LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
Data File : BG046763.D
Acq On : 3 Oct 2020 9:08
Operator : CG/JU
Sample : L4150-09
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7J5

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-BG092920MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG046763.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.534	28	33	39	rBV2	13179	21372	1.19%	0.092%
2	4.192	139	145	151	rVB8	8796	17020	0.95%	0.074%
3	4.291	155	162	170	rVB	29549	49398	2.76%	0.214%
4	4.503	191	198	203	rVB2	42915	62831	3.51%	0.272%
5	4.644	216	222	226	rBV4	17328	24904	1.39%	0.108%
6	4.914	261	268	277	rBV	28069	44985	2.51%	0.195%
7	5.202	311	317	324	rVB4	6956	10283	0.57%	0.044%
8	5.425	348	355	360	rBV8	5710	11154	0.62%	0.048%
9	5.596	380	384	386	rBV4	5823	8877	0.50%	0.038%
10	6.154	469	479	486	rVB8	23821	44389	2.48%	0.192%
11	7.200	652	657	660	rBV5	11012	16719	0.93%	0.072%
12	7.253	660	666	675	rVB	262970	462638	25.86%	2.002%
13	7.429	690	696	706	rBV	176163	305859	17.10%	1.323%
14	7.640	725	732	740	rBV	246440	416709	23.30%	1.803%
15	7.729	744	747	750	rVV4	7461	10902	0.61%	0.047%
16	7.799	754	759	763	rVB4	11766	18911	1.06%	0.082%
17	8.111	803	812	819	rVV	254143	433320	24.23%	1.875%
18	8.163	819	821	827	rVB4	8419	12651	0.71%	0.055%
19	8.704	909	913	917	rVB4	7188	9850	0.55%	0.043%
20	8.804	923	930	941	rVB	316950	532307	29.76%	2.303%
21	8.927	947	951	962	rVB3	20113	38306	2.14%	0.166%
22	9.268	1003	1009	1018	rVB	246481	431791	24.14%	1.868%
23	9.338	1018	1021	1026	rBV6	5455	9056	0.51%	0.039%
24	9.433	1031	1037	1043	rVB2	71968	114086	6.38%	0.494%
25	9.697	1077	1082	1088	rVB3	18807	28834	1.61%	0.125%
26	9.997	1125	1133	1140	rBV	210721	378576	21.16%	1.638%
27	10.326	1183	1189	1191	rBV6	5972	11999	0.67%	0.052%
28	10.425	1202	1206	1211	rBV4	13097	24532	1.37%	0.106%
29	10.531	1216	1224	1232	rBV2	323103	585265	32.72%	2.532%
30	10.925	1281	1291	1297	rBV	341425	625070	34.95%	2.704%
31	10.972	1297	1299	1302	rVV2	15974	22000	1.23%	0.095%
32	11.048	1306	1312	1321	rVB	206066	384746	21.51%	1.665%
33	11.565	1395	1400	1405	rBV7	9198	17672	0.99%	0.076%
34	11.853	1443	1449	1455	rBV7	5741	11054	0.62%	0.048%
35	11.935	1457	1463	1466	rBV4	7167	13063	0.73%	0.057%
36	12.241	1511	1515	1520	rVB6	7908	12990	0.73%	0.056%

LSC Area Percent Report

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Operator : CG/JU
Sample : L4150-09
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7J5

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-BG092920MA.M
Title : SVOA CALIBRATION

37	12.323	1526	1529	1535	rBV4	6375	10493	0.59%	0.045%
38	12.458	1547	1552	1556	rBV5	9551	12902	0.72%	0.056%
39	12.570	1566	1571	1575	rVB3	12128	20822	1.16%	0.090%
40	12.787	1604	1608	1613	rVB5	8833	14721	0.82%	0.064%
41	13.099	1657	1661	1664	rVB3	5504	8490	0.47%	0.037%
42	13.187	1671	1676	1679	rBV3	7052	10940	0.61%	0.047%
43	13.269	1686	1690	1698	rVB7	9537	15540	0.87%	0.067%
44	13.340	1698	1702	1712	rBV6	10699	21354	1.19%	0.092%
45	13.557	1735	1739	1743	rVV4	11370	18887	1.06%	0.082%
46	13.622	1743	1750	1757	rVV2	40665	77345	4.32%	0.335%
47	13.686	1759	1761	1767	rVB5	10427	16669	0.93%	0.072%
48	13.898	1793	1797	1799	rBV4	6885	9482	0.53%	0.041%
49	14.057	1819	1824	1829	rBV6	11590	21816	1.22%	0.094%
50	14.133	1829	1837	1842	rBV	616398	893409	49.95%	3.865%
51	14.180	1842	1845	1849	rVB	40604	54859	3.07%	0.237%
52	14.368	1873	1877	1878	rBV3	9537	9164	0.51%	0.040%
53	14.427	1881	1887	1898	rVB	626986	994337	55.59%	4.302%
54	14.591	1911	1915	1919	rBV2	18268	22399	1.25%	0.097%
55	14.732	1932	1939	1947	rVV	593373	944000	52.78%	4.084%
56	14.797	1948	1950	1957	rVB4	19255	25891	1.45%	0.112%
57	14.903	1960	1968	1978	rBV	298418	491626	27.48%	2.127%
58	15.061	1991	1995	1999	rVV6	5358	9985	0.56%	0.043%
59	15.132	2001	2007	2014	rVB3	43993	69050	3.86%	0.299%
60	15.390	2047	2051	2058	rVB9	16647	27503	1.54%	0.119%
61	15.531	2067	2075	2076	rBV3	13390	21356	1.19%	0.092%
62	15.725	2101	2108	2113	rBV	838530	1268787	70.93%	5.489%
63	15.778	2113	2117	2120	rVV2	15811	21308	1.19%	0.092%
64	15.825	2120	2125	2131	rVB	238569	339068	18.96%	1.467%
65	16.066	2164	2166	2167	rBV2	8477	8276	0.46%	0.036%
66	16.230	2190	2194	2199	rVV8	9044	24533	1.37%	0.106%
67	16.283	2200	2203	2212	rVB9	23986	37078	2.07%	0.160%
68	16.442	2225	2230	2237	rVV10	16784	32877	1.84%	0.142%
69	16.501	2239	2240	2244	rVB4	8395	8811	0.49%	0.038%
70	16.589	2252	2255	2258	rBV4	12756	18669	1.04%	0.081%
71	16.789	2286	2289	2292	rVV4	10295	13702	0.77%	0.059%
72	16.924	2310	2312	2317	rVB6	5872	9470	0.53%	0.041%
73	16.982	2317	2322	2323	rBV4	5492	8338	0.47%	0.036%
74	17.124	2343	2346	2351	rVB4	18630	24369	1.36%	0.105%
75	17.294	2371	2375	2379	rBV2	13510	21663	1.21%	0.094%
76	17.394	2388	2392	2396	rVB6	11882	17475	0.98%	0.076%

