LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG100716\
Data File : BG024220.D
Acq On : 7 Oct 2016 4:47
Operator : UM/SJ
Sample : H5082-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDPJ7

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

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

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.922	181	184	191	rBV	72877	119616	2.26%	0.169%
2	4.057	205	207	210	rBV4	13883	11949	0.23%	0.017%
3	5.203	399	402	405	rBV3	12594	15658	0.30%	0.022%
4	5.373	427	431	436	rVB3	23229	38150	0.72%	0.054%
5	5.685	482	484	487	rVB4	12745	12316	0.23%	0.017%
6	5.761	496	497	499	rBV2	12106	9355	0.18%	0.013%
7	5.920	520	524	526	rBV4	8546	10934	0.21%	0.015%
8	6.225	573	576	579	rBV5	12631	18366	0.35%	0.026%
9	6.255	579	581	583	rVV3	12566	10028	0.19%	0.014%
10	6.343	592	596	598	rBV4	13138	14213	0.27%	0.020%
11	6.625	641	644	647	rVB5	10060	14082	0.27%	0.020%
12	6.654	647	649	650	rBV2	10654	9295	0.18%	0.013%
13	6.866	679	685	688	rVB8	11290	15692	0.30%	0.022%
14	6.913	690	693	695	rBV4	15736	23311	0.44%	0.033%
15	7.101	723	725	728	rVB4	14288	11248	0.21%	0.016%
16	7.424	773	780	794	rVB	849947	1590962	30.00%	2.253%
17	7.612	806	812	819	rBV	570244	978531	18.45%	1.386%
18	7.694	824	826	831	rBV6	11930	10782	0.20%	0.015%
19	7.823	840	848	859	rVB	750652	1363438	25.71%	1.931%
20	8.299	921	929	940	rBV	620314	1134122	21.39%	1.606%
21	8.593	976	979	981	rBV3	9348	13101	0.25%	0.019%
22	8.810	1013	1016	1019	rBV5	9626	9859	0.19%	0.014%
23	8.863	1022	1025	1027	rVB4	11367	15542	0.29%	0.022%
24	8.987	1035	1046	1054	rBV	1015884	1814534	34.22%	2.570%
25	9.469	1121	1128	1135	rBV	793384	1457492	27.49%	2.064%
26	9.527	1136	1138	1145	rVB6	23847	36402	0.69%	0.052%
27	9.668	1161	1162	1166	rBV4	10898	11135	0.21%	0.016%
28	9.874	1195	1197	1200	rVB4	10496	10192	0.19%	0.014%
29	10.197	1244	1252	1259	rBV	667975	1246244	23.50%	1.765%
30	10.373	1278	1282	1285	rBV6	9413	16121	0.30%	0.023%
31	10.732	1335	1343	1352	rVV2	993246	1867160	35.21%	2.644%
32	10.879	1366	1368	1373	rVV6	14213	19041	0.36%	0.027%
33	10.949	1377	1380	1382	rVB4	15837	11389	0.21%	0.016%
34	10.984	1382	1386	1387	rBV4	8566	10022	0.19%	0.014%

LSC Area Percent Report

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Data File : BG024220.D
Acq On : 7 Oct 2016 4:47
Operator : UM/SJ
Sample : H5082-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDPJ7

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

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

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Title : SVOA CALIBRATION

35	11.078	1398	1402	1403	rBV4	11001	15005	0.28%	0.021%
36	11.125	1403	1410	1419	rVV	888804	1658936	31.29%	2.349%
37	11.190	1419	1421	1424	rVB4	14042	13671	0.26%	0.019%
38	11.255	1424	1432	1440	rBV	990046	1860926	35.10%	2.635%
39	11.431	1460	1462	1466	rVB5	12116	13262	0.25%	0.019%
40	11.478	1469	1470	1474	rBV4	8875	12278	0.23%	0.017%
41	11.590	1482	1489	1491	rBV5	17214	26704	0.50%	0.038%
42	12.353	1617	1619	1621	rBV3	10738	13231	0.25%	0.019%
43	12.583	1653	1658	1660	rVB6	6995	12215	0.23%	0.017%
44	12.688	1670	1676	1681	rBV8	29272	53589	1.01%	0.076%
45	12.812	1695	1697	1702	rVB5	10031	11698	0.22%	0.017%
46	13.023	1730	1733	1734	rVB3	11755	9526	0.18%	0.013%
47	13.041	1734	1736	1738	rBV3	7586	9278	0.17%	0.013%
48	13.094	1743	1745	1749	rVB4	6881	9573	0.18%	0.014%
49	13.235	1766	1769	1770	rVB3	13960	10937	0.21%	0.015%
50	13.270	1772	1775	1780	rVV7	12906	22853	0.43%	0.032%
51	13.517	1813	1817	1821	rVB6	16943	20644	0.39%	0.029%
52	13.781	1857	1862	1867	rBV2	73265	110468	2.08%	0.156%
53	13.910	1881	1884	1887	rVV4	6449	10309	0.19%	0.015%
54	14.233	1936	1939	1941	rBV4	11923	16602	0.31%	0.024%
55	14.251	1941	1942	1944	rVV2	14833	12723	0.24%	0.018%
56	14.310	1946	1952	1956	rVV	2049479	3124744	58.93%	4.425%
57	14.357	1956	1960	1968	rVB	1602200	2312385	43.61%	3.275%
58	14.521	1986	1988	1991	rVB4	10183	11841	0.22%	0.017%
59	14.610	1996	2003	2012	rVB	1965659	3289398	62.04%	4.658%
60	14.774	2026	2031	2035	rBV6	23930	41273	0.78%	0.058%
61	14.915	2047	2055	2063	rBV	1536010	2510548	47.35%	3.555%
62	15.074	2075	2082	2096	rBV2	1153152	1787491	33.71%	2.531%
63	15.315	2120	2123	2126	rBV5	11452	15518	0.29%	0.022%
64	15.461	2145	2148	2151	rBV4	13061	19043	0.36%	0.027%
65	15.632	2173	2177	2181	rVB3	44482	56145	1.06%	0.080%
66	15.679	2181	2185	2187	rBV4	15103	18837	0.36%	0.027%
67	15.720	2188	2192	2197	rVB2	55931	71510	1.35%	0.101%
68	15.785	2201	2203	2205	rBV2	9142	9309	0.18%	0.013%
69	15.896	2216	2222	2232	rVB	2746173	4400516	82.99%	6.232%
70	15.996	2232	2239	2250	rBV	980454	1560226	29.43%	2.210%
71	16.131	2259	2262	2267	rVB6	44534	77337	1.46%	0.110%

LSC Area Percent Report

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Acq On : 7 Oct 2016 4:47
Operator : UM/SJ
Sample : H5082-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDPJ7

