LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048241.D
Acq On : 20 Feb 2021 18:04
Operator : CG/JU
Sample : M1412-14
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBGX9

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.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	3.516	25	30	35	rVV3	5804	8460	0.71%	0.057%
2	4.051	117	121	129	rVV6	3589	6777	0.57%	0.046%
3	4.145	132	137	138	rBV3	1858	2284	0.19%	0.015%
4	4.169	138	141	146	rVB2	5195	8562	0.72%	0.058%
5	4.274	153	159	163	rVB3	5923	8645	0.72%	0.059%
6	5.203	312	317	318	rBV2	4630	6099	0.51%	0.041%
7	5.214	318	319	325	rVB3	3818	5408	0.45%	0.037%
8	5.297	331	333	338	rVB3	1673	2673	0.22%	0.018%
9	5.896	433	435	440	rBV	1232	2353	0.20%	0.016%
10	6.284	497	501	504	rBV	2135	2708	0.23%	0.018%
11	6.624	557	559	563	rBV2	2706	3601	0.30%	0.024%
12	6.777	581	585	591	rVB4	6522	11073	0.93%	0.075%
13	7.089	634	638	643	rVB5	7687	11703	0.98%	0.079%
14	7.259	660	667	670	rBV2	13202	22301	1.86%	0.151%
15	7.312	670	676	687	rVV	146106	277116	23.17%	1.878%
16	7.441	691	698	706	rVV	95217	162671	13.60%	1.102%
17	7.518	706	711	722	rVB3	5820	11322	0.95%	0.077%
18	7.659	728	735	742	rBV	140129	243471	20.35%	1.650%
19	7.717	742	745	749	rVB2	4380	6016	0.50%	0.041%
20	8.111	806	812	821	rVB	145926	249392	20.85%	1.690%
21	8.246	828	835	840	rVV3	14720	22548	1.88%	0.153%
22	8.322	844	848	853	rVB2	5200	7519	0.63%	0.051%
23	8.851	931	938	948	rBV2	179300	310500	25.96%	2.104%
24	8.951	948	955	959	rBV4	5079	8416	0.70%	0.057%
25	9.286	1004	1012	1020	rVV	139954	248301	20.76%	1.683%
26	9.433	1033	1037	1038	rBV2	1981	2250	0.19%	0.015%
27	9.656	1073	1075	1081	rVB2	2682	4530	0.38%	0.031%
28	9.838	1103	1106	1111	rVB3	5897	9674	0.81%	0.066%
29	10.009	1128	1135	1143	rVV	130370	230515	19.27%	1.562%
30	10.238	1168	1174	1178	rBV2	8343	14603	1.22%	0.099%
31	10.285	1180	1182	1185	rVV3	3633	4338	0.36%	0.029%
32	10.338	1185	1191	1198	rVV7	9030	17043	1.42%	0.116%
33	10.567	1223	1230	1241	rBV	187306	339696	28.40%	2.302%
34	10.708	1248	1254	1262	rBV3	18115	34428	2.88%	0.233%

LSC Area Percent Report

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Instrument : BNA_G
ClientSampleId : DBGX9

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Title : SVOA CALIBRATION

35	10.931	1284	1292	1298	rBV	196039	346896	29.00%	2.351%
36	10.984	1298	1301	1307	rVB5	10843	17026	1.42%	0.115%
37	11.078	1308	1317	1330	rBV2	83207	169231	14.15%	1.147%
38	11.384	1366	1369	1372	rBV3	2396	2582	0.22%	0.017%
39	11.472	1380	1384	1386	rBV2	3677	5076	0.42%	0.034%
40	11.742	1428	1430	1437	rVB4	3830	4860	0.41%	0.033%
41	12.095	1486	1490	1493	rBV2	1969	2949	0.25%	0.020%
42	12.500	1554	1559	1566	rBV4	4027	9738	0.81%	0.066%
43	12.653	1579	1585	1594	rVB7	5879	10911	0.91%	0.074%
44	12.735	1594	1599	1600	rBV3	2767	3498	0.29%	0.024%
45	12.753	1600	1602	1605	rVV3	2939	3693	0.31%	0.025%
46	12.805	1608	1611	1616	rVV2	2239	3592	0.30%	0.024%
47	12.847	1616	1618	1621	rVB3	2690	2731	0.23%	0.019%
48	13.199	1677	1678	1681	rBV2	2598	2354	0.20%	0.016%
49	13.334	1698	1701	1702	rBV3	2891	2462	0.21%	0.017%
50	13.346	1702	1703	1706	rVV2	4165	3287	0.27%	0.022%
51	13.634	1748	1752	1754	rBV5	6132	8176	0.68%	0.055%
52	13.675	1754	1759	1765	rVB2	34286	55556	4.64%	0.377%
53	13.757	1770	1773	1776	rVB4	2545	3655	0.31%	0.025%
54	13.793	1776	1779	1781	rBV3	2962	3541	0.30%	0.024%
55	13.939	1801	1804	1806	rBV3	2677	3025	0.25%	0.021%
56	13.981	1810	1811	1815	rBV3	2513	2399	0.20%	0.016%
57	14.022	1817	1818	1821	rBV3	3720	3291	0.28%	0.022%
58	14.063	1823	1825	1828	rVV4	2820	3357	0.28%	0.023%
59	14.092	1828	1830	1832	rVV3	2399	3032	0.25%	0.021%
60	14.145	1833	1839	1844	rVV	325311	480739	40.19%	3.258%
61	14.192	1844	1847	1852	rVV	24724	34908	2.92%	0.237%
62	14.257	1855	1858	1860	rBV4	2910	2318	0.19%	0.016%
63	14.368	1872	1877	1878	rBV3	4626	5367	0.45%	0.036%
64	14.439	1880	1889	1898	rVV	360627	566471	47.36%	3.839%
65	14.527	1900	1904	1905	rBV3	3241	3766	0.31%	0.026%
66	14.603	1913	1917	1922	rVB7	8087	15420	1.29%	0.105%
67	14.739	1934	1940	1948	rVB	303641	496413	41.50%	3.364%
68	14.850	1955	1959	1960	rBV4	6763	7447	0.62%	0.050%
69	14.897	1964	1967	1970	rBV5	2621	3818	0.32%	0.026%
70	14.944	1973	1975	1976	rVB2	4154	2252	0.19%	0.015%
71	14.991	1976	1983	1990	rBV	144242	263009	21.99%	1.782%

LSC Area Percent Report

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Sample : M1412-14
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBGX9