LSC Area Percent Report

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Operator : CG/JU
Sample : L4150-09
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7J5

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-BG092920MA.M
Title : SVOA CALIBRATION

77	17.476	2400	2406	2410	rBV	768228	1134855	63.45%	4.910%
78	17.517	2410	2413	2417	rVV	262157	366364	20.48%	1.585%
79	17.576	2417	2423	2434	rVB	871193	1539943	86.09%	6.662%
80	17.658	2434	2437	2439	rBV4	10077	12361	0.69%	0.053%
81	17.870	2467	2473	2478	rBV3	29383	54426	3.04%	0.235%
82	18.134	2516	2518	2525	rVB7	12496	18057	1.01%	0.078%
83	18.310	2546	2548	2551	rBV2	15999	18013	1.01%	0.078%
84	18.357	2551	2556	2559	rVV2	69071	99682	5.57%	0.431%
85	18.398	2559	2563	2567	rVB2	59492	76807	4.29%	0.332%
86	18.545	2582	2588	2591	rBV3	60075	108847	6.09%	0.471%
87	18.845	2634	2639	2642	rBV	43168	65811	3.68%	0.285%
88	18.880	2643	2645	2650	rVB3	33633	41300	2.31%	0.179%
89	19.256	2705	2709	2716	rBV6	48881	91733	5.13%	0.397%
90	19.374	2726	2729	2741	rVB2	82311	168729	9.43%	0.730%
91	19.521	2749	2754	2759	rVV	515397	719786	40.24%	3.114%
92	19.568	2759	2762	2767	rVB3	51230	70355	3.93%	0.304%
93	19.856	2805	2811	2814	rBV	1175435	1788712	100.00%	7.738%
94	21.389	3068	3072	3081	rVB4	293172	503134	28.13%	2.177%
95	21.747	3126	3133	3138	rBV2	796291	1633905	91.35%	7.069%
96	23.980	3508	3513	3520	rBV	150844	396952	22.19%	1.717%
97	24.127	3533	3538	3542	rBV	153263	255212	14.27%	1.104%
98	24.803	3646	3653	3660	rBV	449378	1069340	59.78%	4.626%
99	25.020	3683	3690	3699	rBV2	395695	1096298	61.29%	4.743%
100	26.272	3897	3903	3913	rVB2	306396	882340	49.33%	3.817%

Sum of corrected areas: 23114545

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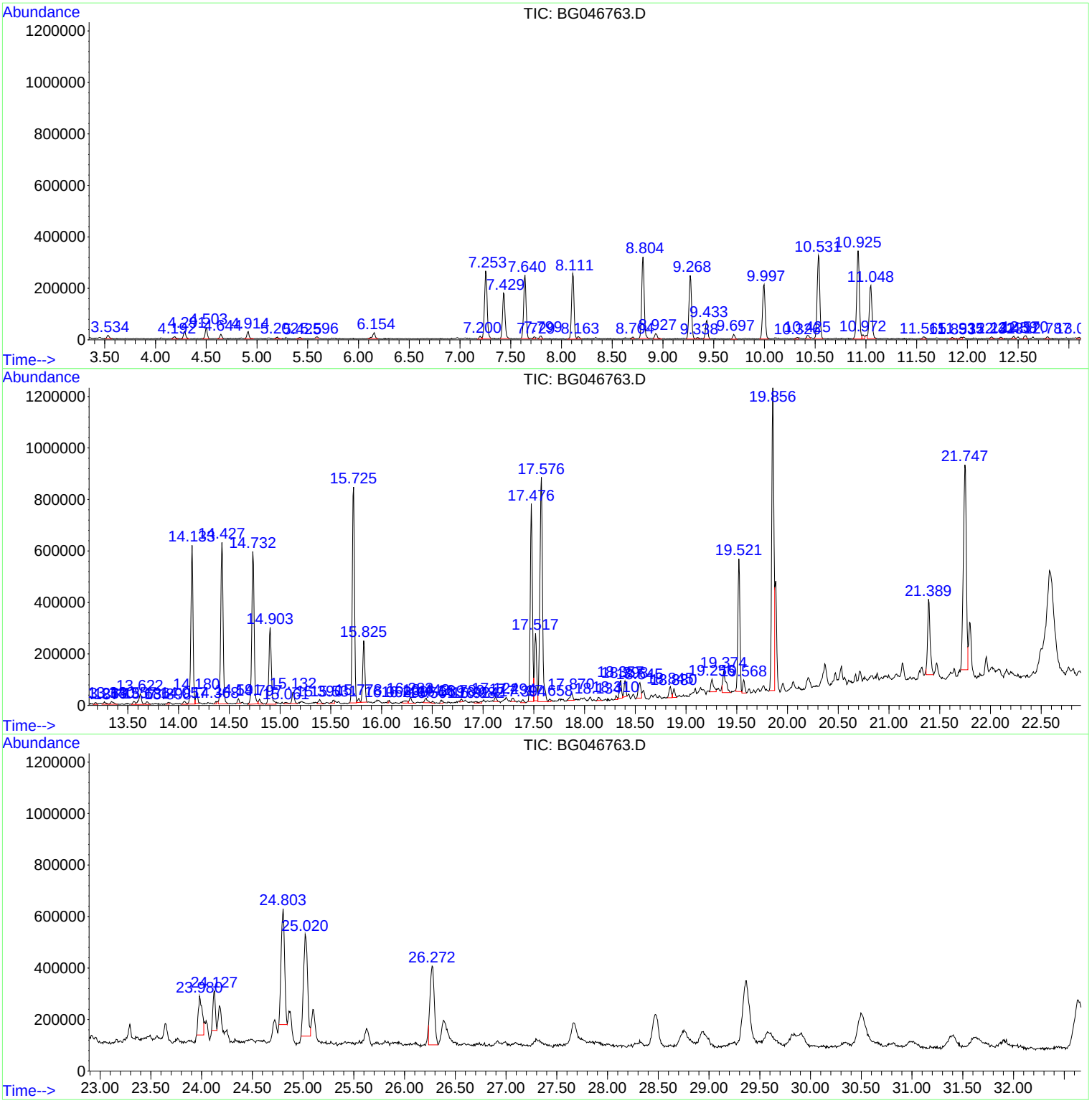
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CB7J5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
Data File : BG046763.D
Acq On : 3 Oct 2020 9:08
Operator : CG/JU
Sample : L4150-09
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7J5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