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

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Peak separation: 5

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Title : SVOA CALIBRATION

72	16.278	2283	2287	2289	rBV5	14909	22777	0.43%	0.032%
73	16.578	2333	2338	2346	rBV2	82944	144330	2.72%	0.204%
74	16.766	2366	2370	2374	rVB7	19871	31047	0.59%	0.044%
75	16.836	2378	2382	2387	rBV8	25827	55674	1.05%	0.079%
76	16.889	2390	2391	2394	rBV3	10686	12815	0.24%	0.018%
77	16.924	2394	2397	2403	rBV8	13359	31292	0.59%	0.044%
78	17.154	2435	2436	2439	rBV3	10448	13252	0.25%	0.019%
79	17.371	2469	2473	2476	rBV	86609	105203	1.98%	0.149%
80	17.653	2514	2521	2529	rVB2	1940100	3094970	58.37%	4.383%
81	17.753	2531	2538	2546	rVV2	3169599	4782471	90.19%	6.773%
82	17.894	2559	2562	2566	rBV	112361	143845	2.71%	0.204%
83	17.941	2566	2570	2575	rVV	238077	315298	5.95%	0.447%
84	17.982	2575	2577	2580	rVB4	16070	12193	0.23%	0.017%
85	18.105	2595	2598	2604	rVB2	57137	69836	1.32%	0.099%
86	18.499	2660	2665	2677	rBV2	360769	560207	10.57%	0.793%
87	19.198	2780	2784	2787	rBV2	155115	177493	3.35%	0.251%
88	19.239	2787	2791	2795	rVV	311864	372539	7.03%	0.528%
89	19.657	2858	2862	2865	rBV4	50347	72484	1.37%	0.103%
90	19.756	2875	2879	2886	rBV2	334717	480933	9.07%	0.681%
91	20.021	2918	2924	2931	rVB	3544187	5302382	100.00%	7.509%
92	20.309	2968	2973	2975	rBV	1535916	1742132	32.86%	2.467%
93	20.344	2975	2979	2985	rVB	3130165	3514579	66.28%	4.977%
94	21.313	3141	3144	3147	rBV	529895	633405	11.95%	0.897%
95	21.355	3147	3151	3156	rVB	1095096	1375251	25.94%	1.948%
96	21.543	3180	3183	3189	rVB4	242661	334719	6.31%	0.474%
97	21.954	3247	3253	3260	rVB	1859150	3326334	62.73%	4.711%
98	22.530	3347	3351	3357	rVB2	437104	624005	11.77%	0.884%
99	25.174	3792	3801	3815	rVB2	1582485	4732435	89.25%	6.702%
100	25.415	3833	3842	3855	rVB3	1052645	3348916	63.16%	4.743%

Sum of corrected areas: 70613648

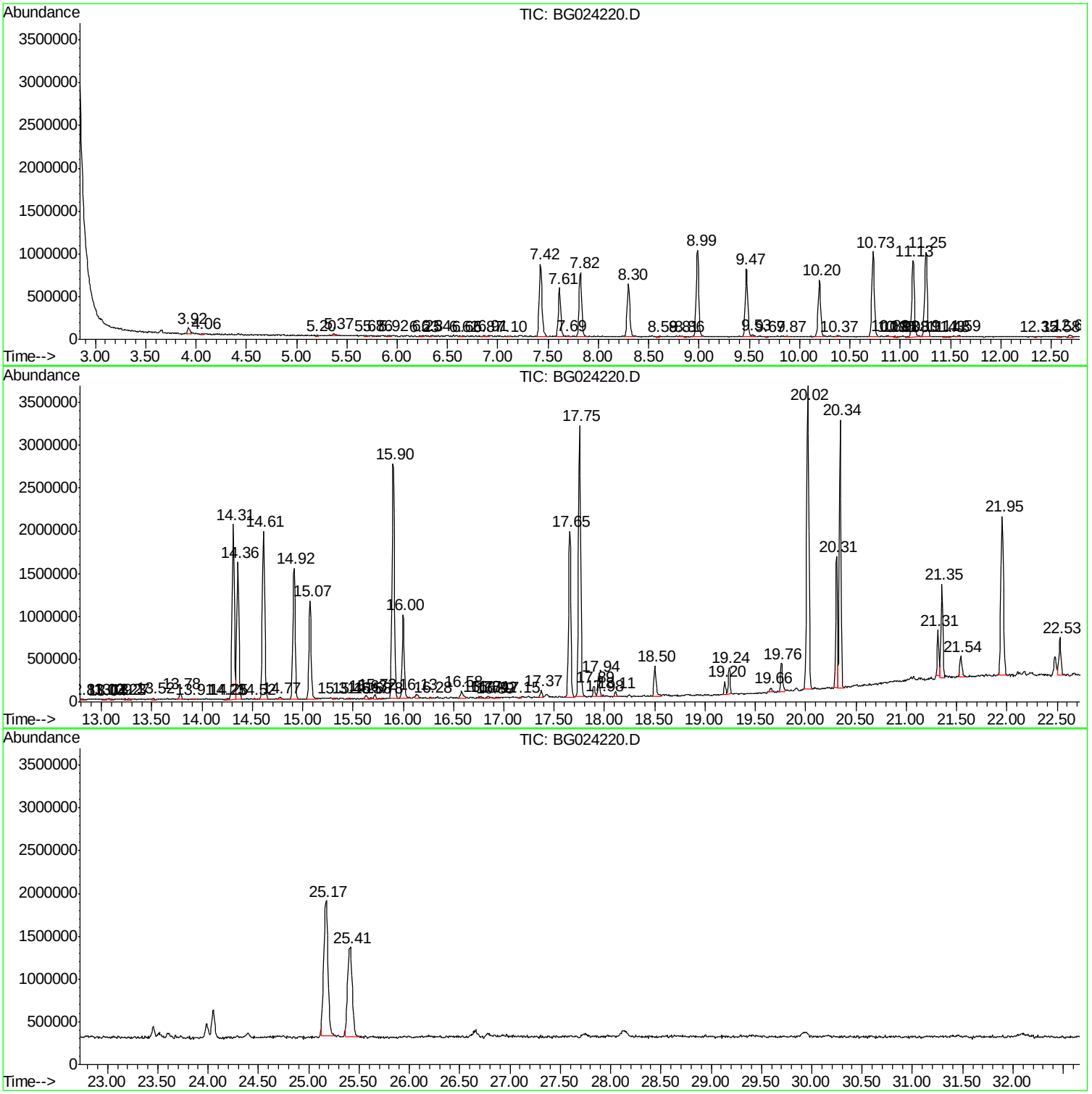
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG100716\
Data File : BG024220.D
Acq On : 7 Oct 2016 4:47
Operator : UM/SJ
Sample : H5082-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDPJ7

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG100716\
Data File : BG024220.D
Acq On : 7 Oct 2016 4:47
Operator : UM/SJ
Sample : H5082-18
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Instrument : BNA_G
ClientSampleId : BDPJ7

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

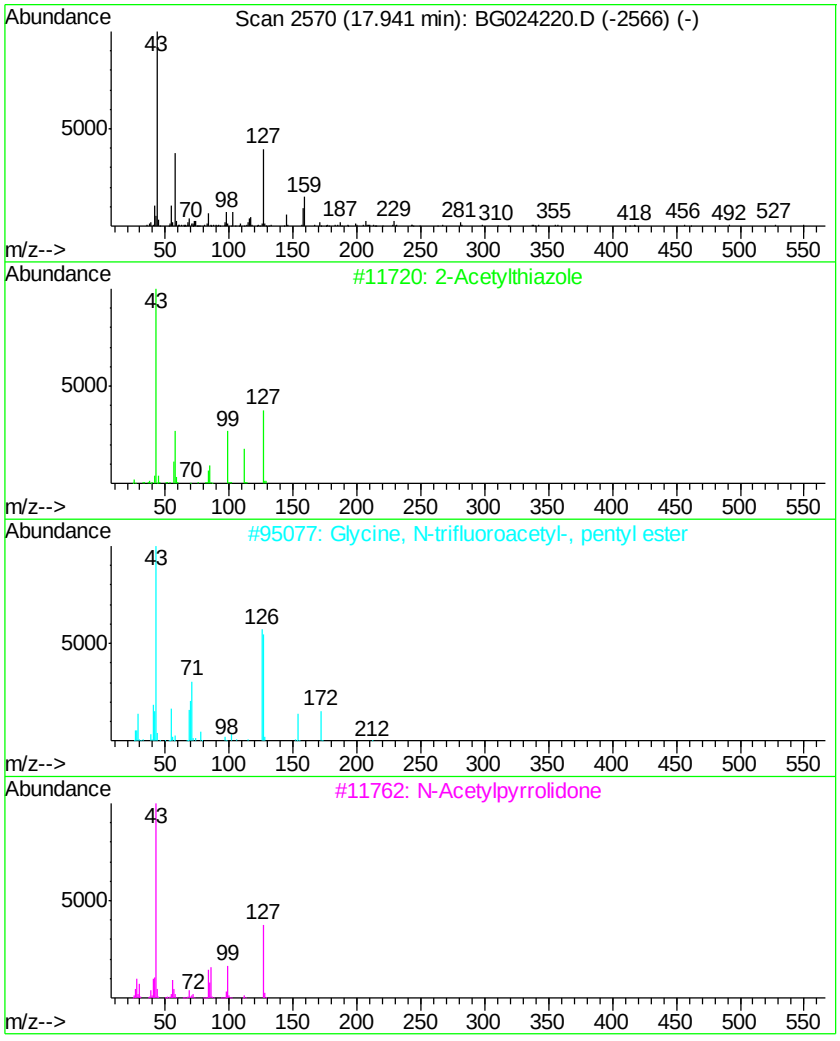
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

