

Data Path : Z:\HPCHEM1\BNA G\DATA\BG083017\
 Data File : BG028567.D
 Acq On : 31 Aug 2017 00:17
 Operator : SJ/JU
 Sample : I4917-09
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B0B51

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.1
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 1

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.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	3.818	35	39	47	rVB5	6626	13370	1.23%	0.102%
2	4.446	141	146	153	rVB4	4513	10055	0.92%	0.077%
3	4.558	154	165	170	rVB7	5185	13661	1.26%	0.104%
4	4.758	190	199	207	rBV	21299	37448	3.44%	0.285%
5	4.875	212	219	232	rVV7	7367	20441	1.88%	0.156%
6	5.157	262	267	276	rVB2	12955	21652	1.99%	0.165%
7	5.445	309	316	321	rBV	4725	8918	0.82%	0.068%
8	5.851	381	385	391	rBV3	6644	9685	0.89%	0.074%
9	5.986	399	408	419	rVB2	27266	67947	6.25%	0.517%
10	6.356	461	471	475	rVB7	4732	14368	1.32%	0.109%
11	6.603	507	513	520	rVV4	2181	6210	0.57%	0.047%
12	6.820	546	550	554	rVV3	3021	5671	0.52%	0.043%
13	6.867	554	558	564	rVV4	4463	6557	0.60%	0.050%
14	6.973	570	576	582	rVB6	4380	9872	0.91%	0.075%
15	7.190	609	613	621	rBV6	2656	6702	0.62%	0.051%
