LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111920\
Data File : BG047369.D
Acq On : 20 Nov 2020 2:07
Operator : CG/JU
Sample : L4710-23
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ0

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-BG111920MA.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	4.154	136	139	145	rVB3	9741	11322	1.41%	0.122%
2	4.265	156	158	159	rBV	2564	2198	0.27%	0.024%
3	4.389	177	179	184	rBV3	4724	5713	0.71%	0.061%
4	5.340	336	341	345	rVB4	5147	8117	1.01%	0.087%
5	5.575	379	381	385	rVB3	3318	3009	0.37%	0.032%
6	6.134	472	476	479	rBV2	2891	2817	0.35%	0.030%
7	6.404	517	522	524	rBV2	1802	3121	0.39%	0.034%
8	6.604	553	556	557	rBV	2437	2428	0.30%	0.026%
9	6.792	583	588	594	rBV4	7719	15234	1.90%	0.164%
10	6.856	597	599	603	rVB3	2643	2511	0.31%	0.027%
11	6.950	609	615	623	rBV	89961	152930	19.04%	1.644%
12	7.033	627	629	634	rVB	1607	2349	0.29%	0.025%
13	7.221	656	661	662	rBV4	7277	10517	1.31%	0.113%
14	7.262	663	668	676	rVB3	71817	125922	15.68%	1.353%
15	7.403	684	692	702	rVB	94337	174896	21.78%	1.880%
16	7.585	717	723	730	rBV	62718	103040	12.83%	1.107%
17	7.720	739	746	753	rBV	144519	241520	30.07%	2.596%
18	7.796	753	759	766	rVV3	86535	154880	19.29%	1.665%
19	7.861	766	770	773	rVV3	3514	4785	0.60%	0.051%
20	7.961	785	787	791	rBV	2669	2360	0.29%	0.025%
21	8.026	796	798	803	rVB3	3082	3439	0.43%	0.037%
22	8.178	818	824	830	rBV3	34571	58533	7.29%	0.629%
23	8.278	834	841	847	rVV	172525	303329	37.77%	3.260%
24	8.407	862	863	868	rVB2	4874	5455	0.68%	0.059%
25	8.501	870	879	884	rVB4	27252	46186	5.75%	0.496%
26	8.595	890	895	900	rVB5	8953	13883	1.73%	0.149%
27	8.660	903	906	911	rVB5	7834	11103	1.38%	0.119%
28	8.736	916	919	924	rBV	1864	2355	0.29%	0.025%
29	8.895	944	946	949	rBV	1853	2462	0.31%	0.026%
30	8.960	951	957	964	rVV3	96912	175396	21.84%	1.885%
31	9.095	974	980	982	rBV3	3081	4833	0.60%	0.052%
32	9.236	1001	1004	1007	rVB2	2295	2421	0.30%	0.026%
33	9.283	1009	1012	1016	rBV	1549	2337	0.29%	0.025%
34	9.436	1032	1038	1046	rVB2	87035	160948	20.04%	1.730%

LSC Area Percent Report

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Acq On : 20 Nov 2020 2:07
Operator : CG/JU
Sample : L4710-23
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ0

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-BG111920MA.M
Title : SVOA CALIBRATION

35	9.518	1050	1052	1054	rBV3	2382	2442	0.30%	0.026%
36	9.859	1108	1110	1113	rVB	3909	3183	0.40%	0.034%
37	10.082	1146	1148	1149	rBV	4010	2849	0.35%	0.031%
38	10.164	1156	1162	1171	rVB3	76298	144751	18.02%	1.556%
39	10.705	1246	1254	1261	rVB	124363	225384	28.07%	2.422%
40	11.104	1314	1322	1329	rBV	221860	423076	52.68%	4.547%
41	11.222	1335	1342	1349	rBV2	82929	147940	18.42%	1.590%
42	11.298	1352	1355	1357	rBV	2768	2355	0.29%	0.025%
43	11.416	1372	1375	1379	rVV4	4511	7001	0.87%	0.075%
44	11.486	1381	1387	1392	rVB3	4204	8907	1.11%	0.096%
45	11.615	1405	1409	1410	rBV	3156	3266	0.41%	0.035%
46	11.862	1447	1451	1453	rBV	2281	3082	0.38%	0.033%
47	11.886	1453	1455	1459	rVB	3183	3794	0.47%	0.041%
48	12.027	1476	1479	1482	rVB	2039	2204	0.27%	0.024%
49	12.086	1486	1489	1492	rVV2	2061	2793	0.35%	0.030%
50	12.127	1494	1496	1498	rVB	2570	2209	0.28%	0.024%
51	12.156	1498	1501	1502	rBV	2751	2466	0.31%	0.027%
52	12.262	1515	1519	1520	rVB3	3222	3518	0.44%	0.038%
53	12.438	1542	1549	1555	rBV3	49156	77453	9.64%	0.832%
54	12.614	1577	1579	1583	rVB	2442	2577	0.32%	0.028%
55	12.661	1583	1587	1589	rBV2	4366	5905	0.74%	0.063%
56	12.814	1611	1613	1618	rVB4	3017	3769	0.47%	0.041%
57	12.955	1633	1637	1640	rBV	3238	5721	0.71%	0.061%
58	13.090	1658	1660	1662	rVB	2904	2226	0.28%	0.024%
59	13.125	1662	1666	1667	rBV2	1987	2899	0.36%	0.031%
60	13.237	1682	1685	1686	rBV2	2583	2744	0.34%	0.029%
61	13.349	1701	1704	1706	rVB2	3206	3441	0.43%	0.037%
62	13.502	1727	1730	1731	rVB	3339	2784	0.35%	0.030%
63	13.519	1731	1733	1739	rBV2	1484	2670	0.33%	0.029%
64	13.643	1750	1754	1761	rVB7	8146	19997	2.49%	0.215%
65	13.719	1764	1767	1771	rBV4	6221	10808	1.35%	0.116%
66	13.772	1771	1776	1782	rVV7	17252	31229	3.89%	0.336%
67	13.831	1782	1786	1792	rVB5	8701	14639	1.82%	0.157%
68	13.907	1792	1799	1804	rBV5	4614	8799	1.10%	0.095%
69	13.989	1809	1813	1815	rBV	6853	9234	1.15%	0.099%
70	14.218	1848	1852	1857	rBV5	17370	26378	3.28%	0.284%
71	14.277	1857	1862	1866	rBV	229440	340480	42.40%	3.659%

LSC Area Percent Report

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Acq On : 20 Nov 2020 2:07
Operator : CG/JU
Sample : L4710-23
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ0

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-BG111920MA.M
Title : SVOA CALIBRATION