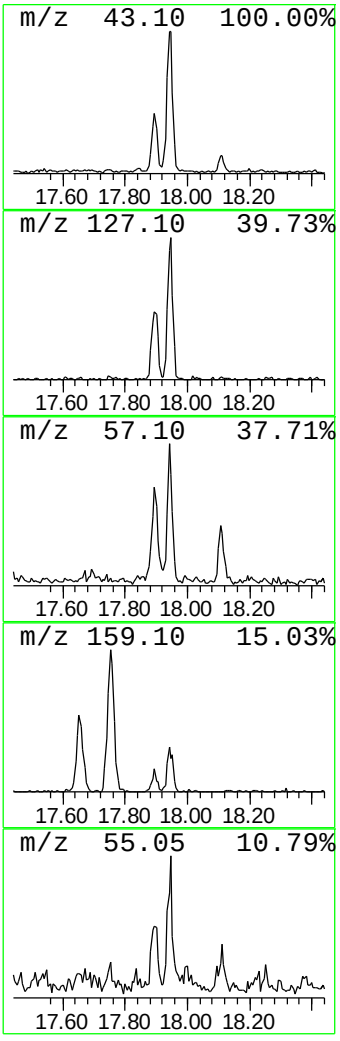
****

Peak Number 2 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.94	2.04 ng/ul	315298	Phenanthrene-d10	17.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Acetylthiazole	127	C5H5NOS	024295-03-2	32
2	Glycine, N-trifluoroacetyl-, pen...	241	C9H14F3NO3	1000328-03-6	23
3	N-Acetylpyrrolidone	127	C6H9NO2	000932-17-2	9
4	Uracil, 6-acetamido-	169	C6H7N3O3	032970-34-6	9
5	4,6(1H)-Pyrimidinedione, 3,4,5,6...	127	C4H5N3O2	004425-67-6	7





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG100716\
Data File : BG024220.D
Acq On : 7 Oct 2016 4:47
Operator : UM/SJ
Sample : H5082-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDPJ7

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

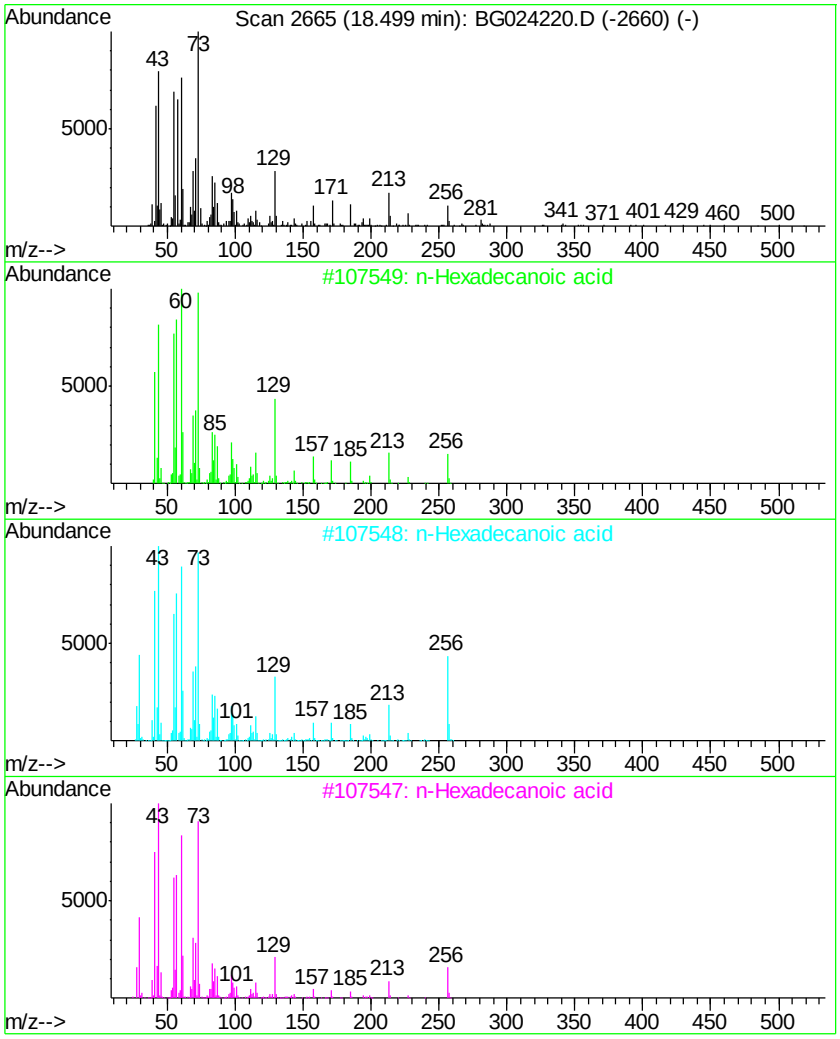
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

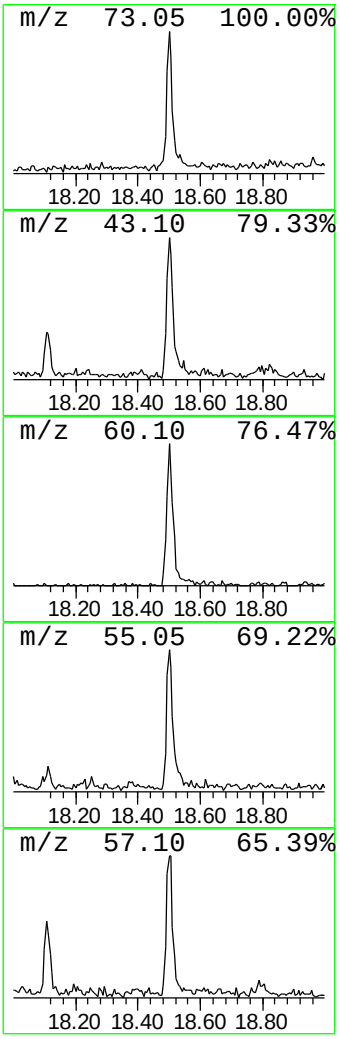
****

Peak Number 3 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.50	3.62 ng/ul	560207	Phenanthrene-d10	17.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
4	Tridecanoic acid	214	C13H26O2	000638-53-9	87
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	76





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG100716\
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Acq On : 7 Oct 2016 4:47
Operator : UM/SJ
Sample : H5082-18
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Instrument : BNA_G
ClientSampleId : BDPJ7

