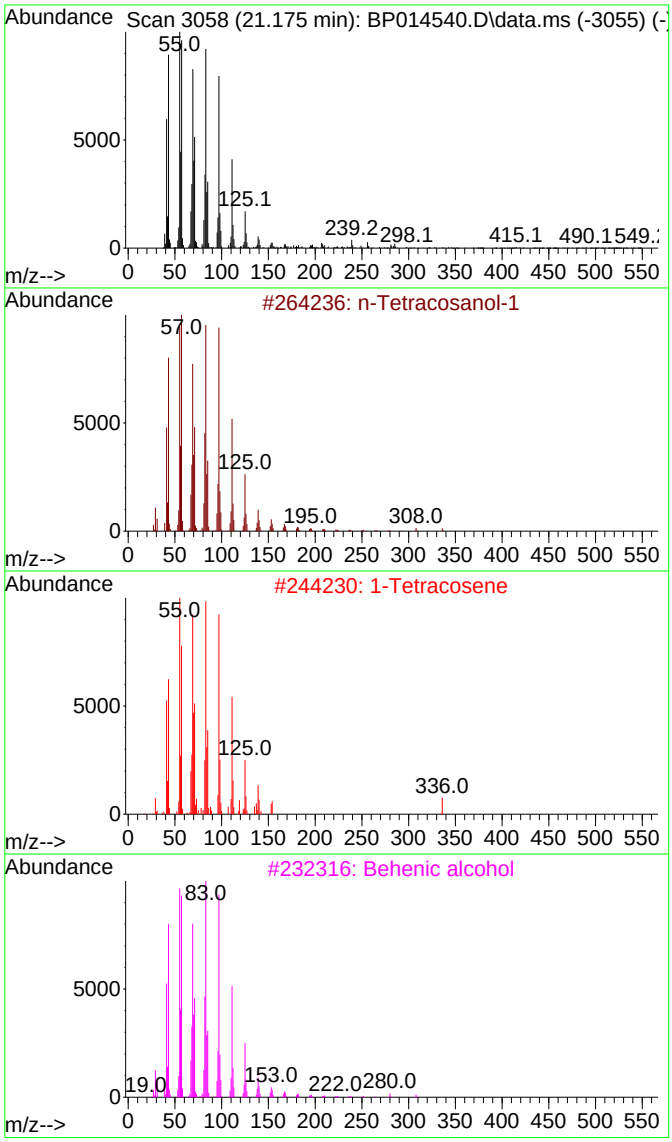
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

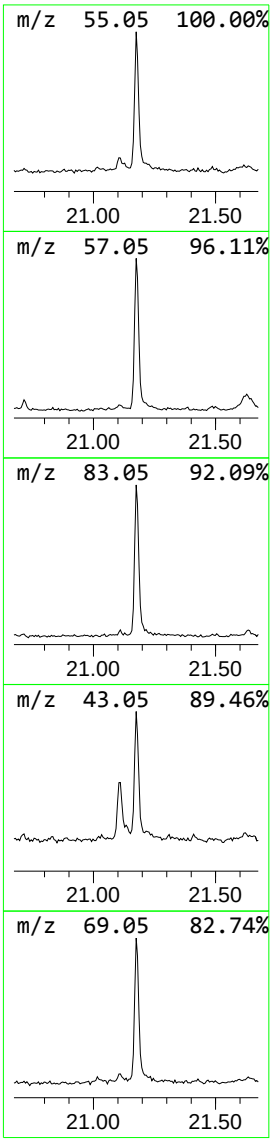
**********

Peak Number 43 n-Tetracosanol-1 Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.175 | 8.86 ng/ul | 1075310 | Chrysene-d12 | 21.504 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Tetracosanol-1 | 354 | C24H50O | 000506-51-4 | 94 |
| 2 | 1-Tetracosene | 336 | C24H48 | 010192-32-2 | 94 |
| 3 | Behenic alcohol | 326 | C22H46O | 000661-19-8 | 94 |
| 4 | 1-Heneicosyl formate | 340 | C22H44O2 | 077899-03-7 | 94 |
| 5 | Heptacos-1-ene | 378 | C27H54 | 015306-27-1 | 94 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
Data File : BP014540.D
Acq On : 04 Apr 2023 22:00
Operator : CG/JU
Sample : 02122-04
Misc :
ALS Vial : 63 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
CB9A6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.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 |
| Paraldehyde | 4.211 | 5.2 | ng/ul | 201035 | 1 | 7.999 | 774195 | 20.0 |
| 1,3-Oxathiolane | 4.593 | 11.6 | ng/ul | 448886 | 1 | 7.999 | 774195 | 20.0 |
| N-Ethylmorpholine | 5.717 | 8.9 | ng/ul | 345609 | 1 | 7.999 | 774195 | 20.0 |
| 1,3-Propanediam... | 6.952 | 3.9 | ng/ul | 150524 | 1 | 7.999 | 774195 | 20.0 |
| unknown-01 | 7.458 | 4.7 | ng/ul | 180467 | 1 | 7.999 | 774195 | 20.0 |
| Tetramethylbuta... | 8.087 | 4.9 | ng/ul | 190880 | 1 | 7.999 | 774195 | 20.0 |
| Benzenamine, 3-... | 9.005 | 180.3 | ng/ul | 6978100 | 1 | 7.999 | 774195 | 20.0 |
| Atrolactic acid | 9.116 | 5.8 | ng/ul | 224968 | 1 | 7.999 | 774195 | 20.0 |
| unknown-02 | 9.528 | 3.3 | ng/ul | 194868 | 2 | 10.822 | 1185840 | 20.0 |
| Hexanoic acid, ... | 9.652 | 5.1 | ng/ul | 302569 | 2 | 10.822 | 1185840 | 20.0 |