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Title : SVOA CALIBRATION

72	15.115	2000	2004	2008	rBV7	9457	14706	1.23%	0.100%
73	15.162	2008	2012	2018	rVB9	7093	13526	1.13%	0.092%
74	15.267	2028	2030	2032	rBV3	5691	4199	0.35%	0.028%
75	15.344	2041	2043	2045	rBV3	3471	3675	0.31%	0.025%
76	15.485	2064	2067	2069	rBV4	7341	7602	0.64%	0.052%
77	15.508	2069	2071	2073	rVV3	4677	4144	0.35%	0.028%
78	15.549	2073	2078	2083	rVB8	14751	26237	2.19%	0.178%
79	15.731	2103	2109	2117	rBV	449529	678365	56.71%	4.597%
80	15.861	2127	2131	2139	rVB2	154054	226359	18.92%	1.534%
81	15.914	2139	2140	2141	rBV	6143	3641	0.30%	0.025%
82	15.972	2146	2150	2155	rVB7	10339	18785	1.57%	0.127%
83	16.319	2208	2209	2211	rBV2	6904	6428	0.54%	0.044%
84	16.789	2287	2289	2291	rBV3	8227	6129	0.51%	0.042%
85	17.488	2402	2408	2415	rBV	381347	570963	47.73%	3.870%
86	17.588	2419	2425	2431	rVV2	451744	694492	58.06%	4.707%
87	18.364	2554	2557	2560	rBV5	35334	41519	3.47%	0.281%
88	18.763	2620	2625	2630	rBV5	79917	143959	12.03%	0.976%
89	19.086	2675	2680	2686	rVB	543467	722159	60.37%	4.894%
90	19.274	2710	2712	2716	rBV3	33828	40076	3.35%	0.272%
91	19.580	2759	2764	2769	rVB	952751	1196189	100.00%	8.107%
92	19.868	2808	2813	2818	rBV	579780	806995	67.46%	5.469%
93	20.232	2872	2875	2880	rVV	271863	312348	26.11%	2.117%
94	20.297	2882	2886	2892	rVB	813153	1060680	88.67%	7.188%
95	20.514	2920	2923	2926	rBV3	82719	92709	7.75%	0.628%
96	20.632	2938	2943	2947	rVB	469691	633596	52.97%	4.294%
97	20.820	2971	2975	2980	rVB	654286	879488	73.52%	5.960%
98	21.766	3131	3136	3145	rVB2	416651	712517	59.57%	4.829%
99	22.183	3204	3207	3212	rVB	134365	185590	15.52%	1.258%
100	24.833	3653	3658	3668	rVB	305237	782439	65.41%	5.303%

Sum of corrected areas: 14755357

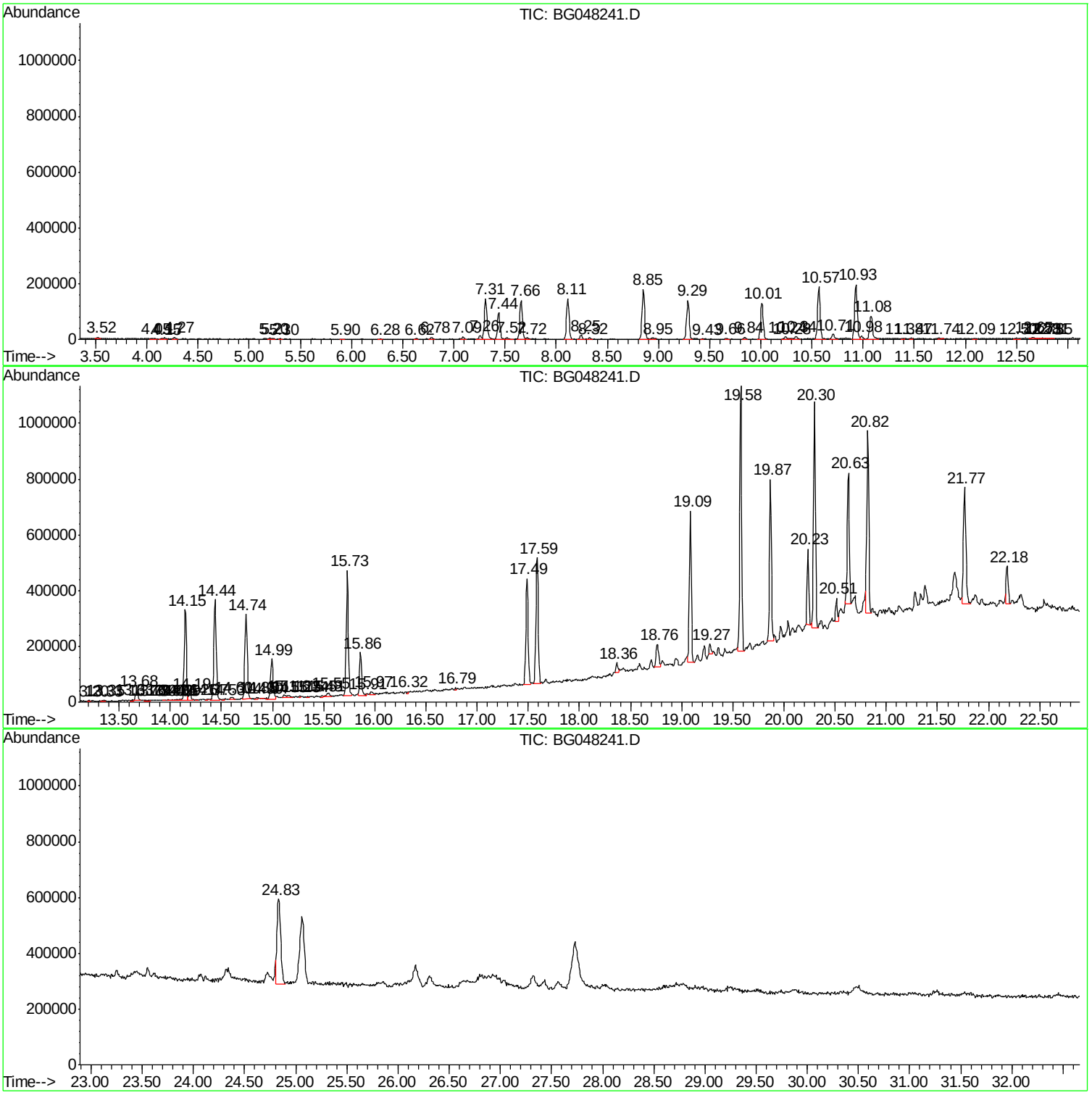
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048241.D
Acq On : 20 Feb 2021 18:04
Operator : CG/JU
Sample : M1412-14
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBGX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048241.D
Acq On : 20 Feb 2021 18:04
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Instrument : BNA_G
ClientSampleId : DBGX9

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Quant Title : SVOA CALIBRATION

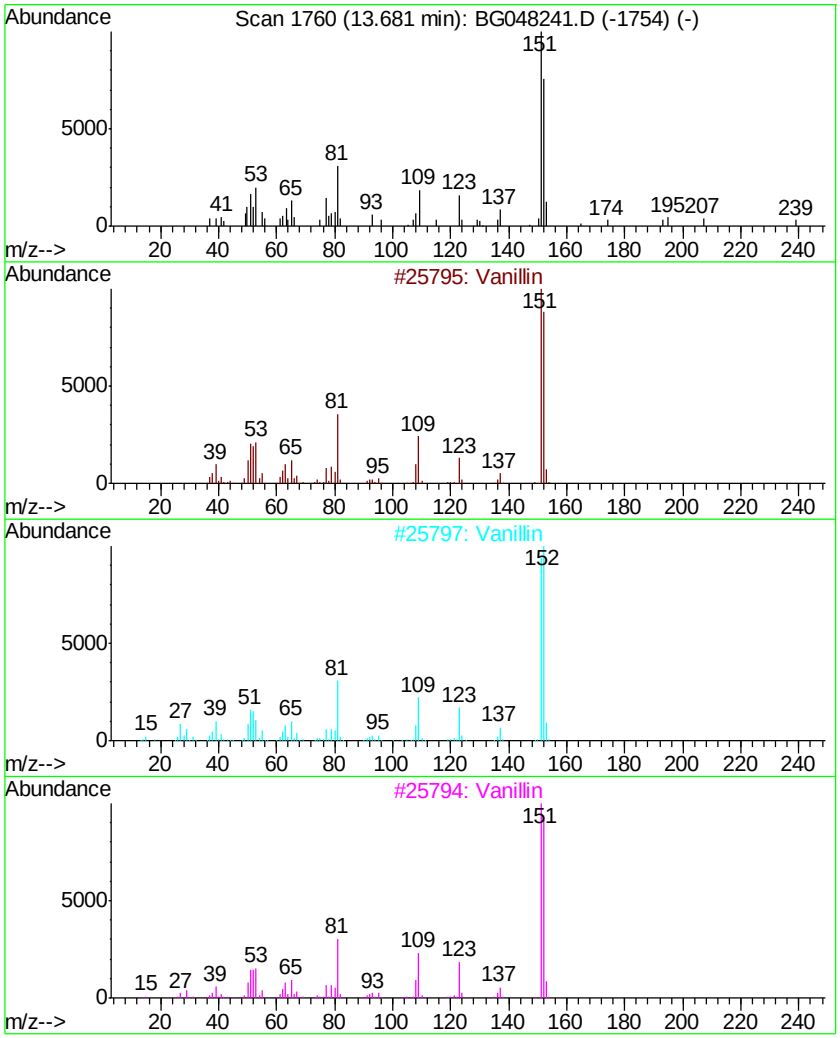
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