16	7.314	627	634	639	rBV2	12053	20254	1.86%	0.154%
17	7.419	647	652	657	rBV2	22601	39428	3.63%	0.300%
18	7.496	657	665	681	rVB	134999	321796	29.60%	2.449%
19	7.654	685	692	701	rBV	99576	177595	16.34%	1.351%
20	7.719	701	703	711	rVB6	4215	5815	0.53%	0.044%
21	7.878	721	730	738	rBV	134404	243407	22.39%	1.852%
22	7.978	742	747	755	rVB	29660	47650	4.38%	0.363%
23	8.136	771	774	780	rBV4	3440	5520	0.51%	0.042%
24	8.289	796	800	803	rBV3	3695	5068	0.47%	0.039%
25	8.342	803	809	823	rVB2	154033	289607	26.64%	2.204%
26	8.448	823	827	834	rBV2	9026	15808	1.45%	0.120%
27	8.718	863	873	879	rBV4	7644	16639	1.53%	0.127%
28	8.818	885	890	895	rBV4	7299	13321	1.23%	0.101%
29	8.877	895	900	909	rVB3	11573	24578	2.26%	0.187%
30	8.971	909	916	921	rBV4	18147	34109	3.14%	0.260%
31	9.041	921	928	940	rVV	163293	314198	28.90%	2.391%
32	9.135	940	944	947	rVV4	4040	5980	0.55%	0.046%
33	9.176	947	951	957	rVV4	6507	10050	0.92%	0.076%
34	9.276	965	968	973	rBV2	4240	6341	0.58%	0.048%

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35	9.347	974	980	984	rVB3	3940	6837	0.63%	0.052%
36	9.441	984	996	1001	rBV2	17372	35132	3.23%	0.267%