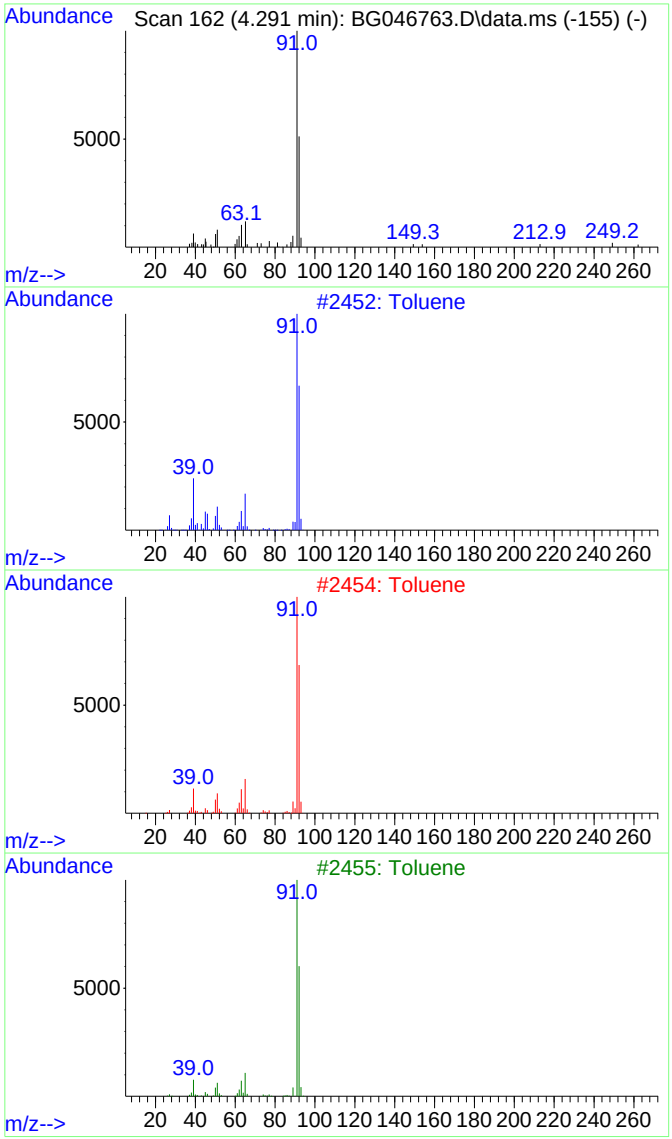
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

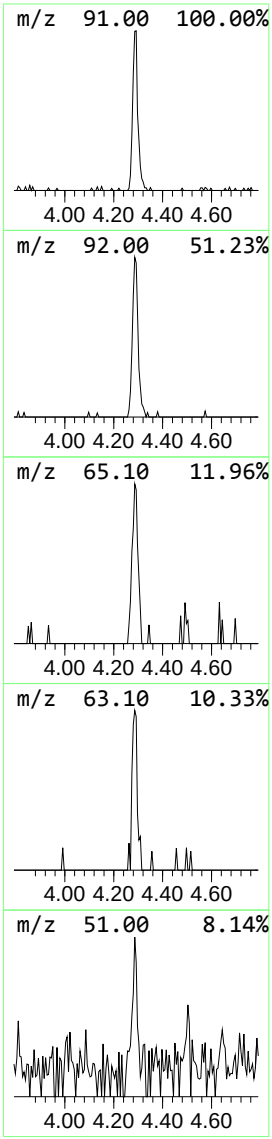
****

Peak Number 1 Toluene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.290	2.28 ng/ul	49398	1,4-Dichlorobenzene-d4	8.111

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene	92	C7H8	000108-88-3	90
2	Toluene	92	C7H8	000108-88-3	90
3	Toluene	92	C7H8	000108-88-3	87
4	Molybdenum, di-.mu.-chlorobis[(1...	532	C20H26Cl2Mo2	035625-66-2	78
5	1,3,5-Cycloheptatriene	92	C7H8	000544-25-2	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
Data File : BG046763.D
Acq On : 3 Oct 2020 9:08
Operator : CG/JU
Sample : L4150-09
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CB7J5

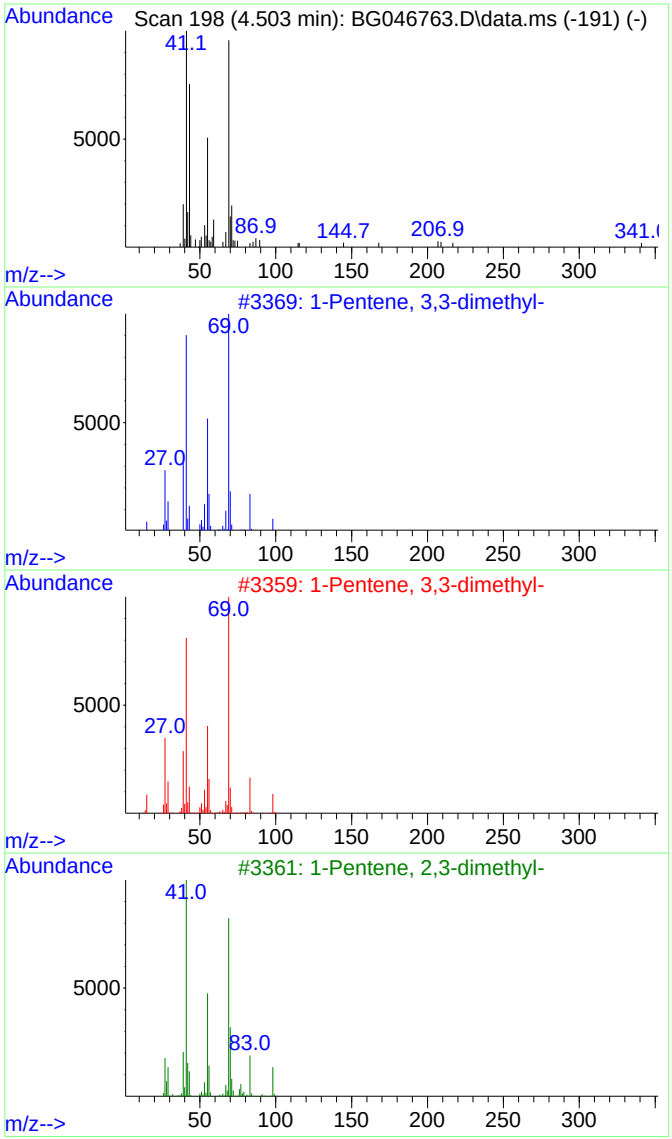
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