| unknown-03 | 10.028 | 4.9 | ng/ul | 289403 | 2 | 10.822 | 1185840 | 20.0 |
| Bicyclo[3.1.1]h... | 10.087 | 5.3 | ng/ul | 314740 | 2 | 10.822 | 1185840 | 20.0 |
| Bicyclo[2.2.1]h... | 10.363 | 8.1 | ng/ul | 481657 | 2 | 10.822 | 1185840 | 20.0 |
| Morpholine, 4-a... | 10.769 | 3.2 | ng/ul | 187865 | 2 | 10.822 | 1185840 | 20.0 |
| unknown-04 | 11.087 | 3.3 | ng/ul | 194790 | 2 | 10.822 | 1185840 | 20.0 |
| (DEL) Alkane: S... | 11.258 | 2.1 | ng/ul | 122827 | 2 | 10.822 | 1185840 | 20.0 |
| unknown-05 | 11.652 | 3.5 | ng/ul | 207256 | 2 | 10.822 | 1185840 | 20.0 |
| unknown-06 | 11.805 | 3.3 | ng/ul | 193440 | 2 | 10.822 | 1185840 | 20.0 |
| Phenol, p-tert-... | 12.169 | 5.7 | ng/ul | 336028 | 2 | 10.822 | 1185840 | 20.0 |
| (1R,2R,4R)-4-(2... | 12.816 | 3.8 | ng/ul | 316109 | 3 | 14.651 | 1664100 | 20.0 |
| 2,4-Heptadiene,... | 13.710 | 3.3 | ng/ul | 270955 | 3 | 14.651 | 1664100 | 20.0 |
| Benzoic acid, p... | 14.481 | 4.6 | ng/ul | 381287 | 3 | 14.651 | 1664100 | 20.0 |
| unknown-07 | 14.581 | 9.1 | ng/ul | 753005 | 3 | 14.651 | 1664100 | 20.0 |
| unknown-08 | 15.910 | 5.7 | ng/ul | 476003 | 3 | 14.651 | 1664100 | 20.0 |
| Benzophenone | 16.016 | 4.5 | ng/ul | 378996 | 3 | 14.651 | 1664100 | 20.0 |
| Benzenesulfonam... | 16.181 | 14.6 | ng/ul | 1682220 | 4 | 17.416 | 2299650 | 20.0 |
| Benzenesulfonam... | 16.692 | 8.7 | ng/ul | 1005400 | 4 | 17.416 | 2299650 | 20.0 |
| n-Hexadecanoic ... | 18.287 | 11.6 | ng/ul | 1330760 | 4 | 17.416 | 2299650 | 20.0 |
| 4-Fluoro-2,3-di... | 19.045 | 4.8 | ng/ul | 549770 | 4 | 17.416 | 2299650 | 20.0 |
| 1H-1,3,2-Benzod... | 21.110 | 9.5 | ng/ul | 1158130 | 5 | 21.504 | 2428230 | 20.0 |
| n-Tetracosanol-1 | 21.175 | 8.9 | ng/ul | 1075310 | 5 | 21.504 | 2428230 | 20.0 |