72	14.324	1866	1870	1873	rVB2	77044	102952	12.82%	1.107%
73	14.412	1881	1885	1893	rVB5	54610	89399	11.13%	0.961%
74	14.583	1907	1914	1921	rVB	227243	372103	46.33%	3.999%
75	14.647	1921	1925	1928	rBV4	3981	6990	0.87%	0.075%
76	14.847	1954	1959	1961	rBV	152610	215114	26.79%	2.312%
77	14.882	1961	1965	1972	rVB	371220	595651	74.17%	6.402%
78	14.947	1973	1976	1982	rBV6	11761	18394	2.29%	0.198%
79	15.035	1986	1991	2001	rVB2	102060	165557	20.62%	1.779%
80	15.123	2004	2006	2007	rBV2	3764	2234	0.28%	0.024%
81	15.153	2007	2011	2015	rVB6	26447	37109	4.62%	0.399%
82	15.276	2029	2032	2036	rVB3	4267	5195	0.65%	0.056%
83	15.393	2047	2052	2060	rBV2	81903	136856	17.04%	1.471%
84	15.764	2111	2115	2123	rVB6	7536	15306	1.91%	0.165%
85	15.869	2127	2133	2141	rVB	294230	473695	58.98%	5.091%
86	15.963	2144	2149	2155	rBV2	73457	119238	14.85%	1.282%
87	16.145	2178	2180	2182	rBV2	2192	2464	0.31%	0.026%
88	16.281	2200	2203	2206	rBV3	8458	11915	1.48%	0.128%
89	16.351	2208	2215	2219	rBV6	17722	31277	3.89%	0.336%
90	16.645	2263	2265	2268	rVB	4519	3671	0.46%	0.039%
91	16.709	2274	2276	2278	rBV3	5792	3688	0.46%	0.040%
92	17.185	2354	2357	2358	rBV3	2622	2664	0.33%	0.029%
93	17.614	2424	2430	2437	rBV	507631	788497	98.18%	8.475%
94	17.714	2441	2447	2453	rVV	349327	516584	64.33%	5.552%
95	18.484	2575	2578	2586	rVB6	37496	62597	7.79%	0.673%
96	19.254	2707	2709	2715	rVB2	53065	54390	6.77%	0.585%
97	19.982	2827	2833	2843	rVB	426568	672293	83.71%	7.226%
98	21.516	3090	3094	3099	rBV2	81528	119859	14.92%	1.288%
99	21.898	3153	3159	3165	rVB2	472906	803078	100.00%	8.631%
100	22.673	3287	3291	3297	rVB3	259273	487991	60.77%	5.245%

Sum of corrected areas: 9304083

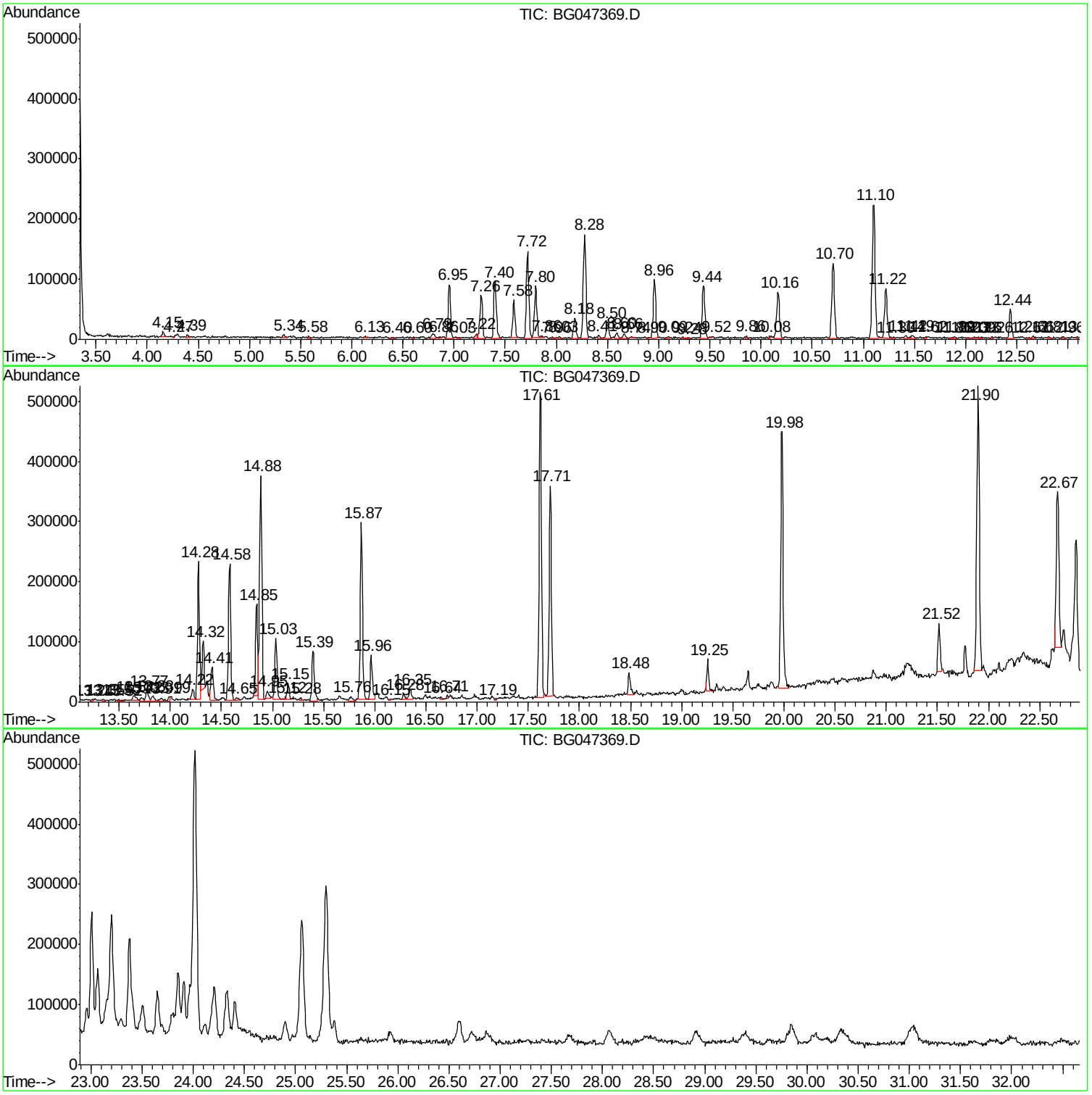
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111920\
Data File : BG047369.D
Acq On : 20 Nov 2020 2:07
Operator : CG/JU
Sample : L4710-23
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.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\BG111920\
Data File : BG047369.D
Acq On : 20 Nov 2020 2:07
Operator : CG/JU
Sample : L4710-23
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

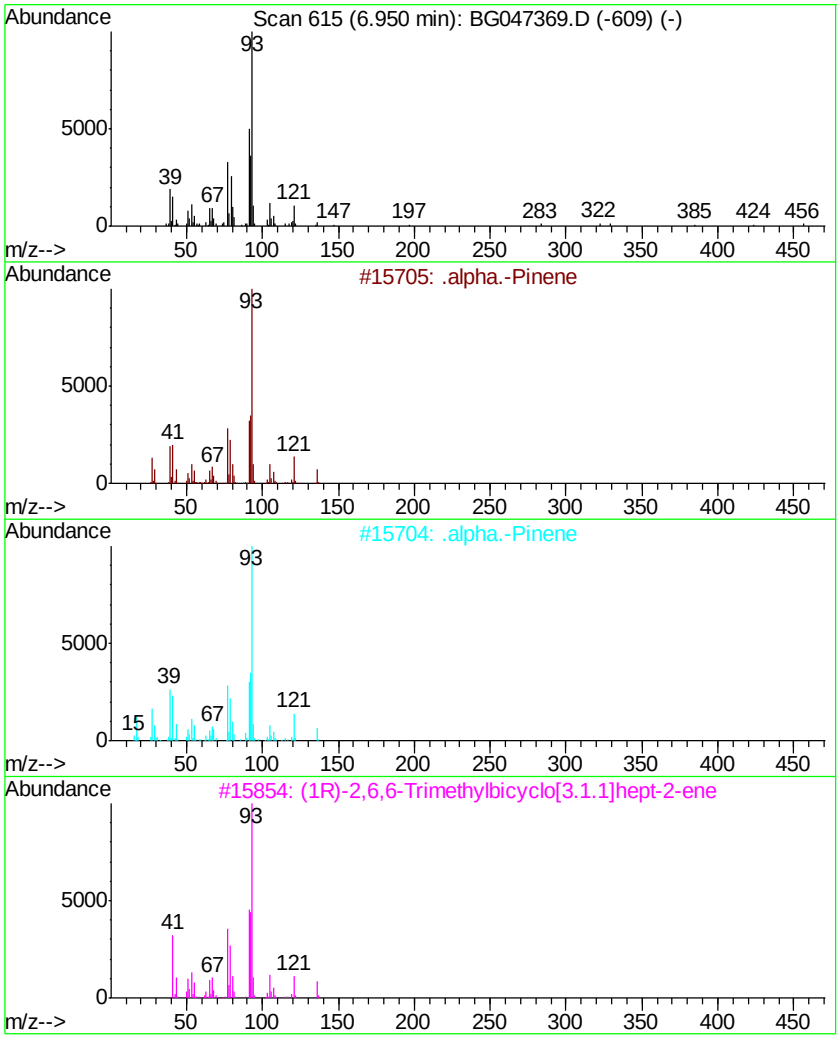
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