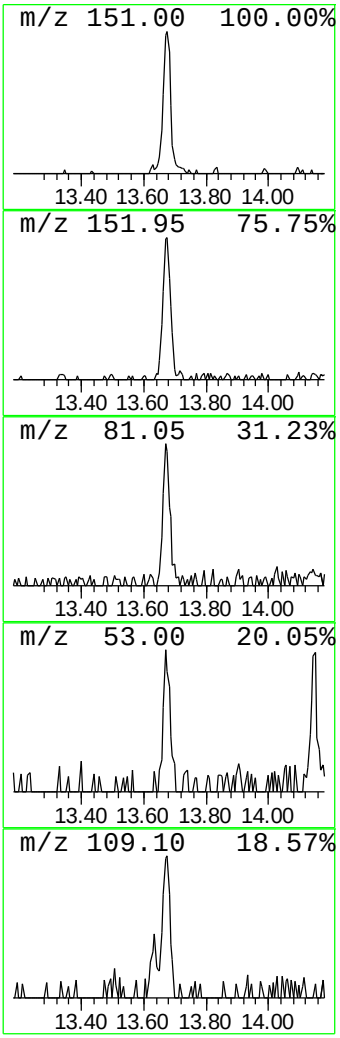
****

Peak Number 1 Vanillin Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.68	2.24 ng/ul	55556	Acenaphthene-d10	14.74

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Vanillin	152	C8H8O3	000121-33-5	97
2	Vanillin	152	C8H8O3	000121-33-5	97
3	Vanillin	152	C8H8O3	000121-33-5	97
4	Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	95
5	Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	93





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048241.D
Acq On : 20 Feb 2021 18:04
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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBGX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

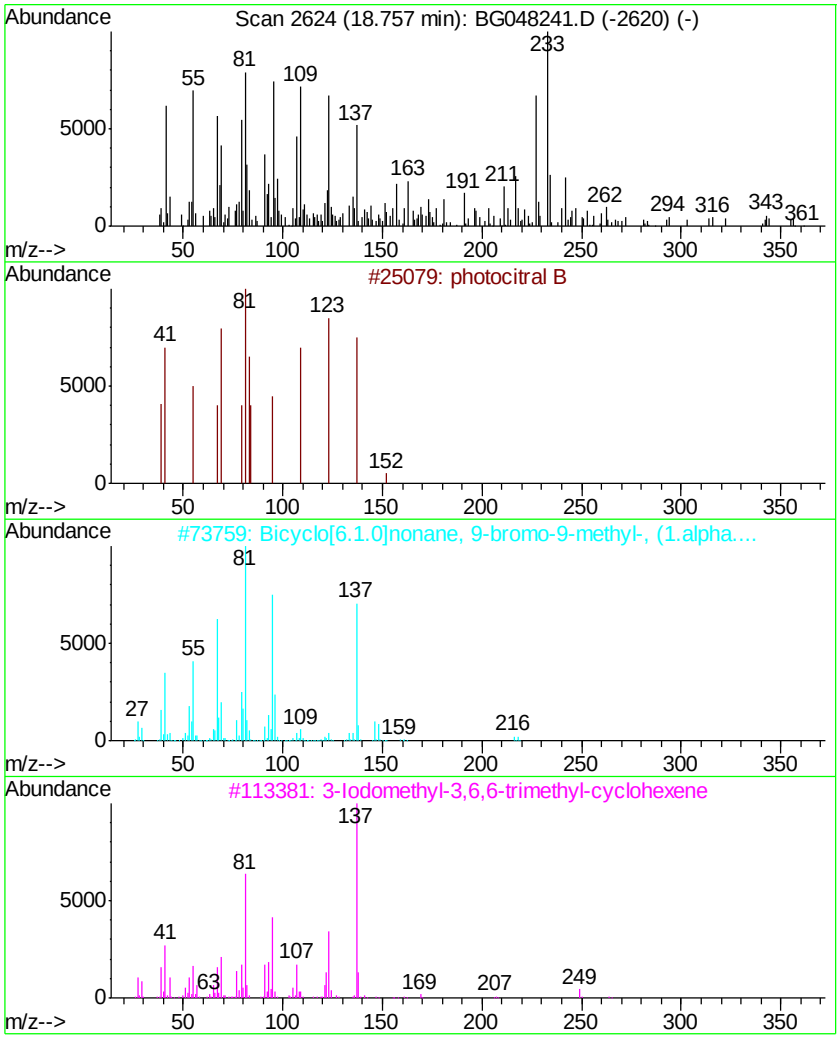
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

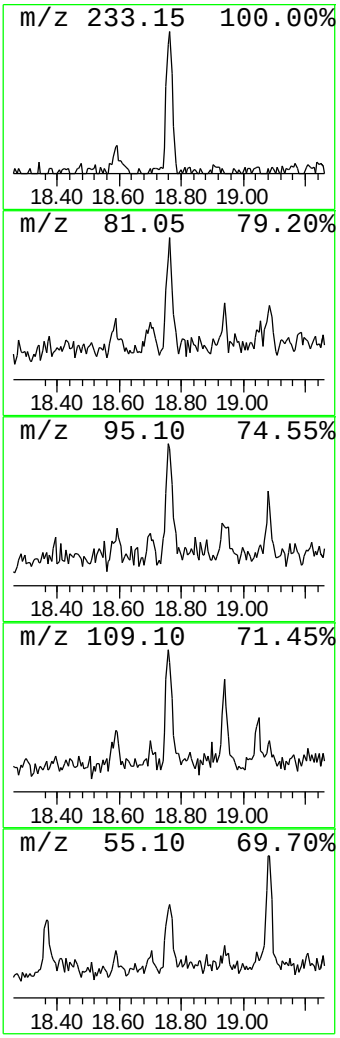
****

Peak Number 2 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.76	5.04 ng/ul	143959	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	photocitral B	152	C10H16O	006040-45-5	49
2	Bicyclo[6.1.0]nonane, 9-bromo-9-...	216	C10H17Br	059474-03-2	46
3	3-Iodomethyl-3,6,6-trimethyl-cyc...	264	C10H17I	1000193-08-2	38
4	3,9-Dimethyltricyclo[4.2.1.1(2,5...	180	C12H20O	1000185-61-6	35
5	Trans, trans-2-ethylbicyclo[4.4....	166	C12H22	066660-37-5	35





Library Search Compound Report

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Quant Title : SVOA CALIBRATION

