

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
 Data File : BM021272.D  
 Acq On : 29 Jun 2019 21:17  
 Operator : HP/JU  
 Sample : K3593-16  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C7BA8

## Integration Parameters: LSCINT.P

Integrator: RTE  
 Smoothing : OFF  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 0.1 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
 Title : SVOA CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 3.016       | 3             | 5           | 24           | rVB      | 2749499        | 3271375       | 12.32%          | 1.377%        |
| 2         | 4.352       | 228           | 232         | 237          | rBV      | 301215         | 382547        | 1.44%           | 0.161%        |
| 3         | 4.404       | 237           | 241         | 246          | rVB      | 340329         | 448545        | 1.69%           | 0.189%        |
| 4         | 5.828       | 479           | 483         | 494          | rVB3     | 111901         | 187228        | 0.71%           | 0.079%        |
| 5         | 6.863       | 653           | 659         | 664          | rBV      | 248504         | 369377        | 1.39%           | 0.155%        |
| 6         | 6.934       | 664           | 671         | 682          | rVB      | 1100481        | 1811353       | 6.82%           | 0.762%        |
| 7         | 7.104       | 695           | 700         | 711          | rVB      | 823770         | 1319524       | 4.97%           | 0.555%        |
| 8         | 7.293       | 726           | 732         | 740          | rVB      | 1155450        | 1802302       | 6.79%           | 0.759%        |
| 9         | 7.757       | 805           | 811         | 818          | rBV      | 1088962        | 1754464       | 6.61%           | 0.738%        |
| 10        | 7.887       | 827           | 833         | 840          | rVB      | 442763         | 653472        | 2.46%           | 0.275%        |
| 11        | 8.057       | 857           | 862         | 871          | rVB      | 304012         | 475433        | 1.79%           | 0.200%        |
| 12        | 8.357       | 903           | 913         | 918          | rBV3     | 101646         | 200085        | 0.75%           | 0.084%        |
| 13        | 8.469       | 925           | 932         | 941          | rVB      | 1168782        | 1923943       | 7.25%           | 0.810%        |
| 14        | 8.922       | 1002          | 1009        | 1017         | rBV      | 1154358        | 1887608       | 7.11%           | 0.795%        |
| 15        | 9.069       | 1030          | 1034        | 1042         | rVV2     | 120196         | 197831        | 0.75%           | 0.083%        |
| 16        | 9.245       | 1057          | 1064        | 1068         | rVV      | 490058         | 773173        | 2.91%           | 0.325%        |
| 17        | 9.287       | 1068          | 1071        | 1080         | rVB      | 177936         | 255957        | 0.96%           | 0.108%        |
| 18        | 9.634       | 1121          | 1130        | 1142         | rBV      | 953524         | 1681821       | 6.33%           | 0.708%        |
| 19        | 9.957       | 1179          | 1185        | 1192         | rBV2     | 82496          | 130733        | 0.49%           | 0.055%        |
| 20        | 10.169      | 1214          | 1221        | 1239         | rVB      | 1578016        | 2668070       | 10.05%          | 1.123%        |
| 21        | 10.545      | 1278          | 1285        | 1291         | rBV      | 1578307        | 2618464       | 9.86%           | 1.102%        |
| 22        | 10.692      | 1304          | 1310        | 1323         | rBV      | 527443         | 981377        | 3.70%           | 0.413%        |
| 23        | 11.222      | 1394          | 1400        | 1407         | rBV2     | 119022         | 195546        | 0.74%           | 0.082%        |
| 24        | 11.486      | 1434          | 1445        | 1451         | rBV2     | 114575         | 192669        | 0.73%           | 0.081%        |
| 25        | 11.775      | 1490          | 1494        | 1502         | rVB2     | 119384         | 193212        | 0.73%           | 0.081%        |
| 26        | 12.663      | 1639          | 1645        | 1651         | rVV3     | 76365          | 158874        | 0.60%           | 0.067%        |
| 27        | 12.745      | 1653          | 1659        | 1669         | rVB3     | 53978          | 125579        | 0.47%           | 0.053%        |
| 28        | 12.845      | 1669          | 1676        | 1681         | rBV4     | 101267         | 154714        | 0.58%           | 0.065%        |
| 29        | 13.045      | 1706          | 1710        | 1717         | rVB      | 78072          | 113782        | 0.43%           | 0.048%        |
| 30        | 13.263      | 1742          | 1747        | 1755         | rVV5     | 59114          | 142101        | 0.54%           | 0.060%        |
| 31        | 13.710      | 1818          | 1823        | 1830         | rVB5     | 81901          | 135774        | 0.51%           | 0.057%        |
| 32        | 13.816      | 1830          | 1841        | 1845         | rBV      | 2754530        | 3893133       | 14.66%          | 1.639%        |
| 33        | 13.863      | 1845          | 1849        | 1855         | rVB      | 1307681        | 1744583       | 6.57%           | 0.734%        |
| 34        | 14.092      | 1882          | 1888        | 1901         | rVB      | 3117891        | 4587858       | 17.28%          | 1.931%        |

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 Stop Thrs : 0

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 Min Area: 0.1 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 14.404 | 1934 | 1941 | 1947 | rVV2 | 2214408 | 3487490 | 13.14% | 1.468% |
| 36 | 14.463 | 1947 | 1951 | 1957 | rVB  | 85530   | 150389  | 0.57%  | 0.063% |
| 37 | 14.621 | 1970 | 1978 | 1989 | rBV  | 802979  | 1422841 | 5.36%  | 0.599% |
| 38 | 15.398 | 2104 | 2110 | 2116 | rVV  | 3970407 | 5666034 | 21.34% | 2.385% |
| 39 | 15.451 | 2116 | 2119 | 2126 | rVB  | 73803   | 104579  | 0.39%  | 0.044% |
| 40 | 15.527 | 2126 | 2132 | 2138 | rBV  | 751464  | 1081211 | 4.07%  | 0.455% |
| 41 | 15.639 | 2145 | 2151 | 2158 | rBV5 | 124208  | 200946  | 0.76%  | 0.085% |
| 42 | 15.963 | 2199 | 2206 | 2214 | rVB  | 235163  | 380668  | 1.43%  | 0.160% |
| 43 | 16.104 | 2225 | 2230 | 2232 | rBV2 | 72840   | 108768  | 0.41%  | 0.046% |
| 44 | 16.268 | 2252 | 2258 | 2262 | rBV3 | 106345  | 151180  | 0.57%  | 0.064% |
| 45 | 16.474 | 2288 | 2293 | 2297 | rVB3 | 82938   | 111079  | 0.42%  | 0.047% |
| 46 | 16.651 | 2319 | 2323 | 2327 | rBV  | 137815  | 198891  | 0.75%  | 0.084% |
| 47 | 16.904 | 2359 | 2366 | 2371 | rBV4 | 80773   | 189248  | 0.71%  | 0.080% |
| 48 | 16.968 | 2374 | 2377 | 2383 | rVB  | 74283   | 108631  | 0.41%  | 0.046% |
| 49 | 17.045 | 2386 | 2390 | 2392 | rBV3 | 88432   | 113253  | 0.43%  | 0.048% |
| 50 | 17.151 | 2402 | 2408 | 2412 | rVV  | 2381076 | 3496117 | 13.17% | 1.472% |
| 51 | 17.192 | 2412 | 2415 | 2419 | rVV  | 1109474 | 1489392 | 5.61%  | 0.627% |
| 52 | 17.251 | 2419 | 2425 | 2434 | rVB  | 3392766 | 5137510 | 19.35% | 2.162% |
| 53 | 17.527 | 2465 | 2472 | 2475 | rBV4 | 136233  | 232001  | 0.87%  | 0.098% |
| 54 | 17.668 | 2492 | 2496 | 2502 | rVB2 | 76272   | 119525  | 0.45%  | 0.050% |
| 55 | 17.927 | 2535 | 2540 | 2545 | rBV  | 177937  | 290089  | 1.09%  | 0.122% |
| 56 | 17.986 | 2545 | 2550 | 2554 | rBV  | 1042155 | 1446927 | 5.45%  | 0.609% |
| 57 | 18.045 | 2554 | 2560 | 2569 | rVB2 | 2044859 | 3215323 | 12.11% | 1.353% |
| 58 | 18.221 | 2585 | 2590 | 2594 | rBV3 | 209737  | 375196  | 1.41%  | 0.158% |
| 59 | 18.345 | 2605 | 2611 | 2616 | rBV3 | 815343  | 1209060 | 4.55%  | 0.509% |
| 60 | 18.527 | 2638 | 2642 | 2646 | rBV  | 130376  | 178697  | 0.67%  | 0.075% |
| 61 | 18.574 | 2646 | 2650 | 2660 | rVV2 | 173964  | 276649  | 1.04%  | 0.116% |
| 62 | 18.774 | 2680 | 2684 | 2687 | rBV2 | 77691   | 107687  | 0.41%  | 0.045% |
| 63 | 19.056 | 2727 | 2732 | 2744 | rVB2 | 736702  | 1156245 | 4.35%  | 0.487% |
| 64 | 19.180 | 2749 | 2753 | 2754 | rBV  | 308914  | 370975  | 1.40%  | 0.156% |
| 65 | 19.209 | 2754 | 2758 | 2764 | rVV  | 3206415 | 4607229 | 17.35% | 1.939% |
| 66 | 19.256 | 2764 | 2766 | 2772 | rVV2 | 195964  | 248641  | 0.94%  | 0.105% |
| 67 | 19.321 | 2773 | 2777 | 2787 | rVV3 | 471586  | 720527  | 2.71%  | 0.303% |
| 68 | 19.545 | 2809 | 2815 | 2818 | rBV  | 3911061 | 5866132 | 22.09% | 2.469% |
| 69 | 19.574 | 2818 | 2820 | 2825 | rVB  | 2181001 | 2493642 | 9.39%  | 1.050% |
| 70 | 19.756 | 2848 | 2851 | 2857 | rVB  | 132984  | 206481  | 0.78%  | 0.087% |
| 71 | 19.892 | 2871 | 2874 | 2881 | rVB3 | 104694  | 225763  | 0.85%  | 0.095% |

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 Peak Location: TOP

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 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |          |          |         |         |
|-----|--------|------|------|------|------|----------|----------|---------|---------|
| 72  | 20.074 | 2898 | 2905 | 2913 | rVB3 | 397874   | 927135   | 3.49%   | 0.390%  |
| 73  | 20.168 | 2918 | 2921 | 2925 | rBV  | 192758   | 255892   | 0.96%   | 0.108%  |
| 74  | 20.815 | 3028 | 3031 | 3036 | rBV  | 235947   | 326192   | 1.23%   | 0.137%  |
| 75  | 21.039 | 3065 | 3069 | 3077 | rVB4 | 1352002  | 1966742  | 7.41%   | 0.828%  |
| 76  | 21.115 | 3079 | 3082 | 3087 | rVB  | 302112   | 373474   | 1.41%   | 0.157%  |
| 77  | 21.333 | 3112 | 3119 | 3123 | rBV2 | 3272624  | 5886913  | 22.17%  | 2.478%  |
| 78  | 21.374 | 3123 | 3126 | 3135 | rVB  | 1414739  | 1824342  | 6.87%   | 0.768%  |
| 79  | 21.474 | 3141 | 3143 | 3152 | rVB3 | 388640   | 597800   | 2.25%   | 0.252%  |
| 80  | 21.915 | 3215 | 3218 | 3220 | rBV  | 787865   | 954657   | 3.60%   | 0.402%  |
| 81  | 22.121 | 3250 | 3253 | 3259 | rVB2 | 451956   | 562967   | 2.12%   | 0.237%  |
| 82  | 22.380 | 3293 | 3297 | 3301 | rBV  | 923086   | 1246441  | 4.69%   | 0.525%  |
| 83  | 22.621 | 3332 | 3338 | 3355 | rVB  | 10546546 | 15586878 | 58.71%  | 6.561%  |
| 84  | 22.903 | 3380 | 3386 | 3389 | rBV  | 2661696  | 5262593  | 19.82%  | 2.215%  |
| 85  | 22.950 | 3389 | 3394 | 3405 | rVB2 | 15456777 | 26550191 | 100.00% | 11.175% |
| 86  | 23.415 | 3468 | 3473 | 3476 | rBV  | 644859   | 1187502  | 4.47%   | 0.500%  |
| 87  | 23.462 | 3476 | 3481 | 3486 | rVV  | 3692131  | 7197730  | 27.11%  | 3.030%  |
| 88  | 23.509 | 3486 | 3489 | 3500 | rVV  | 1619383  | 4880093  | 18.38%  | 2.054%  |
| 89  | 23.609 | 3501 | 3506 | 3512 | rVB2 | 1846776  | 3520155  | 13.26%  | 1.482%  |
| 90  | 23.791 | 3531 | 3537 | 3545 | rVB  | 4555479  | 8009404  | 30.17%  | 3.371%  |
| 91  | 24.127 | 3587 | 3594 | 3602 | rVB  | 4912056  | 10013087 | 37.71%  | 4.215%  |
| 92  | 24.215 | 3604 | 3609 | 3622 | rVB  | 2988540  | 6056993  | 22.81%  | 2.549%  |
| 93  | 24.885 | 3716 | 3723 | 3737 | rVB3 | 1775949  | 4308143  | 16.23%  | 1.813%  |
| 94  | 25.338 | 3794 | 3800 | 3811 | rVB  | 2646356  | 6192967  | 23.33%  | 2.607%  |
| 95  | 25.762 | 3864 | 3872 | 3879 | rBV  | 3293978  | 9277671  | 34.94%  | 3.905%  |
| 96  | 25.915 | 3892 | 3898 | 3911 | rVB3 | 1690558  | 4977403  | 18.75%  | 2.095%  |
| 97  | 26.050 | 3914 | 3921 | 3935 | rVB2 | 1801171  | 4965254  | 18.70%  | 2.090%  |
| 98  | 26.691 | 4021 | 4030 | 4041 | rVB2 | 3649596  | 10685407 | 40.25%  | 4.498%  |
| 99  | 27.468 | 4155 | 4162 | 4173 | rVB4 | 1373151  | 4324957  | 16.29%  | 1.820%  |
| 100 | 28.273 | 4293 | 4299 | 4316 | rVB2 | 1514860  | 5280913  | 19.89%  | 2.223%  |

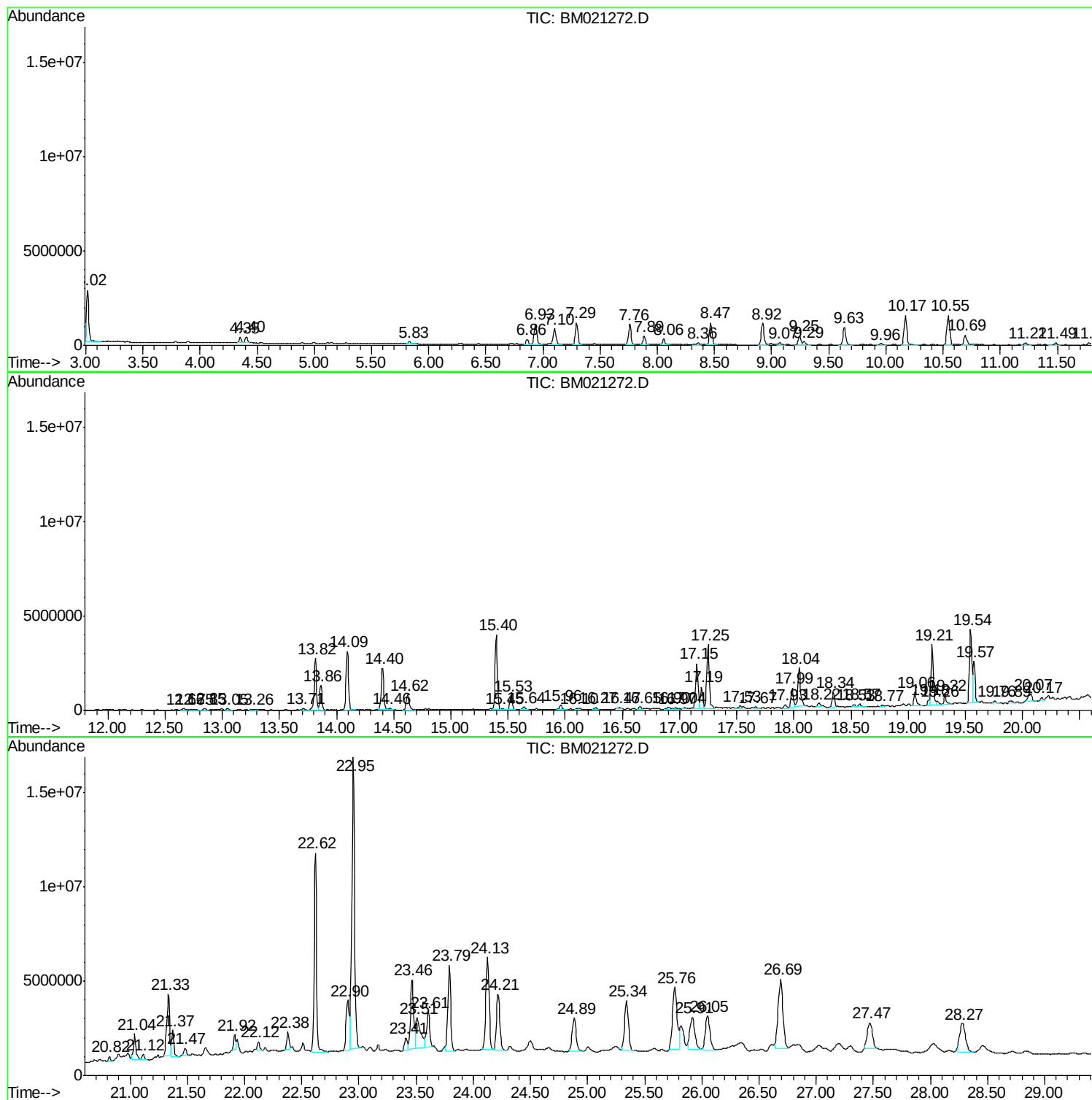
Sum of corrected areas: 237577419

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

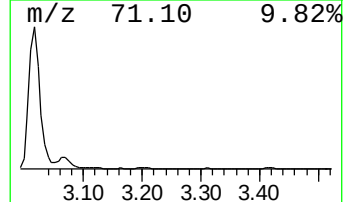
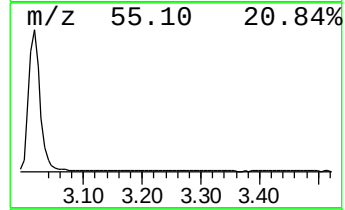
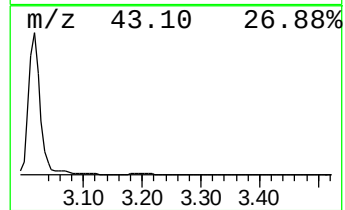
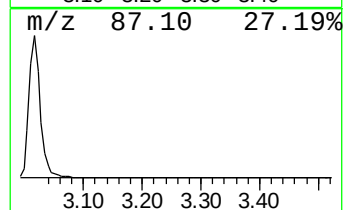
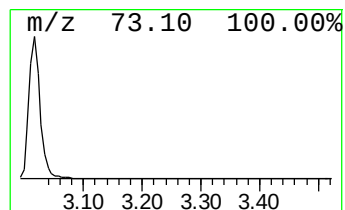
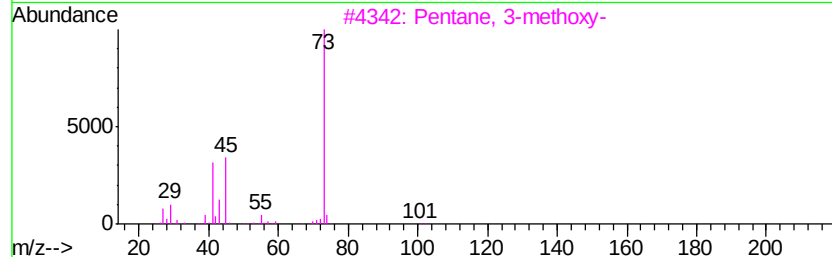
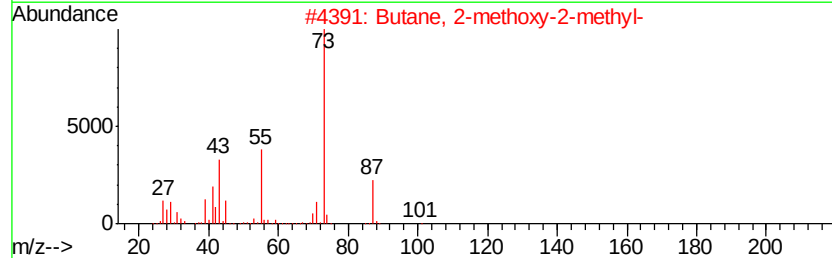
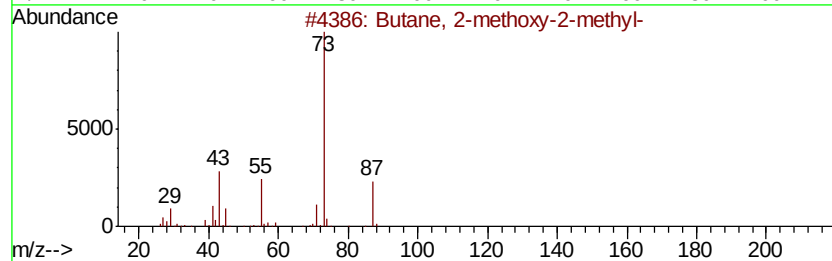
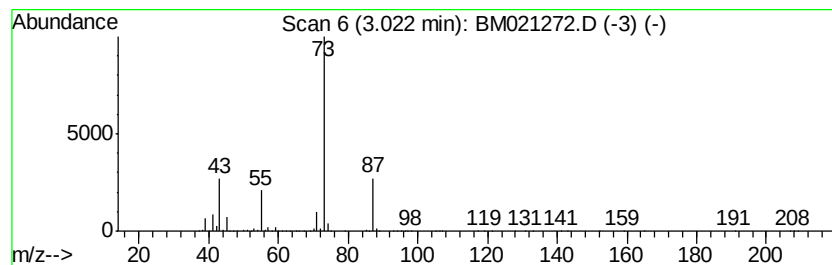
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 6

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 3.02 | 37.29 ng/ul | 3271380 | 1,4-Dichlorobenzene-d4 | 7.76 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 83   |
| 2    |    | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 64   |
| 3    |    | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 17   |
| 4    |    | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |
| 5    |    | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 9    |



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Quant Title : SVOA CALIBRATION

