

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080919\  
 Data File : BF116122.D  
 Acq On : 9 Aug 2019 19:45  
 Operator : HP/JU  
 Sample : K4257-02  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 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\_F\METHODS\8270-BF073019.M

Title : ASP BNA STANDARDS FOR 5 POINT 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         | 2.140       | 3             | 9           | 28           | rVB      | 1647970        | 2459372       | 59.29%          | 5.350%        |
| 2         | 5.475       | 572           | 576         | 611          | rBV      | 2923908        | 3468134       | 83.61%          | 7.545%        |
| 3         | 6.487       | 743           | 748         | 766          | rBV      | 2849415        | 3696465       | 89.11%          | 8.041%        |
| 4         | 6.616       | 766           | 770         | 776          | rVV      | 3721206        | 3541810       | 85.38%          | 7.705%        |
| 5         | 6.840       | 804           | 808         | 815          | rVB      | 861675         | 799245        | 19.27%          | 1.739%        |
| 6         | 6.993       | 830           | 834         | 846          | rBV      | 3002142        | 2986828       | 72.00%          | 6.497%        |
| 7         | 7.398       | 899           | 903         | 927          | rBV      | 2095406        | 2426751       | 58.50%          | 5.279%        |
| 8         | 8.116       | 1022          | 1025        | 1047         | rVB      | 983512         | 1015406       | 24.48%          | 2.209%        |
| 9         | 9.192       | 1203          | 1208        | 1211         | rBV      | 4302076        | 4148221       | 100.00%         | 9.024%        |
| 10        | 9.586       | 1272          | 1275        | 1284         | rVB      | 267559         | 279932        | 6.75%           | 0.609%        |
| 11        | 9.869       | 1319          | 1323        | 1327         | rBV      | 1267870        | 1139122       | 27.46%          | 2.478%        |
| 12        | 9.904       | 1327          | 1329        | 1338         | rVB      | 124679         | 123334        | 2.97%           | 0.268%        |
| 13        | 10.086      | 1353          | 1360        | 1363         | rBV3     | 38420          | 50465         | 1.22%           | 0.110%        |
| 14        | 10.422      | 1414          | 1417        | 1422         | rBV      | 88396          | 96257         | 2.32%           | 0.209%        |
| 15        | 10.663      | 1454          | 1458        | 1471         | rBV      | 2276600        | 2427205       | 58.51%          | 5.280%        |
| 16        | 10.969      | 1504          | 1510        | 1513         | rBV4     | 40106          | 63816         | 1.54%           | 0.139%        |
| 17        | 11.216      | 1548          | 1552        | 1555         | rBV2     | 45915          | 47060         | 1.13%           | 0.102%        |
| 18        | 11.257      | 1555          | 1559        | 1565         | rVB      | 62036          | 61354         | 1.48%           | 0.133%        |
| 19        | 11.357      | 1572          | 1576        | 1578         | rBV      | 1284962        | 1111982       | 26.81%          | 2.419%        |
| 20        | 11.380      | 1578          | 1580        | 1585         | rVV      | 1043648        | 923642        | 22.27%          | 2.009%        |
| 21        | 11.439      | 1585          | 1590        | 1596         | rVB      | 181652         | 245828        | 5.93%           | 0.535%        |
| 22        | 11.598      | 1613          | 1617        | 1623         | rBV      | 57622          | 79514         | 1.92%           | 0.173%        |
| 23        | 11.863      | 1658          | 1662        | 1663         | rBV      | 125790         | 112184        | 2.70%           | 0.244%        |
| 24        | 11.886      | 1663          | 1666        | 1673         | rVV2     | 938377         | 1092245       | 26.33%          | 2.376%        |
| 25        | 11.939      | 1673          | 1675        | 1678         | rVV      | 83070          | 100400        | 2.42%           | 0.218%        |
| 26        | 11.975      | 1678          | 1681        | 1683         | rVV      | 192229         | 219299        | 5.29%           | 0.477%        |
| 27        | 11.998      | 1683          | 1685        | 1691         | rVB      | 90649          | 94704         | 2.28%           | 0.206%        |
| 28        | 12.157      | 1709          | 1712        | 1715         | rBV      | 77649          | 70498         | 1.70%           | 0.153%        |
| 29        | 12.198      | 1715          | 1719        | 1722         | rVV      | 52651          | 54924         | 1.32%           | 0.119%        |
| 30        | 12.410      | 1752          | 1755        | 1758         | rBV      | 68702          | 73700         | 1.78%           | 0.160%        |
| 31        | 12.504      | 1767          | 1771        | 1775         | rVB2     | 35966          | 47498         | 1.15%           | 0.103%        |
| 32        | 12.575      | 1779          | 1783        | 1788         | rBV      | 820669         | 870729        | 20.99%          | 1.894%        |
| 33        | 12.675      | 1796          | 1800        | 1806         | rBV      | 1382384        | 1498844       | 36.13%          | 3.261%        |
| 34        | 12.798      | 1816          | 1821        | 1828         | rVV      | 674762         | 802091        | 19.34%          | 1.745%        |