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

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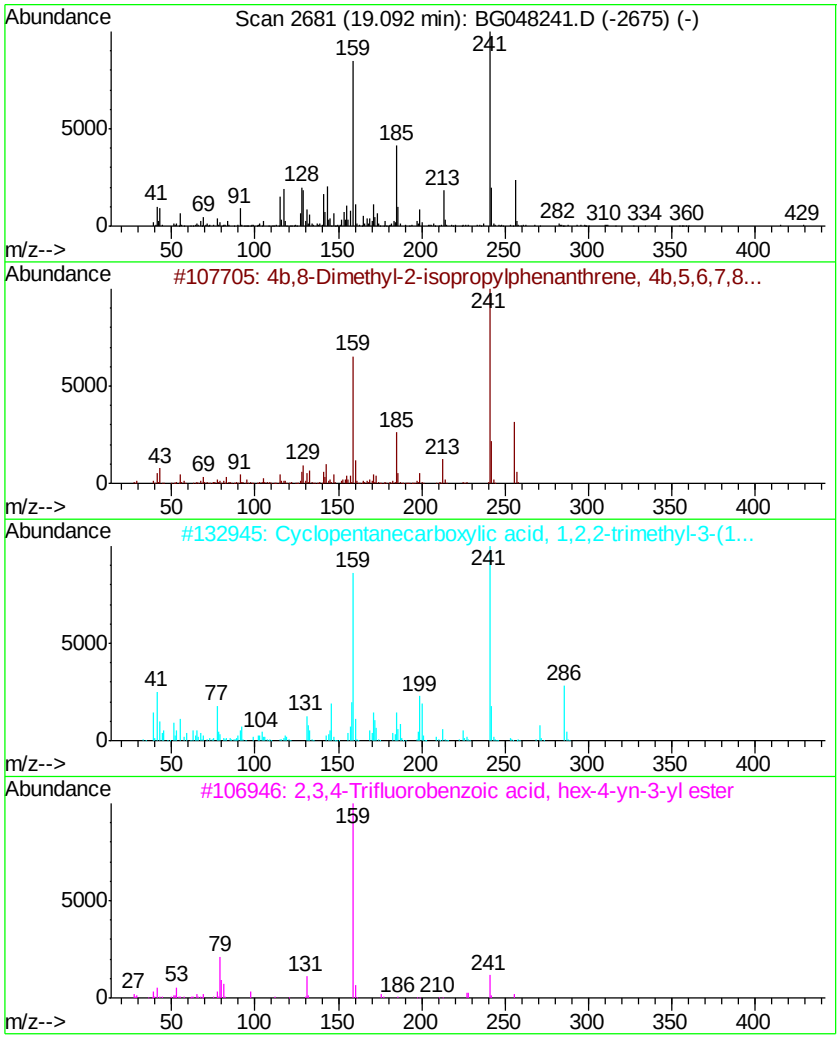
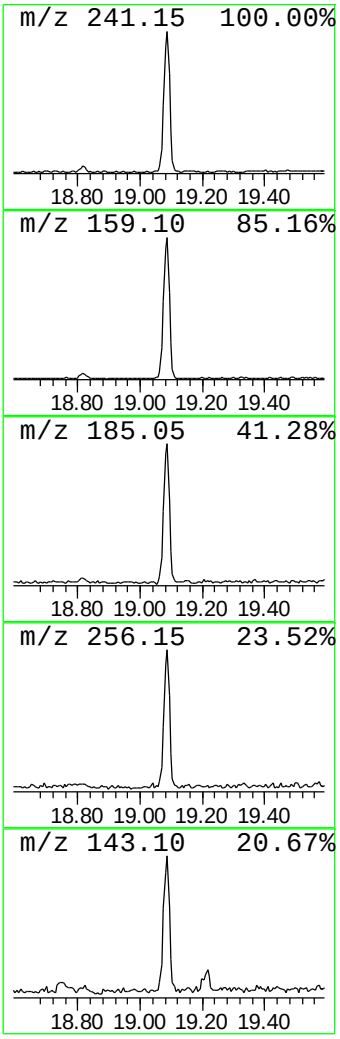
Peak Number 3 4b,8-Dimethyl-2-isopropylph... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.09	25.30 ng/ul	722159	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4b,8-Dimethyl-2-isopropylphenant...	256	C19H28	1000197-14-1	93
2	Cyclopentanecarboxylic acid, 1,2...	286	C17H22N2O2	1000302-75-6	47
3	2,3,4-Trifluorobenzoic acid, hex...	256	C13H11F3O2	1000292-55-4	46
4	1-Phenanthrenecarboxaldehyde, 1,...	284	C20H28O	024035-50-5	38
5	As-Indacene, 1,2,3,6,7,8-hexahyd...	256	C19H28	017465-47-3	27





Library Search Compound Report

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Instrument : BNA_G
ClientSampleId : DBGX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

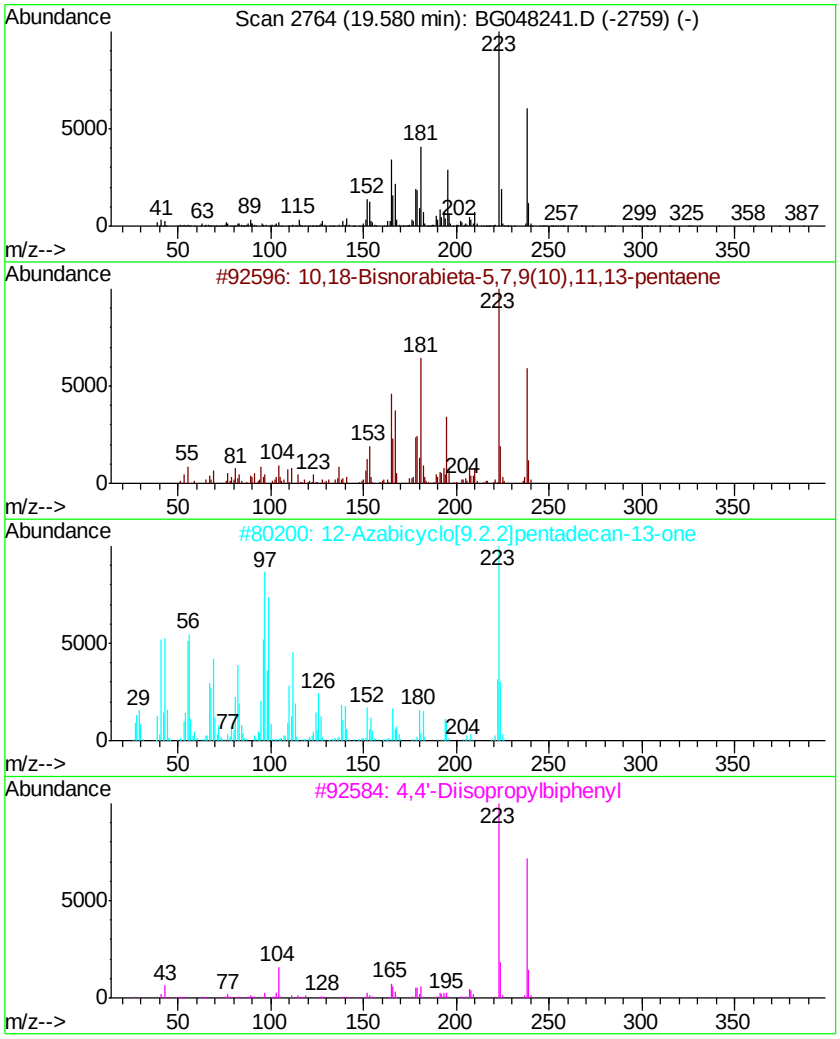
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

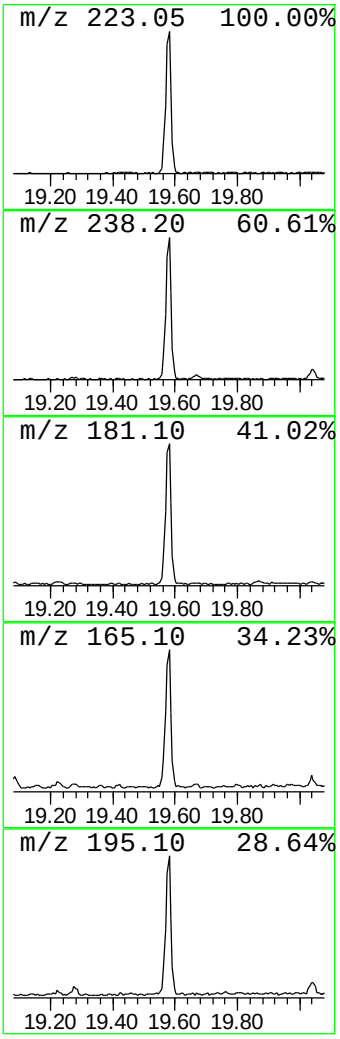
****

Peak Number 4 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.58	41.90 ng/ul	1196190	Phenanthrene-d10	17.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2	12-Azabicyclo[9.2.2]pentadecan-1...	223	C14H25NO	1000191-26-7	64
3	4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	60
4	Anthracene, 1,4-dimethoxy-	238	C16H14O2	013076-29-4	53
5	4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	52