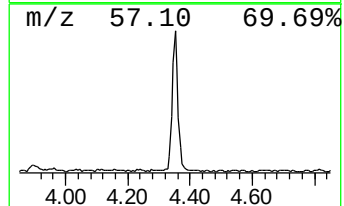
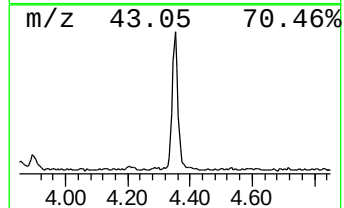
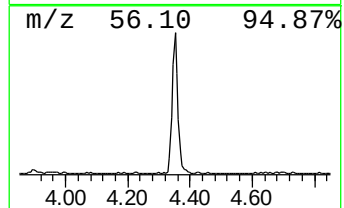
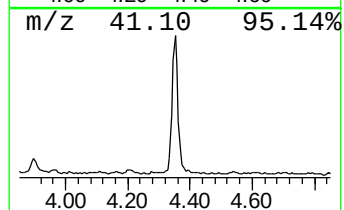
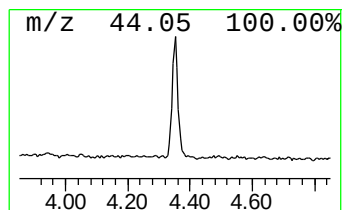
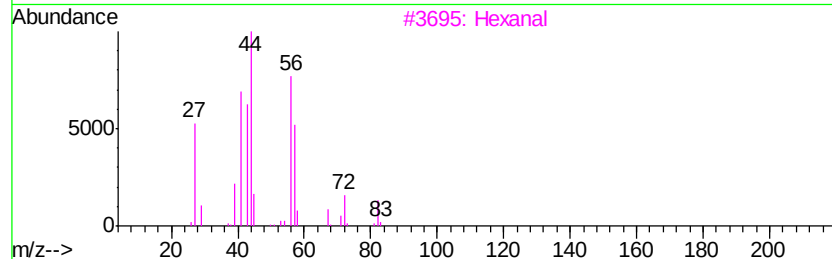
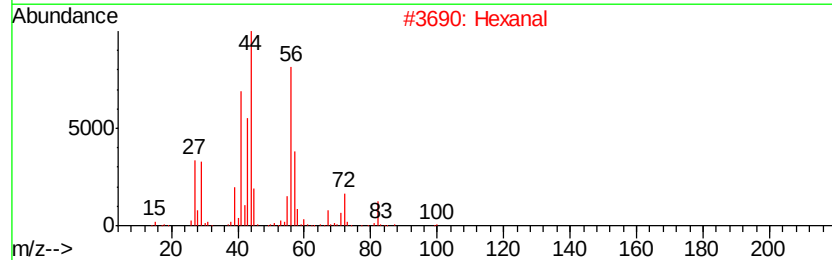
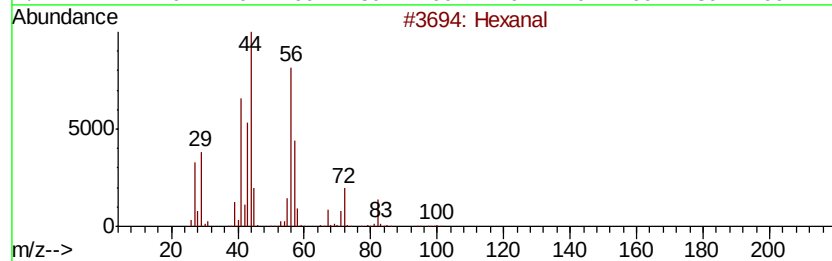
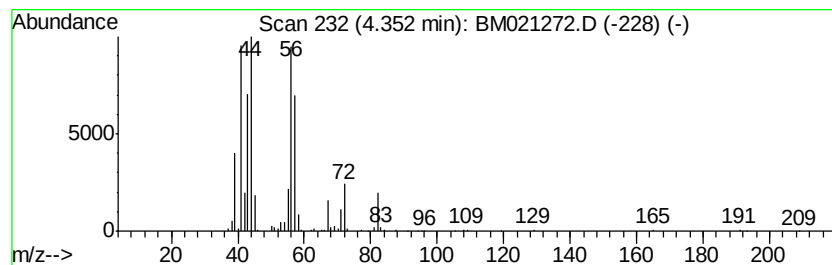
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Hexanal Concentration Rank 22

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 4.35 | 4.36 ng/ul | 382547 | 1,4-Dichlorobenzene-d4 | 7.76 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Hexanal      | 100 | C6H12O  | 000066-25-1 | 64   |
| 2    |    | Hexanal      | 100 | C6H12O  | 000066-25-1 | 59   |
| 3    |    | Hexanal      | 100 | C6H12O  | 000066-25-1 | 53   |
| 4    |    | Hexanal      | 100 | C6H12O  | 000066-25-1 | 50   |
| 5    |    | Butanal      | 72  | C4H8O   | 000123-72-8 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

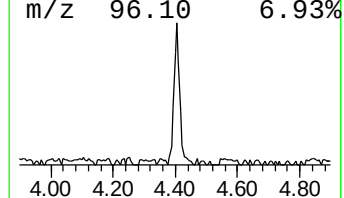
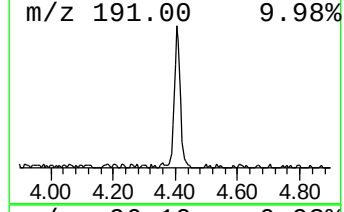
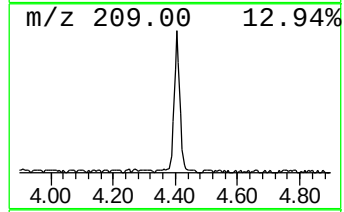
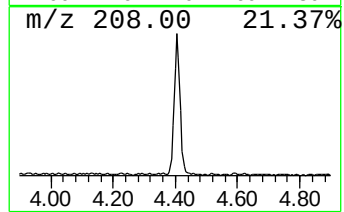
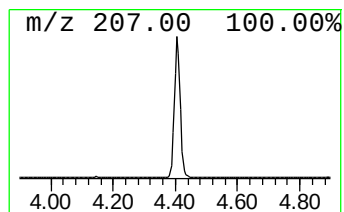
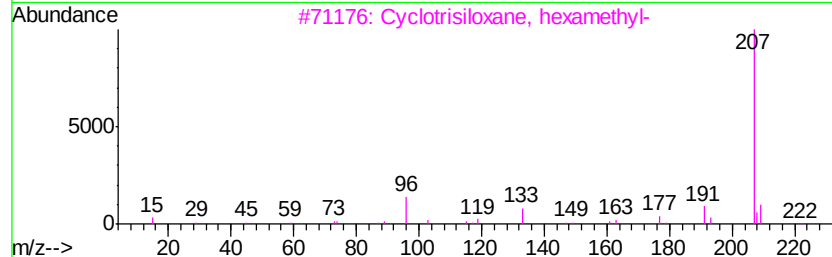
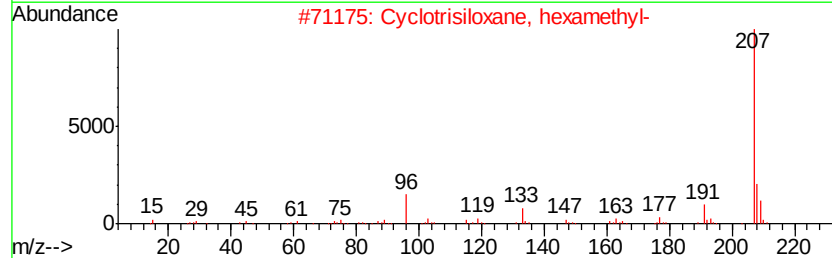
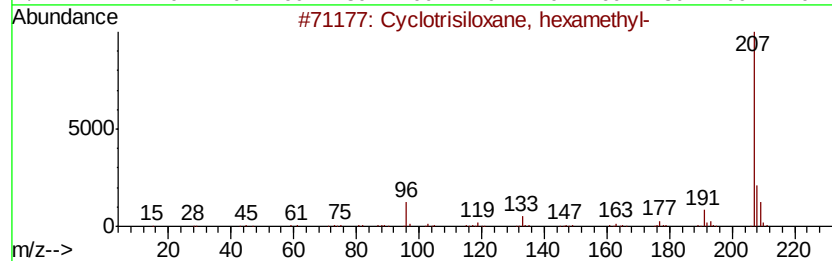
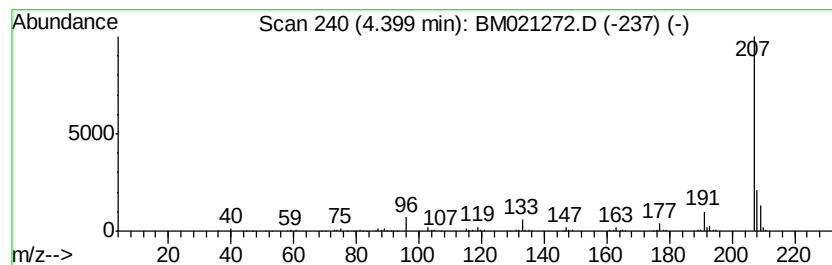
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TIC Integration Parameters: LSCINT.P

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Peak Number 3 Cyclotrisiloxane, hexamethyl- Concentration Rank 21

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 4.40 | 5.11 ng/ul | 448545 | 1,4-Dichlorobenzene-d4 | 7.76 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Cyclotrisiloxane, hexamethyl-       | 222 | C6H18O3Si3 | 000541-05-9  | 91   |
| 2    |    |   | Cyclotrisiloxane, hexamethyl-       | 222 | C6H18O3Si3 | 000541-05-9  | 90   |
| 3    |    |   | Cyclotrisiloxane, hexamethyl-       | 222 | C6H18O3Si3 | 000541-05-9  | 72   |
| 4    |    |   | 2,4,6-Cycloheptatrien-1-one, 3,5... | 250 | C13H22OSi2 | 1000161-21-8 | 64   |
| 5    |    |   | 1,3-Bis(trimethylsilyl)benzene      | 222 | C12H22Si2  | 002060-89-1  | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

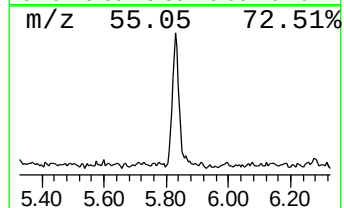
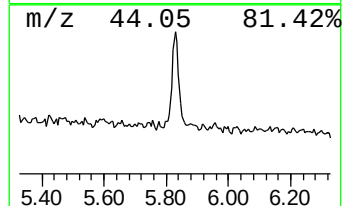
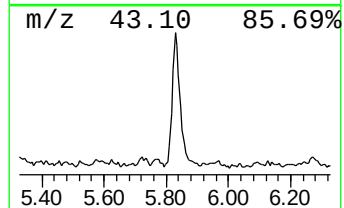
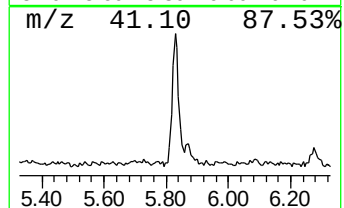
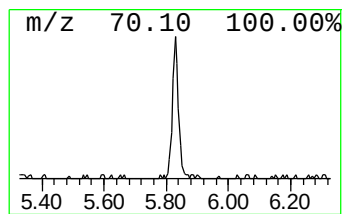
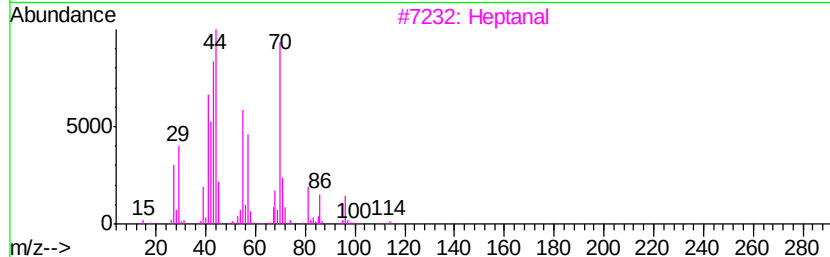
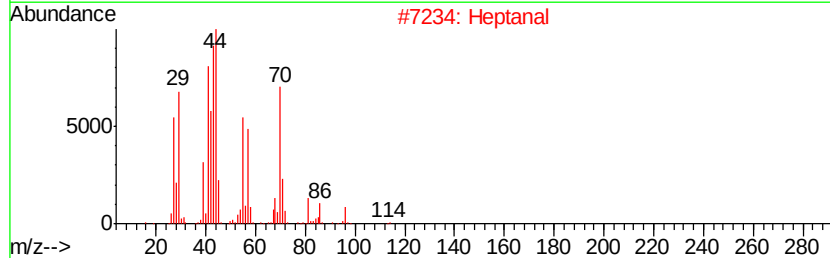
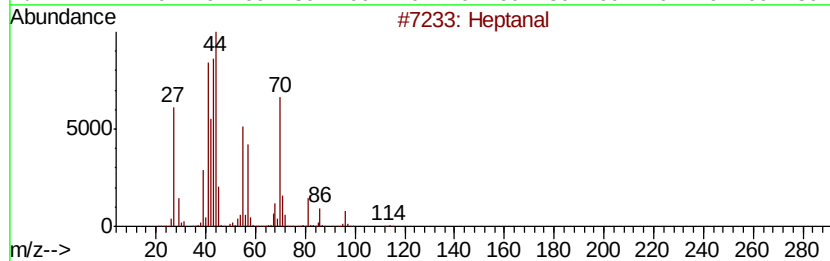
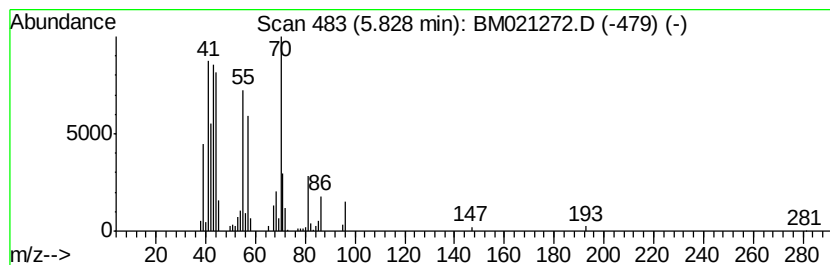
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Heptanal Concentration Rank 32

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 5.83 | 2.13 ng/ul | 187228 | 1,4-Dichlorobenzene-d4 | 7.76 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptanal                            | 114 | C7H14O  | 000111-71-7 | 72   |
| 2    |    | Heptanal                            | 114 | C7H14O  | 000111-71-7 | 64   |
| 3    |    | Heptanal                            | 114 | C7H14O  | 000111-71-7 | 64   |
| 4    |    | Oxirane, 2-(1,1-dimethylethyl)-3... | 114 | C7H14O  | 053897-30-6 | 38   |
| 5    |    | 2-Pentene, (Z)-                     | 70  | C5H10   | 000627-20-3 | 38   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

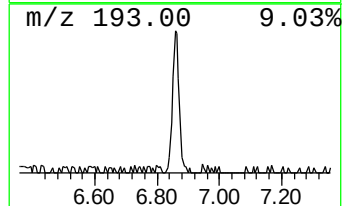
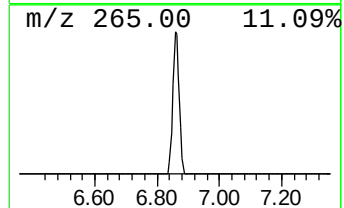
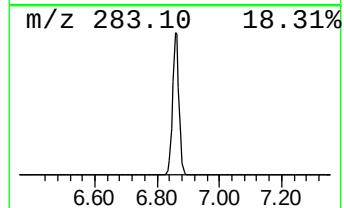
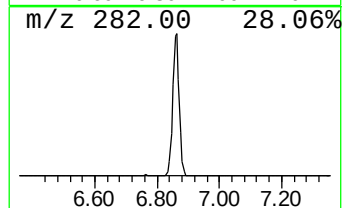
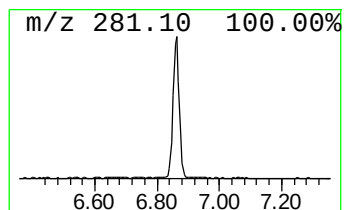
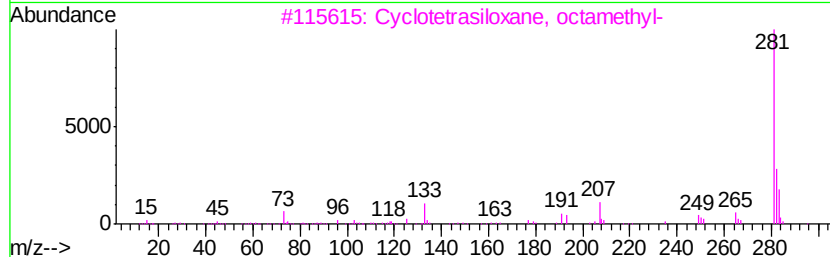
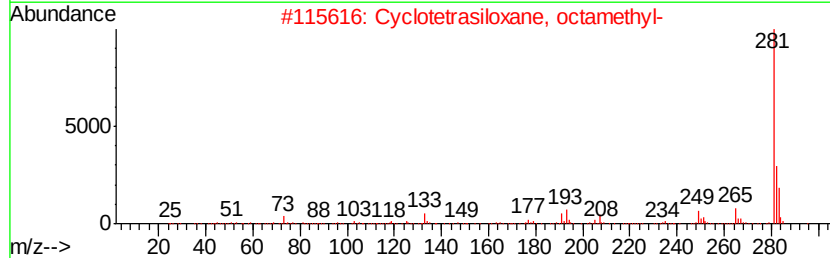
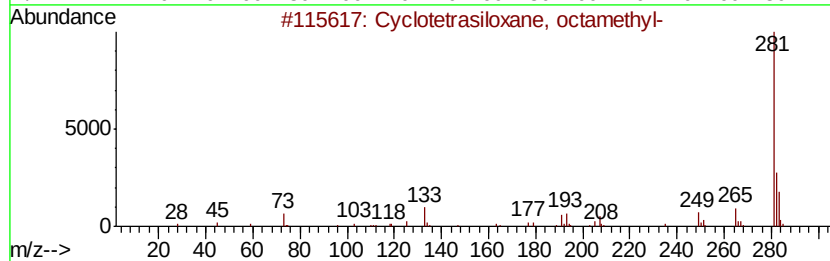
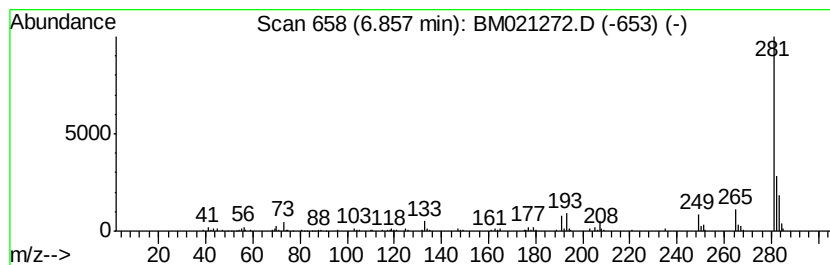
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Cyclotetrasiloxane, octamet... Concentration Rank 24

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.86 | 4.21 ng/ul | 369377 | 1,4-Dichlorobenzene-d4 | 7.76 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Cyclotetrasiloxane, octamethyl-     | 296 | C8H24O4Si4 | 000556-67-2  | 91   |
| 2    |    | Cyclotetrasiloxane, octamethyl-     | 296 | C8H24O4Si4 | 000556-67-2  | 91   |
| 3    |    | Cyclotetrasiloxane, octamethyl-     | 296 | C8H24O4Si4 | 000556-67-2  | 83   |
| 4    |    | Benzofuran-6-ol-3-one, 2-(4-etho... | 310 | C18H14O5   | 1000128-64-6 | 56   |
| 5    |    | 7H-Dibenzo[b,g]carbazole, 7-methyl- | 281 | C21H15N    | 003557-49-1  | 56   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

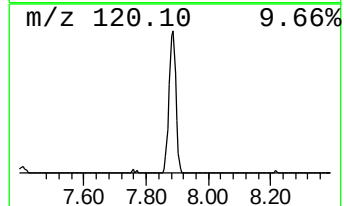
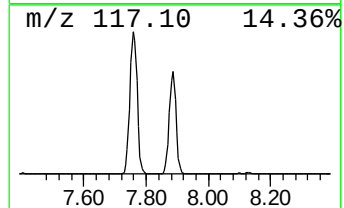
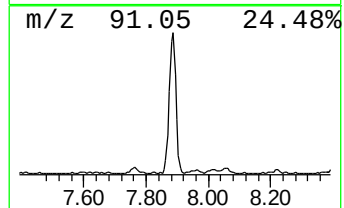
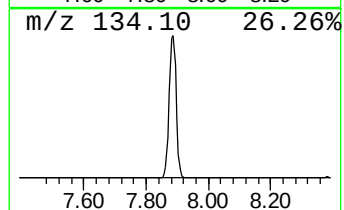
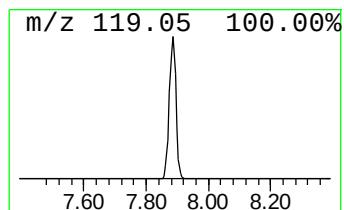
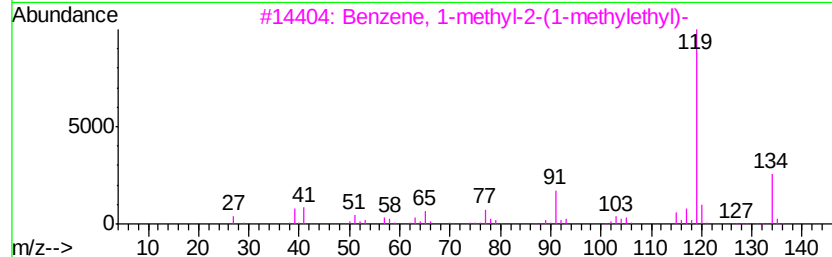
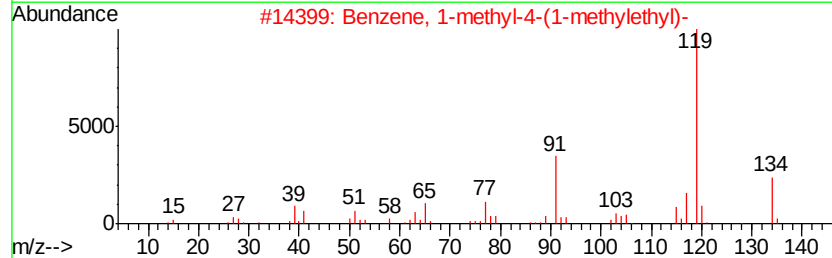
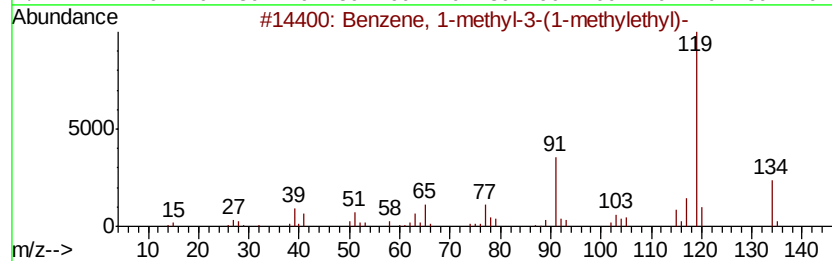
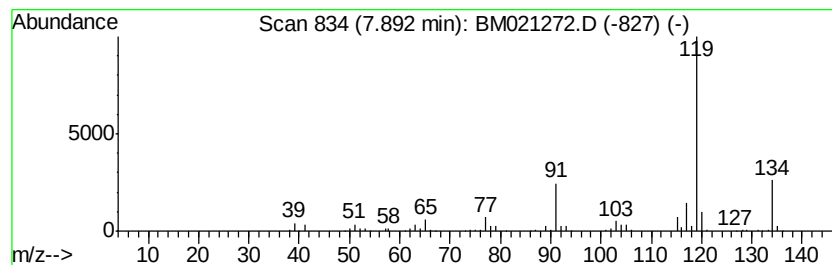
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Benzene, 1-methyl-3-(1-meth... Concentration Rank 15