37	9.511	1001	1008	1018	rVB2	146474	285992	26.31%	2.176%
38	9.605	1018	1024	1027	rBV4	3916	6538	0.60%	0.050%
39	9.681	1035	1037	1043	rVB3	4154	5167	0.48%	0.039%
40	9.969	1078	1086	1092	rBV7	6951	13844	1.27%	0.105%
41	10.022	1092	1095	1100	rBV3	5111	8037	0.74%	0.061%
42	10.234	1121	1131	1142	rBV	114910	230245	21.18%	1.752%
43	10.322	1143	1146	1152	rBV5	4757	8532	0.78%	0.065%
44	10.387	1154	1157	1166	rVB5	6082	14254	1.31%	0.108%
45	10.481	1169	1173	1178	rVB6	3040	5700	0.52%	0.043%
46	10.545	1178	1184	1190	rBV3	10846	19419	1.79%	0.148%
47	10.598	1190	1193	1198	rVB5	4306	5907	0.54%	0.045%
48	10.727	1206	1215	1217	rBV6	2723	6448	0.59%	0.049%
49	10.780	1217	1224	1232	rBV	171451	326326	30.02%	2.483%
50	11.027	1260	1266	1270	rBV3	9840	17582	1.62%	0.134%
51	11.162	1281	1289	1303	rVB	236798	442893	40.74%	3.370%
52	11.291	1303	1311	1327	rVB	115005	245066	22.54%	1.865%
53	11.814	1398	1400	1414	rVB5	2822	7410	0.68%	0.056%
54	12.132	1447	1454	1461	rBV6	3760	9011	0.83%	0.069%
55	12.519	1515	1520	1527	rBV4	3901	6391	0.59%	0.049%
56	12.731	1549	1556	1561	rVB6	4117	9059	0.83%	0.069%
57	12.795	1561	1567	1574	rBV4	4999	10219	0.94%	0.078%