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 BNA\_F  
 ClientSampleId :

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : OFF  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0  
 Filtering: 5  
 Min Area: 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\_F\METHODS\8270-BF073019.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 12.857 | 1828 | 1831 | 1835 | rVB2 | 37262   | 45036   | 1.09%  | 0.098% |
| 36 | 12.939 | 1840 | 1845 | 1849 | rBV  | 3784652 | 4020906 | 96.93% | 8.747% |
| 37 | 13.022 | 1855 | 1859 | 1863 | rBV2 | 50552   | 74650   | 1.80%  | 0.162% |
| 38 | 13.133 | 1870 | 1878 | 1881 | rBV  | 104689  | 161427  | 3.89%  | 0.351% |
| 39 | 13.198 | 1886 | 1889 | 1892 | rVV  | 80835   | 80516   | 1.94%  | 0.175% |
| 40 | 13.239 | 1892 | 1896 | 1903 | rVB3 | 67091   | 100356  | 2.42%  | 0.218% |
| 41 | 13.645 | 1962 | 1965 | 1969 | rVB2 | 61049   | 71801   | 1.73%  | 0.156% |
| 42 | 13.757 | 1980 | 1984 | 1986 | rBV2 | 67882   | 79952   | 1.93%  | 0.174% |
| 43 | 13.857 | 1998 | 2001 | 2010 | rVB  | 127088  | 190102  | 4.58%  | 0.414% |
| 44 | 13.963 | 2016 | 2019 | 2022 | rBV  | 1064089 | 1052354 | 25.37% | 2.289% |
| 45 | 13.998 | 2022 | 2025 | 2028 | rVV2 | 1001416 | 1197995 | 28.88% | 2.606% |
| 46 | 14.027 | 2028 | 2030 | 2038 | rVB  | 282955  | 285571  | 6.88%  | 0.621% |
| 47 | 14.451 | 2099 | 2102 | 2107 | rVB3 | 43969   | 50484   | 1.22%  | 0.110% |
| 48 | 14.751 | 2150 | 2153 | 2156 | rVB  | 58729   | 59345   | 1.43%  | 0.129% |
| 49 | 15.045 | 2199 | 2203 | 2206 | rBV  | 184563  | 286228  | 6.90%  | 0.623% |
| 50 | 15.157 | 2219 | 2222 | 2227 | rBV  | 35472   | 50855   | 1.23%  | 0.111% |
| 51 | 15.345 | 2250 | 2254 | 2260 | rVB  | 104321  | 138103  | 3.33%  | 0.300% |
| 52 | 15.404 | 2260 | 2264 | 2269 | rVV  | 151562  | 197321  | 4.76%  | 0.429% |
| 53 | 15.463 | 2269 | 2274 | 2289 | rVB2 | 696322  | 1068586 | 25.76% | 2.325% |
| 54 | 15.880 | 2342 | 2345 | 2351 | rVB  | 66617   | 99586   | 2.40%  | 0.217% |
| 55 | 16.380 | 2425 | 2430 | 2435 | rVB2 | 50300   | 92096   | 2.22%  | 0.200% |
| 56 | 16.945 | 2521 | 2526 | 2536 | rVB2 | 77305   | 180433  | 4.35%  | 0.393% |
| 57 | 17.127 | 2553 | 2557 | 2563 | rBV7 | 22324   | 42127   | 1.02%  | 0.092% |
| 58 | 17.392 | 2597 | 2602 | 2614 | rVB  | 49271   | 114269  | 2.75%  | 0.249% |