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.89 | 7.45 ng/ul | 653472 | 1,4-Dichlorobenzene-d4 | 7.76 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 95   |
| 2    |    | Benzene, 1-methyl-4-(1-methylethyl)- | 134 | C10H14  | 000099-87-6 | 95   |
| 3    |    | Benzene, 1-methyl-2-(1-methylethyl)- | 134 | C10H14  | 000527-84-4 | 94   |
| 4    |    | Benzene, 1-methyl-2-(1-methylethyl)- | 134 | C10H14  | 000527-84-4 | 94   |
| 5    |    | Benzene, 1-methyl-2-(1-methylethyl)- | 134 | C10H14  | 000527-84-4 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

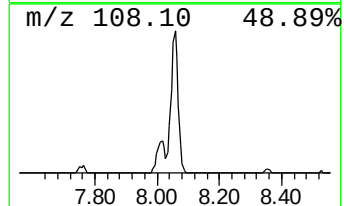
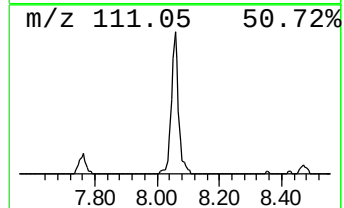
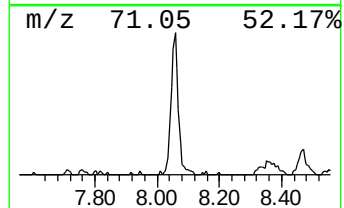
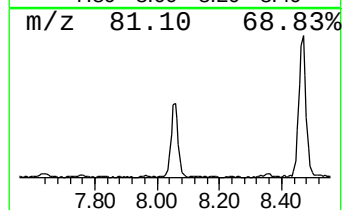
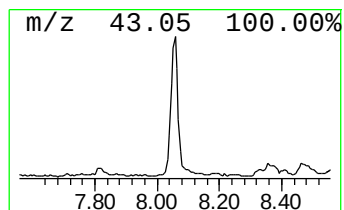
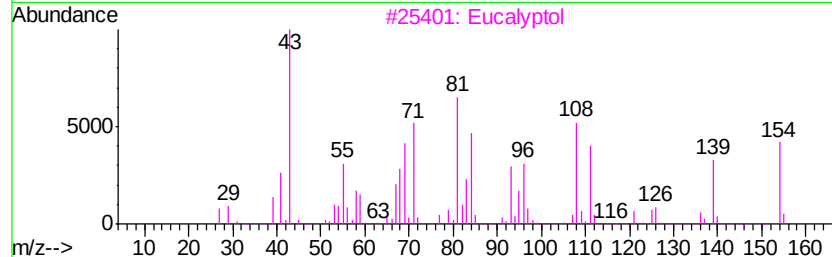
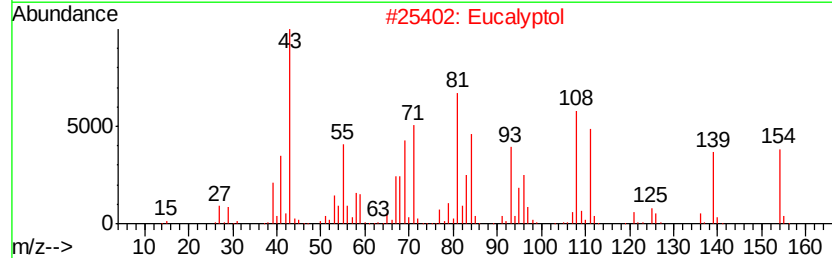
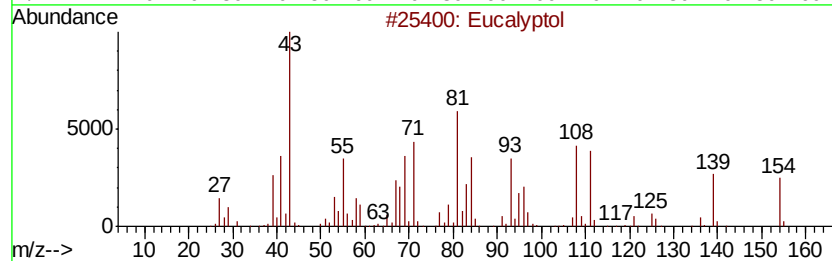
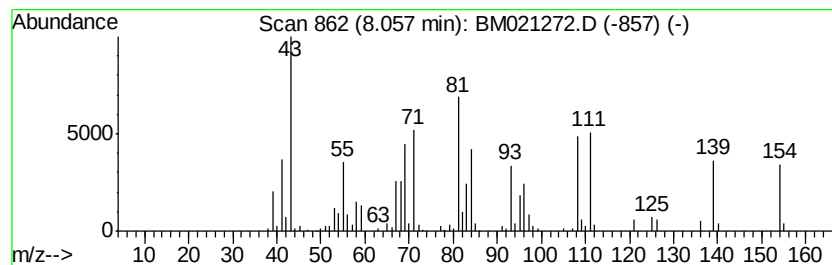
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Eucalyptol Concentration Rank 20

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.06 | 5.42 ng/ul | 475433 | 1,4-Dichlorobenzene-d4 | 7.76 |

| Hit# of | 5          | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|------------|--------------|-----|---------|-------------|------|
| 1       | Eucalyptol |              | 154 | C10H18O | 000470-82-6 | 99   |
| 2       | Eucalyptol |              | 154 | C10H18O | 000470-82-6 | 99   |
| 3       | Eucalyptol |              | 154 | C10H18O | 000470-82-6 | 97   |
| 4       | Eucalyptol |              | 154 | C10H18O | 000470-82-6 | 93   |
| 5       | Eucalyptol |              | 154 | C10H18O | 000470-82-6 | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

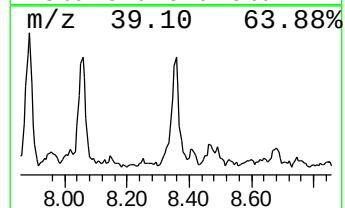
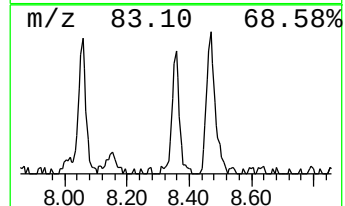
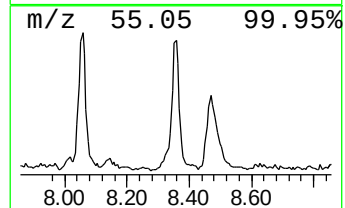
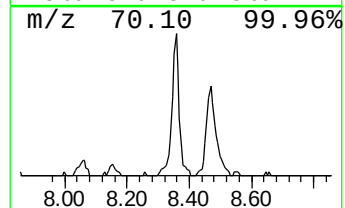
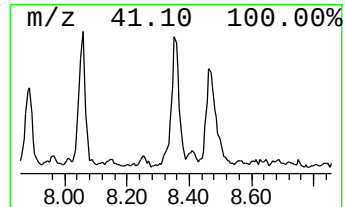
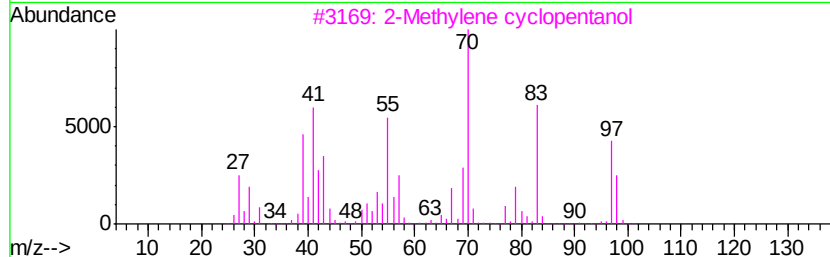
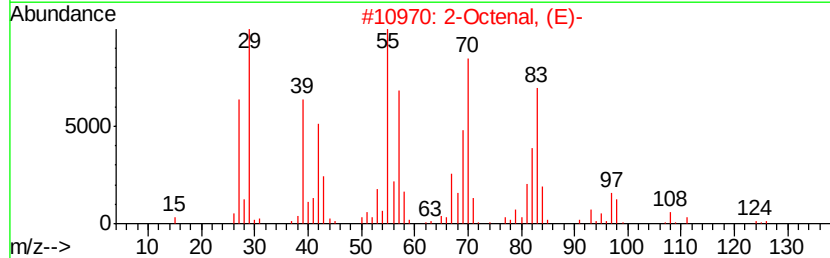
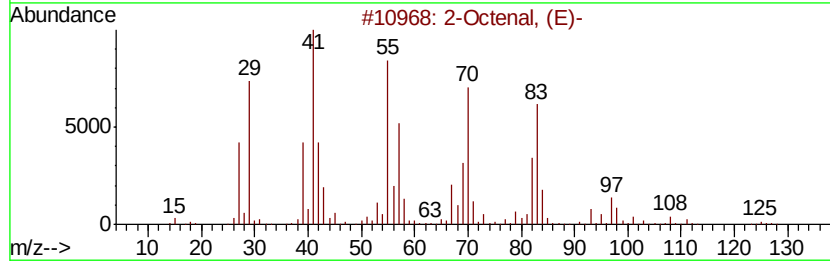
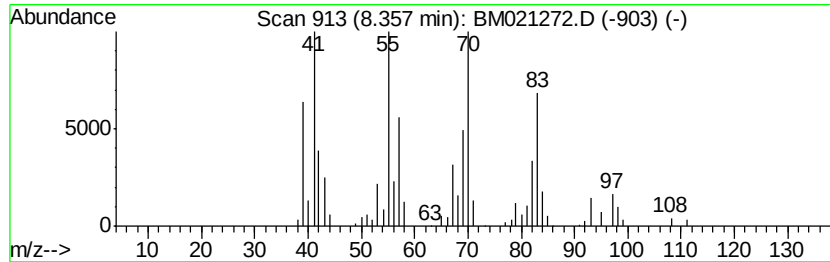
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 2-Octenal, (E)- Concentration Rank 28

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.36 | 2.28 ng/ul | 200085 | 1,4-Dichlorobenzene-d4 | 7.76 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Octenal, (E)-           | 126 | C8H14O  | 002548-87-0 | 64   |
| 2    |    | 2-Octenal, (E)-           | 126 | C8H14O  | 002548-87-0 | 59   |
| 3    |    | 2-Methylene cyclopentanol | 98  | C6H10O  | 020461-31-8 | 47   |
| 4    |    | 2-Pentene, (E)-           | 70  | C5H10   | 000646-04-8 | 43   |
| 5    |    | Butane, 2-cyclopropyl-    | 98  | C7H14   | 005750-02-7 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

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Quant Title : SVOA CALIBRATION

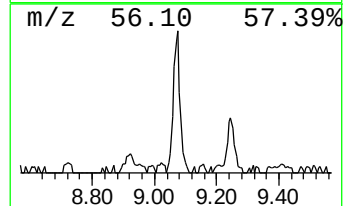
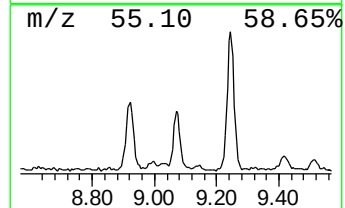
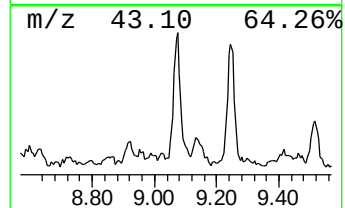
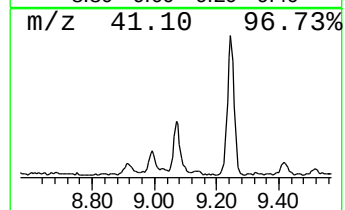
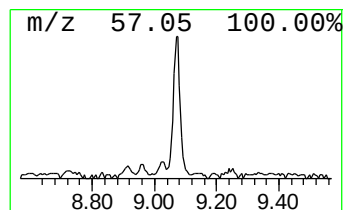
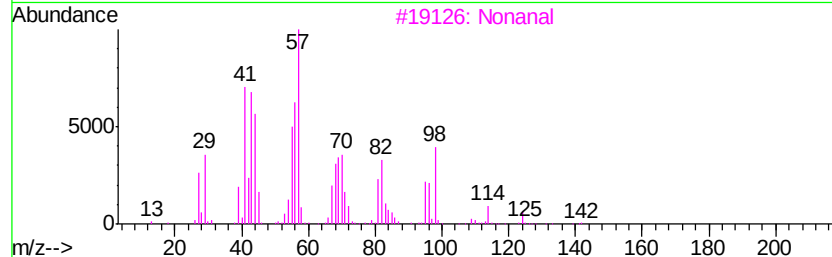
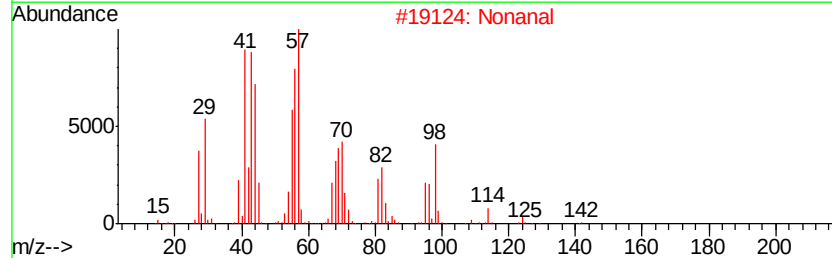
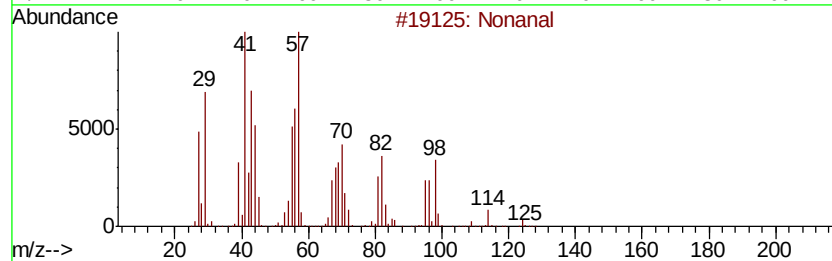
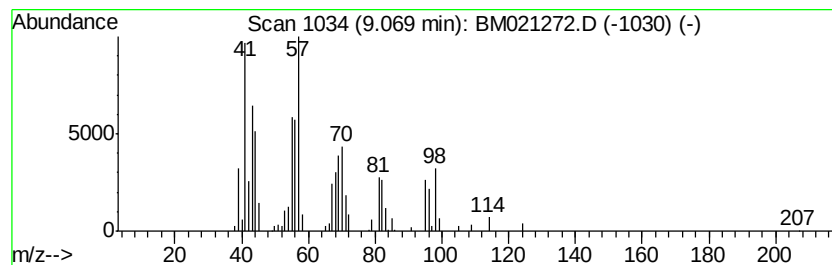
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TIC Integration Parameters: LSCINT.P

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Peak Number 9 Nonanal Concentration Rank 29

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 9.07 | 2.26 ng/ul | 197831 | 1,4-Dichlorobenzene-d4 | 7.76 |

| Hit# of | 5                               | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|---------------------------------|--------------|-----|---------|-------------|------|
| 1       | Nonanal                         |              | 142 | C9H18O  | 000124-19-6 | 91   |
| 2       | Nonanal                         |              | 142 | C9H18O  | 000124-19-6 | 90   |
| 3       | Nonanal                         |              | 142 | C9H18O  | 000124-19-6 | 72   |
| 4       | 2-Propen-1-amine, N-2-propenyl- |              | 97  | C6H11N  | 000124-02-7 | 25   |
| 5       | Formic acid, octyl ester        |              | 158 | C9H18O2 | 000112-32-3 | 22   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

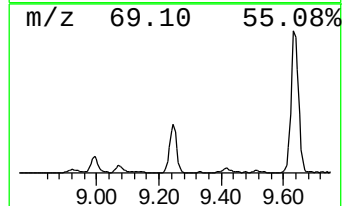
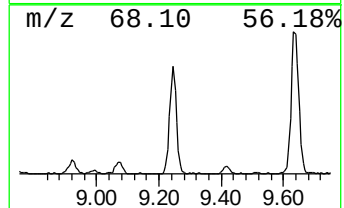
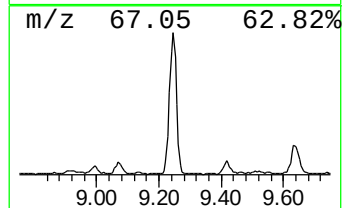
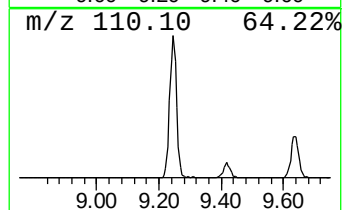
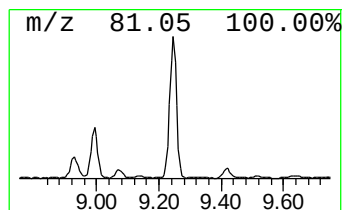
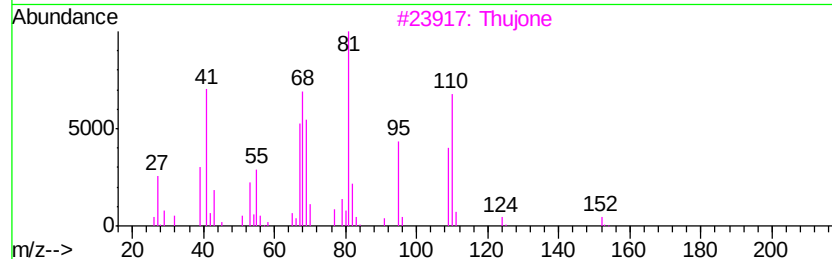
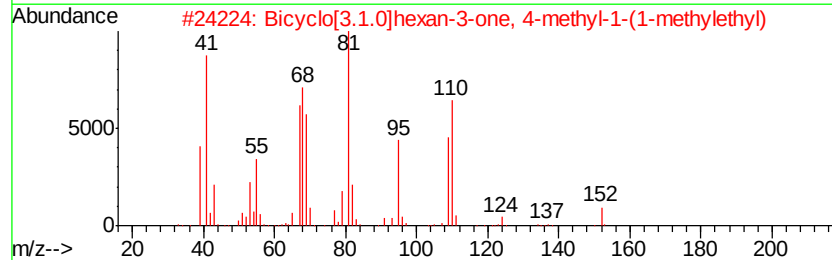
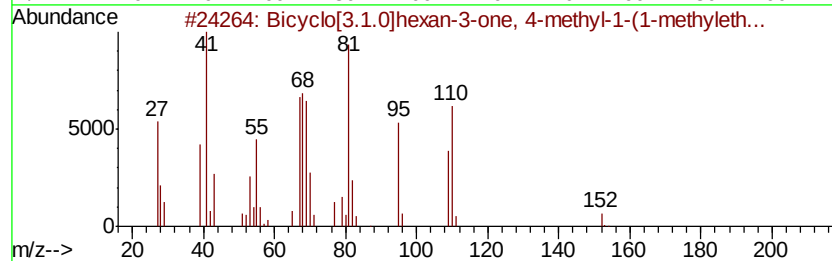
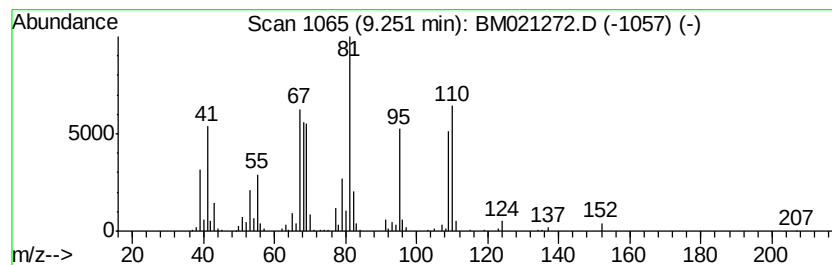
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 Bicyclo[3.1.0]hexan-3-one, ... Concentration Rank 19

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.25 | 5.91 ng/ul | 773173 | Naphthalene-d8   | 10.55 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Bicyclo[3.1.0]hexan-3-one, 4-met... | 152 | C10H16O | 000471-15-8 | 93   |
| 2    |    |   | Bicyclo[3.1.0]hexan-3-one, 4-met... | 152 | C10H16O | 001125-12-8 | 83   |
| 3    |    |   | Thujone                             | 152 | C10H16O | 000546-80-5 | 81   |
| 4    |    |   | Thujone                             | 152 | C10H16O | 000546-80-5 | 81   |
| 5    |    |   | 3,5-Octadiene, (Z,Z)-               | 110 | C8H14   | 007348-80-3 | 74   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

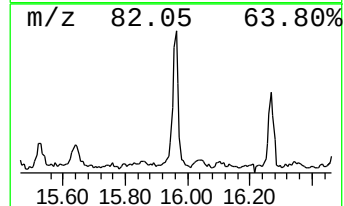
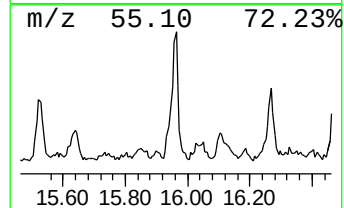
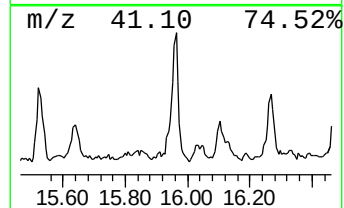
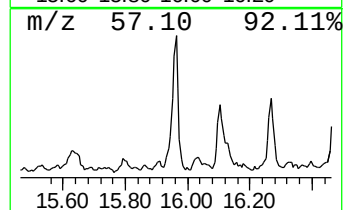
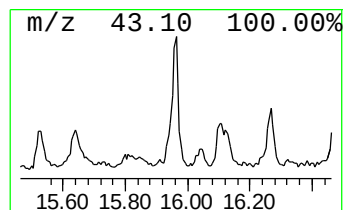
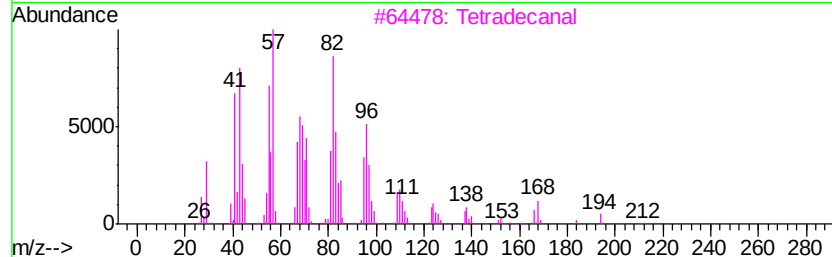
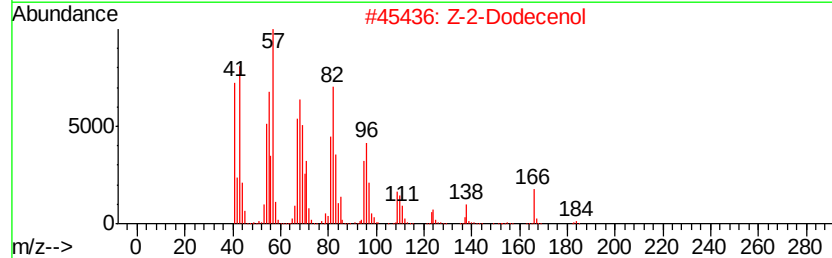
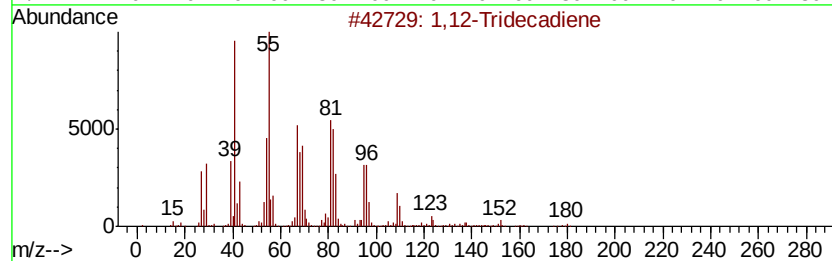
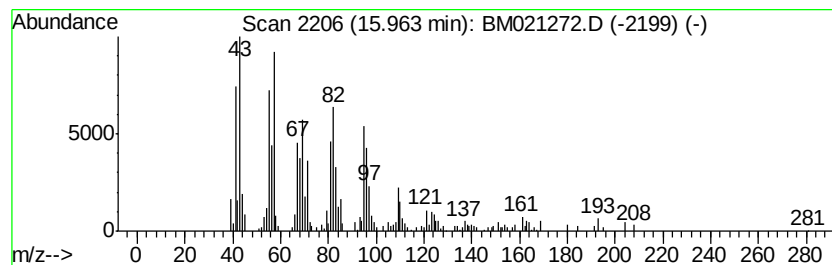
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TIC Integration Parameters: LSCINT.P