58	13.448	1674	1678	1684	rVB5	3782	6588	0.61%	0.050%
59	13.542	1689	1694	1698	rVB3	3373	6708	0.62%	0.051%
60	13.736	1722	1727	1732	rBV5	4484	7014	0.65%	0.053%
61	14.341	1821	1830	1834	rBV	335176	501140	46.10%	3.813%
62	14.388	1834	1838	1849	rVB	117274	181603	16.70%	1.382%
63	14.646	1875	1882	1896	rVB	361716	592764	54.53%	4.511%
64	14.799	1901	1908	1915	rVB3	8112	14092	1.30%	0.107%
65	14.952	1924	1934	1947	rVB2	365140	614736	56.55%	4.678%
66	15.157	1962	1969	1986	rBV3	71157	175856	16.18%	1.338%
67	15.351	1996	2002	2008	rVB8	7203	18852	1.73%	0.143%
68	15.416	2010	2013	2020	rVB5	4069	8086	0.74%	0.062%
69	15.751	2065	2070	2075	rVB3	13566	21721	2.00%	0.165%
70	15.945	2095	2103	2111	rBV	494824	782867	72.01%	5.957%
71	16.068	2118	2124	2131	rBV	127964	206755	19.02%	1.573%

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72	16.156	2133	2139	2141	rBV3	5125	7530	0.69%	0.057%
73	16.250	2152	2155	2159	rVB4	3285	5446	0.50%	0.041%
74	16.609	2212	2216	2225	rBV3	18337	35883	3.30%	0.273%
75	16.803	2245	2249	2258	rVB7	3856	9090	0.84%	0.069%
76	16.891	2258	2264	2269	rBV5	3008	7124	0.66%	0.054%
77	17.061	2291	2293	2299	rVB5	3835	7148	0.66%	0.054%