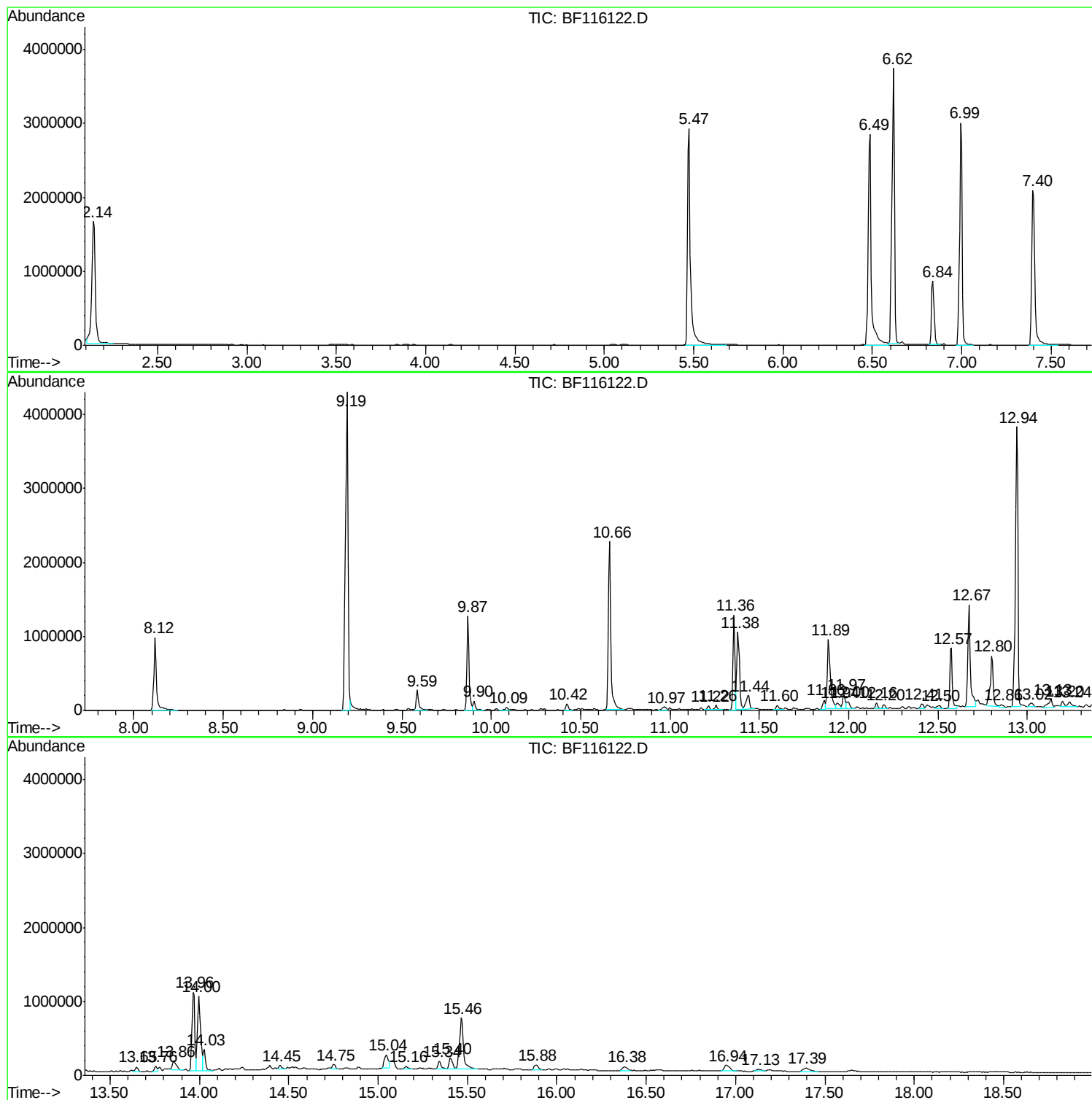
Sum of corrected areas: 45968958

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ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF080919\  
 Data File : BF116122.D  
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 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
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 TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
 Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 2.14 | 61.54 ng | 2459370 | 1,4-Dichlorobenzene-d4 | 6.84 |

| Hit# of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|---------|---|-----------------------------|-----|---------|-------------|------|
| 1       |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 83   |
| 2       |   | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 39   |
| 3       |   | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 25   |
| 4       |   | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 23   |
| 5       |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 17   |