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Peak Number 11 1,12-Tridecadiene Concentration Rank 30

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.96 | 2.18 ng/ul | 380668 | Phenanthrene-d10 | 17.15 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | 1,12-Tridecadiene    | 180 | C13H24  | 021964-48-7 | 76   |
| 2    |    |   | Z-2-Dodecenol        | 184 | C12H24O | 069064-36-4 | 55   |
| 3    |    |   | Tetradecanal         | 212 | C14H28O | 000124-25-4 | 53   |
| 4    |    |   | Oxirane, tetradecyl- | 240 | C16H32O | 007320-37-8 | 53   |
| 5    |    |   | Pentadecanal-        | 226 | C15H30O | 002765-11-9 | 52   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

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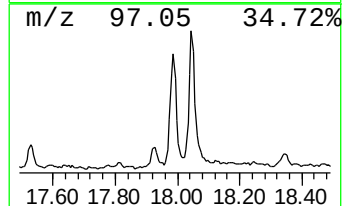
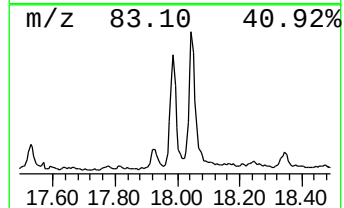
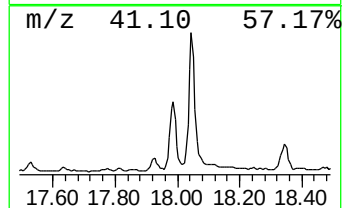
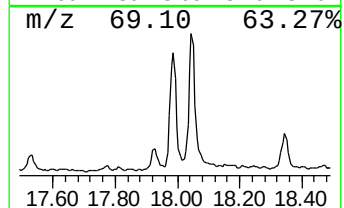
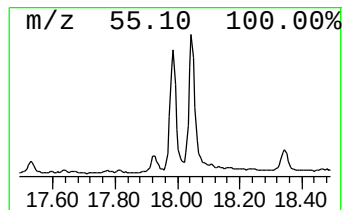
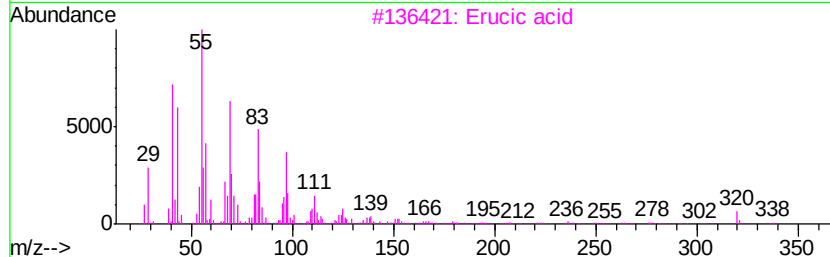
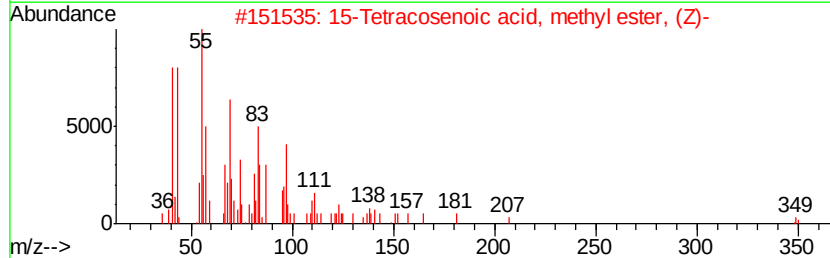
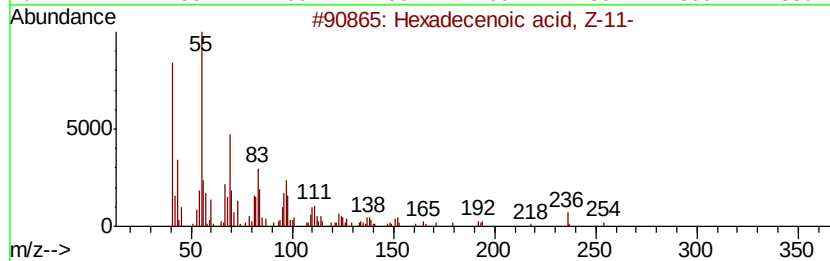
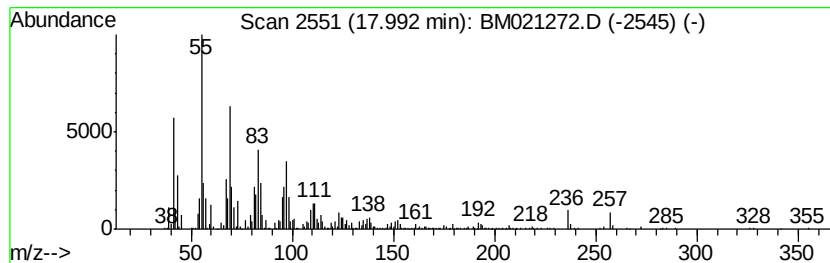
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 Hexadecenoic acid, Z-11- Concentration Rank 14

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 17.99 | 8.28 ng/ul | 1446930 | Phenanthrene-d10 | 17.15 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Hexadecenoic acid, Z-11-            | 254 | C16H30O2 | 002416-20-8  | 92   |
| 2    |    |   | 15-Tetracosenoic acid, methyl es... | 380 | C25H48O2 | 002733-88-2  | 81   |
| 3    |    |   | Erucic acid                         | 338 | C22H42O2 | 000112-86-7  | 70   |
| 4    |    |   | E-9-Tetradecenoic acid              | 226 | C14H26O2 | 1000131-35-8 | 53   |
| 5    |    |   | 1-Decene                            | 140 | C10H20   | 000872-05-9  | 53   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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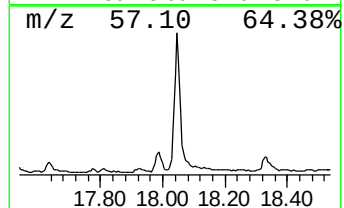
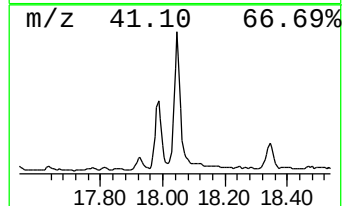
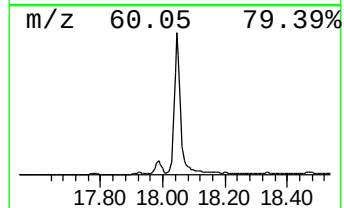
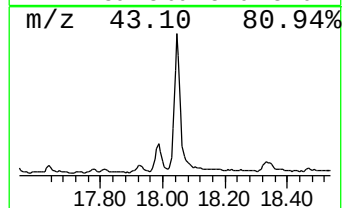
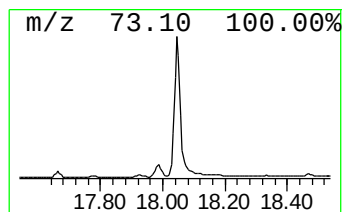
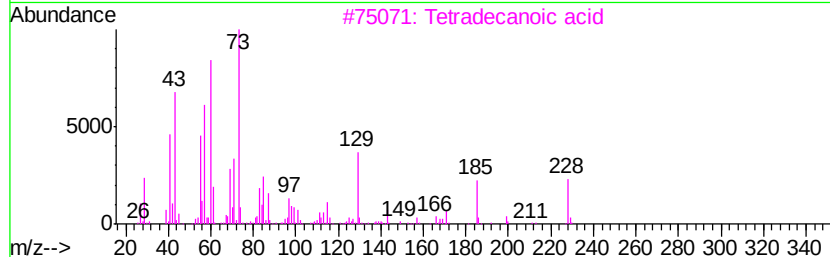
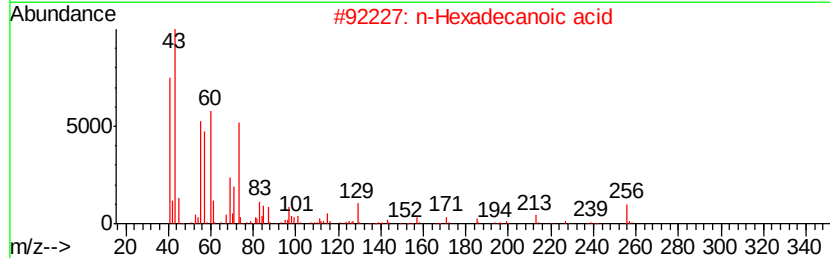
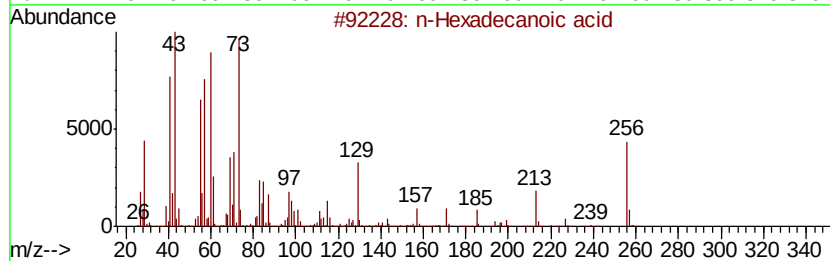
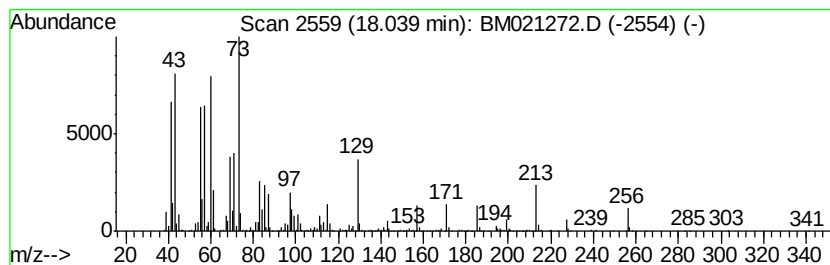
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 n-Hexadecanoic acid Concentration Rank 13

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 18.04 | 18.39 ng/ul | 3215320 | Phenanthrene-d10 | 17.15 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98   |
| 2    |    | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 97   |
| 3    |    | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 94   |
| 4    |    | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 90   |
| 5    |    | n-Decanoic acid     | 172 | C10H20O2 | 000334-48-5 | 76   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Title : SVOA CALIBRATION

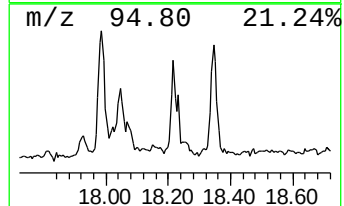
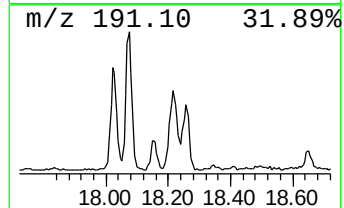
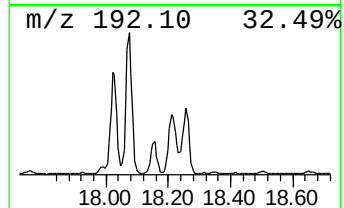
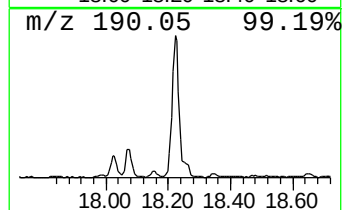
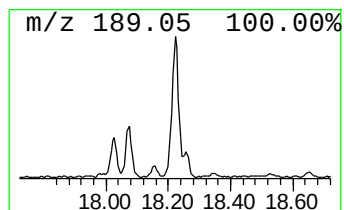
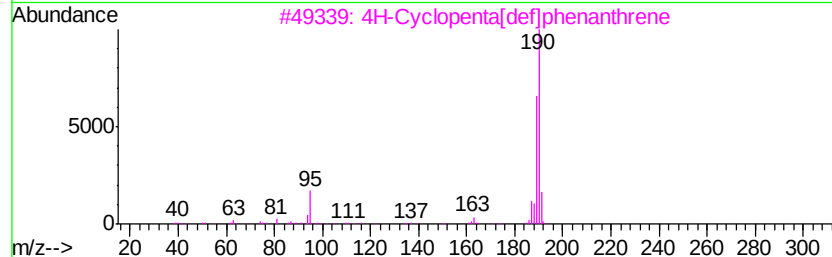
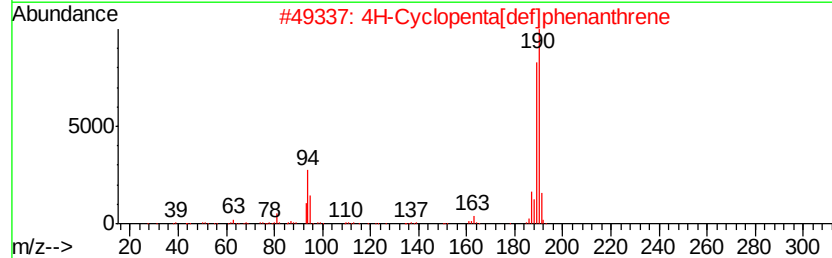
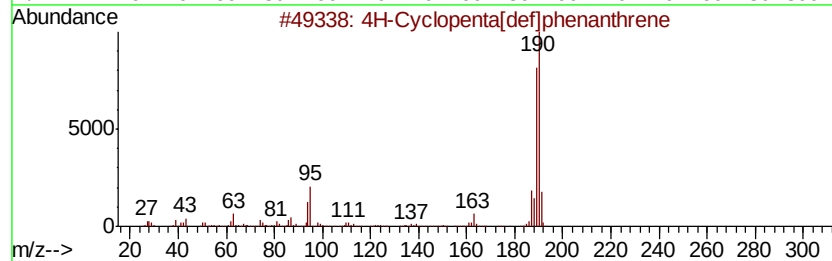
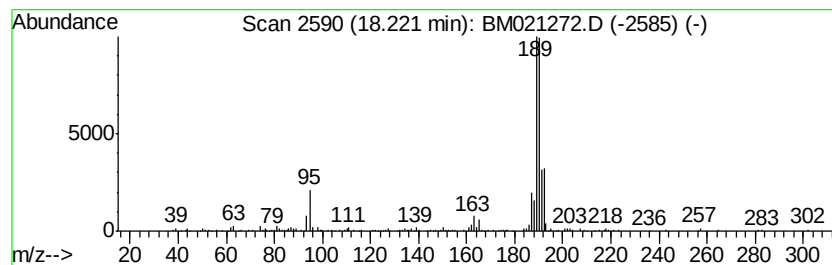
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 4H-Cyclopenta[def]phenanthrene Concentration Rank 31

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.22 | 2.15 ng/ul | 375196 | Phenanthrene-d10 | 17.15 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10    | 000203-64-5 | 90   |
| 2    |    | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10    | 000203-64-5 | 87   |
| 3    |    | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10    | 000203-64-5 | 50   |
| 4    |    | Pyrazole-4-carboxaldehyde, 3-(4-... | 190 | C10H7FN2O | 306936-57-2 | 42   |
| 5    |    | 2-Naphthaldehyde, 1,2,3,4-tetra...  | 190 | C12H14O2  | 002472-02-8 | 18   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

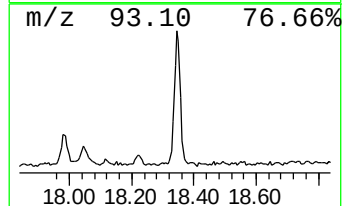
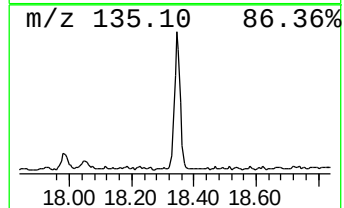
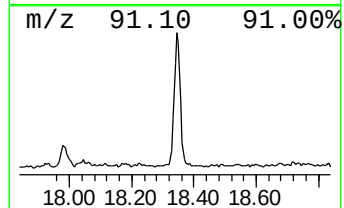
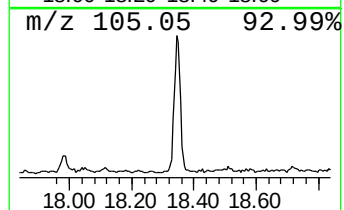
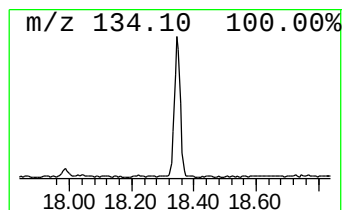
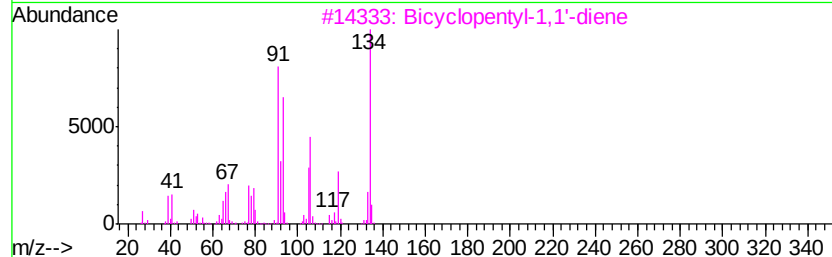
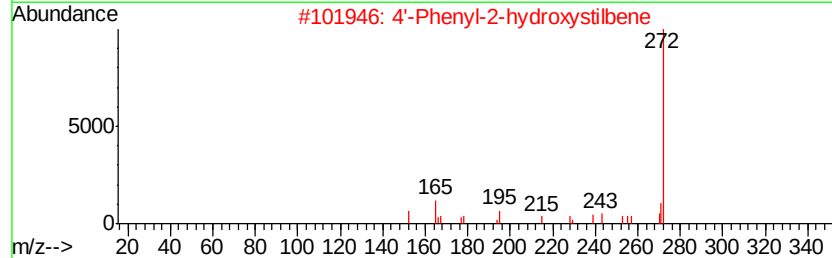
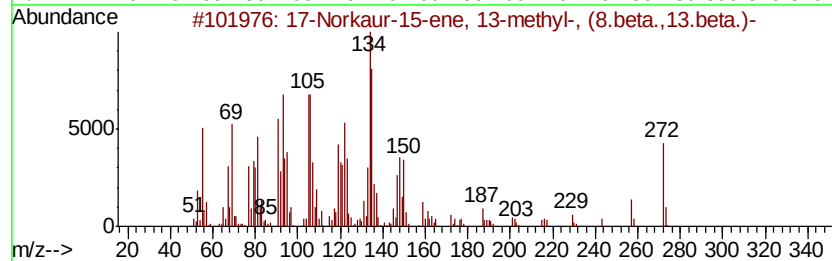
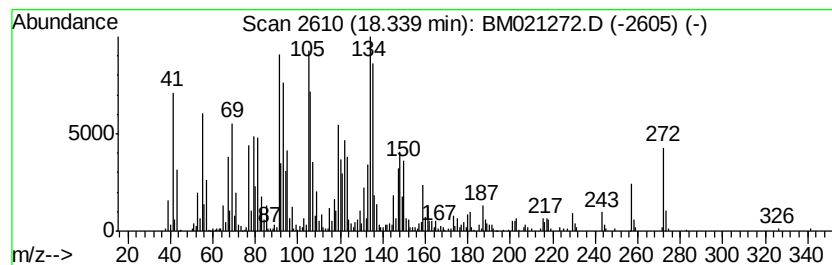
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TIC Integration Parameters: LSCINT.P

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Peak Number 15 17-Norkaur-15-ene, 13-methy... Concentration Rank 16

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 18.34 | 6.92 ng/ul | 1209060 | Phenanthrene-d10 | 17.15 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#         | Qual |
|------|----|--------------------------------------|-----|---------|--------------|------|
| 1    | 5  | 17-Norkaur-15-ene, 13-methyl-, (...) | 272 | C20H32  | 003564-54-3  | 99   |
| 2    |    | 4'-Phenyl-2-hydroxystilbene          | 272 | C20H16O | 1000148-46-6 | 72   |
| 3    |    | Bicyclopentyl-1,1'-diene             | 134 | C10H14  | 000934-02-1  | 35   |
| 4    |    | (E,E)-7,11,15-Trimethyl-3-methyl...  | 272 | C20H32  | 070901-63-2  | 35   |
| 5    |    | N-Benzylformamide                    | 135 | C8H9NO  | 006343-54-0  | 30   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