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

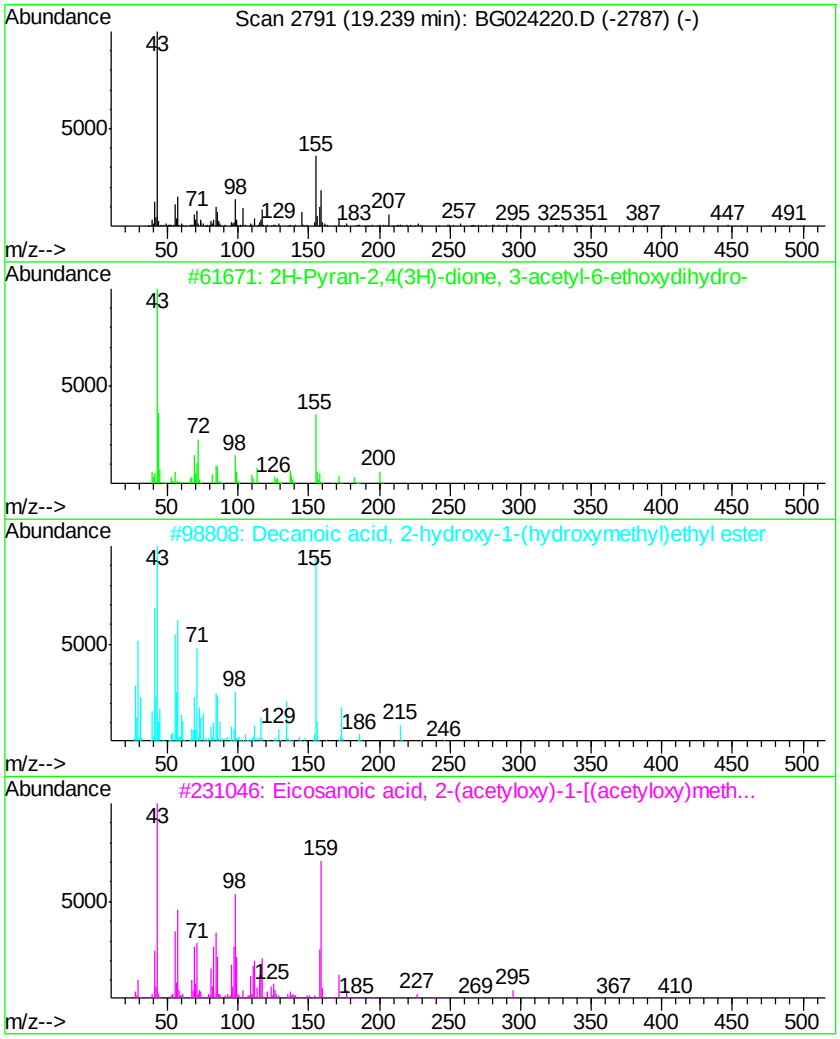
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

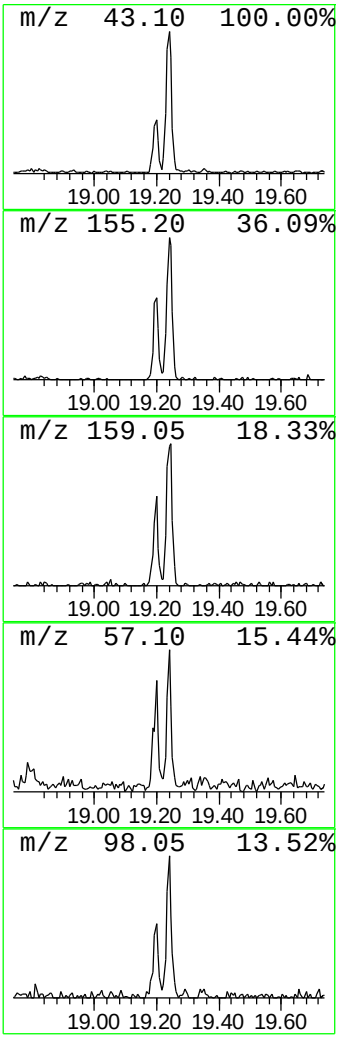
****

Peak Number 4 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.24	2.41 ng/ul	372539	Phenanthrene-d10	17.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran-2,4(3H)-dione, 3-acetyl...	200	C9H12O5	1000318-74-8	25
2	Decanoic acid, 2-hydroxy-1-(hydr...	246	C13H26O4	003376-48-5	23
3	Eicosanoic acid, 2-(acetyloxy)-1...	470	C27H50O6	055429-68-0	16
4	Eicosanoic acid, 2,3-bis(acetylo...	470	C27H50O6	055429-67-9	12
5	Octadecanoic acid, 2-hydroxy-1-(...	358	C21H42O4	000621-61-4	10





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG100716\
Data File : BG024220.D
Acq On : 7 Oct 2016 4:47
Operator : UM/SJ
Sample : H5082-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDPJ7

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

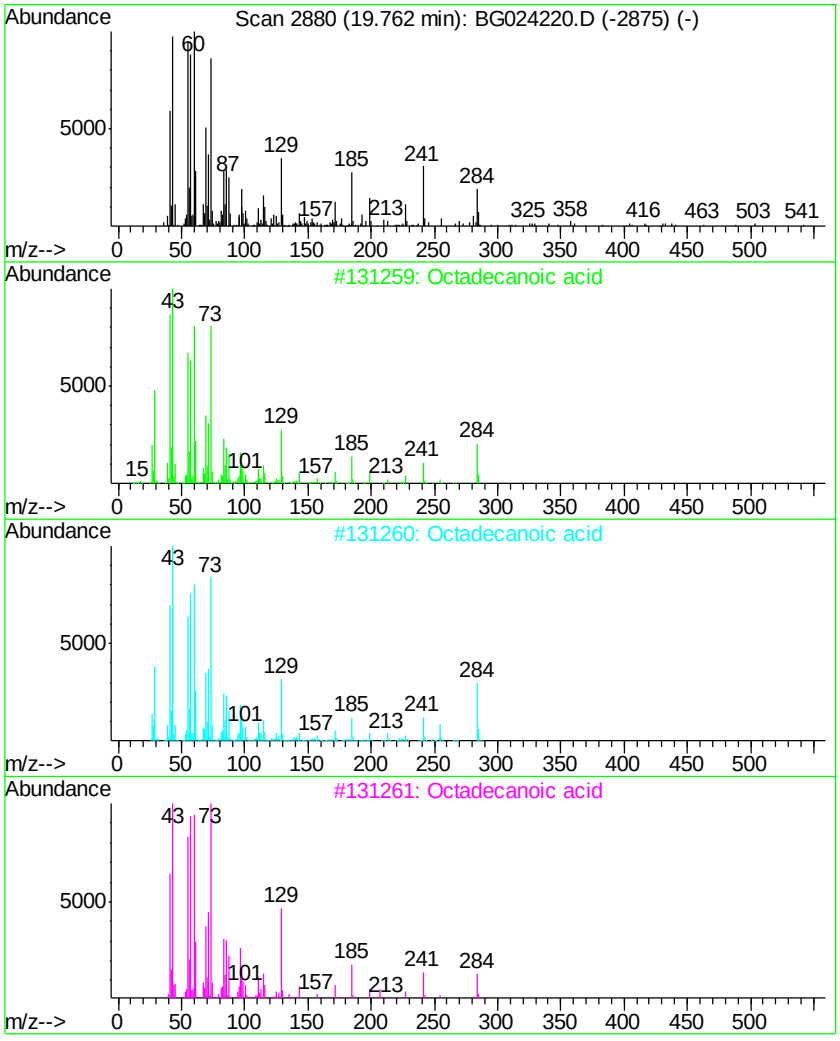
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