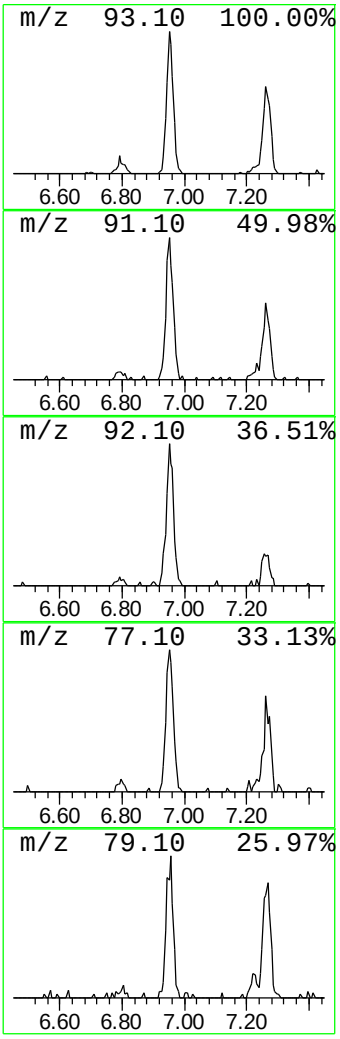
****

Peak Number 1 .alpha.-Pinene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.95	10.08 ng/ul	152930	1,4-Dichlorobenzene-d4	8.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Pinene	136	C10H16	000080-56-8	94
2	.alpha.-Pinene	136	C10H16	000080-56-8	91
3	(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	91
4	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	90
5	3,5-Methanocyclopentapyrazole, 3...	164	C10H16N2	087143-58-6	90





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111920\
Data File : BG047369.D
Acq On : 20 Nov 2020 2:07
Operator : CG/JU
Sample : L4710-23
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

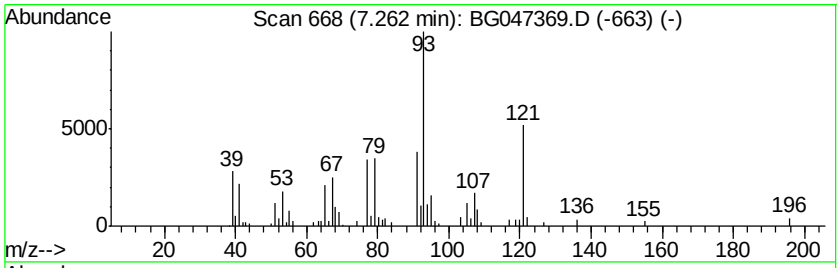
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

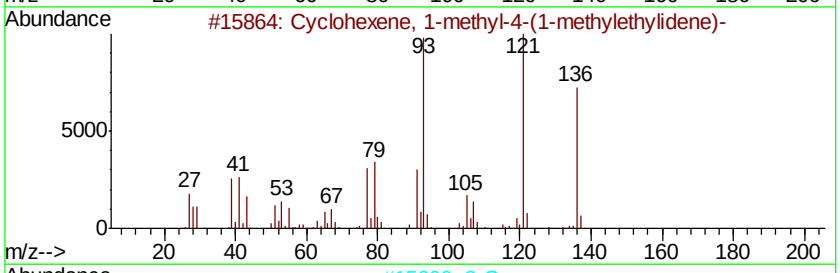
****

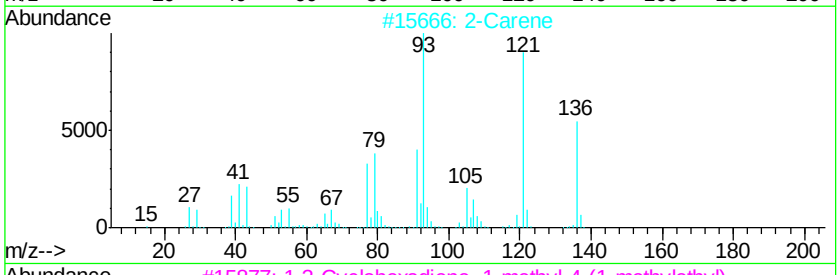
Peak Number 2 2-Carene Concentration Rank 4

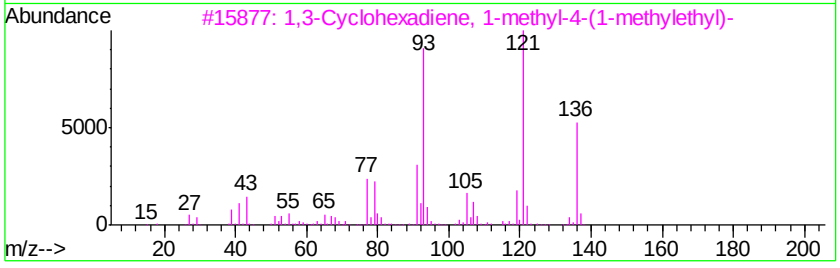
R.T.	EstConc	Area	Relative to ISTD	R.T.
7.26	8.30 ng/ul	125922	1,4-Dichlorobenzene-d4	8.28

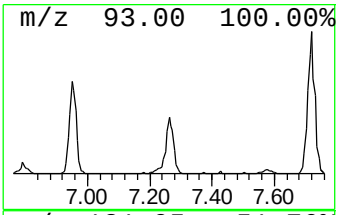
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	000586-62-9	87
2	2-Carene	136	C10H16	000554-61-0	81
3	1,3-Cyclohexadiene, 1-methyl-4-(...	136	C10H16	000099-86-5	81
4	Bicyclo[2.2.1]hept-2-ene, 1,7,7-...	136	C10H16	000464-17-5	74
5	2-Carene	136	C10H16	000554-61-0	74

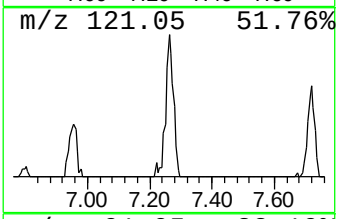


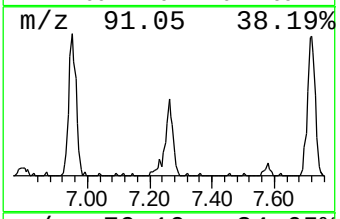


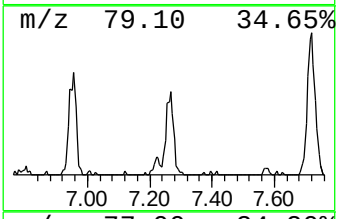


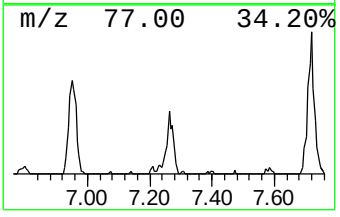












Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111920\
Data File : BG047369.D
Acq On : 20 Nov 2020 2:07
Operator : CG/JU
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Misc :
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Instrument : BNA_G
ClientSampleId : C0AZ0