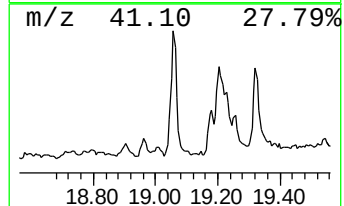
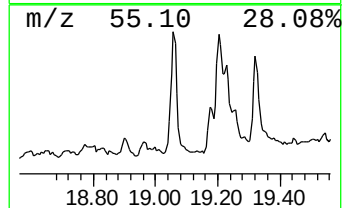
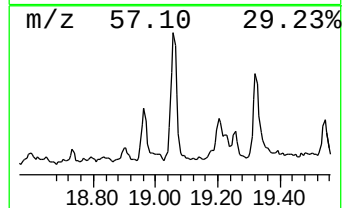
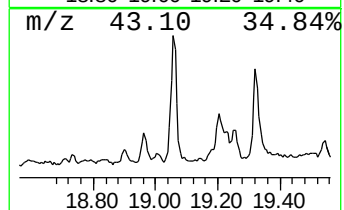
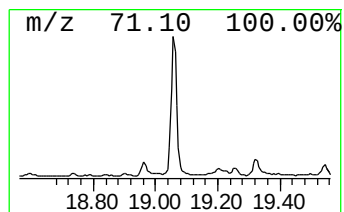
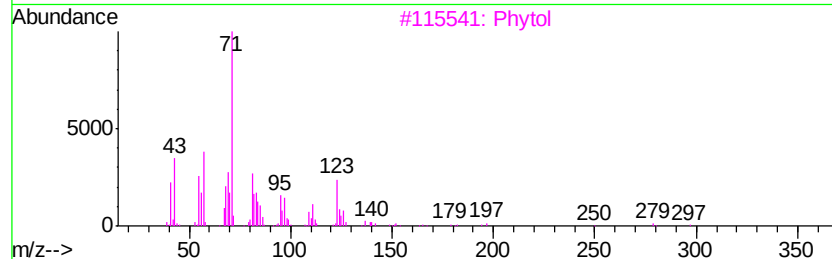
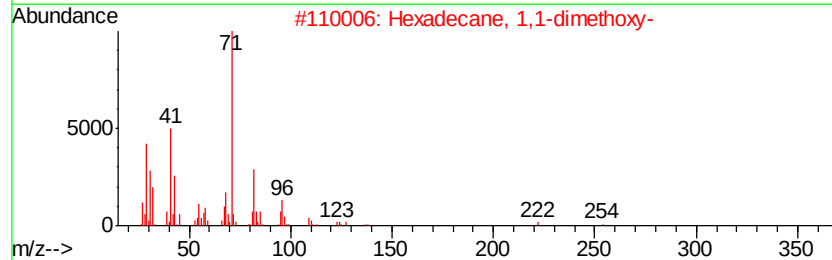
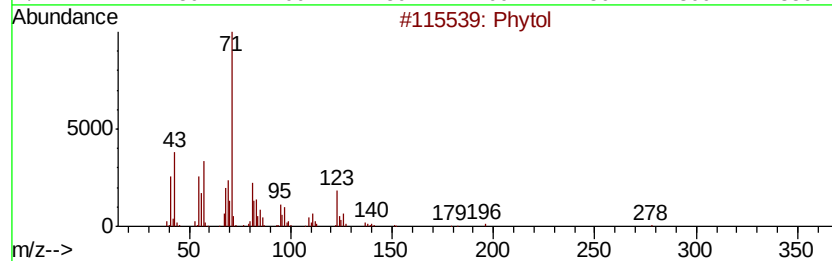
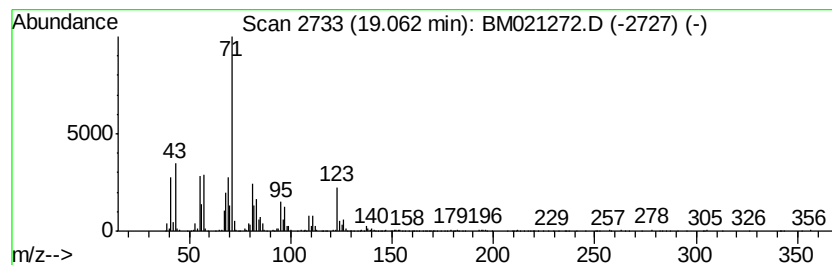
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Phytol Concentration Rank 18

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 19.06 | 6.61 ng/ul | 1156250 | Phenanthrene-d10 | 17.15 |

| Hit# | of                         | 5 | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|------|----------------------------|---|--------------|-----|----------|-------------|------|
| 1    | Phytol                     |   |              | 296 | C20H40O  | 000150-86-7 | 91   |
| 2    | Hexadecane, 1,1-dimethoxy- |   |              | 286 | C18H38O2 | 002791-29-9 | 59   |
| 3    | Phytol                     |   |              | 296 | C20H40O  | 000150-86-7 | 58   |
| 4    | Phytol                     |   |              | 296 | C20H40O  | 000150-86-7 | 58   |
| 5    | Phytol                     |   |              | 296 | C20H40O  | 000150-86-7 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

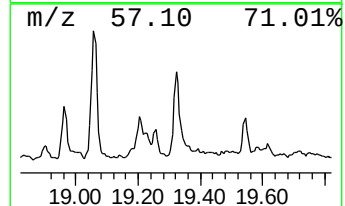
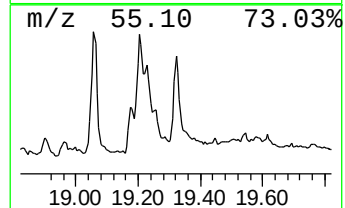
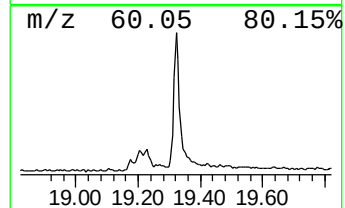
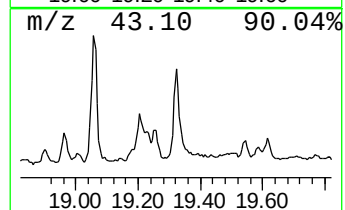
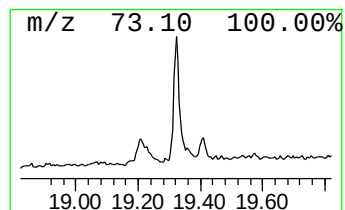
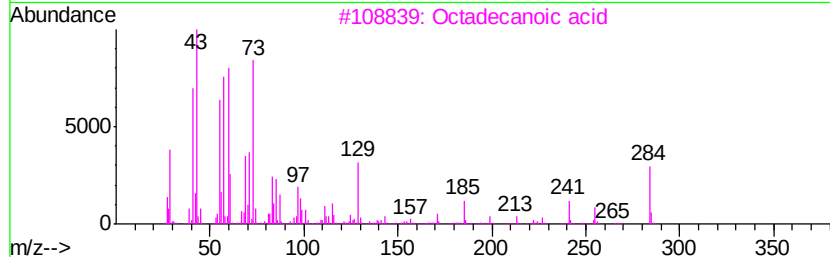
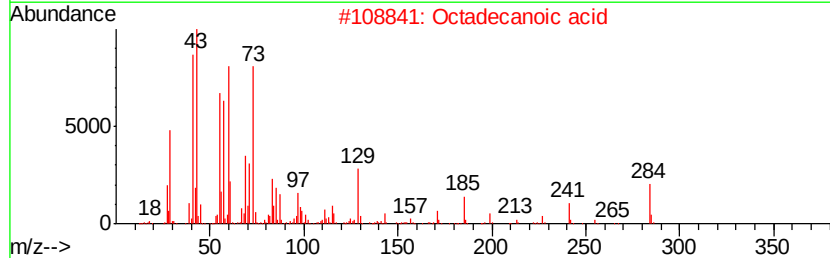
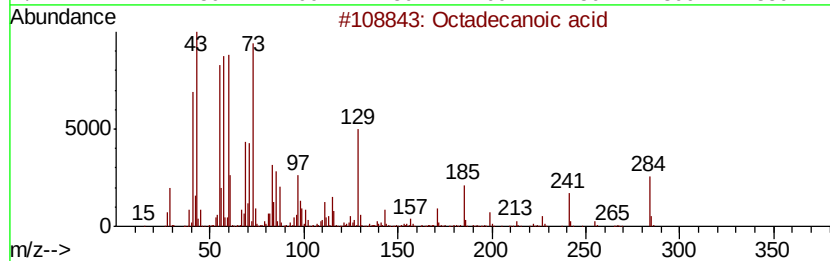
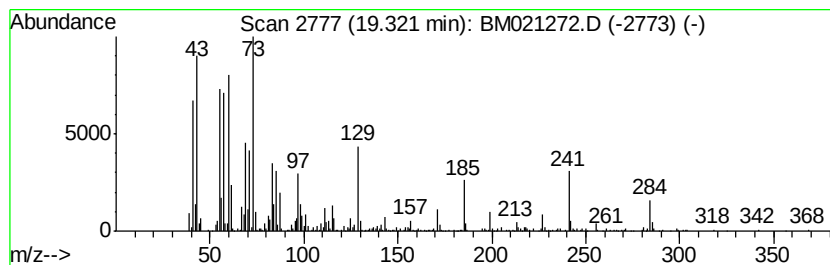
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Octadecanoic acid Concentration Rank 27

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.32 | 2.45 ng/ul | 720527 | Chrysene-d12     | 21.34 |

| 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 | 94   |
| 4    |    |   | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 93   |
| 5    |    |   | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

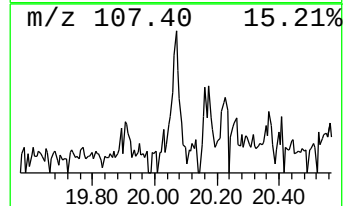
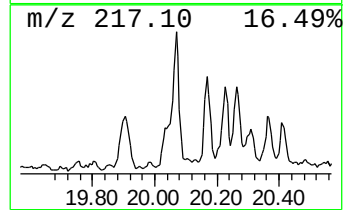
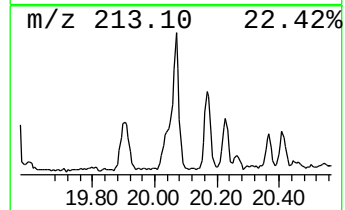
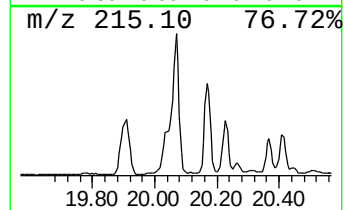
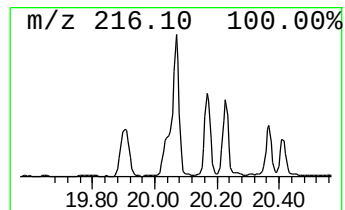
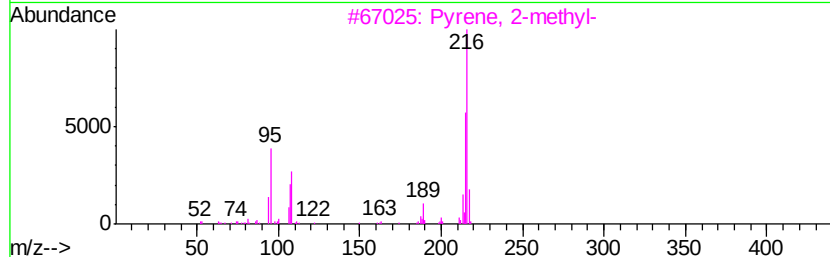
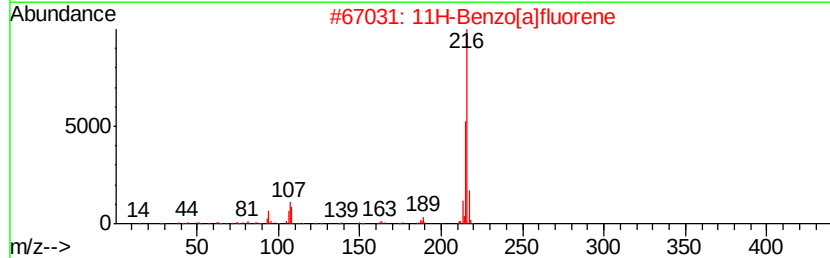
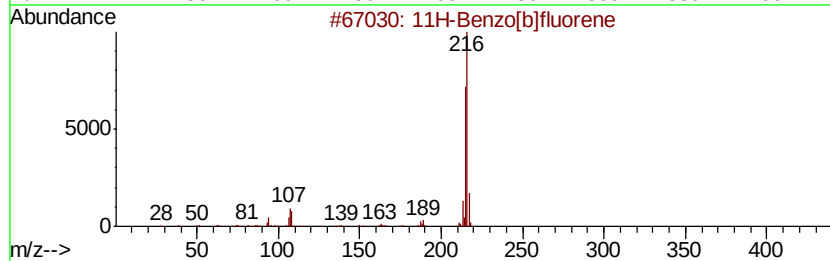
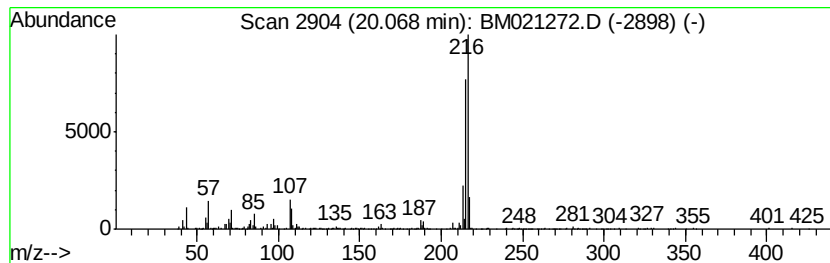
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 11H-Benzo[b]fluorene Concentration Rank 26

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.07 | 3.15 ng/ul | 927135 | Chrysene-d12     | 21.34 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | 11H-Benzo[b]fluorene | 216 | C17H12  | 000243-17-4 | 76   |
| 2    |    |   | 11H-Benzo[a]fluorene | 216 | C17H12  | 000238-84-6 | 76   |
| 3    |    |   | Pvrene, 2-methyl-    | 216 | C17H12  | 003442-78-2 | 76   |
| 4    |    |   | 11H-Benzo[a]fluorene | 216 | C17H12  | 000238-84-6 | 70   |
| 5    |    |   | 11H-Benzo[b]fluorene | 216 | C17H12  | 000243-17-4 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

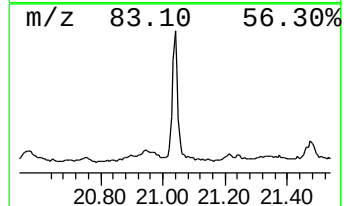
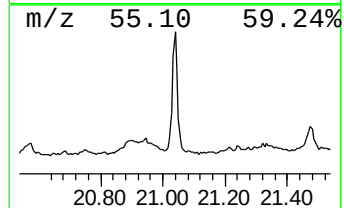
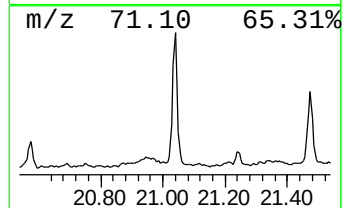
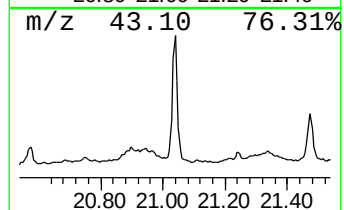
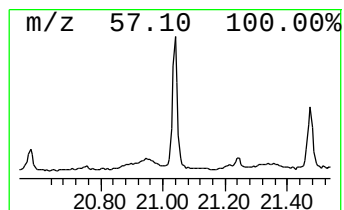
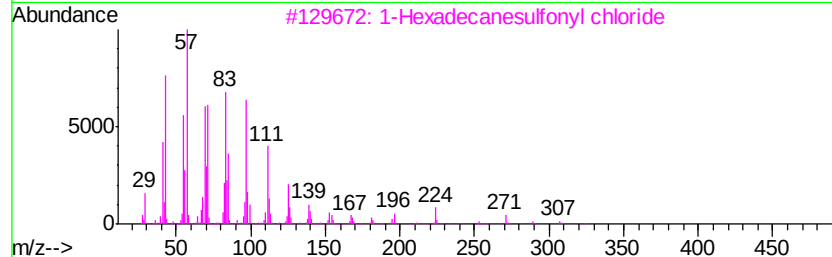
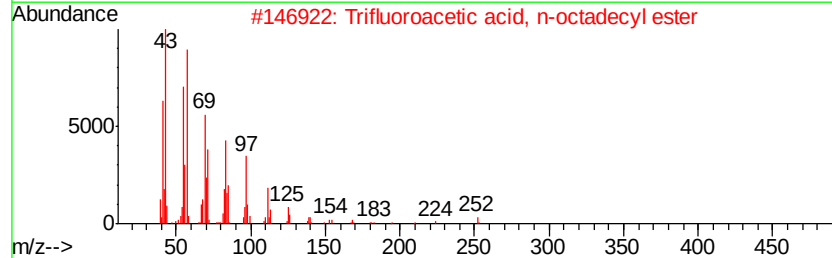
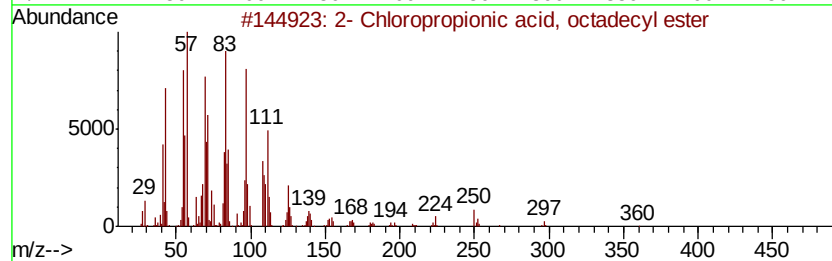
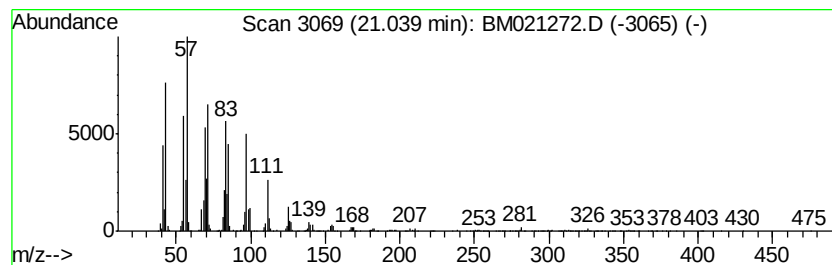
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 2- Chloropropionic acid, oc... Concentration Rank 17

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 21.04 | 6.68 ng/ul | 1966740 | Chrysene-d12     | 21.34 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | 2- Chloropropionic acid, octadec... | 360 | C21H41ClO2  | 088104-31-8  | 89   |
| 2    |    |   | Trifluoroacetic acid, n-octadecy... | 366 | C20H37F3O2  | 079392-43-1  | 87   |
| 3    |    |   | 1-Hexadecanesulfonyl chloride       | 324 | C16H33ClO2S | 038775-38-1  | 87   |
| 4    |    |   | Pentafluoropropionic acid, hepta... | 402 | C20H35F5O2  | 1000283-04-2 | 76   |
| 5    |    |   | 9-Tricosene, (Z)-                   | 322 | C23H46      | 027519-02-4  | 76   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

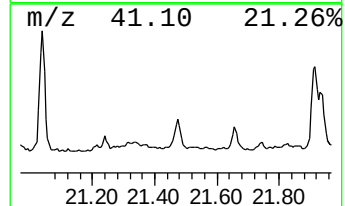
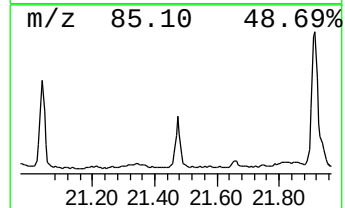
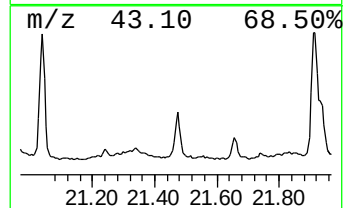
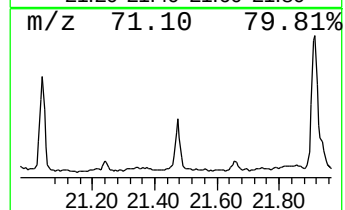
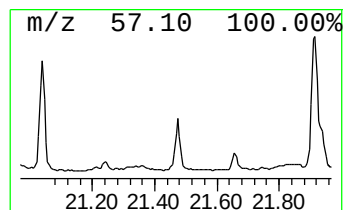
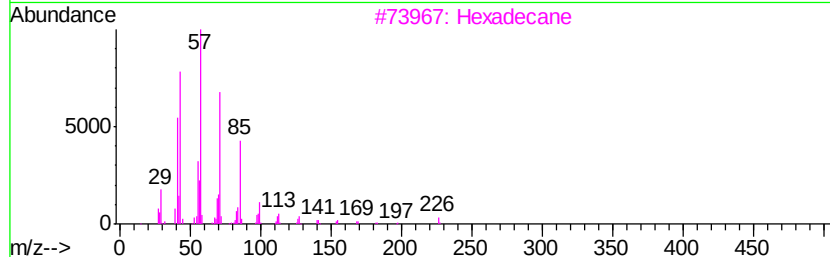
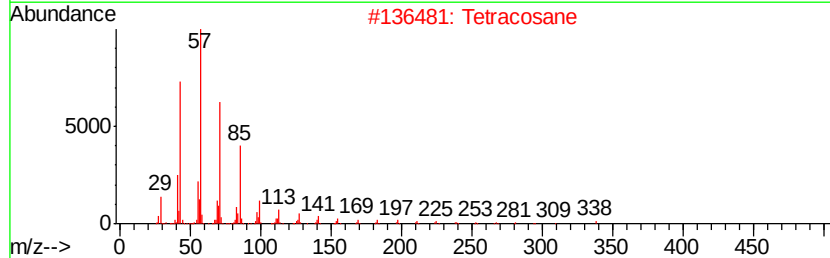
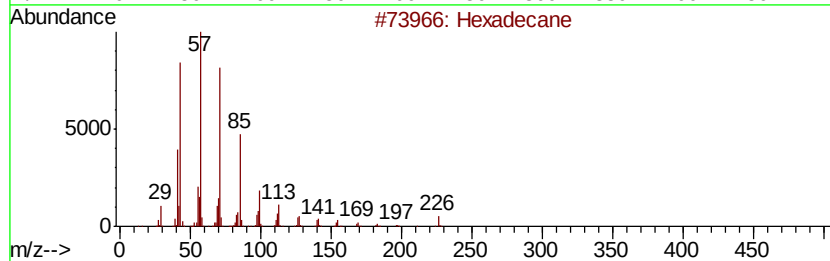
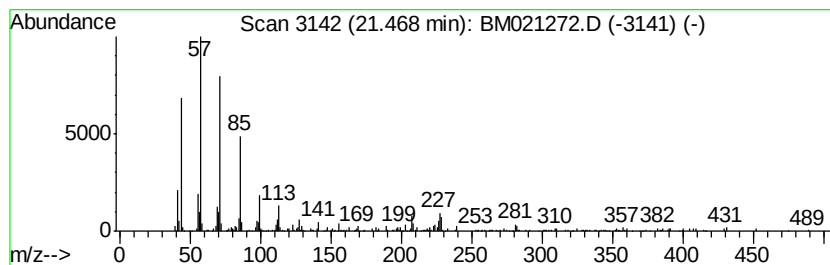
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 33

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.47 | 2.03 ng/ul | 597800 | Chrysene-d12     | 21.34 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 83   |
| 2    |    |   | Tetracosane  | 338 | C24H50  | 000646-31-1 | 78   |
| 3    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 76   |
| 4    |    |   | Tetracosane  | 338 | C24H50  | 000646-31-1 | 62   |
| 5    |    |   | Pentacosane  | 352 | C25H52  | 000629-99-2 | 58   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