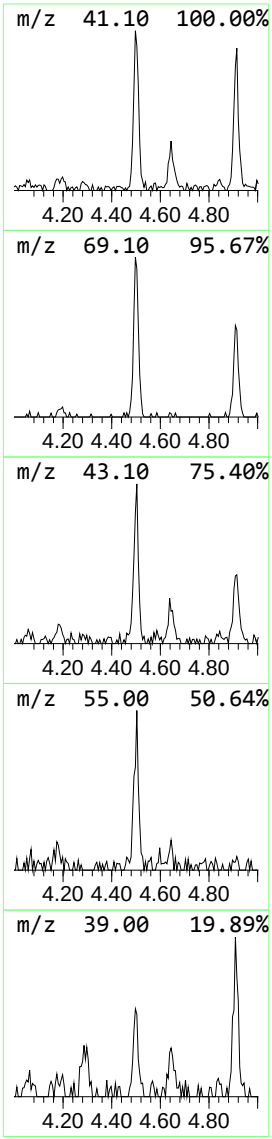
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.500	2.90 ng/ul	62831	1,4-Dichlorobenzene-d4	8.111

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene, 3,3-dimethyl-	98	C7H14	003404-73-7	38
2	1-Pentene, 3,3-dimethyl-	98	C7H14	003404-73-7	38
3	1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	38
4	2-Butene, 1-bromo-3-methyl-	148	C5H9Br	000870-63-3	37
5	1-Pentene, 3-ethyl-	98	C7H14	004038-04-4	37





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
Data File : BG046763.D
Acq On : 3 Oct 2020 9:08
Operator : CG/JU
Sample : L4150-09
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7J5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

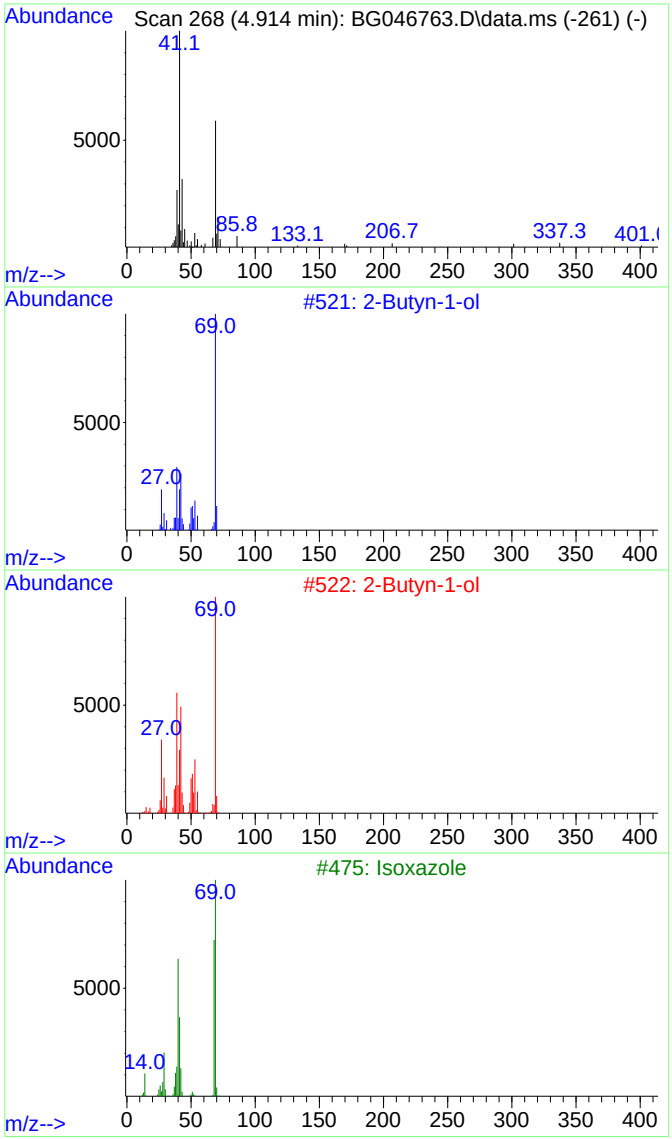
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

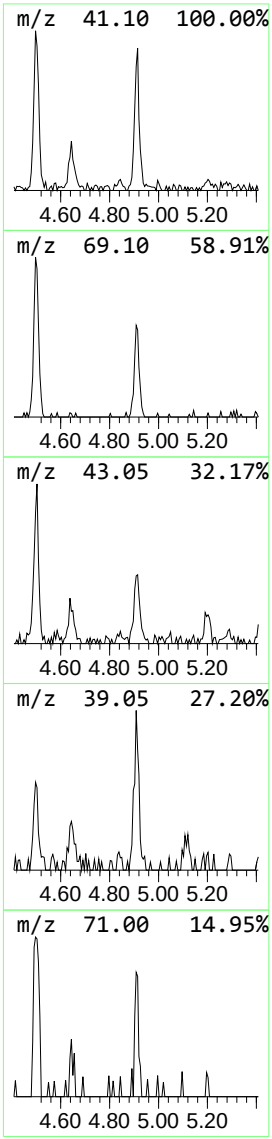
****

Peak Number 3 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.910	2.08 ng/ul	44985	1,4-Dichlorobenzene-d4	8.111