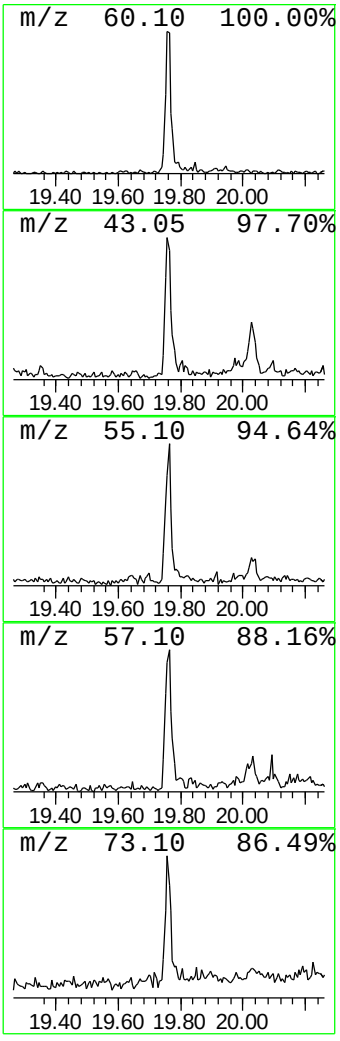
****

Peak Number 5 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.76	3.11 ng/ul	480933	Phenanthrene-d10	17.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	98
2	Octadecanoic acid	284	C18H36O2	000057-11-4	96
3	Octadecanoic acid	284	C18H36O2	000057-11-4	95
4	Octadecanoic acid	284	C18H36O2	000057-11-4	93
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	93





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG100716\
Data File : BG024220.D
Acq On : 7 Oct 2016 4:47
Operator : UM/SJ
Sample : H5082-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDPJ7

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

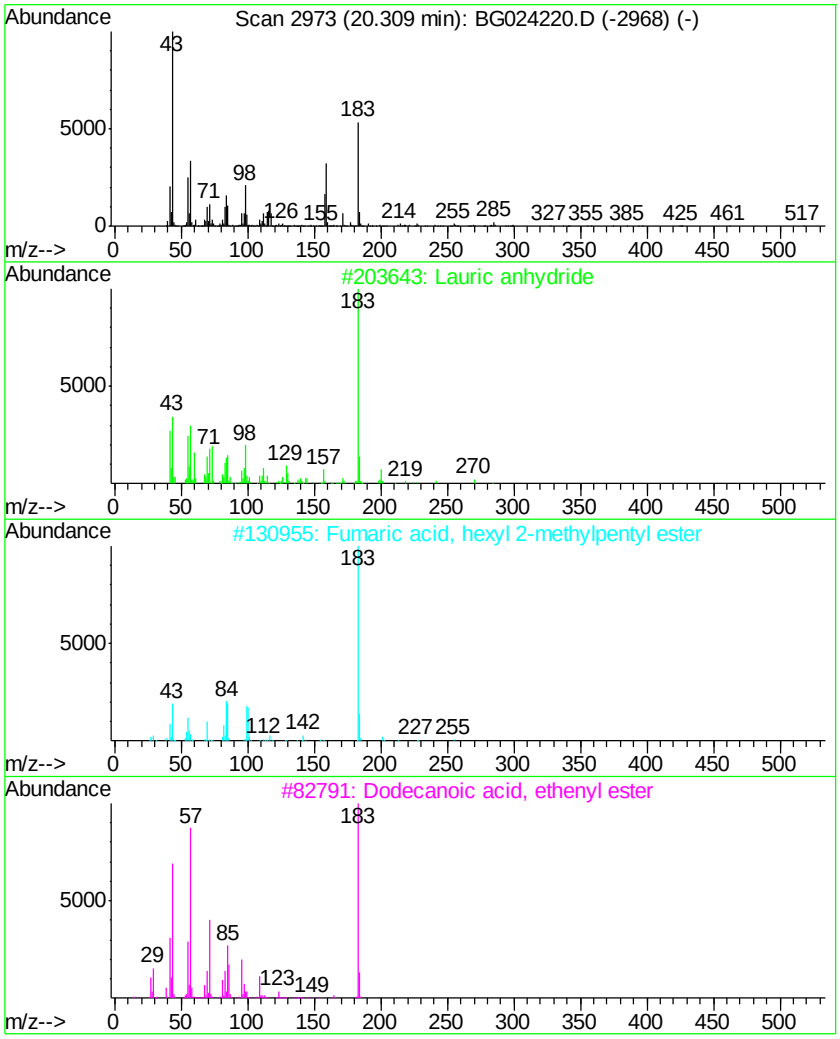
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

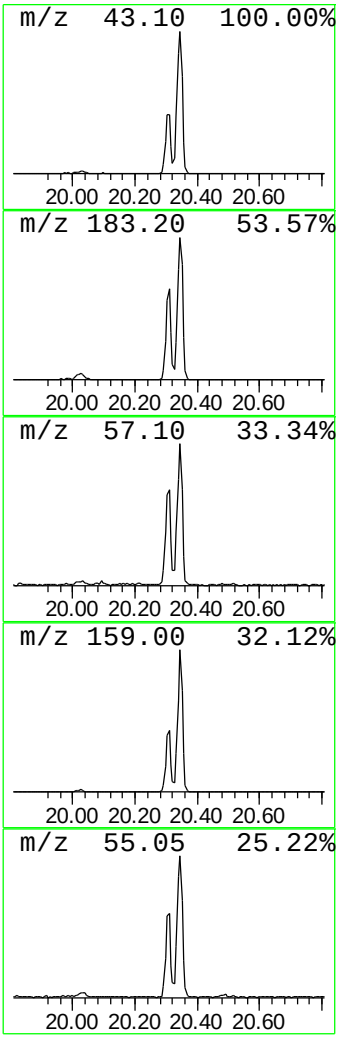
****

Peak Number 6 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.31	10.47 ng/ul	1742130	Chrysene-d12	21.95

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Lauric anhydride	382	C24H46O3	000645-66-9	43
2	Fumaric acid, hexyl 2-methylpent...	284	C16H28O4	1000348-72-5	38
3	Dodecanoic acid, ethenyl ester	226	C14H26O2	002146-71-6	35
4	Fumaric acid, 4-heptyl hexyl ester	298	C17H30O4	1000348-51-5	32
5	Fumaric acid, cyclobutyl isohexy...	254	C14H22O4	1000345-03-5	32





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG100716\
Data File : BG024220.D
Acq On : 7 Oct 2016 4:47
Operator : UM/SJ
Sample : H5082-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDPJ7