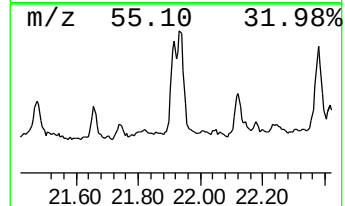
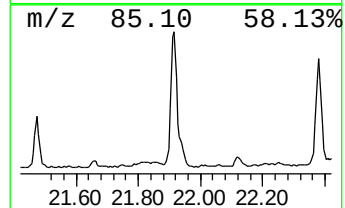
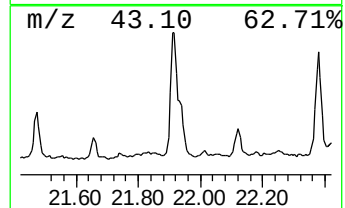
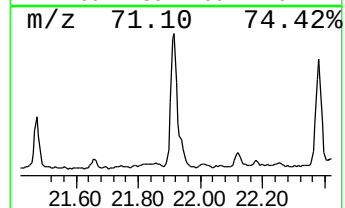
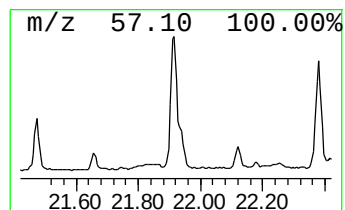
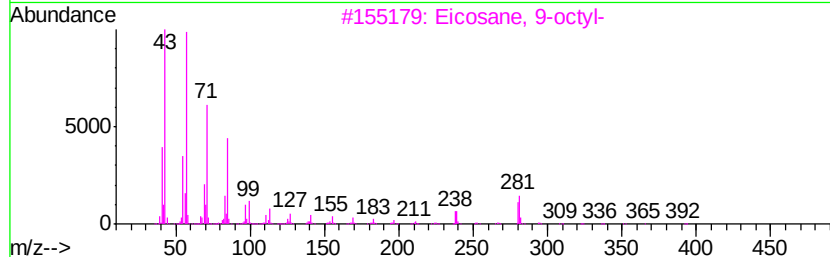
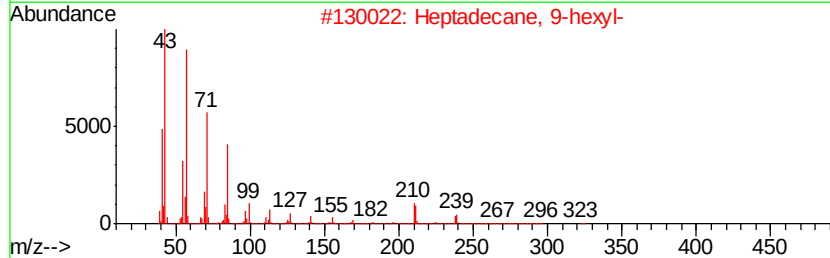
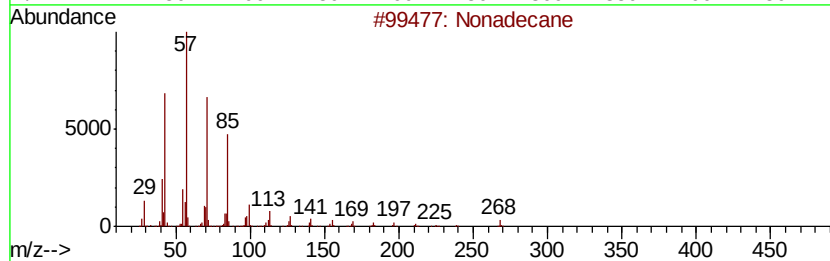
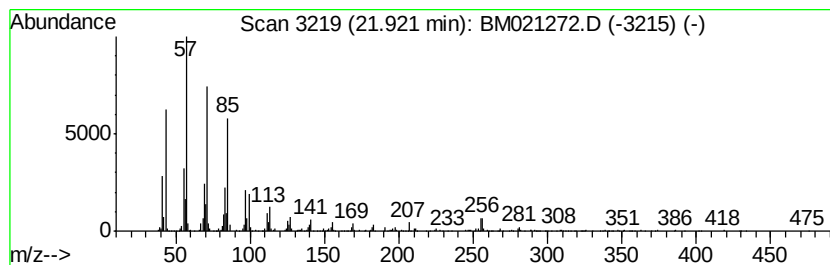
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 25

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.92 | 3.24 ng/ul | 954657 | Chrysene-d12     | 21.34 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Nonadecane            | 268 | C19H40  | 000629-92-5 | 94   |
| 2    |    | Heptadecane, 9-hexyl- | 324 | C23H48  | 055124-79-3 | 94   |
| 3    |    | Eicosane, 9-octyl-    | 394 | C28H58  | 013475-77-9 | 94   |
| 4    |    | Heneicosane           | 296 | C21H44  | 000629-94-7 | 91   |
| 5    |    | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

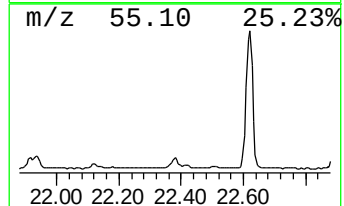
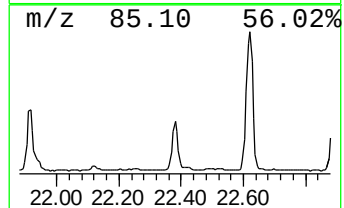
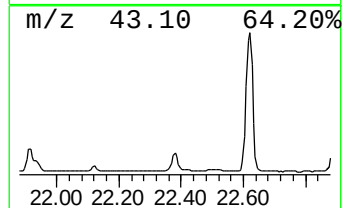
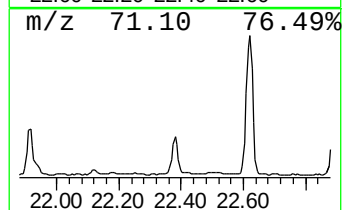
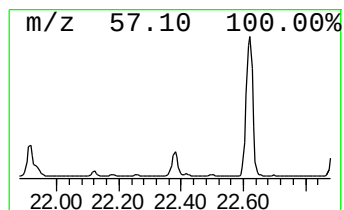
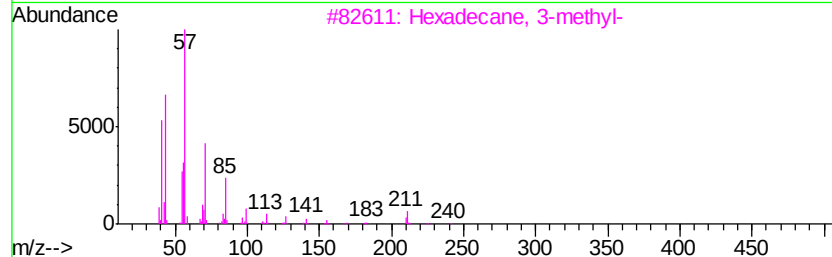
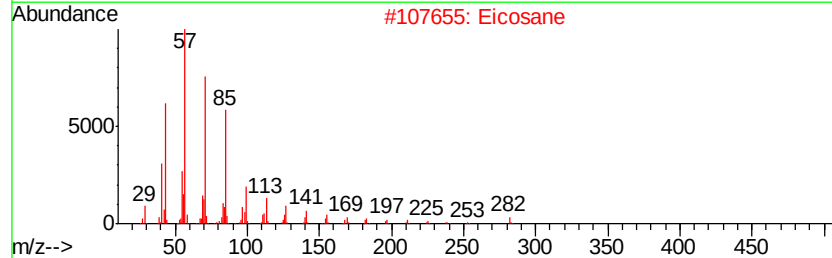
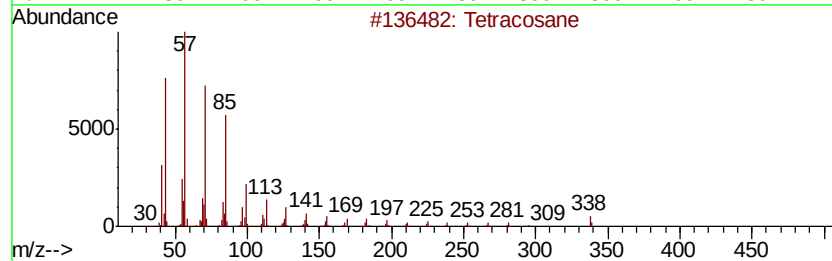
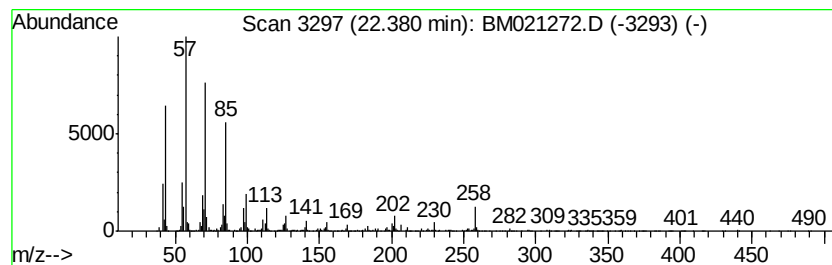
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 23

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 22.38 | 4.23 ng/ul | 1246440 | Chrysene-d12     | 21.34 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Tetracosane           | 338 | C24H50  | 000646-31-1 | 98   |
| 2    |    |   | Eicosane              | 282 | C20H42  | 000112-95-8 | 93   |
| 3    |    |   | Hexadecane, 3-methyl- | 240 | C17H36  | 006418-43-5 | 90   |
| 4    |    |   | Heneicosane           | 296 | C21H44  | 000629-94-7 | 90   |
| 5    |    |   | triacontane           | 422 | C30H62  | 000638-68-6 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

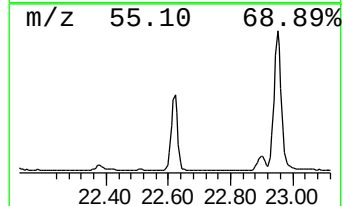
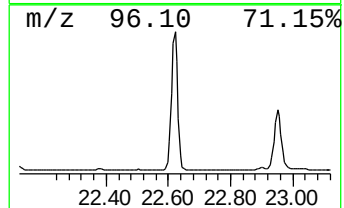
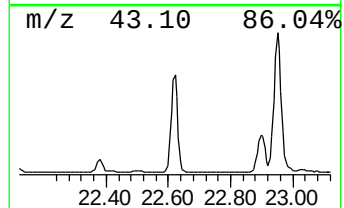
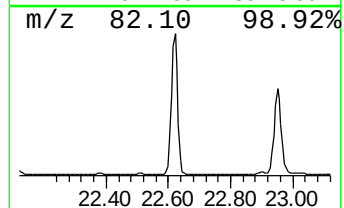
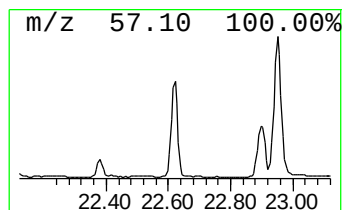
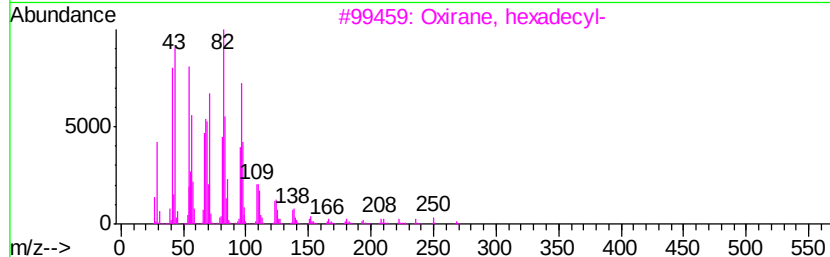
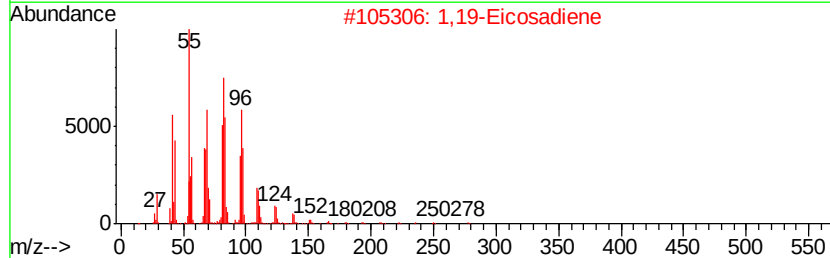
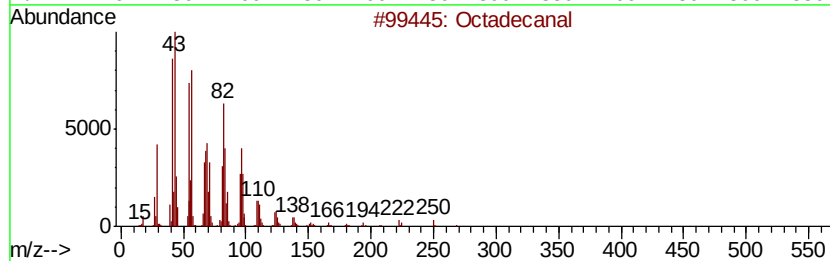
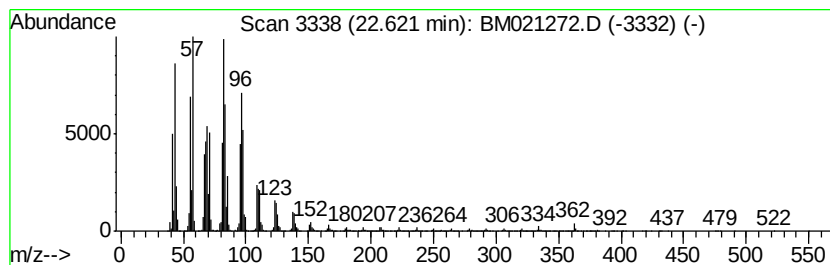
Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 23 Octadecanal Concentration Rank 1

| R.T.    | EstConc              | Area         | Relative to ISTD | R.T.      |
|---------|----------------------|--------------|------------------|-----------|
| 22.62   | 88.56 ng/ul          | 15586900     | Perylene-d12     | 23.61     |
| Hit# of | 5                    | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Octadecanal          | 268 C18H36O  | 000638-66-4      | 96        |
| 2       | 1,19-Eicosadiene     | 278 C20H38   | 014811-95-1      | 94        |
| 3       | Oxirane, hexadecyl-  | 268 C18H36O  | 007390-81-0      | 91        |
| 4       | Oxirane, heptadecyl- | 282 C19H38O  | 067860-04-2      | 91        |
| 5       | Oxirane, hexadecyl-  | 268 C18H36O  | 007390-81-0      | 91        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

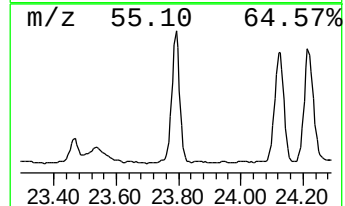
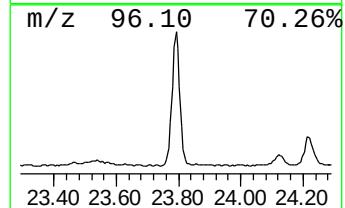
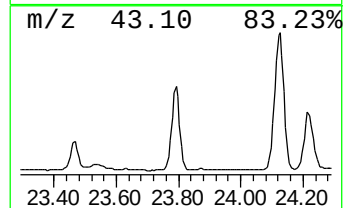
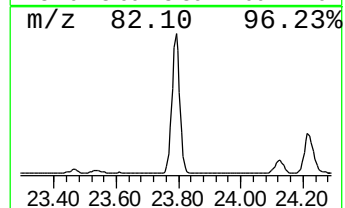
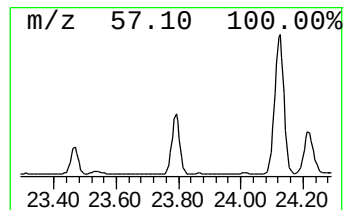
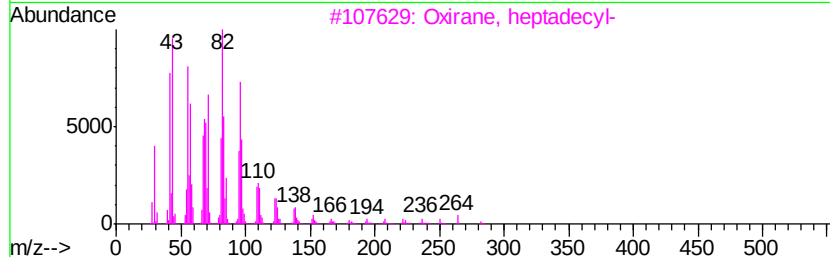
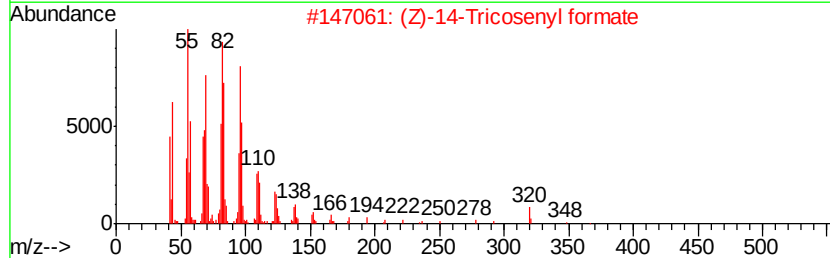
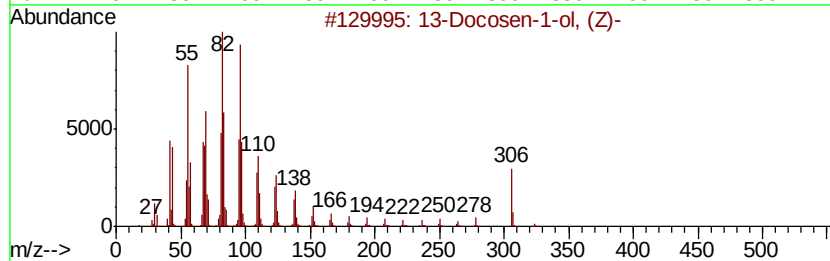
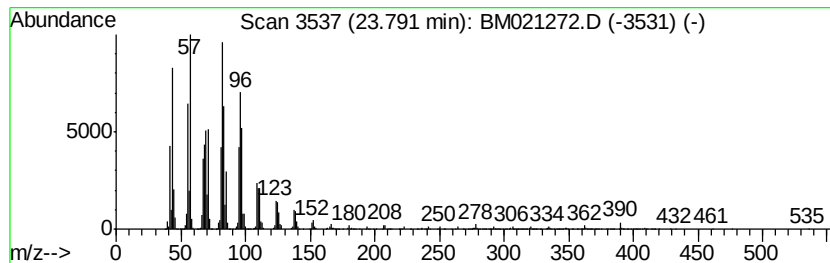
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 13-Docosen-1-ol, (Z)- Concentration Rank 5

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 23.79 | 45.51 ng/ul | 8009400 | Perylene-d12     | 23.61 |

| Hit# | of | Tentative ID              | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------------|-----|----------|-------------|------|
| 1    | 5  | 13-Docosen-1-ol, (Z)-     | 324 | C22H44O  | 000629-98-1 | 98   |
| 2    |    | (Z)-14-Tricosenyl formate | 366 | C24H46O2 | 077899-10-6 | 91   |
| 3    |    | Oxirane, heptadecyl-      | 282 | C19H38O  | 067860-04-2 | 91   |
| 4    |    | Oxirane, hexadecyl-       | 268 | C18H36O  | 007390-81-0 | 91   |
| 5    |    | Octadecanal               | 268 | C18H36O  | 000638-66-4 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

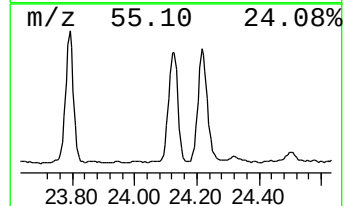
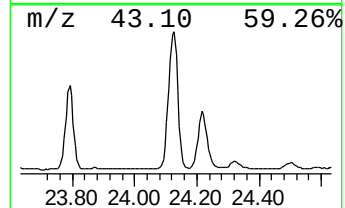
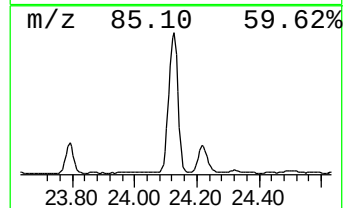
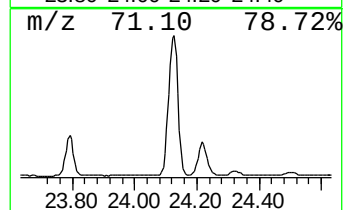
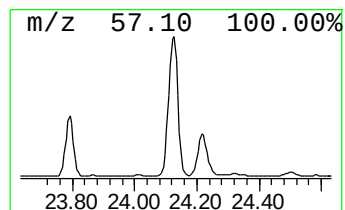
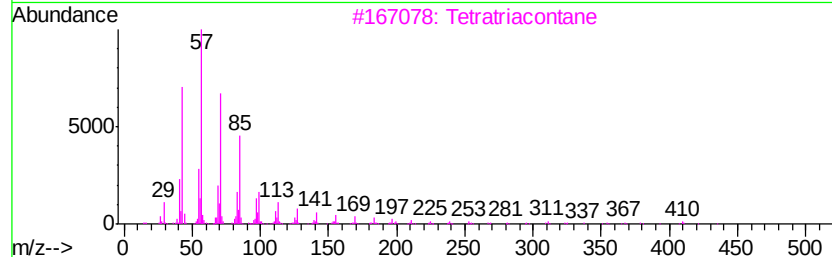
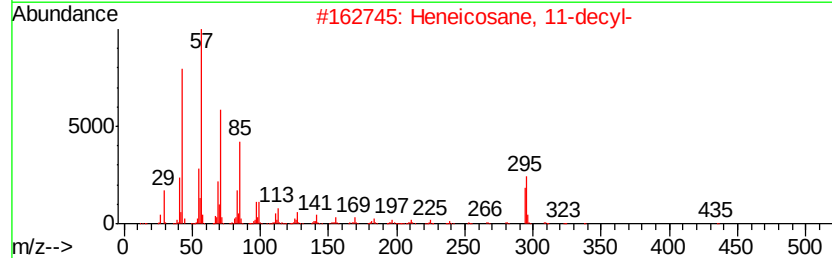
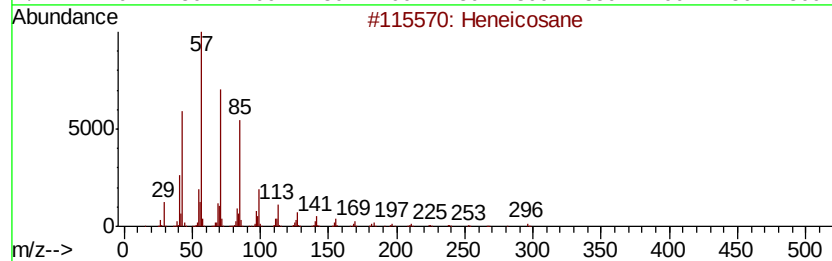
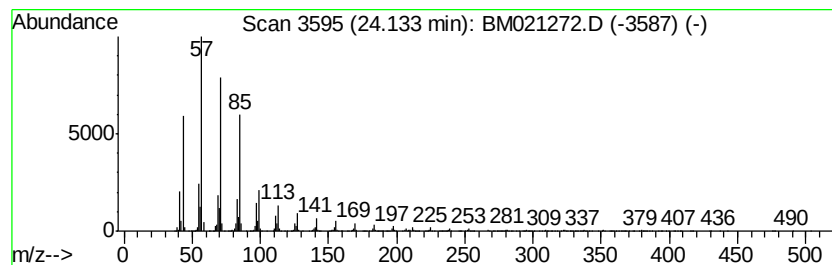
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 24.13 | 56.89 ng/ul | 10013100 | Perylene-d12     | 23.61 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Heneicosane            | 296 | C21H44  | 000629-94-7 | 97   |
| 2    |    | Heneicosane, 11-decyl- | 437 | C31H64  | 055320-06-4 | 95   |
| 3    |    | Tetratriacontane       | 479 | C34H70  | 014167-59-0 | 94   |
| 4    |    | Eicosane               | 282 | C20H42  | 000112-95-8 | 94   |
| 5    |    | Eicosane, 9-octyl-     | 394 | C28H58  | 013475-77-9 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