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 Peak Number 2 3-Chloro-6-fluoro-pyrazine Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.62 | 88.63 ng | 3541810 | 1,4-Dichlorobenzene-d4 | 6.84 |

| Hit# of | 5 | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|---------|---|----------------------------------|-----|-----------|--------------|------|
| 1       |   | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2       |   | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 3       |   | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 12   |
| 4       |   | 4-Chloro-2-fluoro-pyrimidine     | 132 | C4H2ClFN2 | 051422-00-5  | 9    |
| 5       |   | 5-Fluoro-2-chloropyrimidine      | 132 | C4H2ClFN2 | 062802-42-0  | 9    |

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 Peak Number 3 1,4-Dichlorobenzene-D4 Concentration Rank 2

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.99 | 74.74 ng | 2986830 | 1,4-Dichlorobenzene-d4 | 6.84 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|---------|---|-------------------------------------|-----|-----------|-------------|------|
| 1       |   | 1,4-Dichlorobenzene-D4              | 150 | C6D4Cl2   | 003855-82-1 | 94   |
| 2       |   | 1,2-Dichlorobenzene-D4              | 150 | C6D4Cl2   | 002199-69-1 | 94   |
| 3       |   | 9-Borabicyclo[3.3.1]nonane, 9-et... | 150 | C10H19B   | 052102-17-7 | 38   |
| 4       |   | 2-Chloro-4,6-difluoro-pyrimidine    | 150 | C4HClF2N2 | 038953-30-9 | 38   |
| 5       |   | 2-Bromo-2'-nitroacetophenone        | 243 | C8H6BrNO3 | 006851-99-6 | 12   |

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 Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|------|---------|------|------------------|------|
|------|---------|------|------------------|------|

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 Sample : K4257-02  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF073019.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

| 11.86   | 2.02 ng | 112184                  | Phenanthrene-d10 | 11.36          |
|---------|---------|-------------------------|------------------|----------------|
| Hit# of | 5       | Tentative ID            | MW MolForm       | CAS# Qual      |
| 1       |         | Phenanthrene, 2-methyl- | 192 C15H12       | 002531-84-2 93 |
| 2       |         | Phenanthrene, 1-methyl- | 192 C15H12       | 000832-69-9 93 |
| 3       |         | Anthracene, 9-methyl-   | 192 C15H12       | 000779-02-2 93 |
| 4       |         | Anthracene, 1-methyl-   | 192 C15H12       | 000610-48-0 93 |
| 5       |         | Anthracene, 2-methyl-   | 192 C15H12       | 000613-12-7 93 |

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 Peak Number 5 n-Hexadecanoic acid Concentration Rank 5

| R.T.    | EstConc  | Area                | Relative to ISTD | R.T.           |
|---------|----------|---------------------|------------------|----------------|
| 11.89   | 19.65 ng | 1092250             | Phenanthrene-d10 | 11.36          |
| Hit# of | 5        | Tentative ID        | MW MolForm       | CAS# Qual      |
| 1       |          | n-Hexadecanoic acid | 256 C16H32O2     | 000057-10-3 99 |
| 2       |          | Tridecanoic acid    | 214 C13H26O2     | 000638-53-9 92 |
| 3       |          | Pentadecanoic acid  | 242 C15H30O2     | 001002-84-2 81 |
| 4       |          | Tetradecanoic acid  | 228 C14H28O2     | 000544-63-8 58 |
| 5       |          | Dodecanoic acid     | 200 C12H24O2     | 000143-07-7 58 |

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 Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.           |
|---------|---------|-------------------------------------|------------------|----------------|
| 11.97   | 3.94 ng | 219299                              | Phenanthrene-d10 | 11.36          |
| Hit# of | 5       | Tentative ID                        | MW MolForm       | CAS# Qual      |
| 1       |         | 4H-Cyclopenta[def]phenanthrene      | 190 C15H10       | 000203-64-5 92 |
| 2       |         | Methyl diselenide                   | 190 C2H6Se2      | 007101-31-7 43 |
| 3       |         | 2,2'-Bis(4,5-dimethylimidazole)     | 190 C10H14N4     | 069286-06-2 37 |
| 4       |         | 1,4-Naphthalenedione, 2,5-dihydr... | 190 C10H6O4      | 004923-55-1 25 |
| 5       |         | Megastigmatrienone                  | 190 C13H18O      | 038818-55-2 18 |