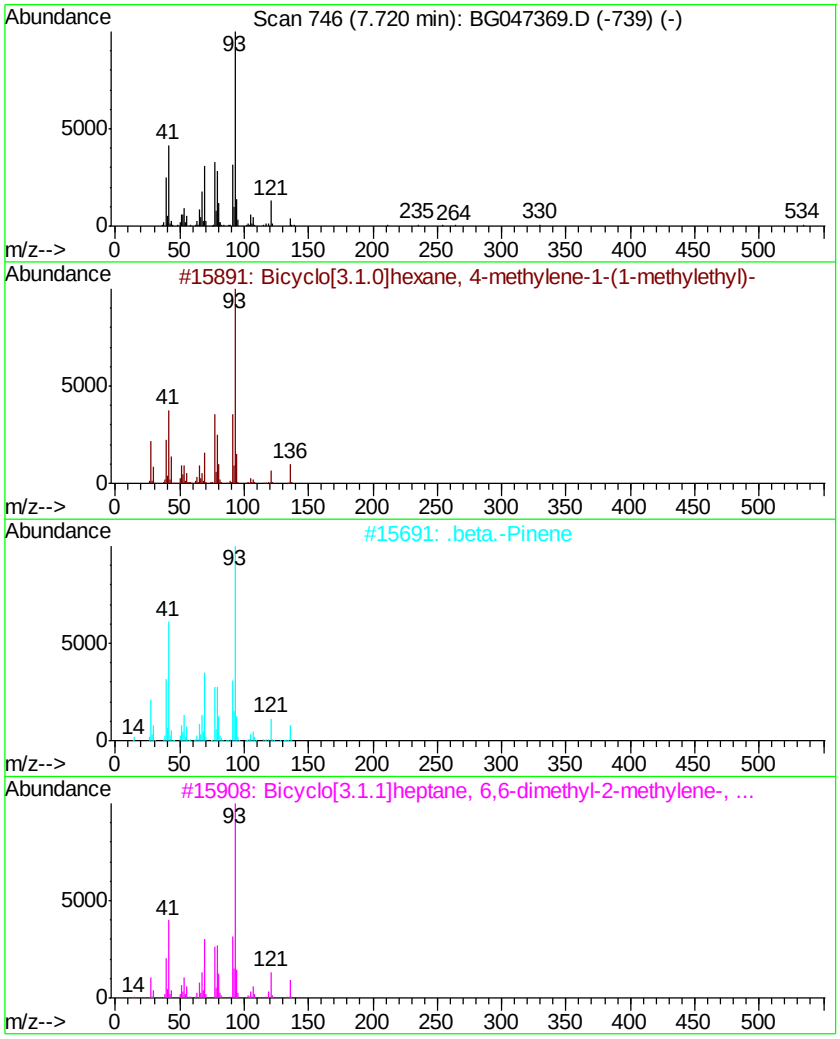
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

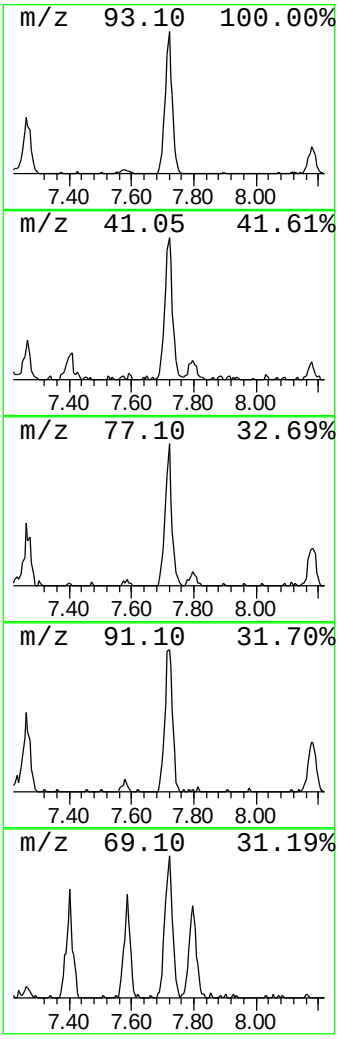
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Bicyclo[3.1.0]hexane, 4-met... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.72	15.92 ng/ul	241520	1,4-Dichlorobenzene-d4	8.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[3.1.0]hexane, 4-methylen...	136	C10H16	003387-41-5	91
2	.beta.-Pinene	136	C10H16	000127-91-3	87
3	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	87
4	.alpha.-Pinene	136	C10H16	000080-56-8	86
5	Cyclohexene, 4-methylene-1-(1-me...	136	C10H16	000099-84-3	86





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Misc :
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Instrument : BNA_G
ClientSampleId : C0AZ0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

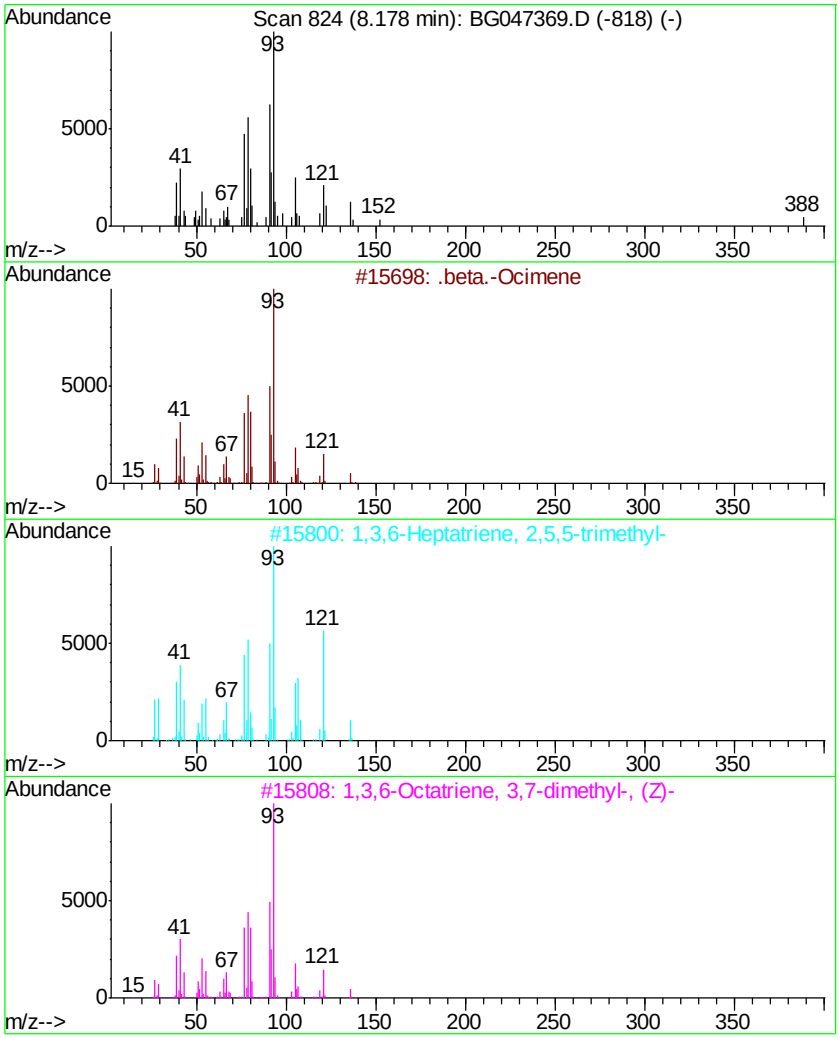
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