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Quant Title : SVOA CALIBRATION

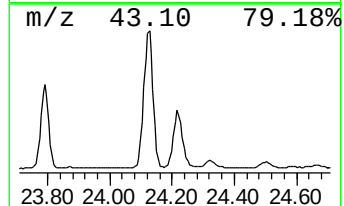
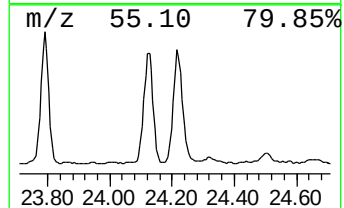
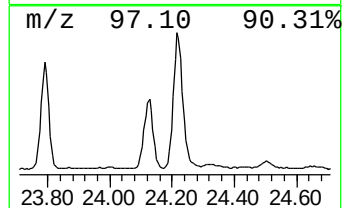
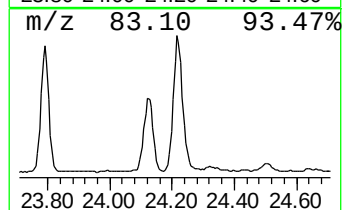
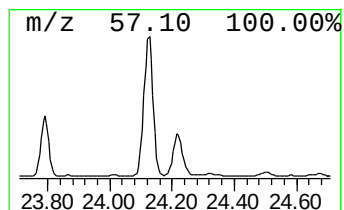
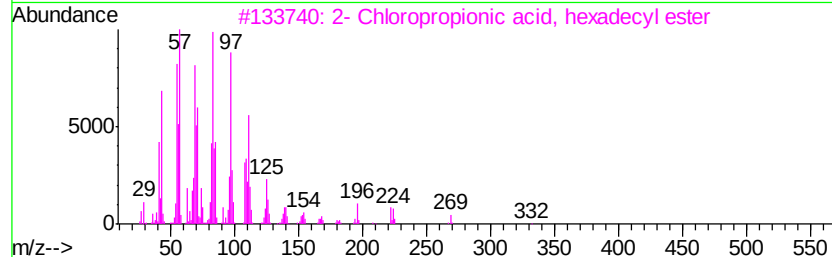
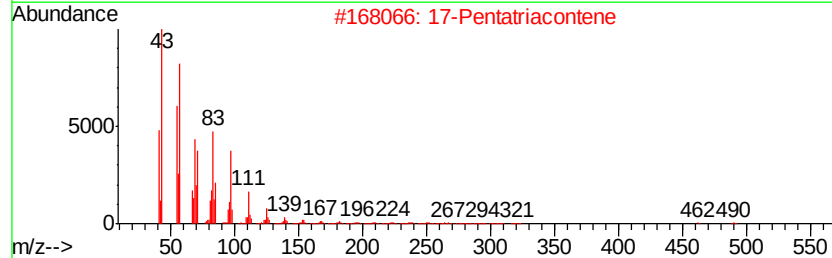
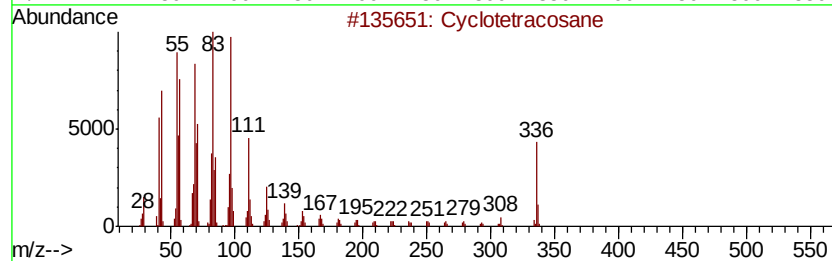
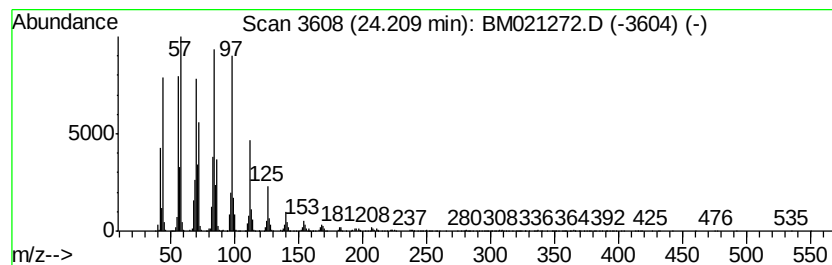
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 26 (DEL) Alkane: Cyclic24.21 Concentration Rank 8

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 24.21 | 34.41 ng/ul | 6056990 | Perylene-d12     | 23.61 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Cyclotetracosane                    | 336 | C24H48      | 000297-03-0  | 91   |
| 2    |    | 17-Pentatriacontene                 | 491 | C35H70      | 006971-40-0  | 90   |
| 3    |    | 2- Chloropropionic acid, hexadec... | 332 | C19H37ClO2  | 086711-81-1  | 83   |
| 4    |    | 2- Chloropropionic acid, octadec... | 360 | C21H41ClO2  | 088104-31-8  | 81   |
| 5    |    | Dichloroacetic acid, heptadecyl ... | 366 | C19H36Cl2O2 | 1000282-98-2 | 81   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

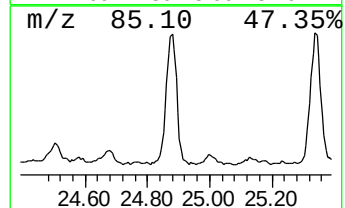
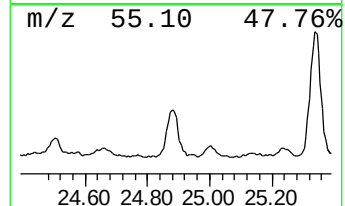
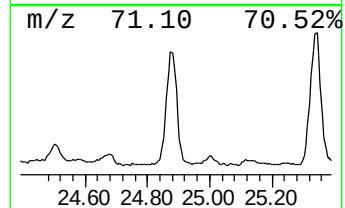
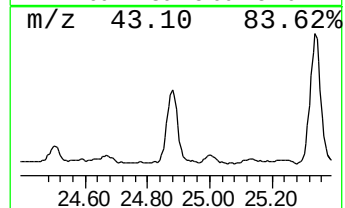
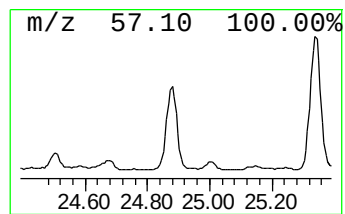
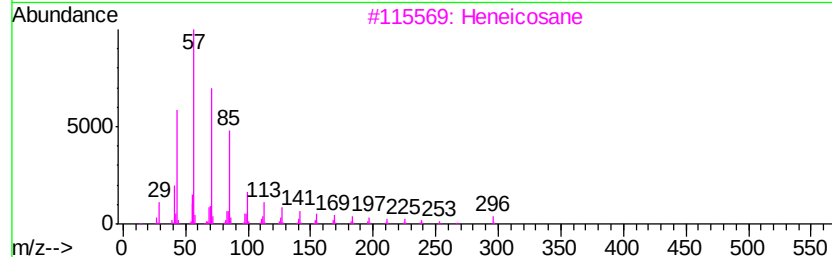
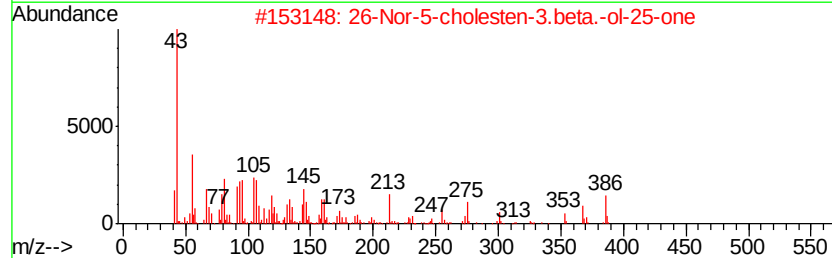
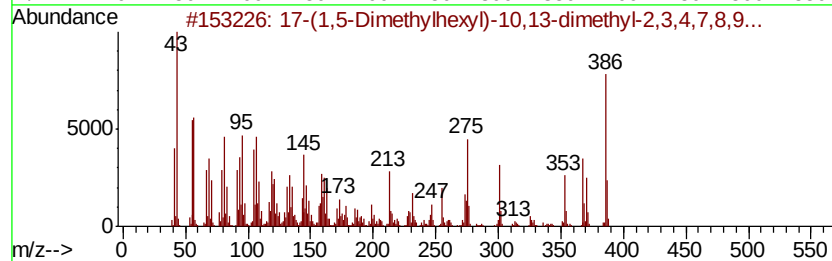
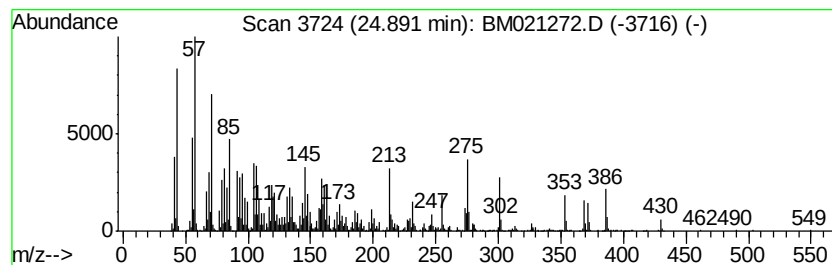
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 27 17-(1,5-Dimethylhexyl)-10,1... Concentration Rank 12

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 24.89 | 24.48 ng/ul | 4308140 | Perylene-d12     | 23.61 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 17-(1,5-Dimethylhexyl)-10,13-dim... | 386 | C27H46O  | 1000210-38-4 | 95   |
| 2    |    | 26-Nor-5-cholesten-3.beta.-ol-25... | 386 | C26H42O2 | 007494-34-0  | 91   |
| 3    |    | Heneicosane                         | 296 | C21H44   | 000629-94-7  | 90   |
| 4    |    | Octacosane                          | 394 | C28H58   | 000630-02-4  | 87   |
| 5    |    | Tridecane, 6-propyl-                | 226 | C16H34   | 055045-10-8  | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

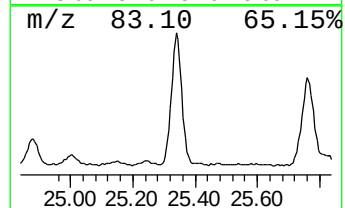
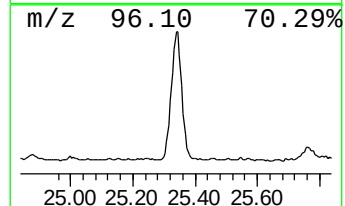
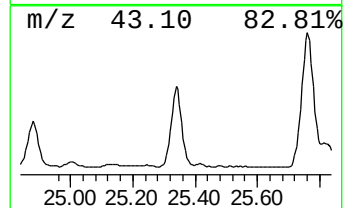
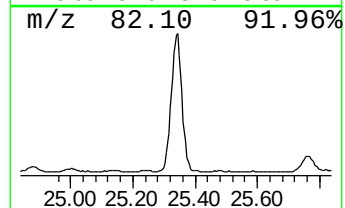
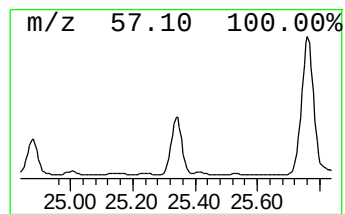
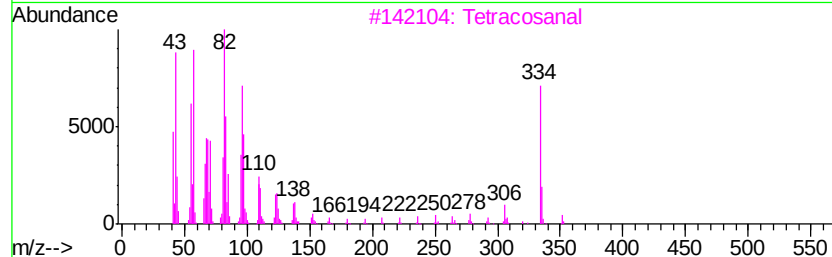
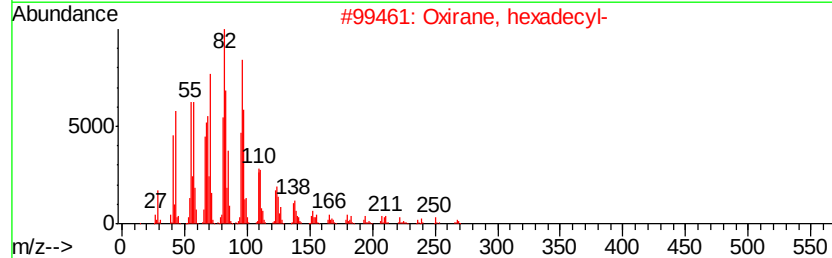
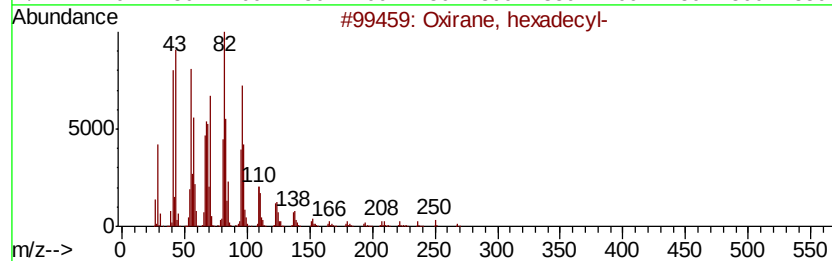
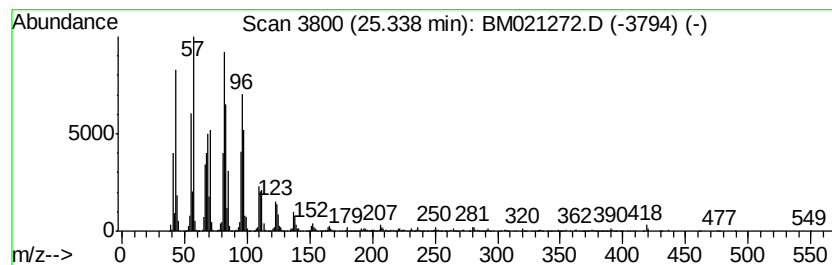
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 28 Oxirane, hexadecyl- Concentration Rank 7

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 25.34 | 35.19 ng/ul | 6192970 | Perylene-d12     | 23.61 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | Oxirane, hexadecyl- | 268 | C18H36O | 007390-81-0 | 96   |
| 2    |    | Oxirane, hexadecyl- | 268 | C18H36O | 007390-81-0 | 91   |
| 3    |    | Tetracosanal        | 352 | C24H48O | 057866-08-7 | 91   |
| 4    |    | Octadecanal         | 268 | C18H36O | 000638-66-4 | 86   |
| 5    |    | 1,13-Tetradecadiene | 194 | C14H26  | 021964-49-8 | 83   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

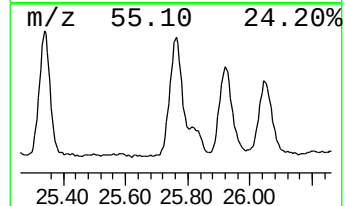
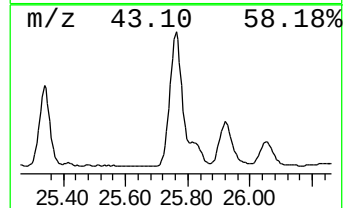
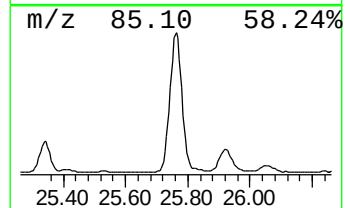
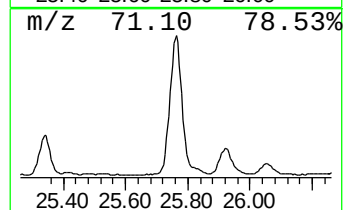
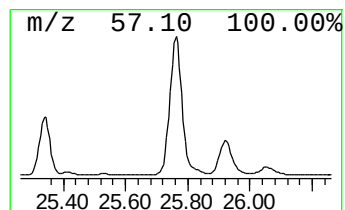
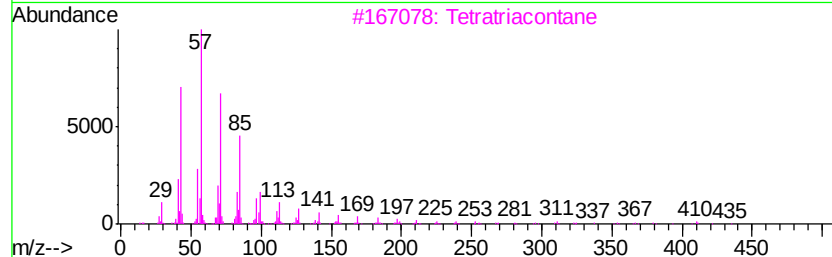
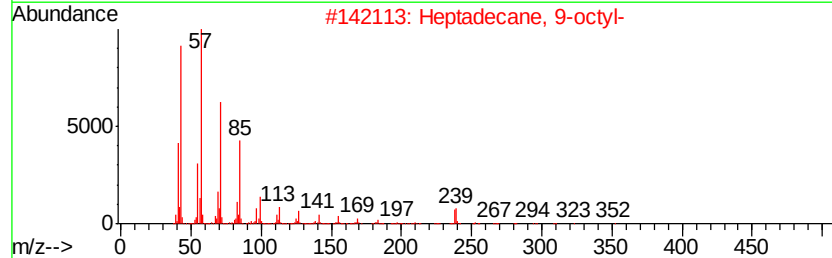
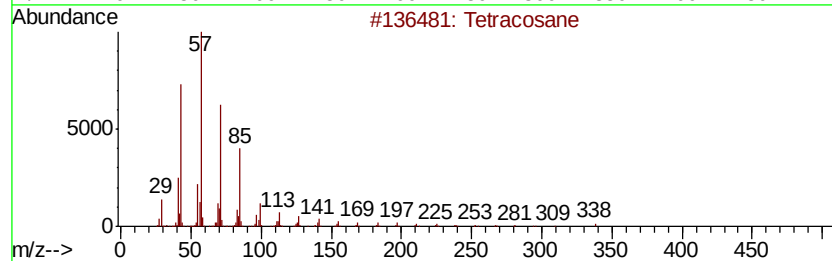
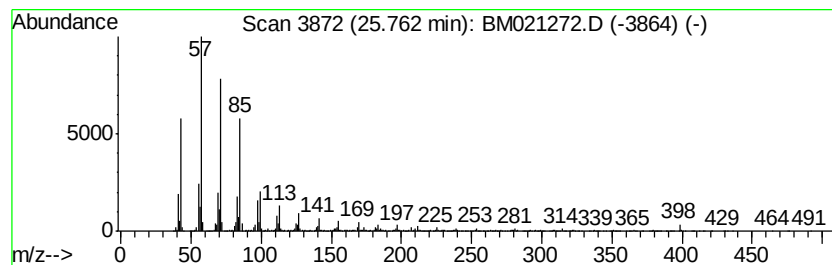
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 25.76 | 52.71 ng/ul | 9277670 | Perylene-d12     | 23.61 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Tetracosane           | 338 | C24H50  | 000646-31-1 | 97   |
| 2    |    | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 95   |
| 3    |    | Tetratriacontane      | 479 | C34H70  | 014167-59-0 | 94   |
| 4    |    | Docosane, 11-butyl-   | 366 | C26H54  | 013475-76-8 | 94   |
| 5    |    | Triacontane           | 422 | C30H62  | 000638-68-6 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

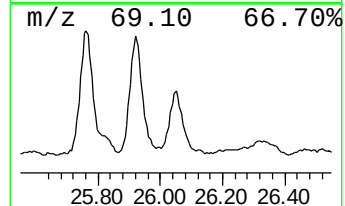
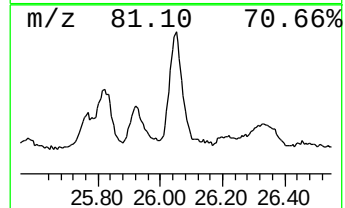
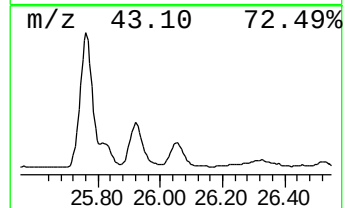
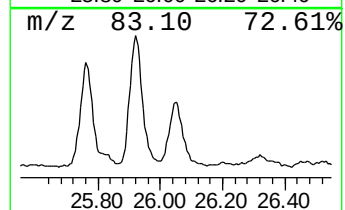
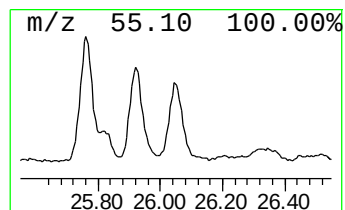
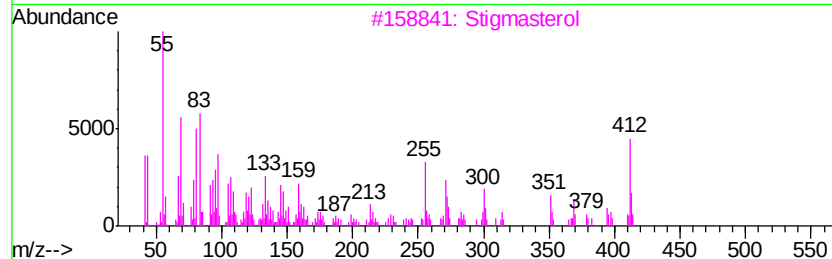
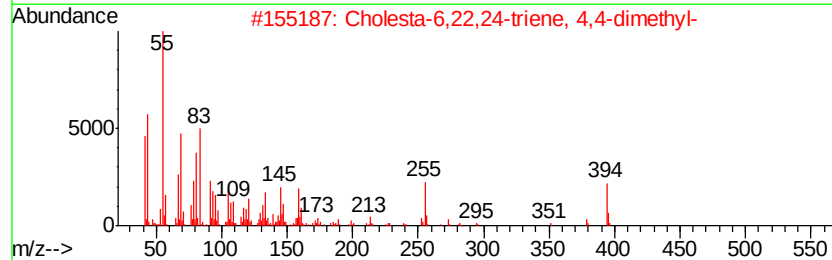
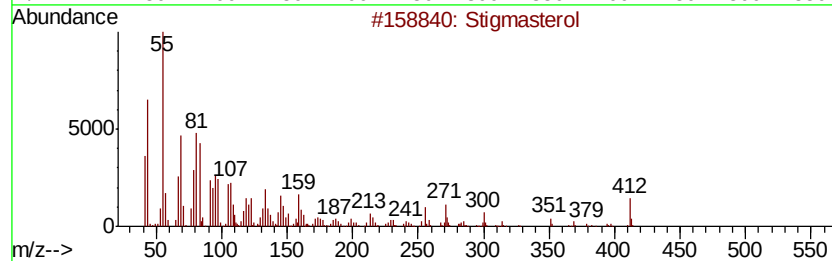
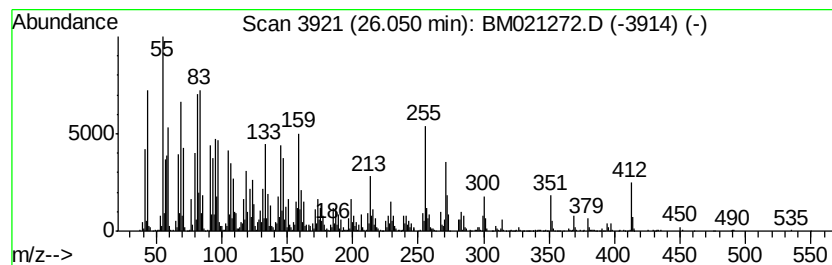
Instrument :  
BNA\_M  
ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 30 Stigmasterol Concentration Rank 10