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048241.D
Acq On : 20 Feb 2021 18:04
Operator : CG/JU
Sample : M1412-14
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBGX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

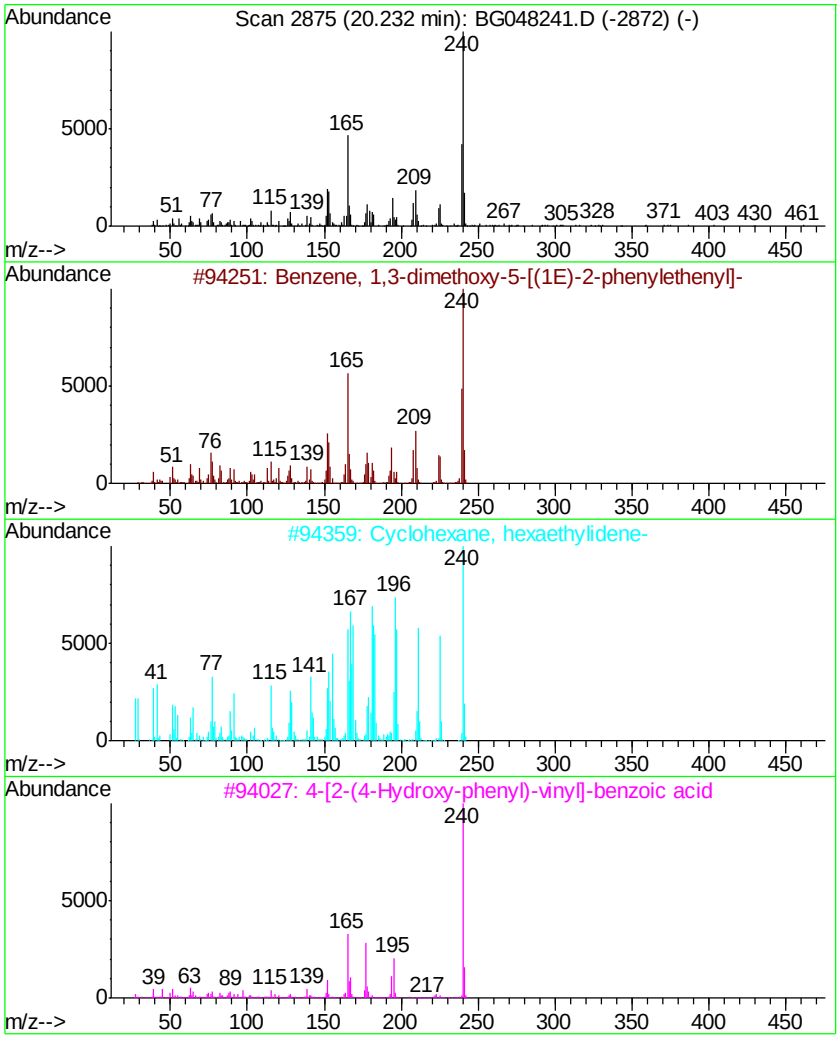
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

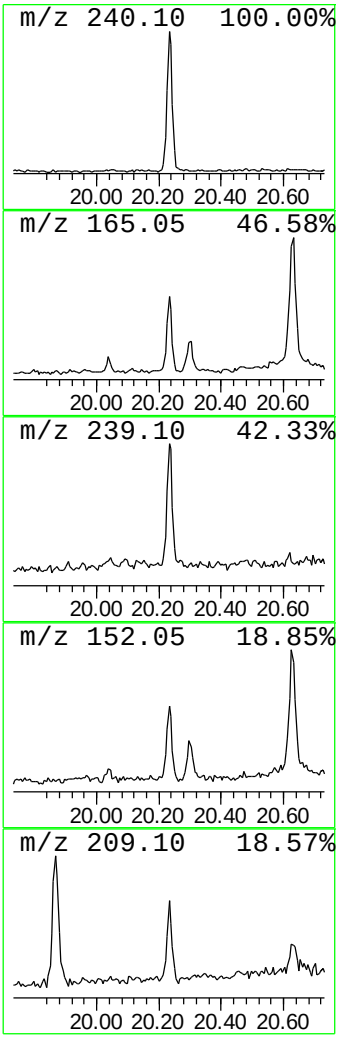
****

Peak Number 5 Benzene, 1,3-dimethoxy-5-[(... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.23	8.77 ng/ul	312348	Chrysene-d12	21.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethoxy-5-[(1E)-2...	240	C16H16O2	021956-56-9	98
2	Cyclohexane, hexaethylidene-	240	C18H24	001482-93-5	91
3	4-[2-(4-Hydroxy-phenyl)-vinyl-b...	240	C15H12O3	152027-61-7	58
4	[1,1'-Biphenyl-4,4'-diamine, 3,...	240	C16H20N2	054827-17-7	53
5	4,6,2',6'-Tetramethyl-biphenyl-2...	240	C16H20N2	004445-46-9	46





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048241.D
Acq On : 20 Feb 2021 18:04
Operator : CG/JU
Sample : M1412-14
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBGX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