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 Peak Number 7 Octadecanoic acid Concentration Rank 4

| R.T.    | EstConc  | Area         | Relative to ISTD | R.T.      |
|---------|----------|--------------|------------------|-----------|
| 12.67   | 26.96 ng | 1498840      | Phenanthrene-d10 | 11.36     |
| Hit# of | 5        | Tentative ID | MW MolForm       | CAS# Qual |

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TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

|   |                    |     |          |             |    |
|---|--------------------|-----|----------|-------------|----|
| 1 | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 99 |
| 2 | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 93 |
| 3 | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 87 |
| 4 | Tridecanoic acid   | 214 | C13H26O2 | 000638-53-9 | 74 |
| 5 | n-Decanoic acid    | 172 | C10H20O2 | 000334-48-5 | 62 |

\*\*\*\*\*  
 Peak Number 8 11H-Benzo[a]fluorene Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.13 | 2.69 ng | 161427 | Chrysene-d12     | 14.00 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 94   |
| 2    |    |   | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 94   |
| 3    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 91   |
| 4    |    |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 90   |
| 5    |    |   | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 87   |

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 Peak Number 9 11H-Benzo[a]fluoren-11-one Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.86 | 3.17 ng | 190102 | Chrysene-d12     | 14.00 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 11H-Benzo[a]fluoren-11-one          | 230 | C17H10O  | 000479-79-8 | 91   |
| 2    |    |   | Dibenzo[c,h][2,6]naphthyridine      | 230 | C16H10N2 | 000218-30-4 | 50   |
| 3    |    |   | 6H-Benz[de]anthracen-6-one          | 230 | C17H10O  | 080252-14-8 | 45   |
| 4    |    |   | 7H-Benz[de]anthracen-7-one          | 230 | C17H10O  | 000082-05-3 | 43   |
| 5    |    |   | (E)-.alpha.,.alpha.'-Dicyanostil... | 230 | C16H10N2 | 002450-55-7 | 38   |

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 Peak Number 10 Benzo[e]pyrene Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.34 | 2.58 ng | 138103 | Perylene-d12     | 15.46 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Benzo[e]pyrene       | 252 | C20H12  | 000192-97-2 | 98   |
| 2    |    |   | Perylene             | 252 | C20H12  | 000198-55-0 | 97   |
| 3    |    |   | Benzo[k]fluoranthene | 252 | C20H12  | 000207-08-9 | 97   |

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TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

|                            |            |                |
|----------------------------|------------|----------------|
| 4 Benz[e]acephenanthrylene | 252 C20H12 | 000205-99-2 95 |
| 5 Benzo[a]pyrene           | 252 C20H12 | 000050-32-8 94 |

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| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| Butane, 2-methoxy... | 2.14  | 61.5    | ng    | 2459370  | 1                     | 6.84  | 799245  | 20.0 |
| 3-Chloro-6-fluoro... | 6.62  | 88.6    | ng    | 3541810  | 1                     | 6.84  | 799245  | 20.0 |
| 1,4-Dichlorobenze... | 6.99  | 74.7    | ng    | 2986830  | 1                     | 6.84  | 799245  | 20.0 |
| Phenanthrene, 2-m... | 11.86 | 2.0     | ng    | 112184   | 4                     | 11.36 | 1111980 | 20.0 |
| n-Hexadecanoic acid  | 11.89 | 19.6    | ng    | 1092250  | 4                     | 11.36 | 1111980 | 20.0 |
| 4H-Cyclopenta[def... | 11.97 | 3.9     | ng    | 219299   | 4                     | 11.36 | 1111980 | 20.0 |
| Octadecanoic acid    | 12.67 | 27.0    | ng    | 1498840  | 4                     | 11.36 | 1111980 | 20.0 |
| 11H-Benzo[a]fluorene | 13.13 | 2.7     | ng    | 161427   | 5                     | 14.00 | 1198000 | 20.0 |
| 11H-Benzo[a]fluor... | 13.86 | 3.2     | ng    | 190102   | 5                     | 14.00 | 1198000 | 20.0 |
| Benzo[e]pyrene       | 15.34 | 2.6     | ng    | 138103   | 6                     | 15.46 | 1068590 | 20.0 |