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butyn-1-ol	70	C4H6O	000764-01-2	10
2	2-Butyn-1-ol	70	C4H6O	000764-01-2	9
3	Isoxazole	69	C3H3NO	000288-14-2	9
4	(Aminomethyl)cyclopropane	71	C4H9N	002516-47-4	9
5	4-Cyclopropylcarbonyloxytridecane	268	C17H32O2	1000245-58-5	9





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Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7J5

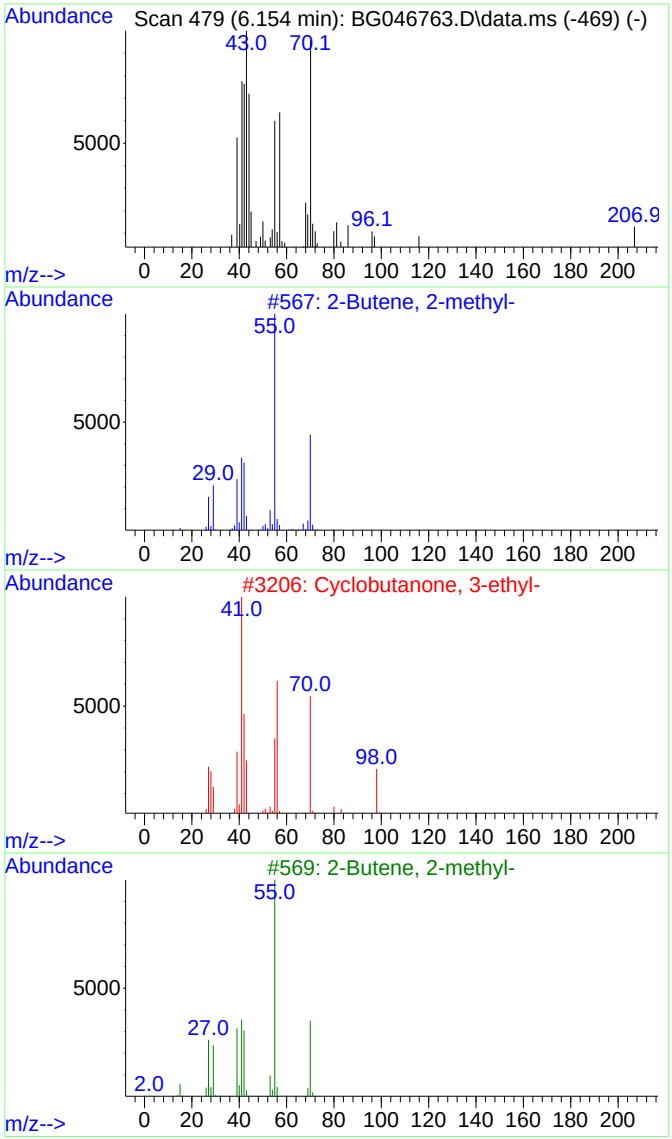
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

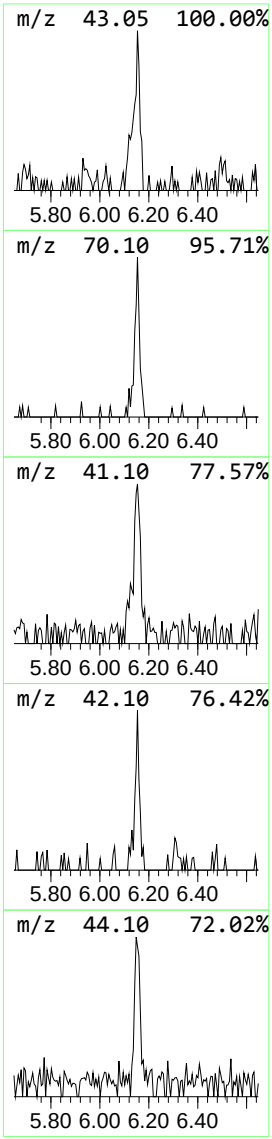
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.150	2.05 ng/ul	44389	1,4-Dichlorobenzene-d4	8.111

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butene, 2-methyl-	70	C5H10	000513-35-9	43
2	Cyclobutanone, 3-ethyl-	98	C6H10O	056335-73-0	43
3	2-Butene, 2-methyl-	70	C5H10	000513-35-9	38
4	2-Butene, 2-methyl-	70	C5H10	000513-35-9	38
5	Pentane, 2-chloro-	106	C5H11Cl	000625-29-6	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100220\
Data File : BG046763.D
Acq On : 3 Oct 2020 9:08
Operator : CG/JU
Sample : L4150-09
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7J5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

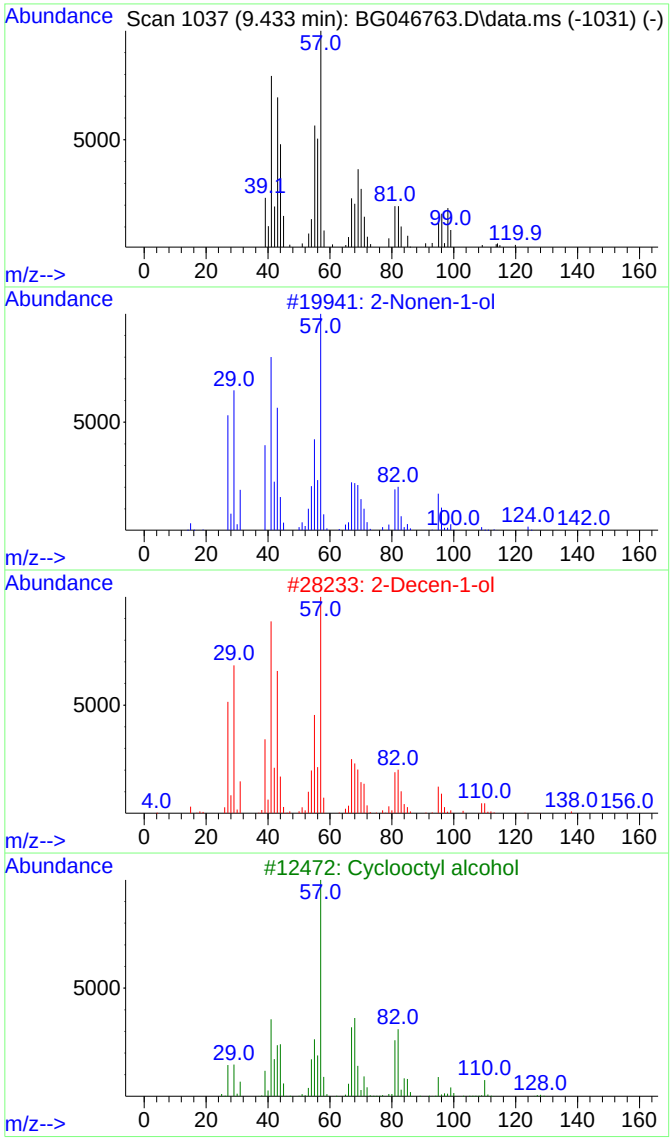
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