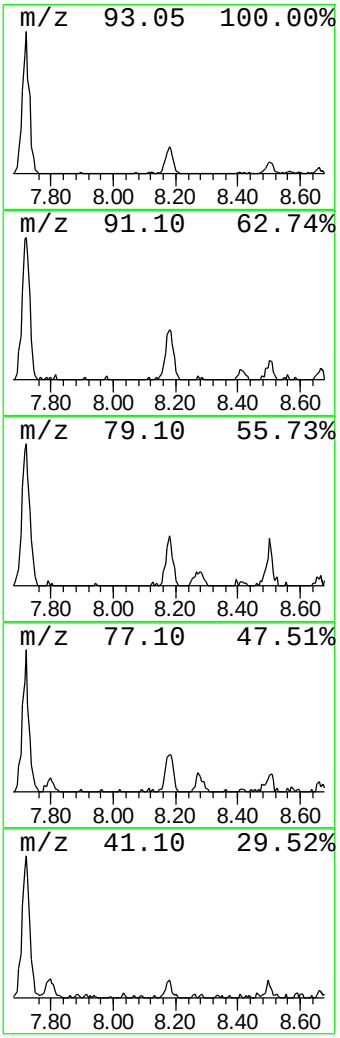
****

Peak Number 4 .beta.-Ocimene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.18	3.86 ng/ul	58533	1,4-Dichlorobenzene-d4	8.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-Ocimene	136	C10H16	013877-91-3	80
2	1,3,6-Heptatriene, 2,5,5-trimethyl-	136	C10H16	029548-02-5	80
3	1,3,6-Octatriene, 3,7-dimethyl-,...	136	C10H16	003338-55-4	80
4	.gamma.-Terpinene	136	C10H16	000099-85-4	68
5	1,3,6-Heptatriene, 2,5,5-trimethyl-	136	C10H16	029548-02-5	68





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Data File : BG047369.D
Acq On : 20 Nov 2020 2:07
Operator : CG/JU
Sample : L4710-23
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

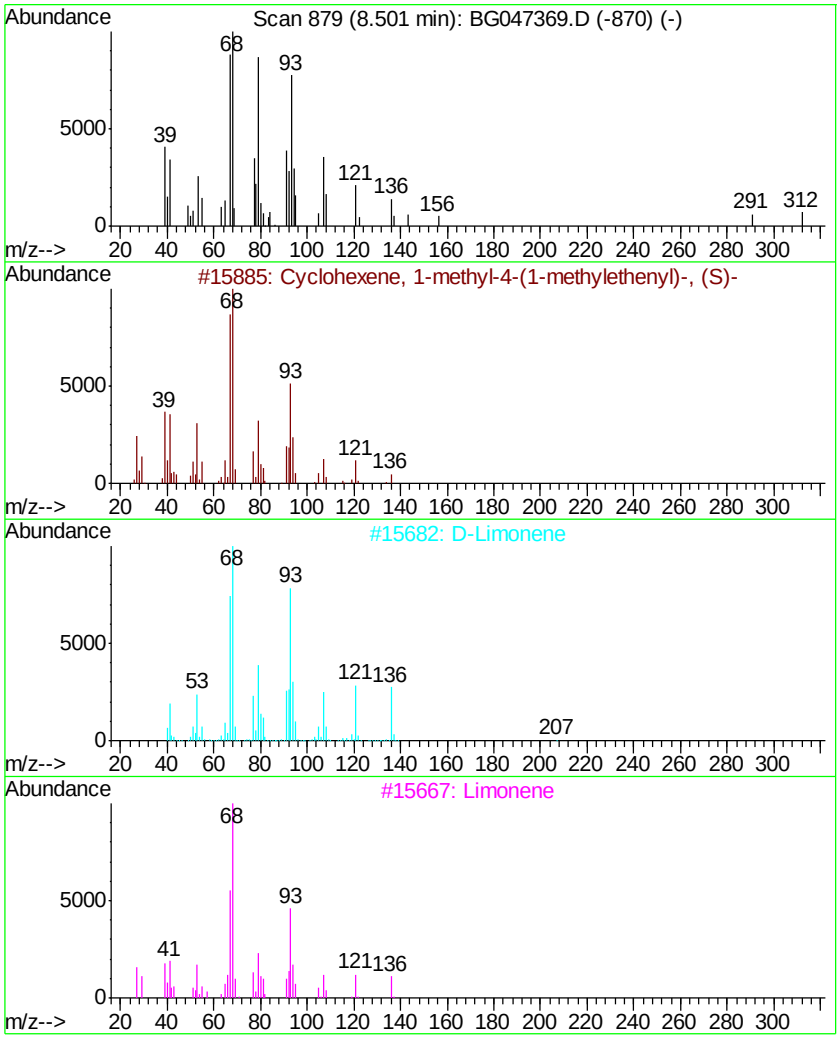
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

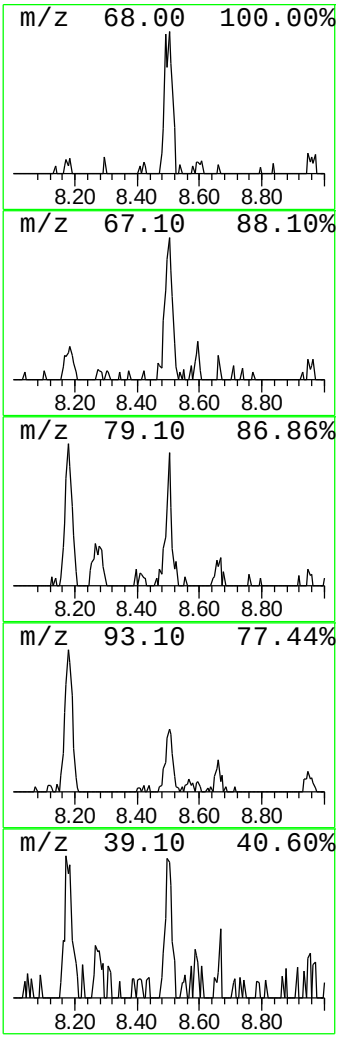
****

Peak Number 5 Cyclohexene, 1-methyl-4-(1-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.50	3.05 ng/ul	46186	1,4-Dichlorobenzene-d4	8.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	005989-54-8	76
2	D-Limonene	136	C10H16	005989-27-5	74
3	Limonene	136	C10H16	000138-86-3	62
4	D-Limonene	136	C10H16	005989-27-5	58
5	D-Limonene	136	C10H16	005989-27-5	52





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111920\
Data File : BG047369.D
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Sample : L4710-23
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

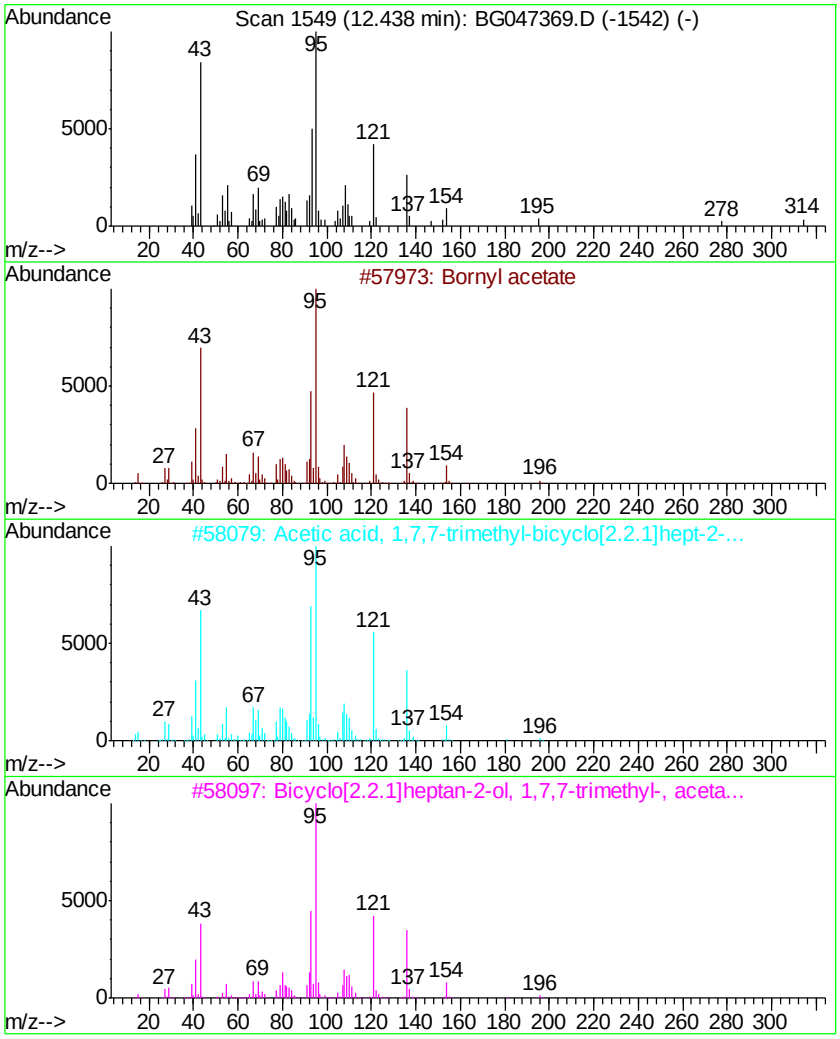
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