78	17.114	2299	2302	2306	rBV4	3984	6510	0.60%	0.050%
79	17.408	2347	2352	2356	rBV4	11682	16784	1.54%	0.128%
80	17.572	2376	2380	2386	rVB7	4775	10129	0.93%	0.077%
81	17.701	2396	2402	2413	rBV2	455530	743746	68.41%	5.660%
82	17.801	2413	2419	2428	rVV	552782	880982	81.04%	6.704%
83	17.878	2428	2432	2438	rVB8	6870	11759	1.08%	0.089%
84	18.142	2474	2477	2484	rVB7	9072	13043	1.20%	0.099%
85	18.289	2497	2502	2503	rBV5	3623	5361	0.49%	0.041%
86	18.318	2505	2507	2513	rVB7	6923	9506	0.87%	0.072%
87	18.424	2521	2525	2528	rBV5	4575	5426	0.50%	0.041%
88	18.548	2541	2546	2555	rBV	103497	176961	16.28%	1.347%
89	18.759	2578	2582	2586	rVB5	4915	9994	0.92%	0.076%
90	18.924	2606	2610	2614	rBV7	9667	17092	1.57%	0.130%
91	19.470	2697	2703	2704	rBV6	10654	19786	1.82%	0.151%
92	19.805	2757	2760	2770	rVB7	33186	51672	4.75%	0.393%
93	19.946	2780	2784	2794	rVB	166758	214839	19.76%	1.635%
94	20.081	2801	2807	2817	rBV	703459	1087126	100.00%	8.272%
95	20.281	2837	2841	2846	rBV2	67183	96782	8.90%	0.736%
96	20.992	2959	2962	2966	rVB2	72087	85390	7.85%	0.650%
97	21.603	3062	3066	3070	rBV5	52601	81955	7.54%	0.624%
98	22.038	3133	3140	3158	rVB	481379	904645	83.21%	6.884%