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

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

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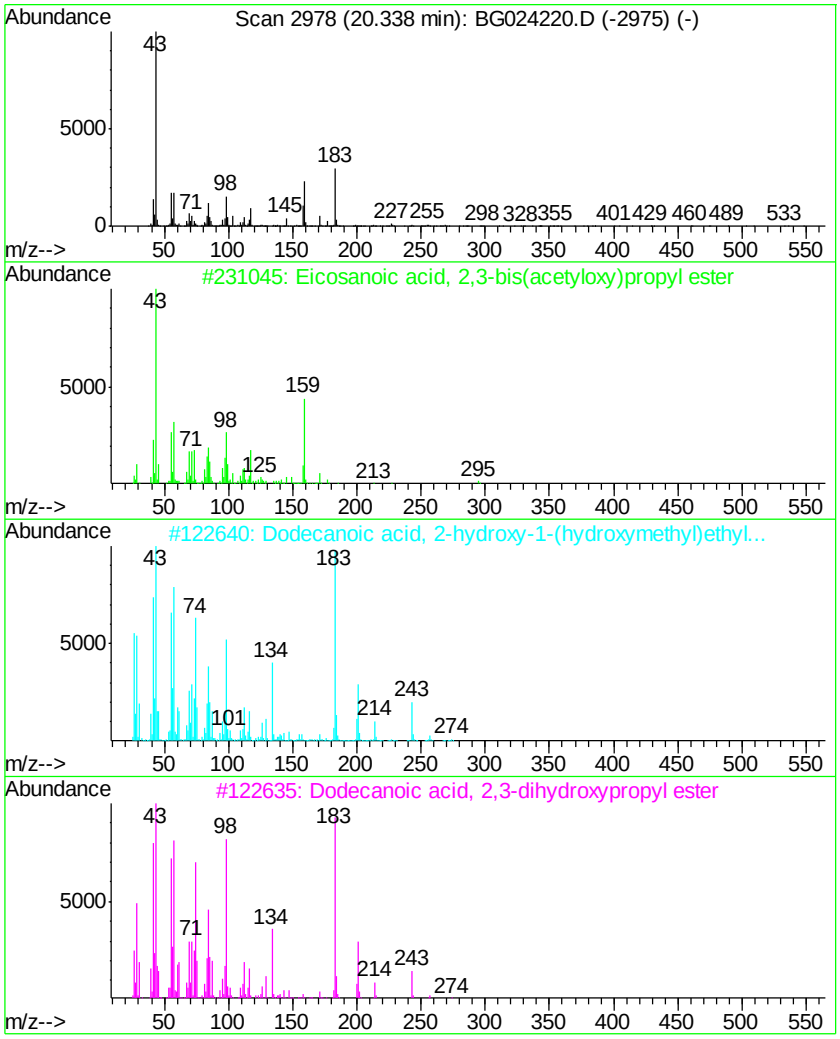
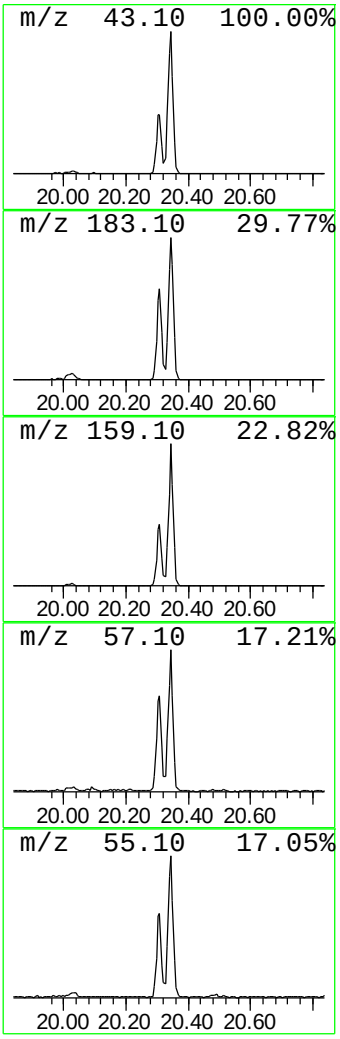
Peak Number 7 unknown-04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.34	21.13 ng/ul	3514580	Chrysene-d12	21.95

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosanoic acid, 2,3-bis(acetylo...	470	C27H50O6	055429-67-9	33
2	Dodecanoic acid, 2-hydroxy-1-(hy...	274	C15H30O4	001678-45-1	32
3	Dodecanoic acid, 2,3-dihydroxypr...	274	C15H30O4	000142-18-7	25
4	9-Octadecenoic acid (Z)-, 2,3-bi...	440	C25H44O6	055401-64-4	23
5	Eicosanoic acid, 2-(acetyloxy)-1...	470	C27H50O6	055429-68-0	10





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG100716\
Data File : BG024220.D
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Operator : UM/SJ
Sample : H5082-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDPJ7

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

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

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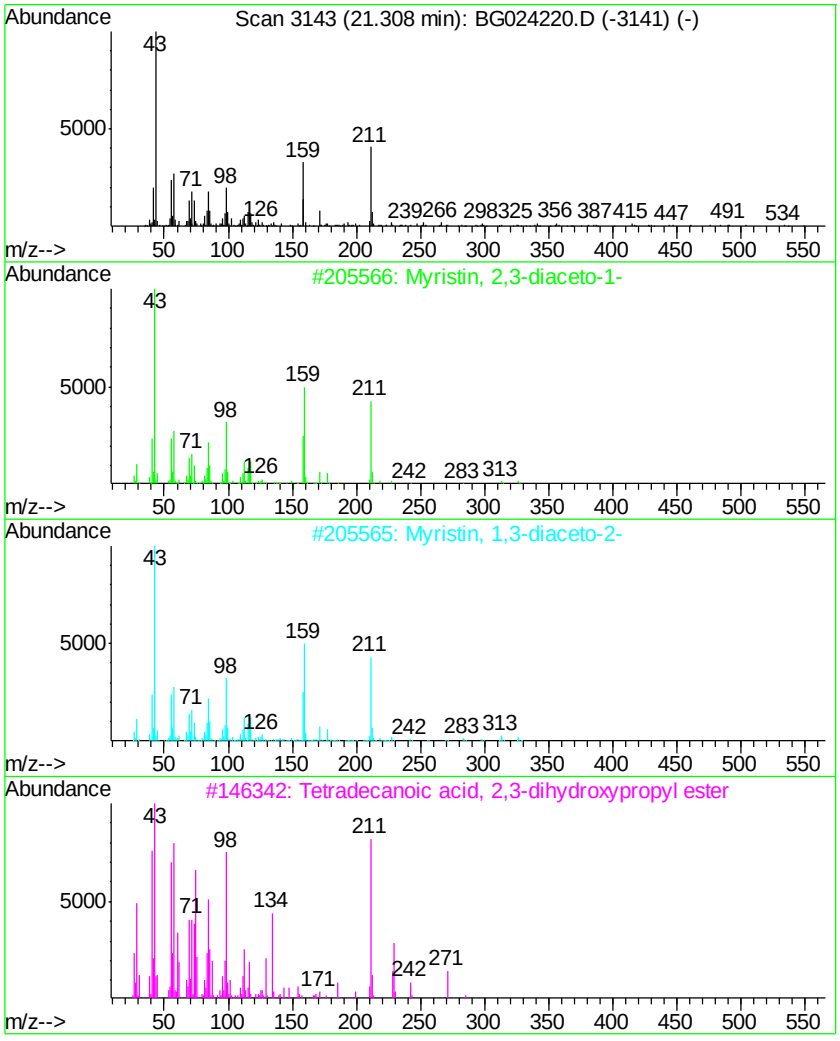
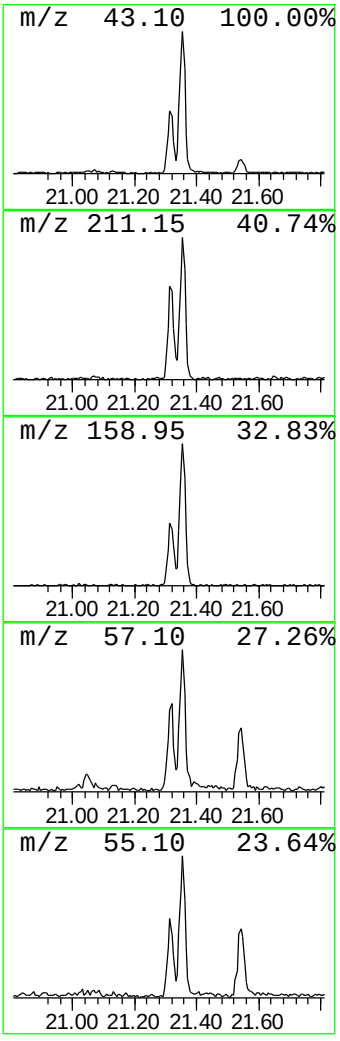
Peak Number 8 Myristin, 2,3-diaceto-1- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.31	3.81 ng/ul	633405	Chrysene-d12	21.95