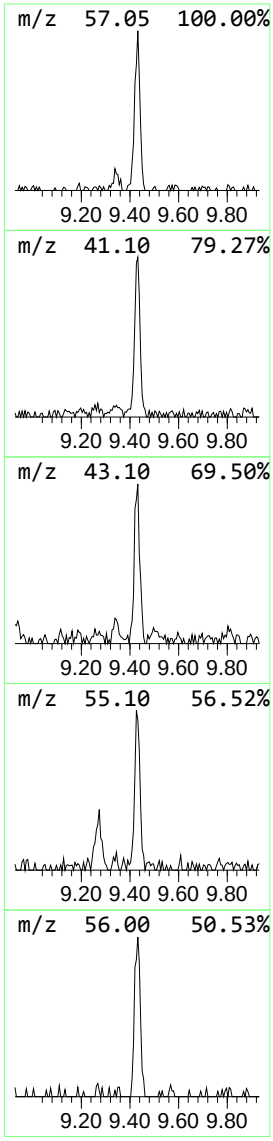
****

Peak Number 5 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.430	5.27 ng/ul	114086	1,4-Dichlorobenzene-d4	8.111

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Nonen-1-ol	142	C9H18O	022104-79-6	27
2	2-Decen-1-ol	156	C10H20O	022104-80-9	27
3	Cyclooctyl alcohol	128	C8H16O	000696-71-9	27
4	Heptane, 2-chloro-	134	C7H15Cl	001001-89-4	22
5	Cyclohexanol	100	C6H12O	000108-93-0	22





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Data File : BG046763.D
Acq On : 3 Oct 2020 9:08
Operator : CG/JU
Sample : L4150-09
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7J5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

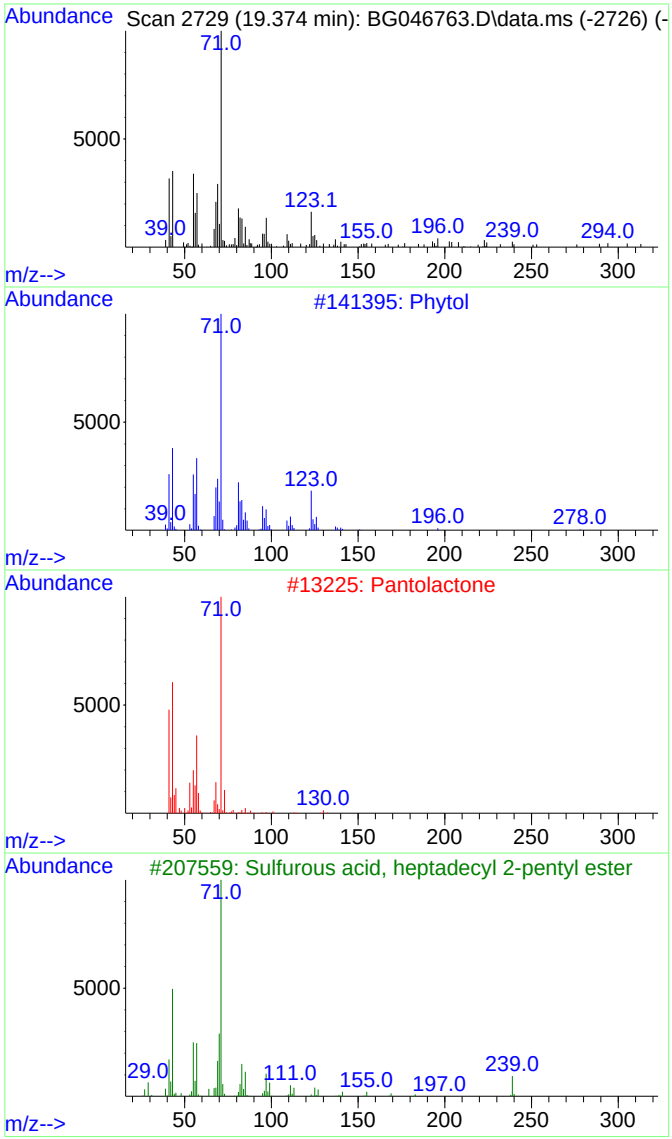
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

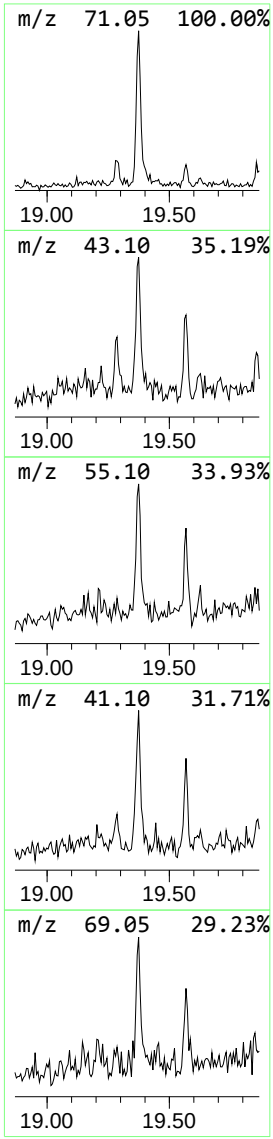
****

Peak Number 6 Phytol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.370	2.97 ng/ul	168729	Phenanthrene-d10	17.476

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phytol	296	C20H40O	000150-86-7	64
2	Pantolactone	130	C6H10O3	000599-04-2	53
3	Sulfurous acid, heptadecyl 2-pen...	390	C22H46O3S	1000309-16-7	50
4	Cyclooctane, methoxy-	142	C9H18O	013213-32-6	49
5	Isobutyric acid, 2,2,2-trichloro...	218	C6H9Cl3O2	1000354-63-4	43





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Sample : L4150-09
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CB7J5