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

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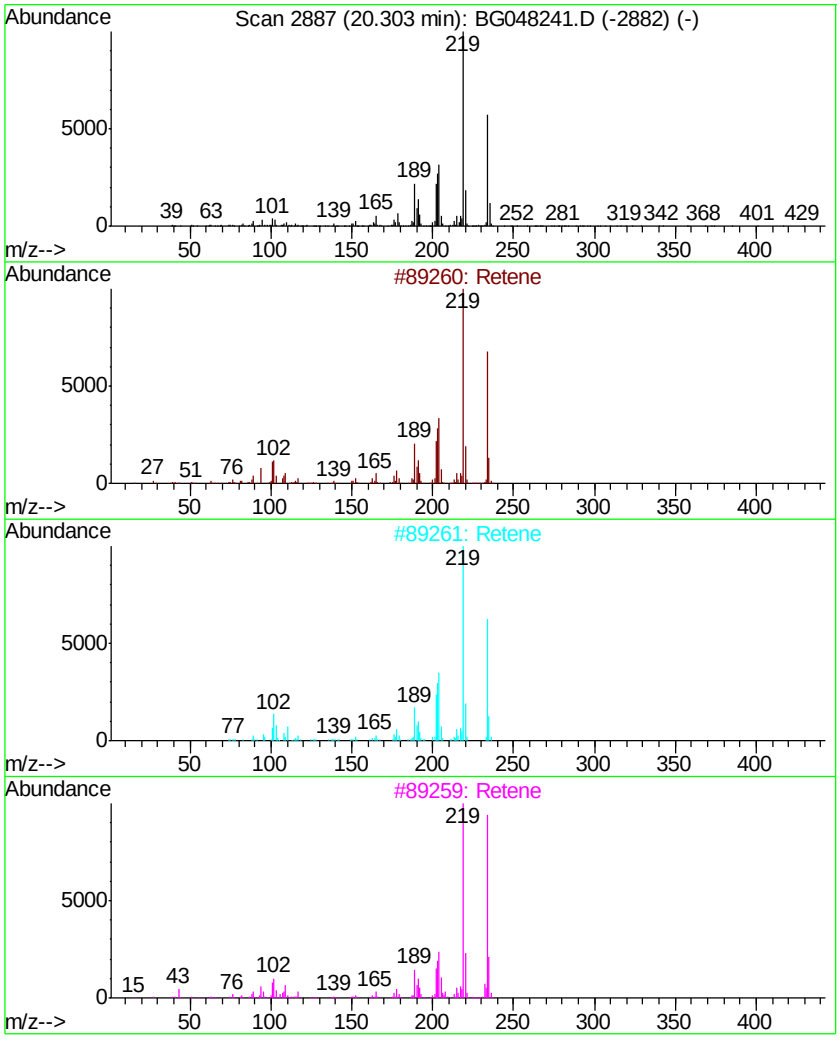
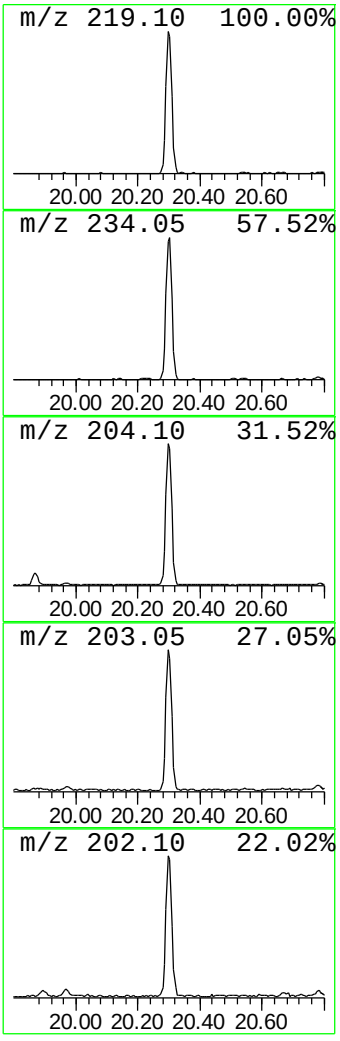
Peak Number 6 Retene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.30	29.77 ng/ul	1060680	Chrysene-d12	21.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Retene	234	C18H18	000483-65-8	98
2	Retene	234	C18H18	000483-65-8	98
3	Retene	234	C18H18	000483-65-8	95
4	2-Isopropyl-10-methylphenanthrene	234	C18H18	066552-97-4	95
5	Phenanthrene, 2,4,5,7-tetramethyl-	234	C18H18	007396-38-5	68





Library Search Compound Report

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Data File : BG048241.D
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Operator : CG/JU
Sample : M1412-14
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBGX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

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

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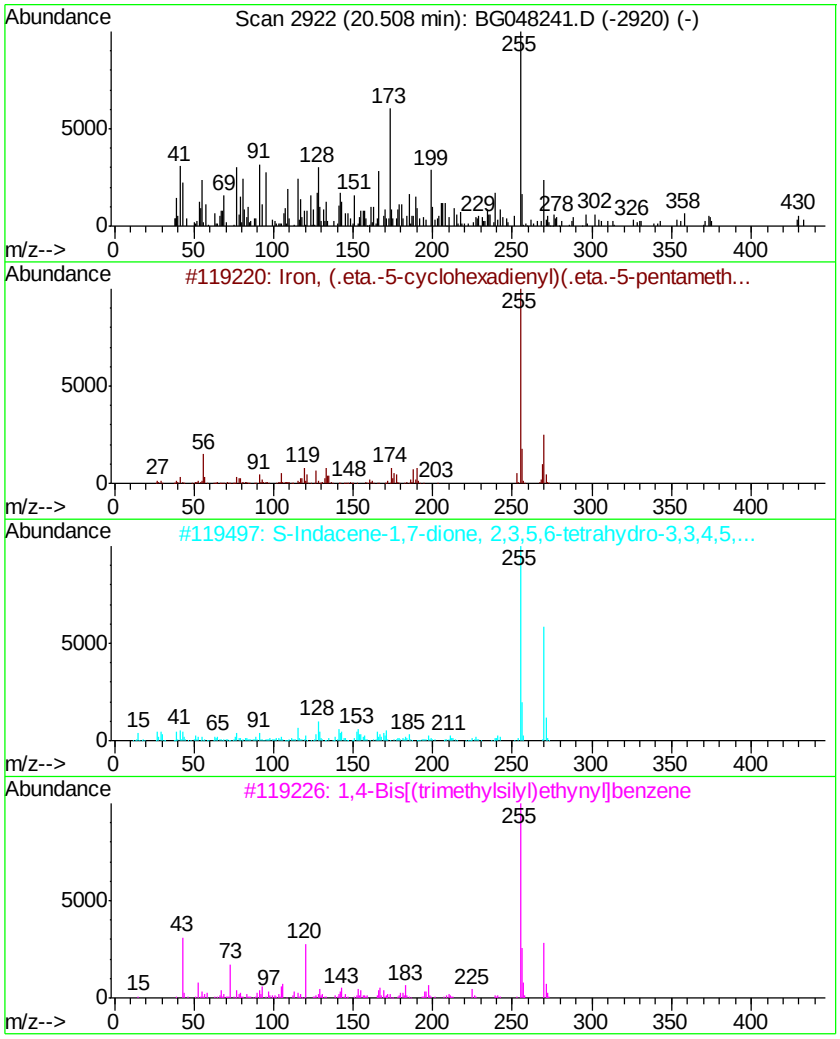
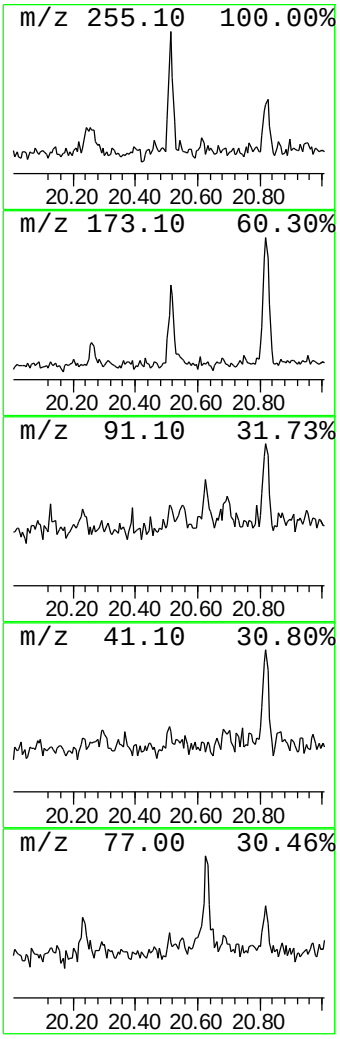
Peak Number 7 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.51	2.60 ng/ul	92709	Chrysene-d12	21.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Iron, (.eta.-5-cyclohexadienyl)(...	270	C16H22Fe	1000162-99-3	30
2	S-Indacene-1,7-dione, 2,3,5,6-te...	270	C18H22O2	055591-16-7	25
3	1,4-Bis[(trimethylsilyl)ethynyl]...	270	C16H22Si2	073392-23-1	25
4	5-tert-Butyl-2,2'-dimethoxy-biph...	270	C18H22O2	1000318-04-7	22
5	4-(2-Hydroxyethyl)-5-oxo-3-pheny...	204	C11H12N2O2	010244-77-6	18





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
Data File : BG048241.D
Acq On : 20 Feb 2021 18:04
Operator : CG/JU
Sample : M1412-14
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBGX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