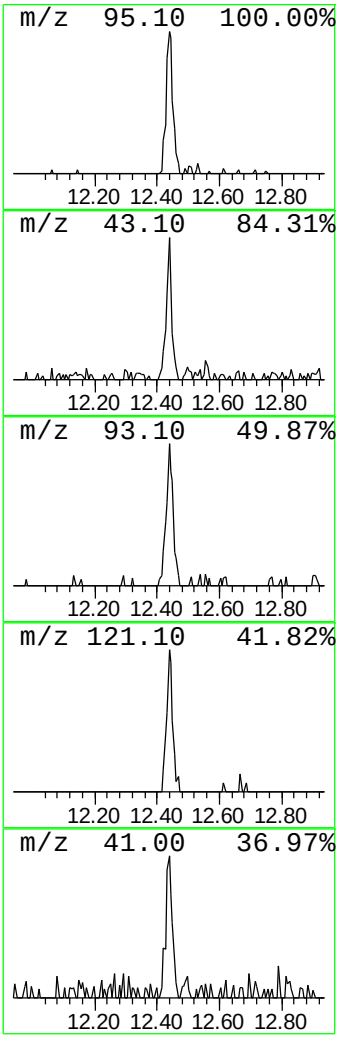
****

Peak Number 6 Bornyl acetate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.44	3.66 ng/ul	77453	Naphthalene-d8	11.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bornyl acetate	196	C12H20O2	000076-49-3	86
2	Acetic acid, 1,7,7-trimethyl-bic...	196	C12H20O2	092618-89-8	80
3	Bicyclo[2.2.1]heptan-2-ol, 1,7,7...	196	C12H20O2	005655-61-8	80
4	Isobornyl formate	182	C11H18O2	001200-67-5	80
5	Bornyl acetate	196	C12H20O2	000076-49-3	78





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Data File : BG047369.D
Acq On : 20 Nov 2020 2:07
Operator : CG/JU
Sample : L4710-23
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

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

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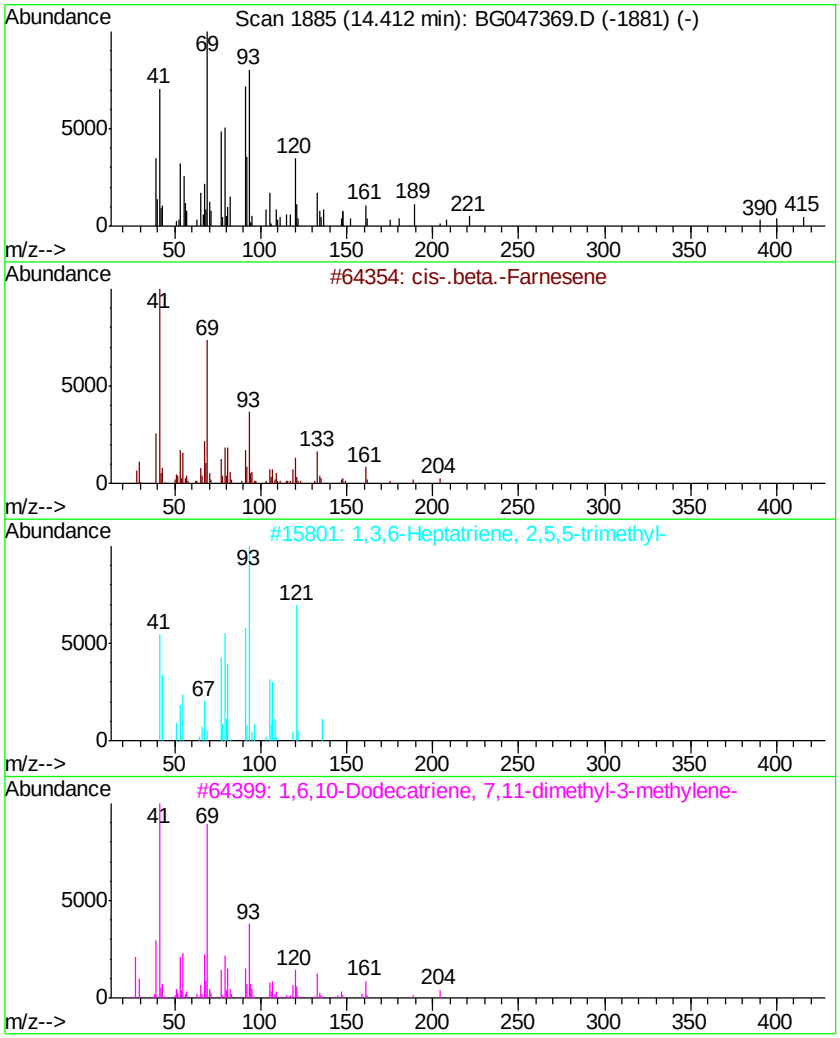
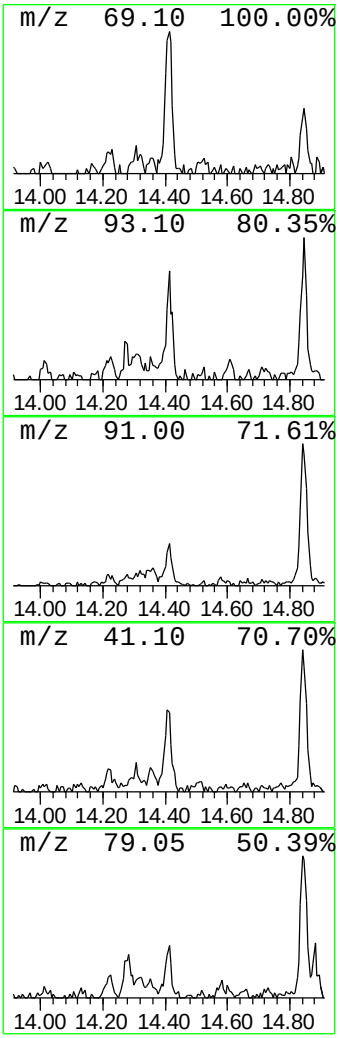
Peak Number 7 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.41	3.00 ng/ul	89399	Acenaphthene-d10	14.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	cis-.beta.-Farnesene	204	C15H24	028973-97-9	40
2	1,3,6-Heptatriene, 2,5,5-trimethyl-	136	C10H16	029548-02-5	38
3	1,6,10-Dodecatriene, 7,11-dimeth...	204	C15H24	077129-48-7	38
4	Bicyclo[3.1.0]hex-2-ene, 2-methy...	136	C10H16	002867-05-2	35
5	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	016022-04-1	35