99	25.310	3687	3697	3715	rVB2	327636	1021689	93.98%	7.774%
100	25.557	3730	3739	3752	rVB2	287532	883638	81.28%	6.724%

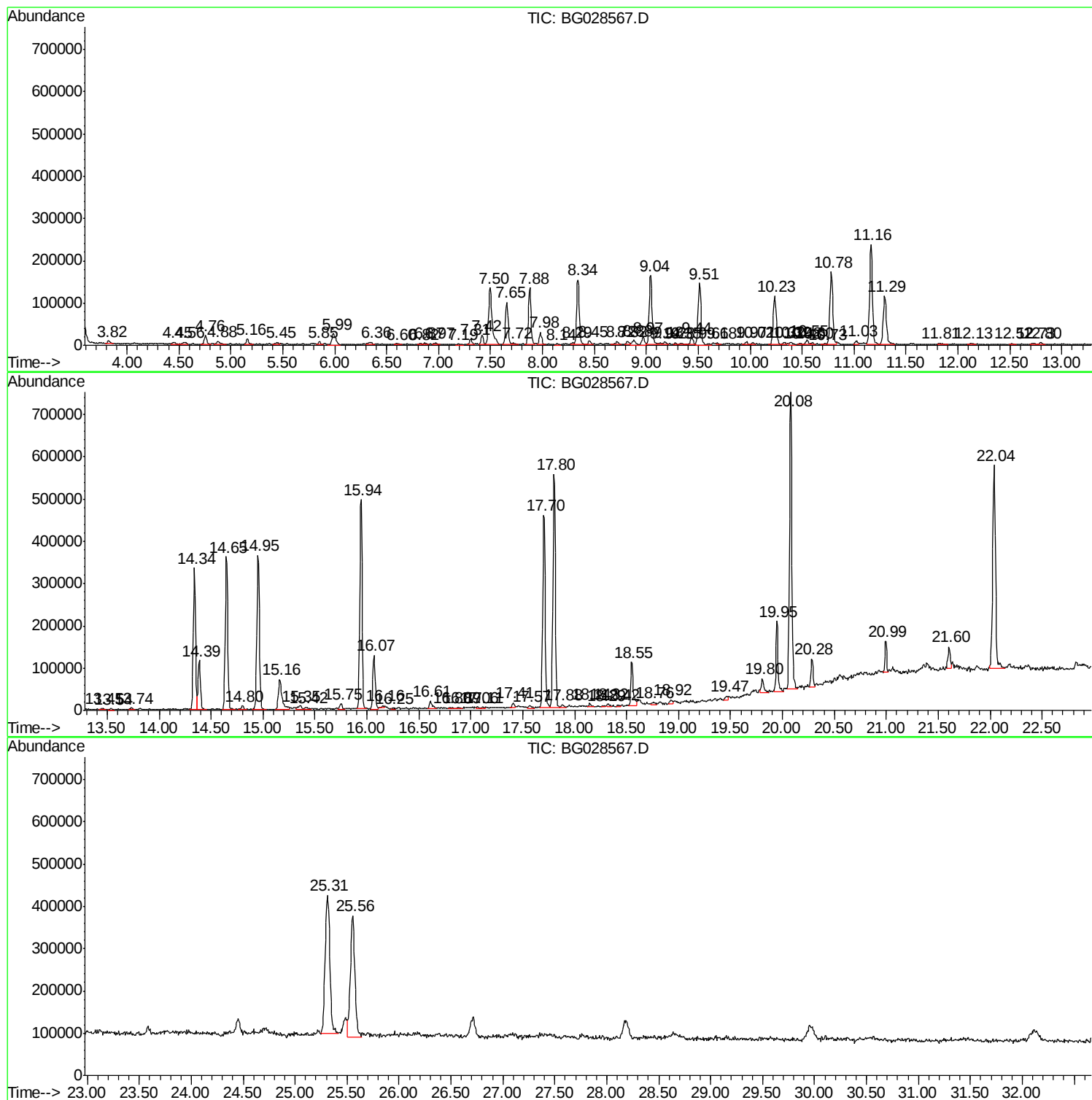
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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



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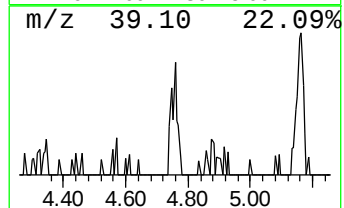
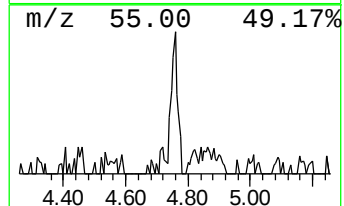
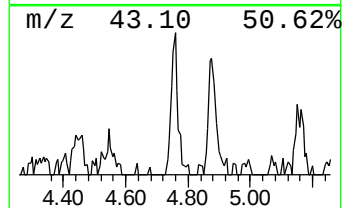
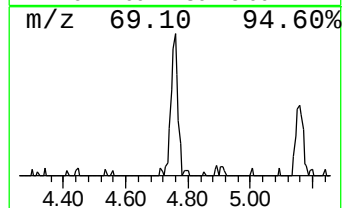
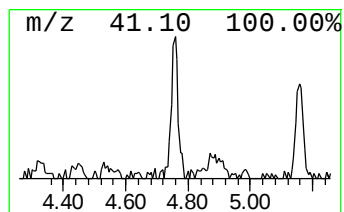
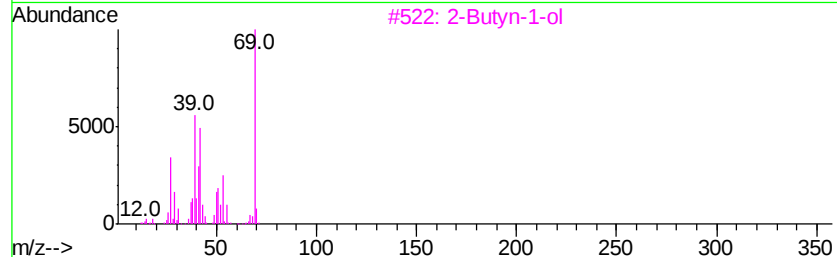
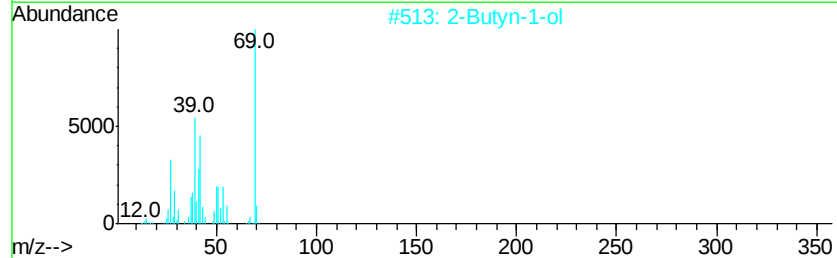
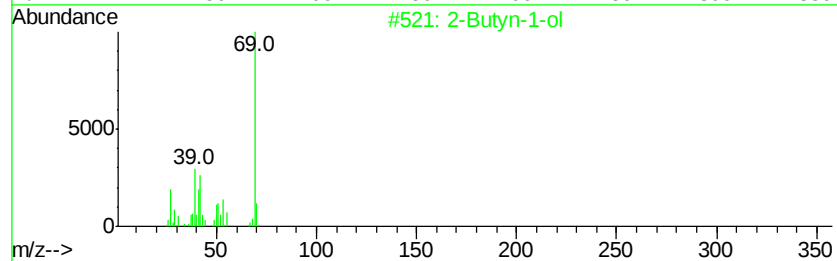
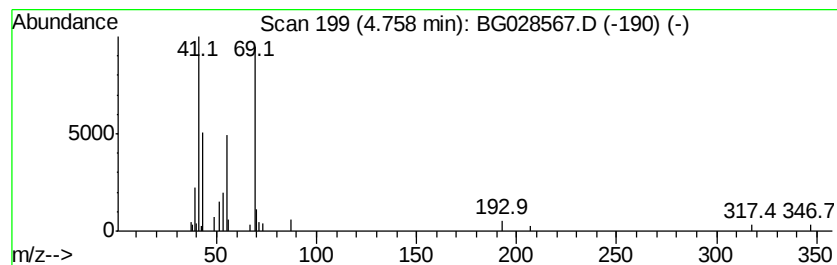
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Butyn-1-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.76	2.59 ng/ul	37448	1,4-Dichlorobenzene-d4	8.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Butyn-1-ol	70	C4H6O	000764-01-2	50
2		2-Butyn-1-ol	70	C4H6O	000764-01-2	40
3		2-Butyn-1-ol	70	C4H6O	000764-01-2	38
4		Methyl 2-hydroxydecanoate	202	C11H22O3	071271-24-4	38
5		2-Pentene, 4-bromo-	148	C5H9Br	001809-26-3	9