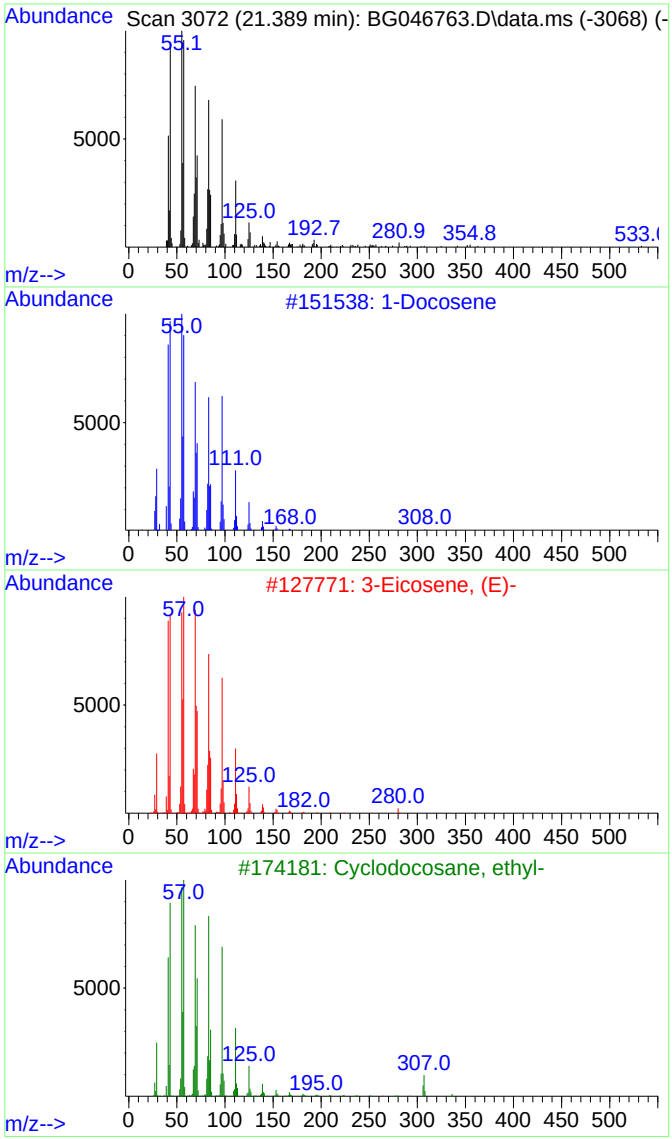
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

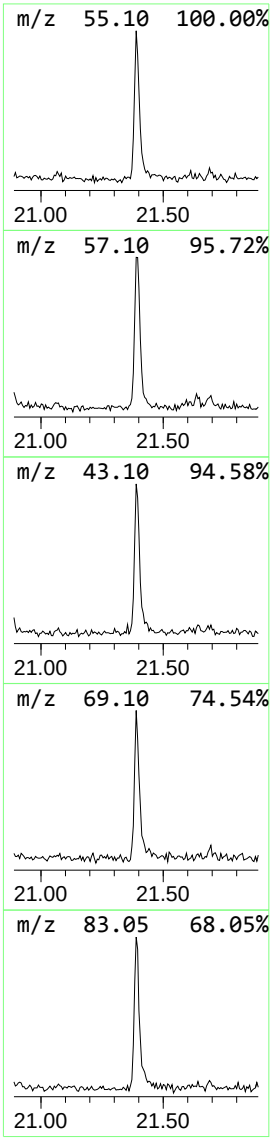
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.390	6.16 ng/ul	503134	Chrysene-d12	21.753

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	92
2	3-Eicosene, (E)-	280	C20H40	074685-33-9	91
3	Cyclodocosane, ethyl-	336	C24H48	1000151-22-6	91
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	91
5	Behenic alcohol	326	C22H46O	000661-19-8	91





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Data File : BG046763.D
Acq On : 3 Oct 2020 9:08
Operator : CG/JU
Sample : L4150-09
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7J5

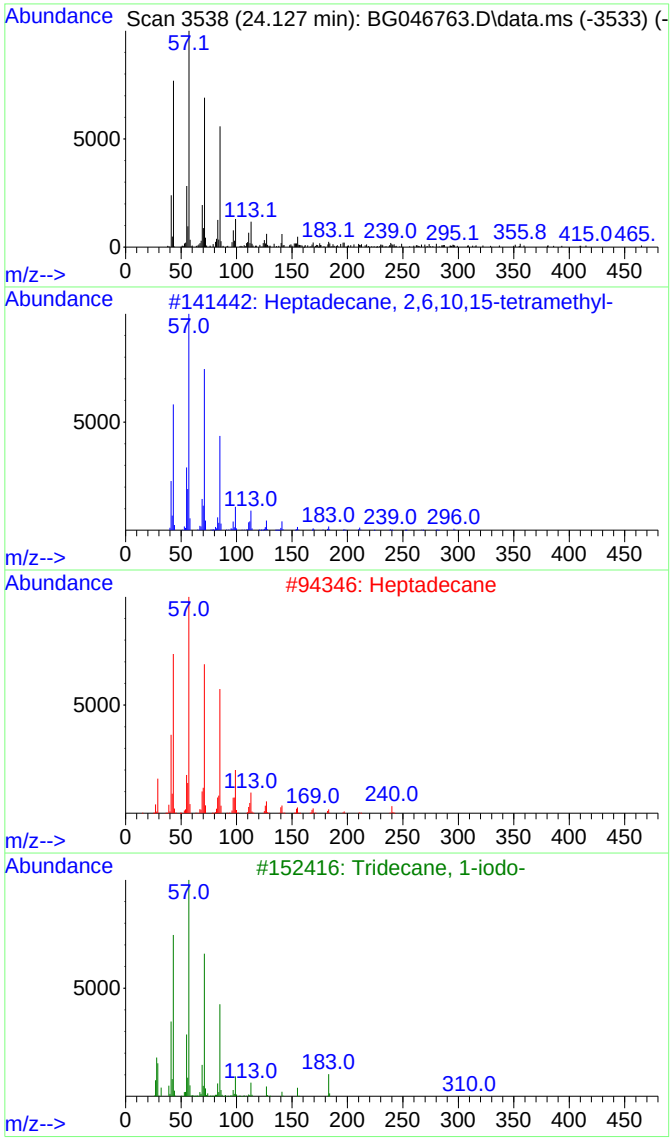
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