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.            |
|---------|-------------------------------------|--------------|------------------|-----------------|
| 26.05   | 28.21 ng/ul                         | 4965250      | Perylene-d12     | 23.61           |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual       |
| 1       | Stigmasterol                        |              | 412 C29H48O      | 000083-48-7 99  |
| 2       | Cholesta-6,22,24-triene, 4,4-dim... |              | 394 C29H46       | 1000128-66-9 87 |
| 3       | Stigmasterol                        |              | 412 C29H48O      | 000083-48-7 74  |
| 4       | Stigmasterol                        |              | 412 C29H48O      | 000083-48-7 58  |
| 5       | Stigmasta-5,22-dien-3-ol, acetat... |              | 454 C31H50O2     | 004651-48-3 58  |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

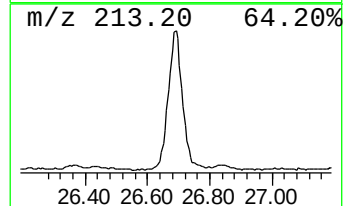
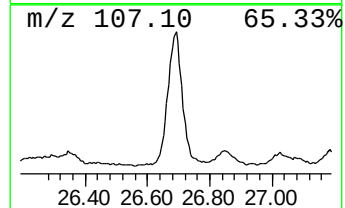
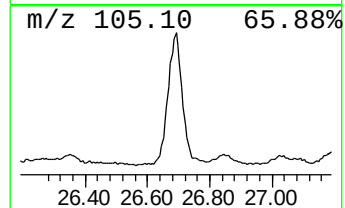
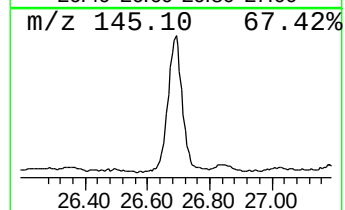
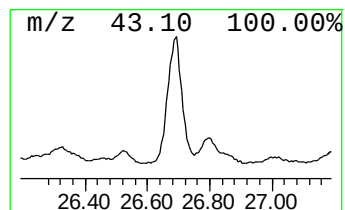
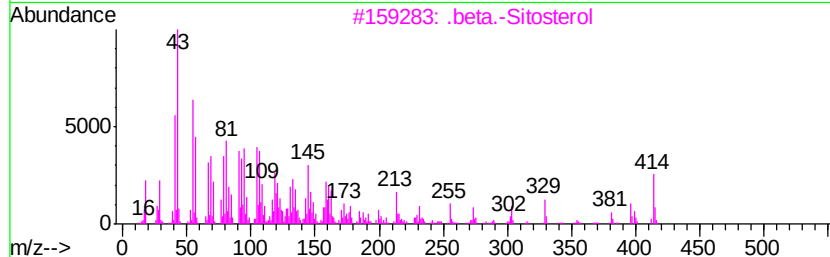
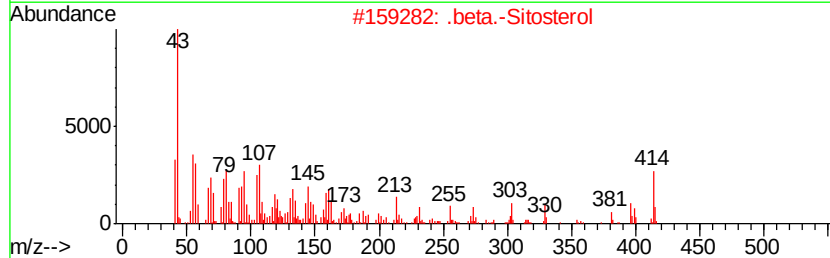
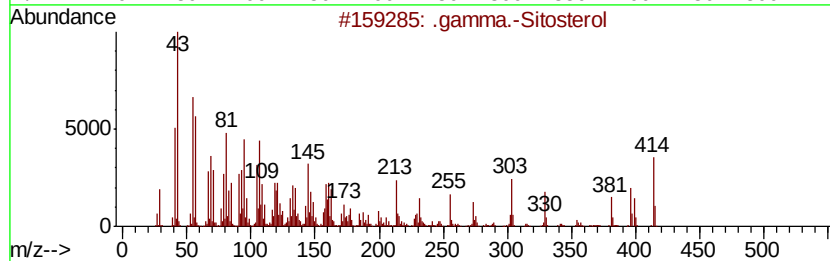
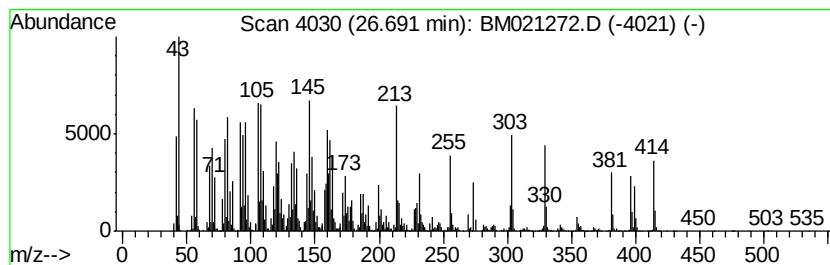
Instrument :  
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ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 31 .gamma.-Sitosterol Concentration Rank 2

| R.T.    | EstConc     | Area                                | Relative to ISTD | R.T.            |
|---------|-------------|-------------------------------------|------------------|-----------------|
| 26.69   | 60.71 ng/ul | 10685400                            | Perylene-d12     | 23.61           |
| Hit# of | 5           | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |             | .gamma.-Sitosterol                  | 414 C29H50O      | 000083-47-6 99  |
| 2       |             | .beta.-Sitosterol                   | 414 C29H50O      | 000083-46-5 96  |
| 3       |             | .beta.-Sitosterol                   | 414 C29H50O      | 000083-46-5 55  |
| 4       |             | 22,26-Oxido-4,17-cholestadien-3.... | 414 C27H42O3     | 1000252-80-4 25 |
| 5       |             | Stigmastan-3-en-6-ol                | 414 C29H50O      | 1000214-19-7 15 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

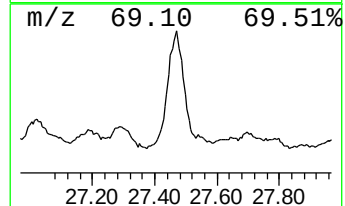
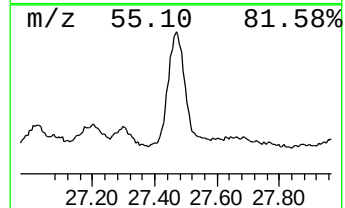
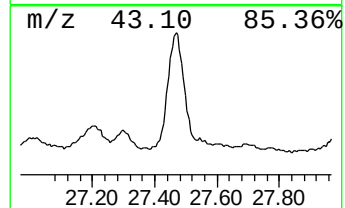
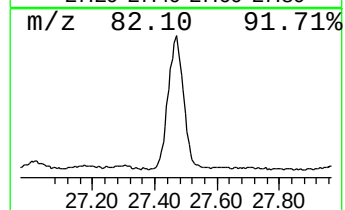
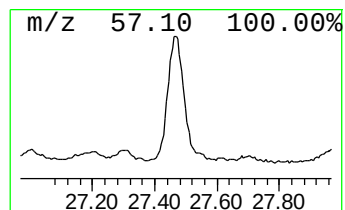
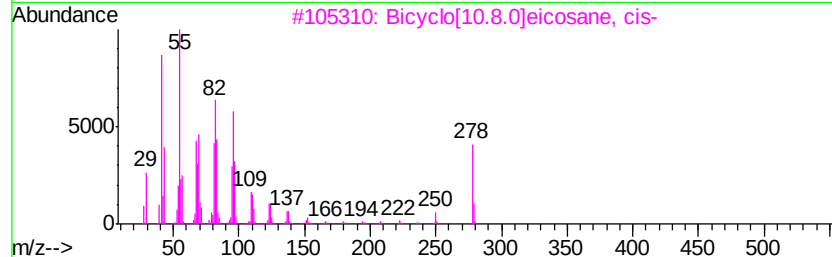
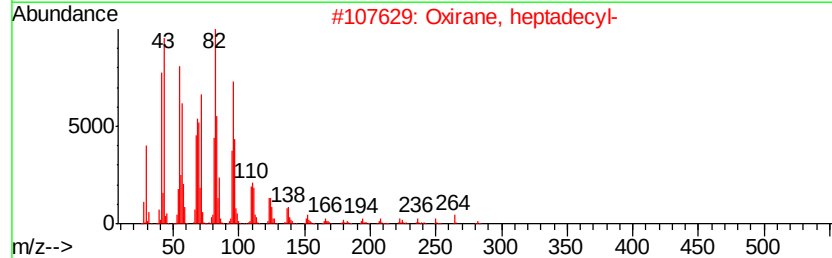
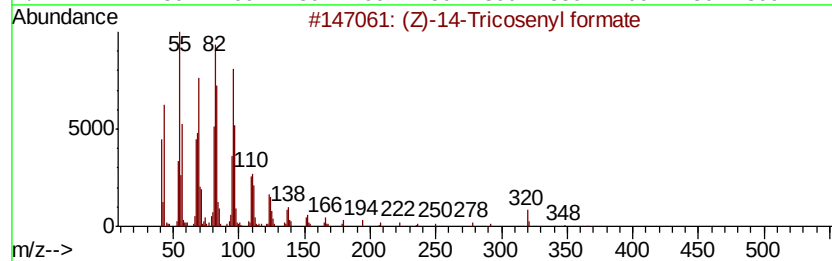
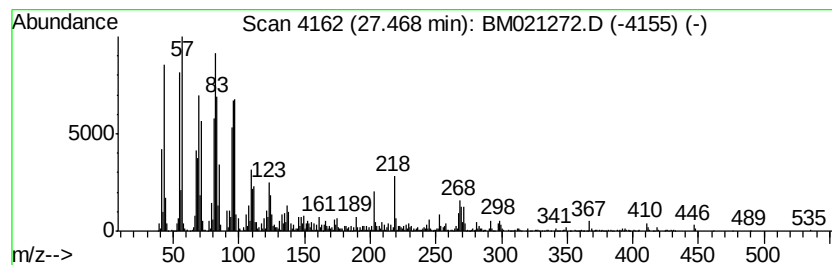
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 32 (Z)-14-Tricosenyl formate Concentration Rank 11

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 27.47 | 24.57 ng/ul | 4324960 | Perylene-d12     | 23.61 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | (Z)-14-Tricosenyl formate           | 366 | C24H46O2 | 077899-10-6  | 93   |
| 2    |    | Oxirane, heptadecyl-                | 282 | C19H38O  | 067860-04-2  | 78   |
| 3    |    | Bicyclo[10.8.0]eicosane, cis-       | 278 | C20H38   | 1000155-82-2 | 70   |
| 4    |    | 2-Octadecyl-propane-1,3-diol        | 328 | C21H44O2 | 005337-61-1  | 55   |
| 5    |    | 6,11-Undecadiene, 1-acetoxy-3,7-... | 252 | C16H28O2 | 1000150-66-0 | 55   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062919\  
Data File : BM021272.D  
Acq On : 29 Jun 2019 21:17  
Operator : HP/JU  
Sample : K3593-16  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

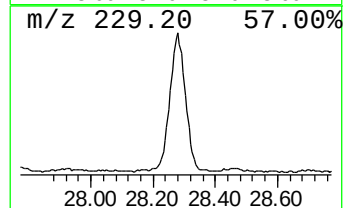
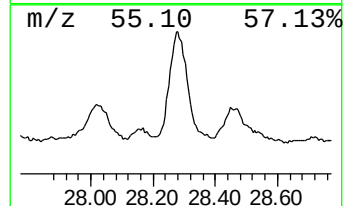
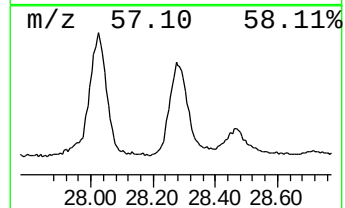
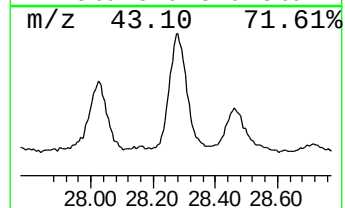
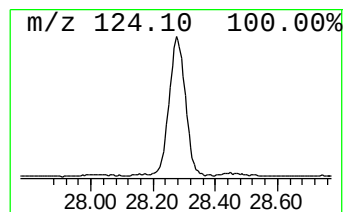
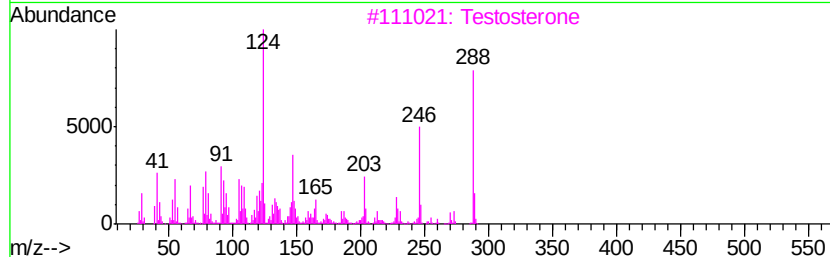
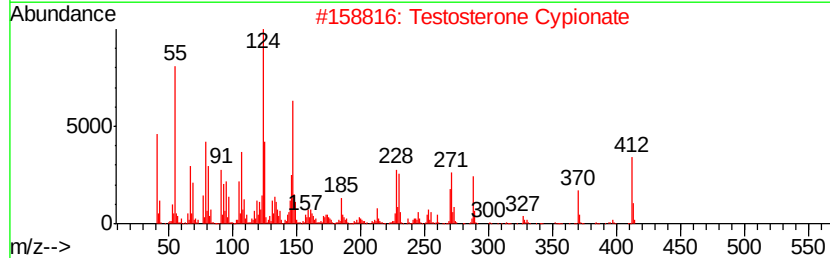
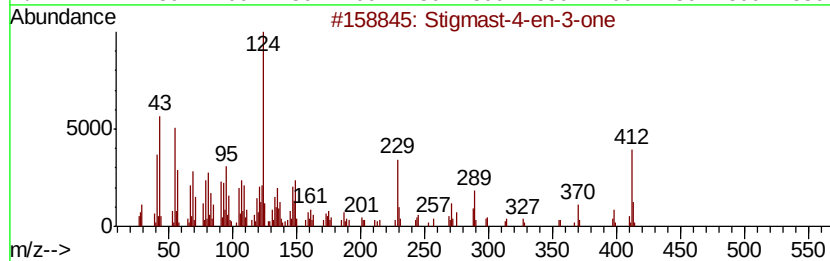
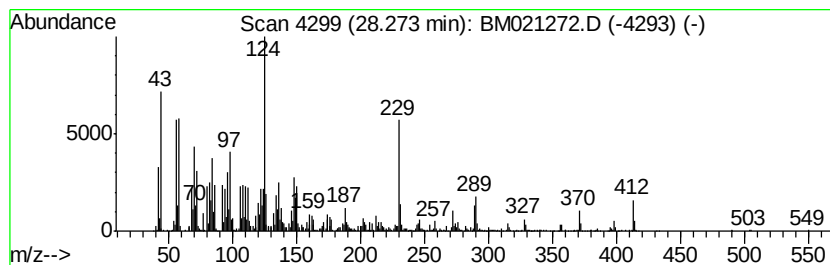
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ClientSampleId :  
C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 33 Stigmast-4-en-3-one Concentration Rank 9

| R.T.    | EstConc     | Area                                | Relative to ISTD | R.T.      |             |      |
|---------|-------------|-------------------------------------|------------------|-----------|-------------|------|
| 28.27   | 30.00 ng/ul | 5280910                             | Perylene-d12     | 23.61     |             |      |
| Hit# of | 5           | Tentative ID                        | MW               | MolForm   | CAS#        | Qual |
| 1       |             | Stigmast-4-en-3-one                 | 412              | C29H48O   | 001058-61-3 | 89   |
| 2       |             | Testosterone Cypionate              | 412              | C27H40O3  | 000058-20-8 | 72   |
| 3       |             | Testosterone                        | 288              | C19H28O2  | 000058-22-0 | 38   |
| 4       |             | Androst-4-en-3-one, 17-hydroxy-,... | 288              | C19H28O2  | 000604-39-7 | 38   |
| 5       |             | Atropine                            | 289              | C17H23NO3 | 000051-55-8 | 30   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM062919\  
 Data File : BM021272.D  
 Acq On : 29 Jun 2019 21:17  
 Operator : HP/JU  
 Sample : K3593-16  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C7BA8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM061419MA.M  
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| Butane, 2-methoxy... | 3.02  | 37.3    | ng/ul | 3271380  | 1                     | 7.76  | 1754460 | 20.0 |
| Hexanal              | 4.35  | 4.4     | ng/ul | 382547   | 1                     | 7.76  | 1754460 | 20.0 |
| Cyclotrisiloxane,... | 4.40  | 5.1     | ng/ul | 448545   | 1                     | 7.76  | 1754460 | 20.0 |
| Heptanal             | 5.83  | 2.1     | ng/ul | 187228   | 1                     | 7.76  | 1754460 | 20.0 |
| Cyclotetrasiloxan... | 6.86  | 4.2     | ng/ul | 369377   | 1                     | 7.76  | 1754460 | 20.0 |
| Benzene, 1-methyl... | 7.89  | 7.5     | ng/ul | 653472   | 1                     | 7.76  | 1754460 | 20.0 |
| Eucalyptol           | 8.06  | 5.4     | ng/ul | 475433   | 1                     | 7.76  | 1754460 | 20.0 |
| 2-Octenal, (E)-      | 8.36  | 2.3     | ng/ul | 200085   | 1                     | 7.76  | 1754460 | 20.0 |
| Nonanal              | 9.07  | 2.3     | ng/ul | 197831   | 1                     | 7.76  | 1754460 | 20.0 |
| Bicyclo[3.1.0]hex... | 9.25  | 5.9     | ng/ul | 773173   | 2                     | 10.55 | 2618460 | 20.0 |
| 1,12-Tridecadiene    | 15.96 | 2.2     | ng/ul | 380668   | 4                     | 17.15 | 3496120 | 20.0 |
| Hexadecenoic acid... | 17.99 | 8.3     | ng/ul | 1446930  | 4                     | 17.15 | 3496120 | 20.0 |
| n-Hexadecanoic acid  | 18.04 | 18.4    | ng/ul | 3215320  | 4                     | 17.15 | 3496120 | 20.0 |
| 4H-Cyclopenta[def... | 18.22 | 2.1     | ng/ul | 375196   | 4                     | 17.15 | 3496120 | 20.0 |
| 17-Norkaur-15-ene... | 18.34 | 6.9     | ng/ul | 1209060  | 4                     | 17.15 | 3496120 | 20.0 |
| Phytol               | 19.06 | 6.6     | ng/ul | 1156250  | 4                     | 17.15 | 3496120 | 20.0 |
| Octadecanoic acid    | 19.32 | 2.5     | ng/ul | 720527   | 5                     | 21.34 | 5886910 | 20.0 |
| 11H-Benzo[b]fluorene | 20.07 | 3.1     | ng/ul | 927135   | 5                     | 21.34 | 5886910 | 20.0 |
| 2-Chloropropioni...  | 21.04 | 6.7     | ng/ul | 1966740  | 5                     | 21.34 | 5886910 | 20.0 |
| (DEL) Alkane: Str... | 21.47 | 2.0     | ng/ul | 597800   | 5                     | 21.34 | 5886910 | 20.0 |
| (DEL) Alkane: Str... | 21.92 | 3.2     | ng/ul | 954657   | 5                     | 21.34 | 5886910 | 20.0 |
| (DEL) Alkane: Str... | 22.38 | 4.2     | ng/ul | 1246440  | 5                     | 21.34 | 5886910 | 20.0 |
| Octadecanal          | 22.62 | 88.6    | ng/ul | 15586900 | 6                     | 23.61 | 3520160 | 20.0 |
| 13-Docosen-1-ol, ... | 23.79 | 45.5    | ng/ul | 8009400  | 6                     | 23.61 | 3520160 | 20.0 |
| (DEL) Alkane: Str... | 24.13 | 56.9    | ng/ul | 10013100 | 6                     | 23.61 | 3520160 | 20.0 |
| (DEL) Alkane: Cyc... | 24.21 | 34.4    | ng/ul | 6056990  | 6                     | 23.61 | 3520160 | 20.0 |
| 17-(1,5-Dimethylh... | 24.89 | 24.5    | ng/ul | 4308140  | 6                     | 23.61 | 3520160 | 20.0 |
| Oxirane, hexadecyl-  | 25.34 | 35.2    | ng/ul | 6192970  | 6                     | 23.61 | 3520160 | 20.0 |
| (DEL) Alkane: Str... | 25.76 | 52.7    | ng/ul | 9277670  | 6                     | 23.61 | 3520160 | 20.0 |
| Stigmasterol         | 26.05 | 28.2    | ng/ul | 4965250  | 6                     | 23.61 | 3520160 | 20.0 |
| .gamma.-Sitosterol   | 26.69 | 60.7    | ng/ul | 10685400 | 6                     | 23.61 | 3520160 | 20.0 |
| (Z)-14-Tricosenyl... | 27.47 | 24.6    | ng/ul | 4324960  | 6                     | 23.61 | 3520160 | 20.0 |
| Stigmast-4-en-3-one  | 28.27 | 30.0    | ng/ul | 5280910  | 6                     | 23.61 | 3520160 | 20.0 |