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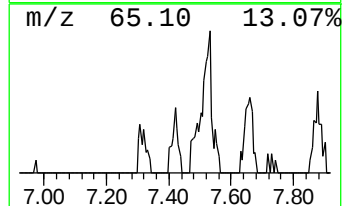
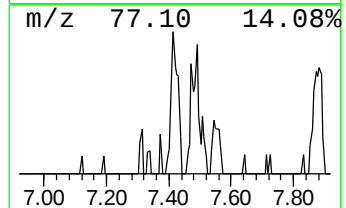
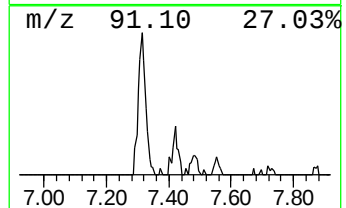
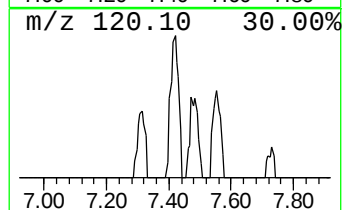
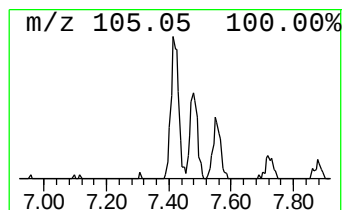
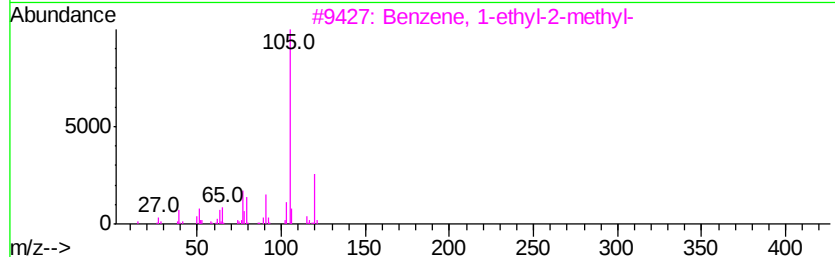
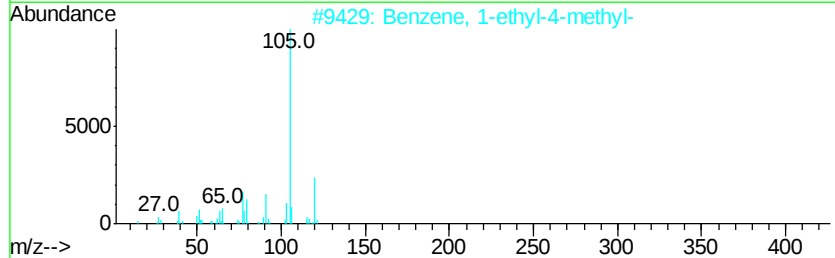
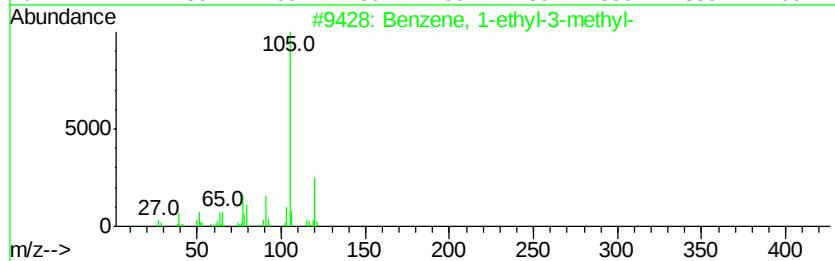
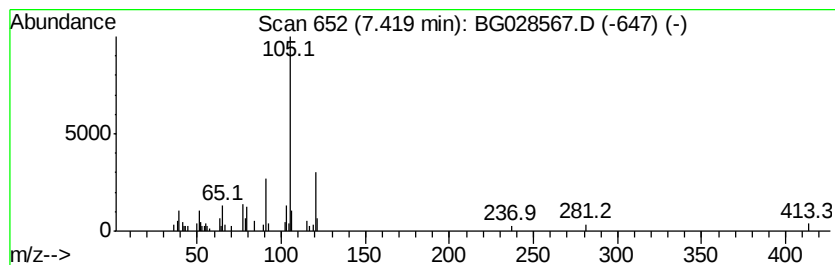
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethyl-3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.42	2.72 ng/ul	39428	1,4-Dichlorobenzene-d4	8.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	76
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	76
3		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	76
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	72
5		Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	72



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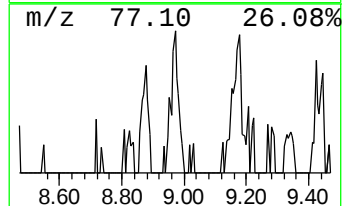
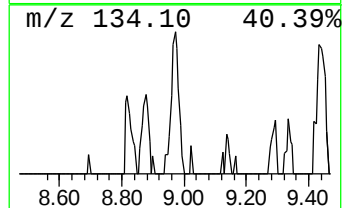
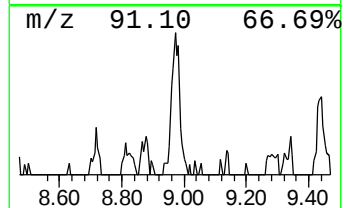
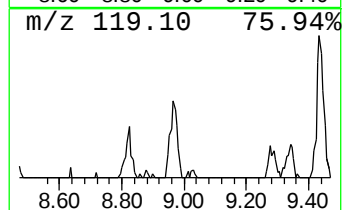
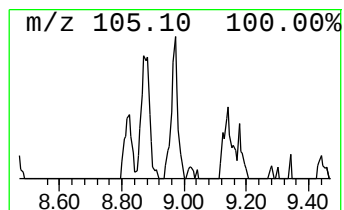
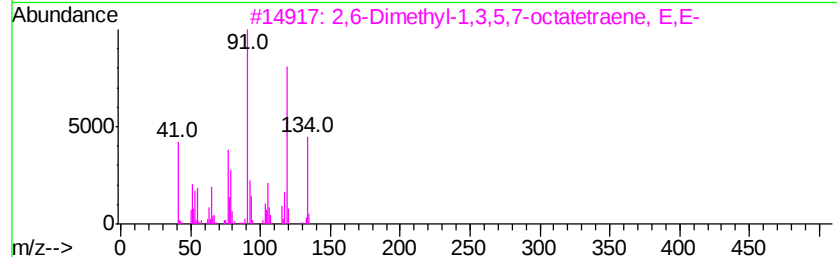