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Myristin, 2,3-diaceto-1-	386	C21H38O6	014473-55-3	76
2	Myristin, 1,3-diaceto-2-	386	C21H38O6	014290-23-4	76
3	Tetradecanoic acid, 2,3-dihydrox...	302	C17H34O4	000589-68-4	27
4	Tetradecanoic acid, 2-hydroxy-1-...	302	C17H34O4	003443-83-2	14
5	2-(4-Benzylphenyl)propan-2-ol	226	C16H18O	1000201-97-7	12





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Data File : BG024220.D
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Operator : UM/SJ
Sample : H5082-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDPJ7

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

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

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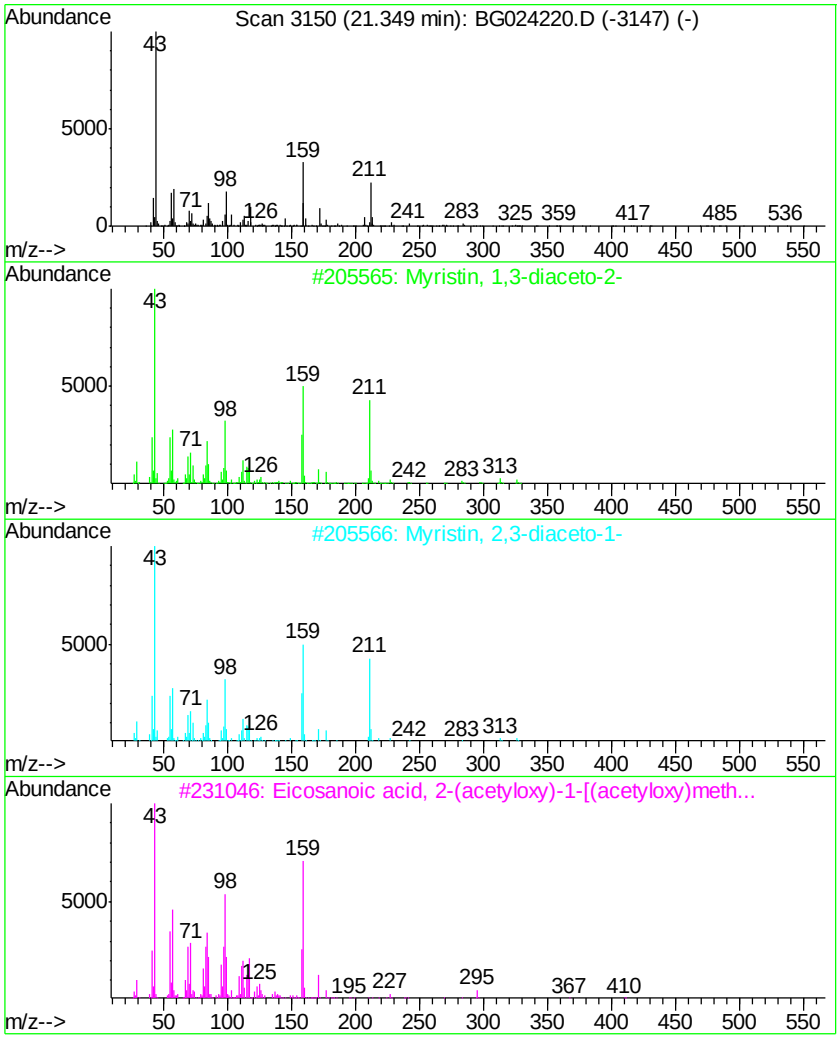
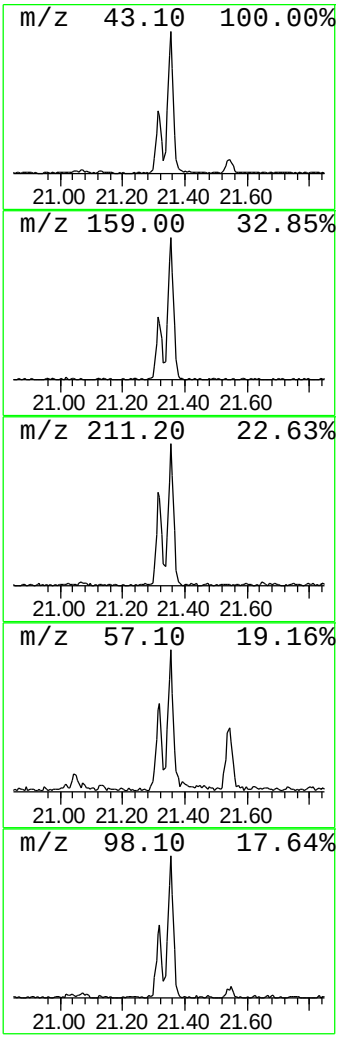
Peak Number 9 Myristin, 1,3-diaceto-2- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.35	8.27 ng/ul	1375250	Chrysene-d12	21.95

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Myristin, 1,3-diaceto-2-	386	C21H38O6	014290-23-4	72
2	Myristin, 2,3-diaceto-1-	386	C21H38O6	014473-55-3	56
3	Eicosanoic acid, 2-(acetyloxy)-1...	470	C27H50O6	055429-68-0	27
4	Tetradecanoic acid, 2,3-dihydrox...	302	C17H34O4	000589-68-4	22
5	Tetradecanoic acid, 2-hydroxy-1-...	302	C17H34O4	003443-83-2	12





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Sample : H5082-18
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDPJ7

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

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

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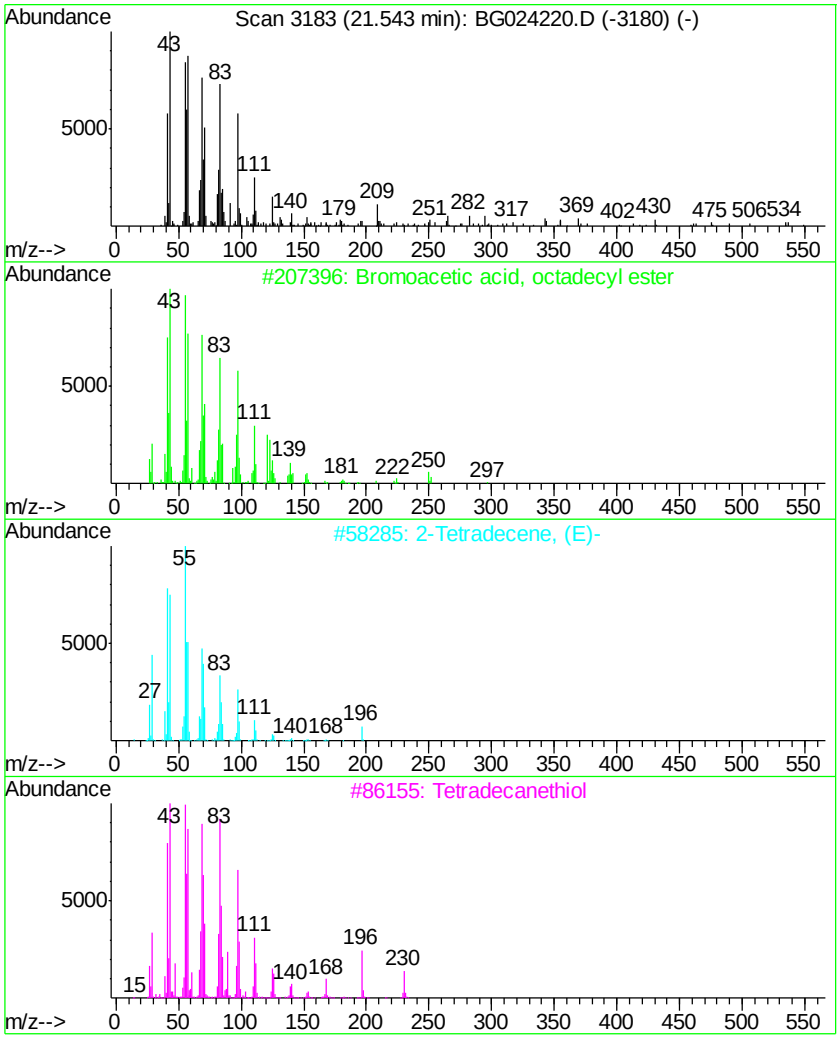
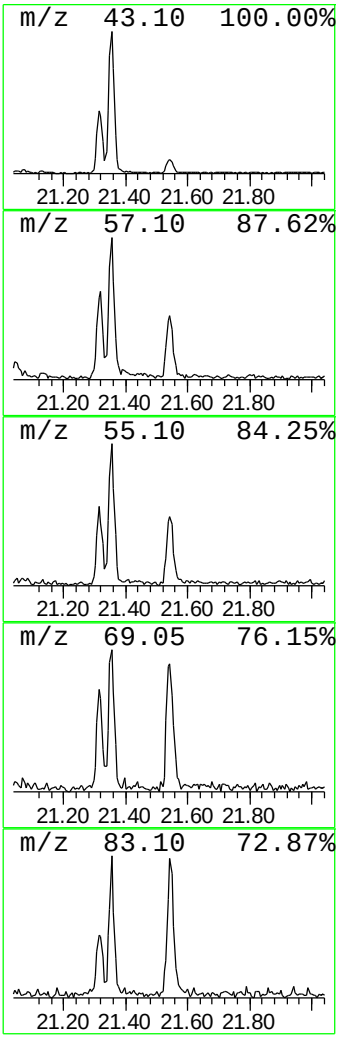
Peak Number 10 Bromoacetic acid, octadecyl... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.54	2.01 ng/ul	334719	Chrysene-d12	21.95