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

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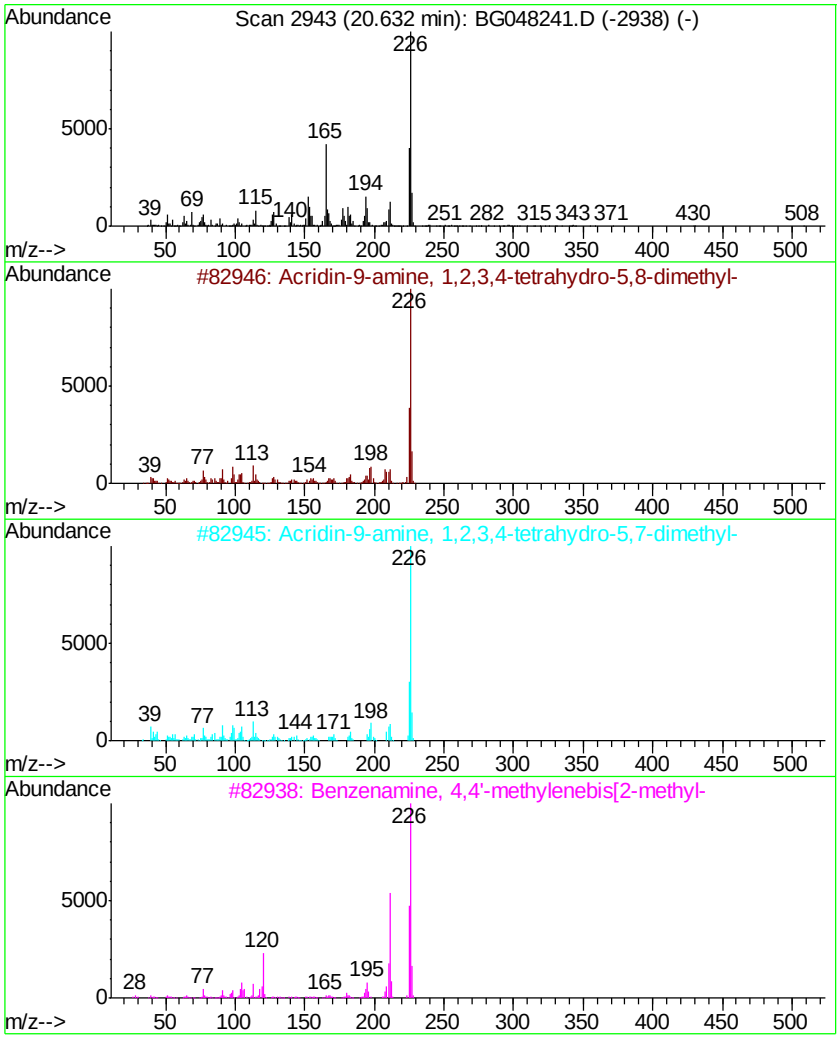
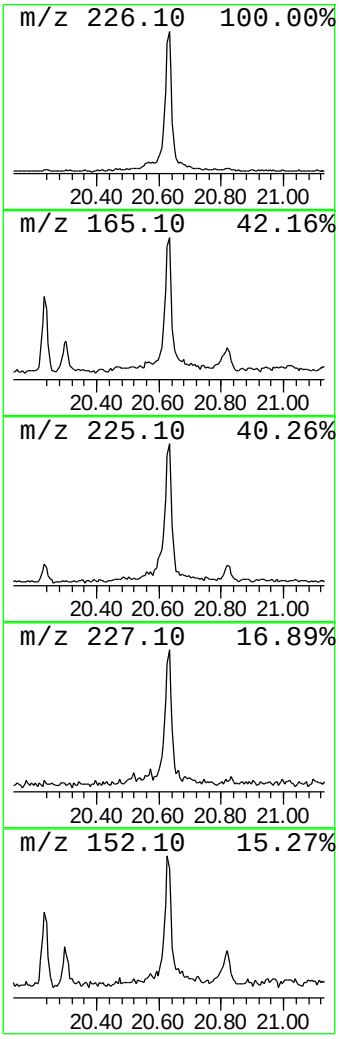
Peak Number 8 Acridin-9-amine, 1,2,3,4-te... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.63	17.78 ng/ul	633596	Chrysene-d12	21.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acridin-9-amine, 1,2,3,4-tetrahy...	226	C15H18N2	297758-19-1	64
2	Acridin-9-amine, 1,2,3,4-tetrahy...	226	C15H18N2	1000300-57-6	58
3	Benzenamine, 4,4'-methylenebis[2...	226	C15H18N2	000838-88-0	50
4	4'-Methoxy-2-hydroxystilbene	226	C15H14O2	1000147-93-1	50
5	9H-Xanthen-9-one, 2-methoxy-	226	C14H10O3	001214-20-6	47





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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBGX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

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

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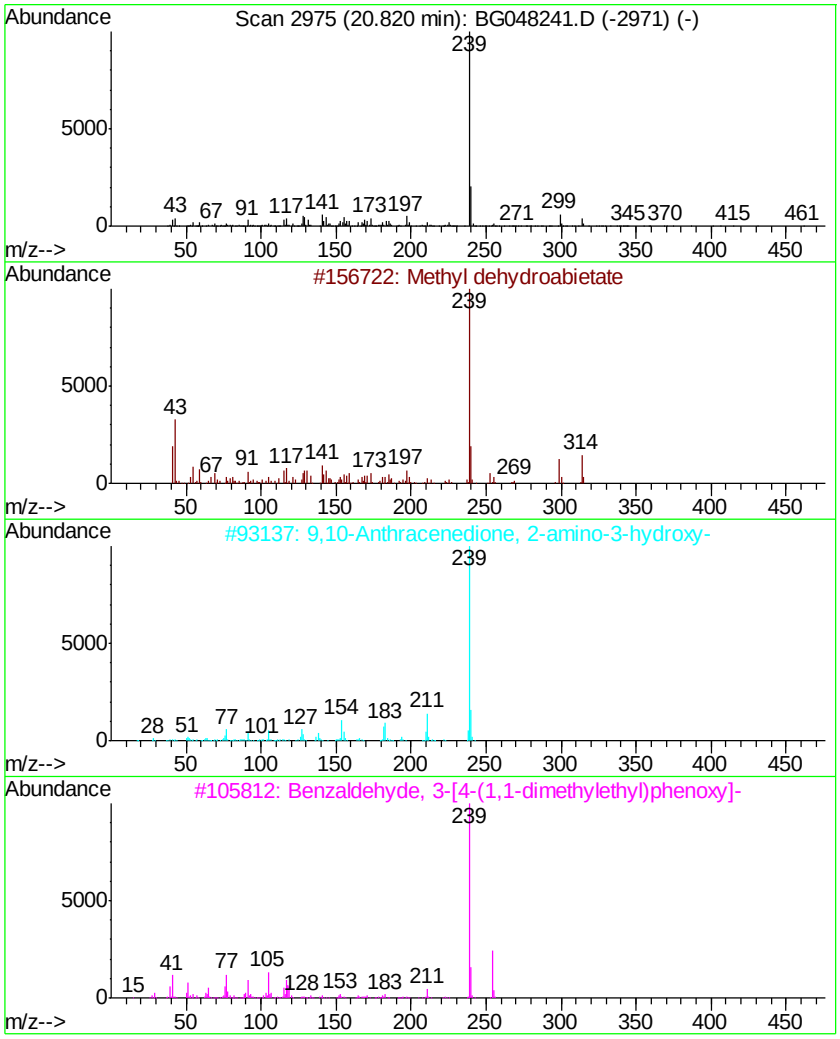
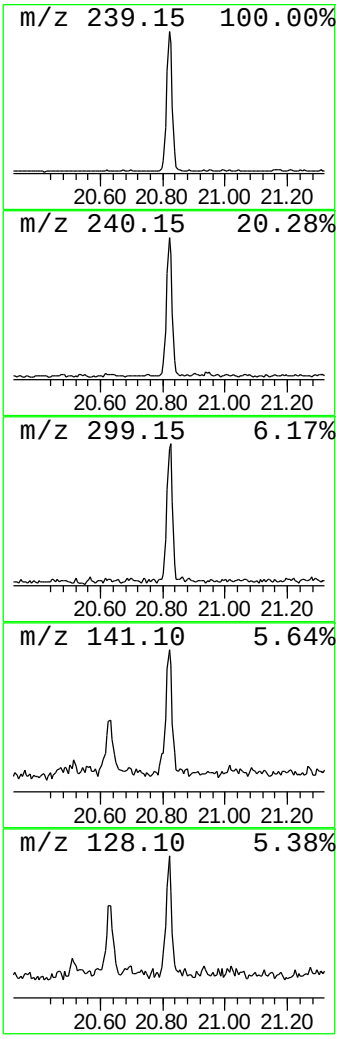
Peak Number 9 Methyl dehydroabietate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.82	24.69 ng/ul	879488	Chrysene-d12	21.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl dehydroabietate	314	C21H30O2	001235-74-1	94
2	9,10-Anthracenedione, 2-amino-3-...	239	C14H9NO3	000117-77-1	64
3	Benzaldehyde, 3-[4-(1,1-dimethyl...	254	C17H18O2	069770-23-6	59
4	Methyl dehydroabietate	314	C21H30O2	001235-74-1	56
5	Imidazolo[2,1-b]thiazole, 5-(3-i...	239	C13H9N3S	327097-37-0	45