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Data File : BG047369.D
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Operator : CG/JU
Sample : L4710-23
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

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

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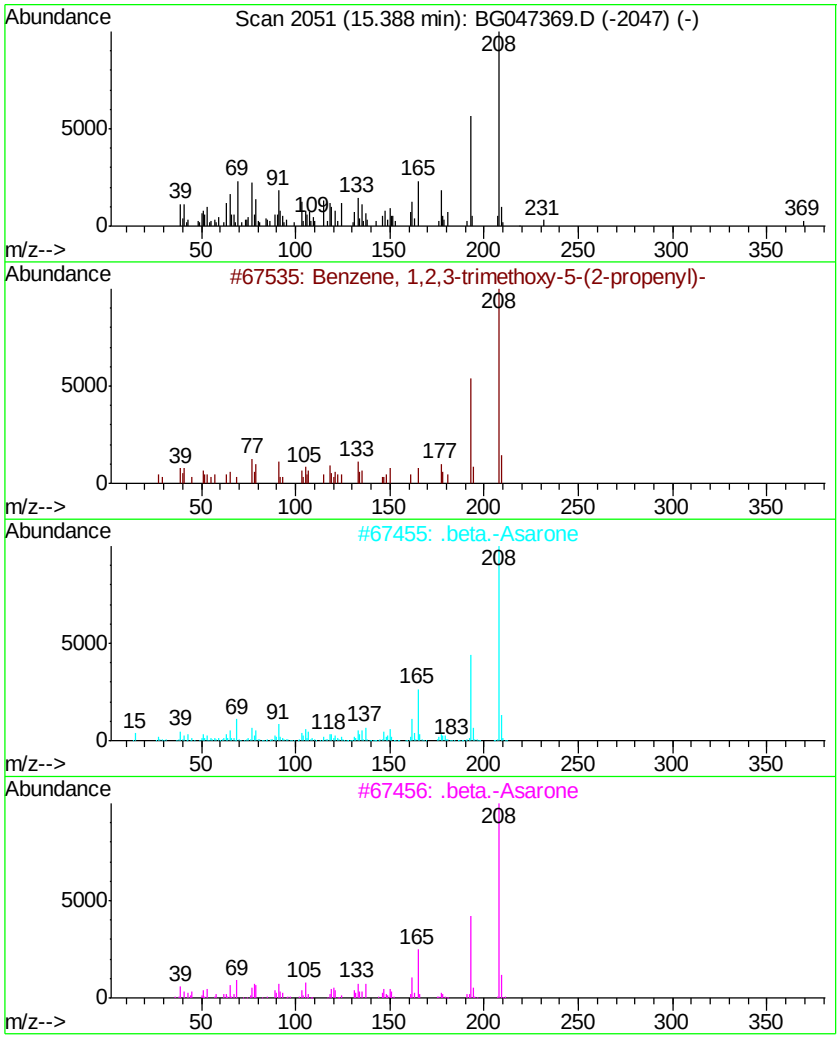
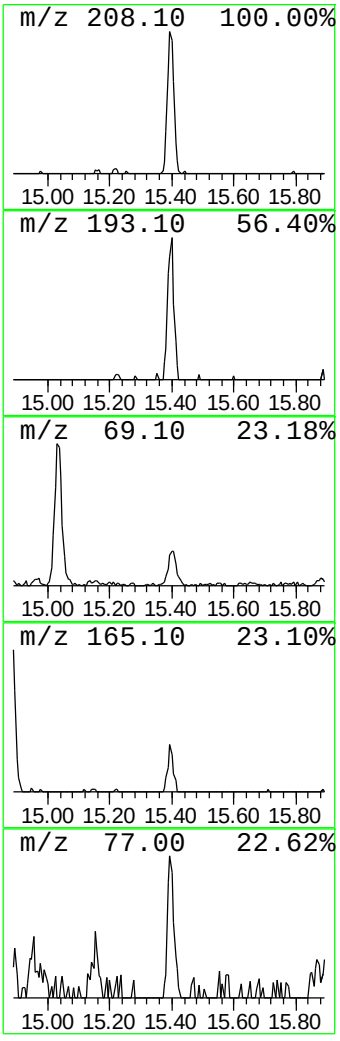
Peak Number 8 Benzene, 1,2,3-trimethoxy-5... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.39	4.60 ng/ul	136856	Acenaphthene-d10	14.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethoxy-5-(2-p...	208	C12H16O3	000487-11-6	93
2	.beta.-Asarone	208	C12H16O3	005273-86-9	86
3	.beta.-Asarone	208	C12H16O3	005273-86-9	83
4	Benzene, 1,2,3-trimethoxy-5-(2-p...	208	C12H16O3	000487-11-6	70
5	Asarone	208	C12H16O3	002883-98-9	70





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Data File : BG047369.D
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Operator : CG/JU
Sample : L4710-23
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

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

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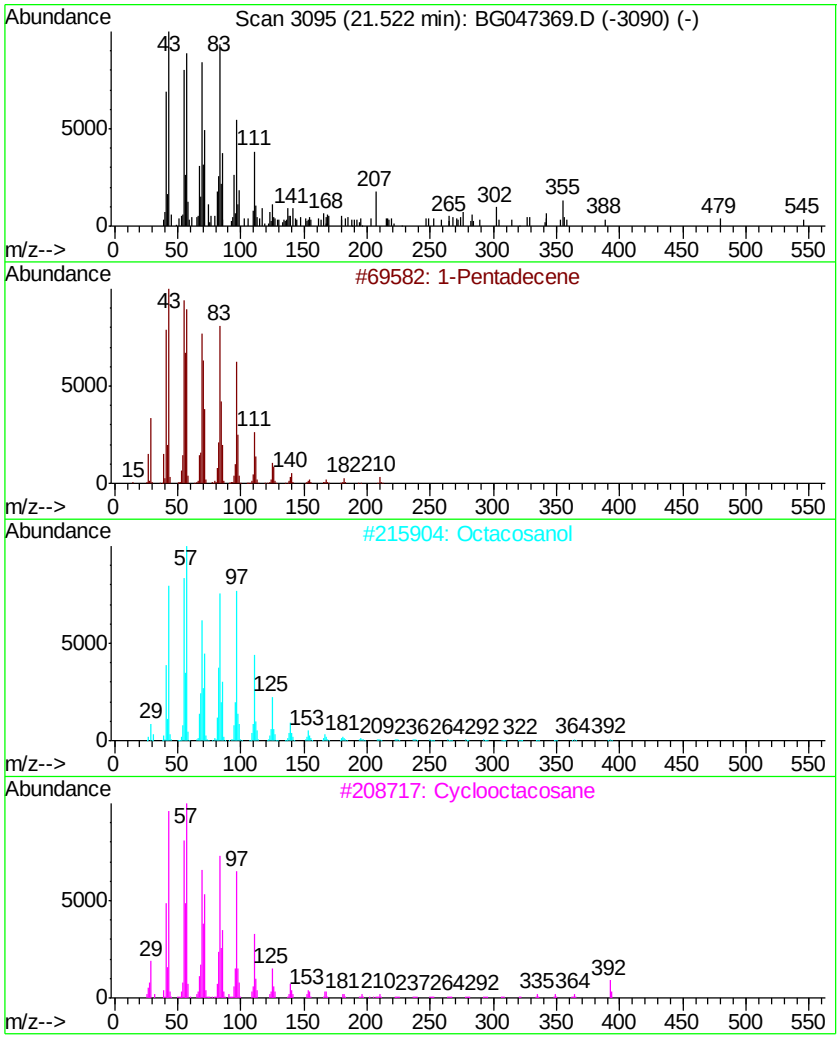
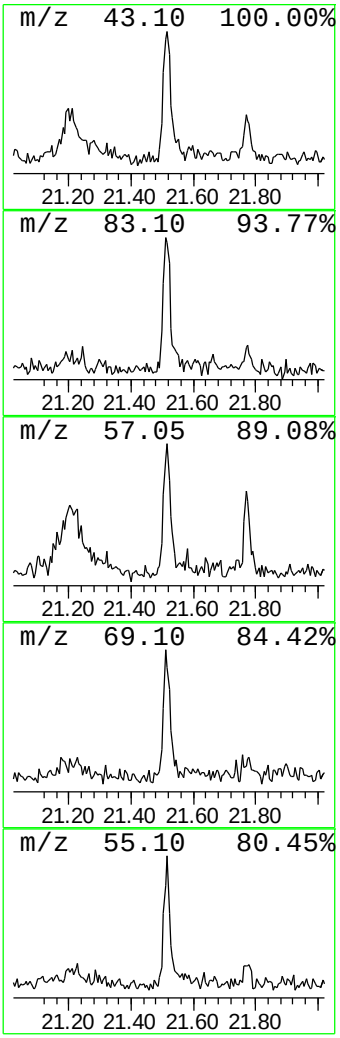
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.52	2.98 ng/ul	119859	Chrysene-d12	21.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentadecene	210	C15H30	013360-61-7	91
2	Octacosanol	410	C28H58O	000557-61-9	90
3	Cyclooctacosane	392	C28H56	000297-24-5	90
4	Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	90
5	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	90