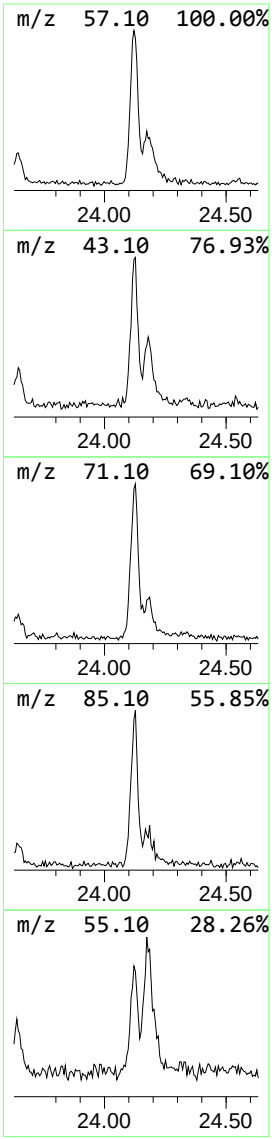
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.130	4.66 ng/ul	255212	Perylene-d12	25.020

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87
2	Heptadecane	240	C17H36	000629-78-7	87
3	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	87
4	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	87
5	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	86





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Data File : BG046763.D
Acq On : 3 Oct 2020 9:08
Operator : CG/JU
Sample : L4150-09
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7J5

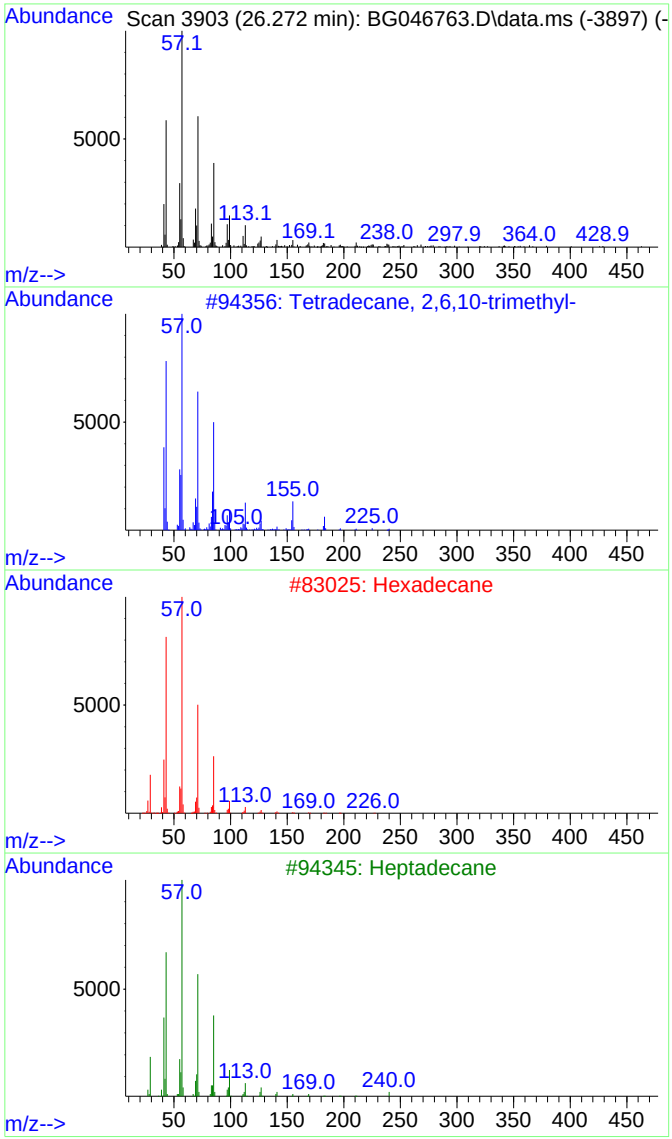
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
Quant Title : SVOA CALIBRATION

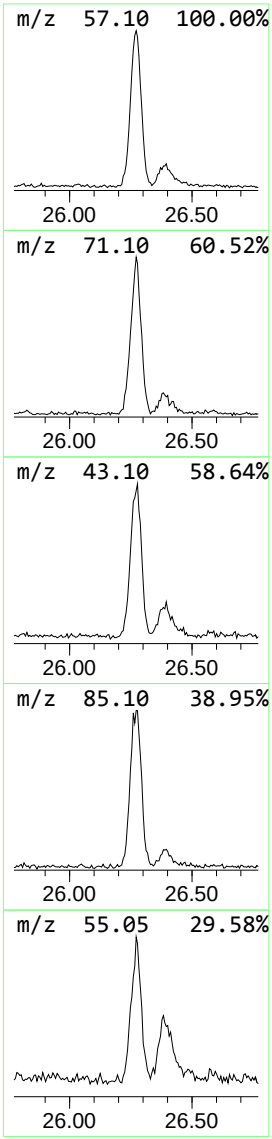
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.270	16.10 ng/ul	882340	Perylene-d12	25.020

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	93
2	Hexadecane	226	C16H34	000544-76-3	91
3	Heptadecane	240	C17H36	000629-78-7	91
4	Octacosane	394	C28H58	000630-02-4	90
5	Hexatriacontane	507	C36H74	000630-06-8	90





Tentatively Identified Compound (LSC) summary

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Data File : BG046763.D
Acq On : 3 Oct 2020 9:08
Operator : CG/JU
Sample : L4150-09
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : CB7J5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG092920MA.M
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
Toluene	4.290	2.3	ng/ul	49398	1	8.111	433320	20.0
(DEL) Alkane: S...	4.500	2.9	ng/ul	62831	1	8.111	433320	20.0
unknown-01	4.910	2.1	ng/ul	44985	1	8.111	433320	20.0
(DEL) Alkane: S...	6.150	2.0	ng/ul	44389	1	8.111	433320	20.0
unknown-02	9.430	5.3	ng/ul	114086	1	8.111	433320	20.0
Phytol	19.370	3.0	ng/ul	168729	4	17.476	1134860	20.0
(DEL) Alkane: S...	21.390	6.2	ng/ul	503134	5	21.753	1633910	20.0
(DEL) Alkane: S...	24.130	4.7	ng/ul	255212	6	25.020	1096300	20.0
(DEL) Alkane: S...	26.270	16.1	ng/ul	882340	6	25.020	1096300	20.0