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
2	2-Tetradecene, (E)-	196	C14H28	035953-54-9	87
3	Tetradecanethiol	230	C14H30S	002079-95-0	87
4	Behenic alcohol	326	C22H46O	000661-19-8	83
5	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	83





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Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : BDPJ7

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG0100716.M
Quant Title : SVOA CALIBRATION

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

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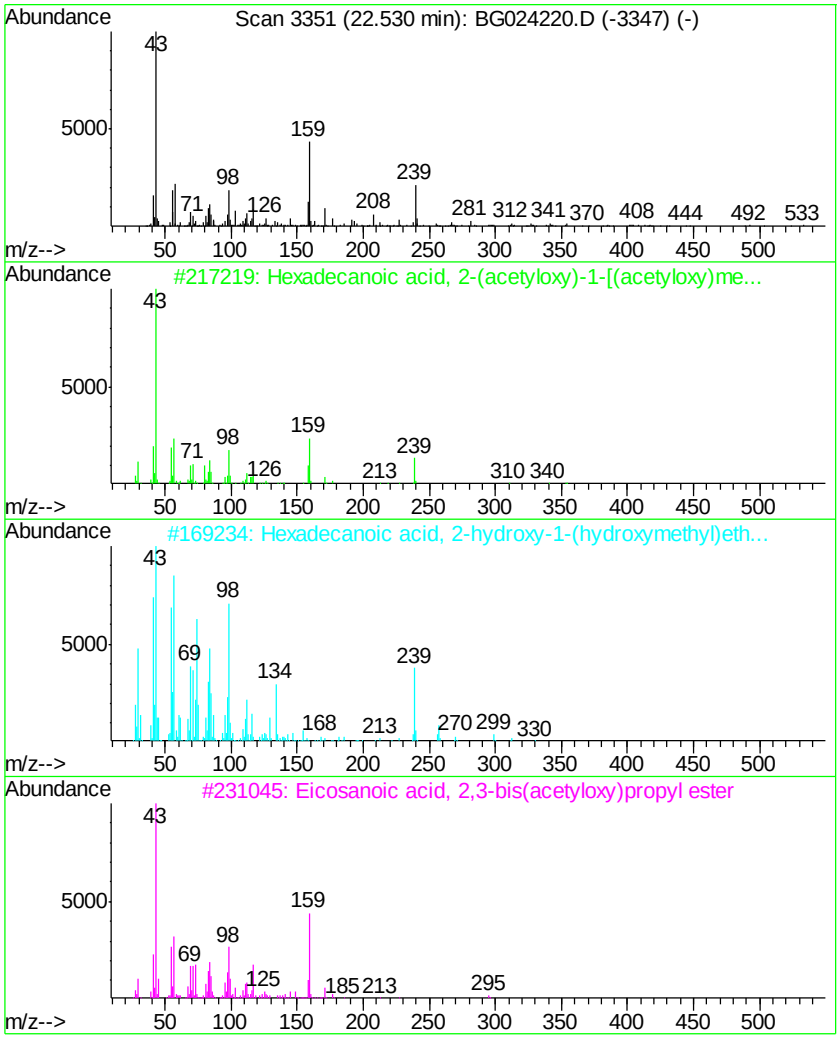
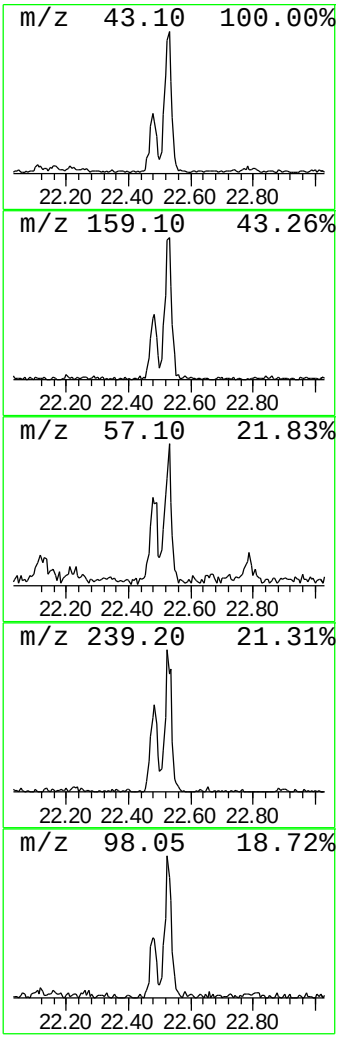
Peak Number 11 unknown-05 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.53	3.75 ng/ul	624005	Chrysene-d12	21.95

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, 2-(acetyloxy)...	414	C23H42O6	055268-69-4	25
2	Hexadecanoic acid, 2-hydroxy-1-(...	330	C19H38O4	023470-00-0	12
3	Eicosanoic acid, 2,3-bis(acetylo...	470	C27H50O6	055429-67-9	12
4	Octadecanoic acid, 2-(acetyloxy)...	442	C25H46O6	055401-62-2	10
5	4,5,6-Trimethyltetrahydro-1,3-ox...	159	C7H13NOS	085333-98-8	9





Tentatively Identified Compound (LSC) summary

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Quant Title : SVOA CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown-01	17.94	2.0	ng/ul	315298	4	17.65	3094970	20.0
n-Hexadecanoic acid	18.50	3.6	ng/ul	560207	4	17.65	3094970	20.0
unknown-02	19.24	2.4	ng/ul	372539	4	17.65	3094970	20.0
Octadecanoic acid	19.76	3.1	ng/ul	480933	4	17.65	3094970	20.0
unknown-03	20.31	10.5	ng/ul	1742130	5	21.95	3326330	20.0
unknown-04	20.34	21.1	ng/ul	3514580	5	21.95	3326330	20.0
Myristin, 2,3-dia...	21.31	3.8	ng/ul	633405	5	21.95	3326330	20.0
Myristin, 1,3-dia...	21.35	8.3	ng/ul	1375250	5	21.95	3326330	20.0
Bromoacetic acid,...	21.54	2.0	ng/ul	334719	5	21.95	3326330	20.0
unknown-05	22.53	3.8	ng/ul	624005	5	21.95	3326330	20.0