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Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

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

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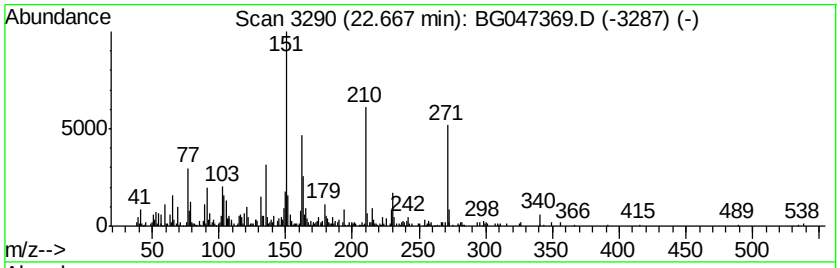
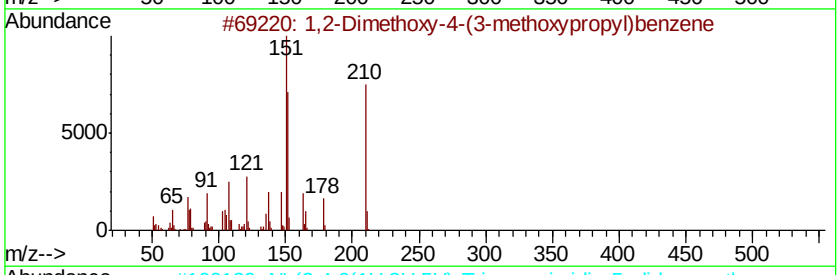
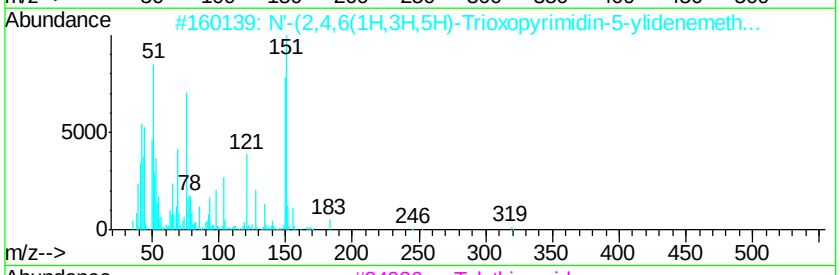
Peak Number 10 unknown-02 Concentration Rank 2

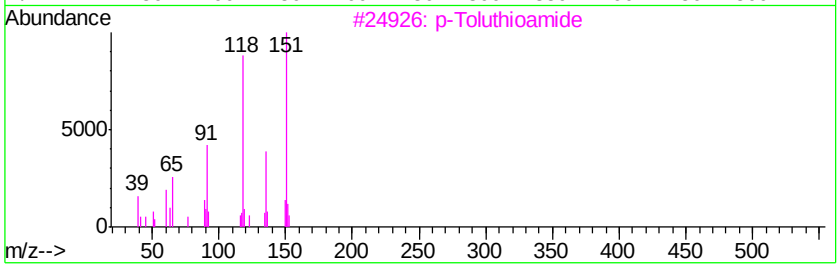
R.T.	EstConc	Area	Relative to ISTD	R.T.
22.67	12.15 ng/ul	487991	Chrysene-d12	21.90

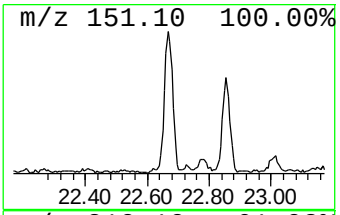
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Dimethoxy-4-(3-methoxypropyl...	210	C12H18O3	1000313-34-9	41
2	N'-(2,4,6(1H,3H,5H)-Trioxopyrimi...	319	C12H9N5O6	1000225-72-5	25
3	p-Toluthioamide	151	C8H9NS	002362-62-1	20
4	1-(2,4-Dimethoxyphenyl)-propan-2...	194	C11H14O3	000831-29-8	18
5	Silane, (3,3-dimethylbut-2-yloxy...	208	C12H20OSi	1000163-31-6	15

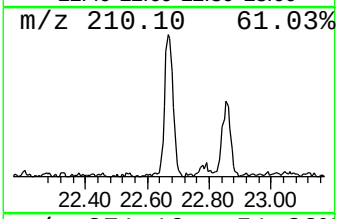


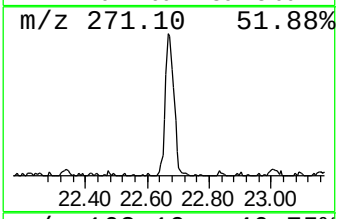


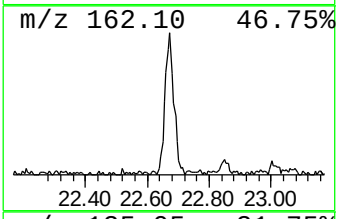


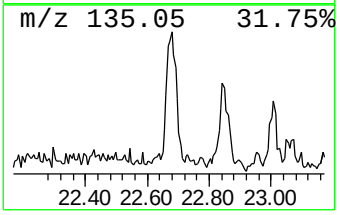












Tentatively Identified Compound (LSC) summary

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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
.alpha.-Pinene	6.95	10.1	ng/ul	152930	1	8.28	303329	20.0
2-Carene	7.26	8.3	ng/ul	125922	1	8.28	303329	20.0
Bicyclo[3.1.0]hex...	7.72	15.9	ng/ul	241520	1	8.28	303329	20.0
.beta.-Ocimene	8.18	3.9	ng/ul	58533	1	8.28	303329	20.0
Cyclohexene, 1-me...	8.50	3.0	ng/ul	46186	1	8.28	303329	20.0
Bornyl acetate	12.44	3.7	ng/ul	77453	2	11.10	423076	20.0
unknown-01	14.41	3.0	ng/ul	89399	3	14.88	595651	20.0
Benzene, 1,2,3-tr...	15.39	4.6	ng/ul	136856	3	14.88	595651	20.0
(DEL) Alkane: Str...	21.52	3.0	ng/ul	119859	5	21.90	803078	20.0
unknown-02	22.67	12.2	ng/ul	487991	5	21.90	803078	20.0