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021921\
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Acq On : 20 Feb 2021 18:04
Operator : CG/JU
Sample : M1412-14
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : DBGX9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG021321MA.M
Quant Title : SVOA CALIBRATION

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

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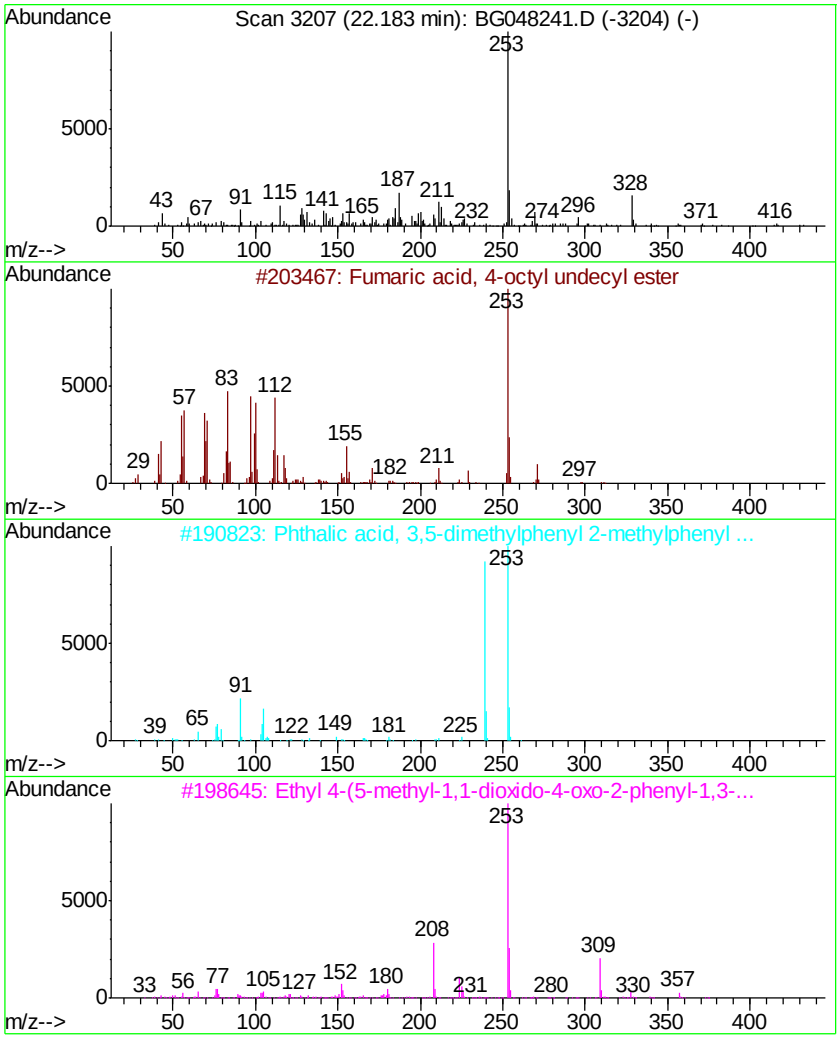
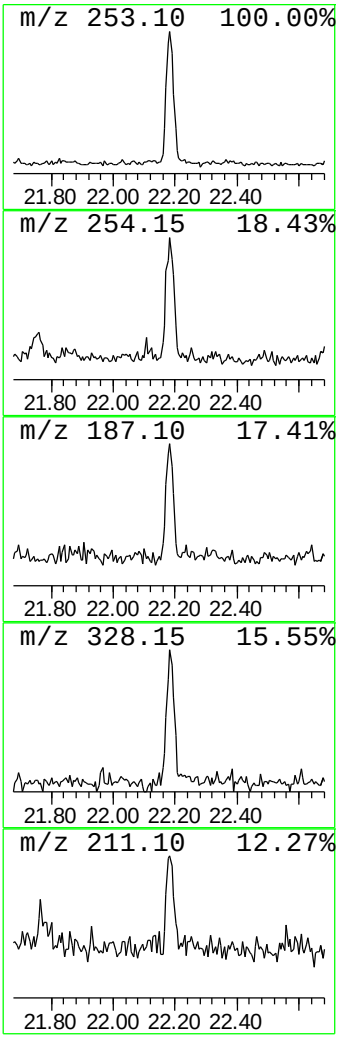
Peak Number 10 Fumaric acid, 4-octyl undec... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.18	5.21 ng/ul	185590	Chrysene-d12	21.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fumaric acid, 4-octyl undecyl ester	382	C23H42O4	1000339-22-2	53
2	Phthalic acid, 3,5-dimethylphenyl...	360	C23H20O4	1000315-59-4	50
3	Ethyl 4-(5-methyl-1,1-dioxido-4-...	373	C19H19NO5S	1000294-64-6	50
4	7-Oxodehydroabietic acid, methyl...	328	C21H28O3	110936-78-2	49
5	Benzaldehyde, 4-((4-(dimethylami...	253	C15H15N3O	039208-00-9	47





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG021921\
Data File : BG048241.D
Acq On : 20 Feb 2021 18:04
Operator : CG/JU
Sample : M1412-14
Misc :
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Instrument : BNA_G
ClientSampleId : DBGX9

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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
Vanillin	13.68	2.2	ng/ul	55556	3	14.74	496413	20.0
unknown-01	18.76	5.0	ng/ul	143959	4	17.49	570963	20.0
4b,8-Dimethyl-2-i...	19.09	25.3	ng/ul	722159	4	17.49	570963	20.0
10,18-Bisnorabiet...	19.58	41.9	ng/ul	1196190	4	17.49	570963	20.0
Benzene, 1,3-dime...	20.23	8.8	ng/ul	312348	5	21.77	712517	20.0
Retene	20.30	29.8	ng/ul	1060680	5	21.77	712517	20.0
unknown-02	20.51	2.6	ng/ul	92709	5	21.77	712517	20.0
Acridin-9-amine, ...	20.63	17.8	ng/ul	633596	5	21.77	712517	20.0
Methyl dehydroabi...	20.82	24.7	ng/ul	879488	5	21.77	712517	20.0
Fumaric acid, 4-o...	22.18	5.2	ng/ul	185590	5	21.77	712517	20.0