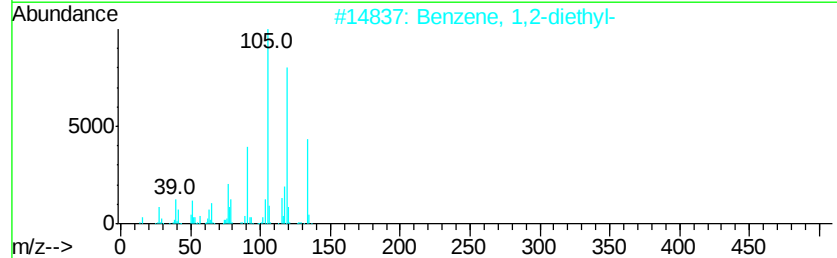
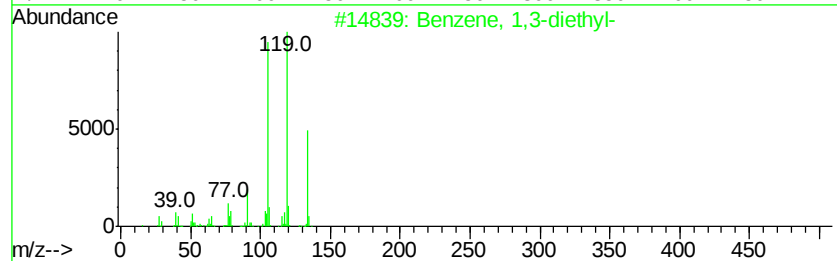
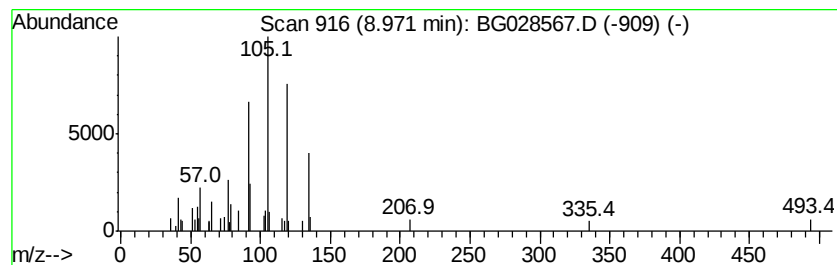
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Peak Number 5 Benzene, 1,3-diethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.97	2.36 ng/ul	34109	1,4-Dichlorobenzene-d4	8.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	64
2		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	64
3		2,6-Dimethyl-1,3,5,7-octatetraen...	134	C10H14	000460-01-5	64
4		1,3,8-p-Menthatriene	134	C10H14	018368-95-1	53
5		3a,6-Methano-3aH-indene, 2,3,4,5...	134	C10H14	098640-10-9	52



Data Path : Z:\HPCHEM1\BNA G\DATA\BG083017\
Data File : BG028567.D
Acq On : 31 Aug 2017 00:17
Operator : SJ/JU
Sample : I4917-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B0B51

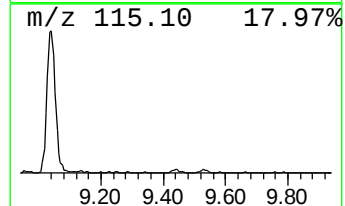
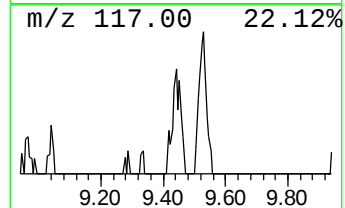
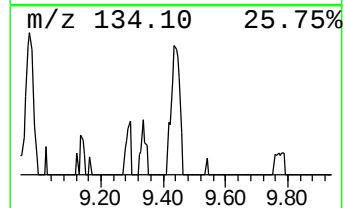
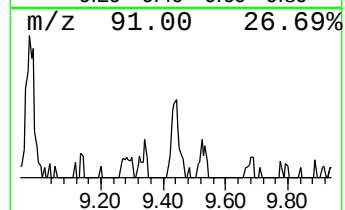
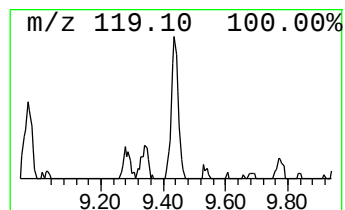
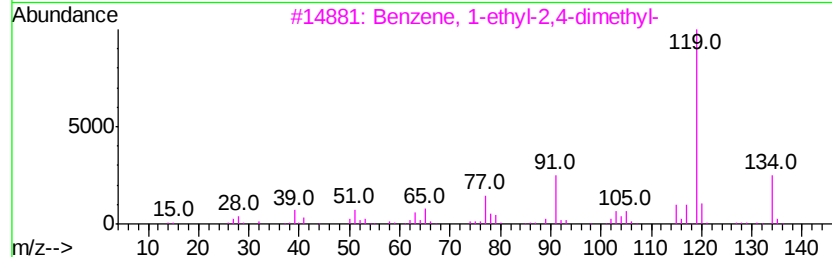
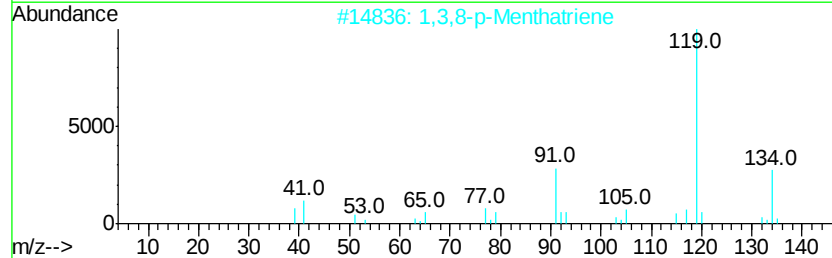
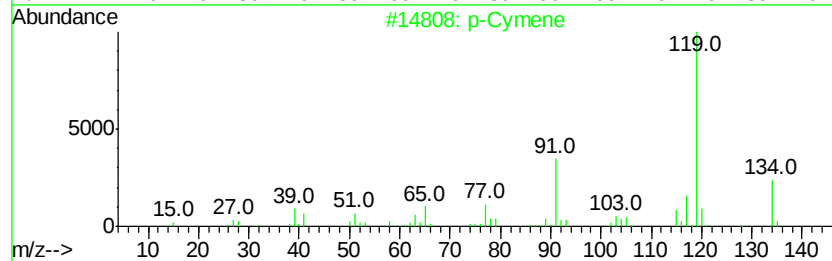
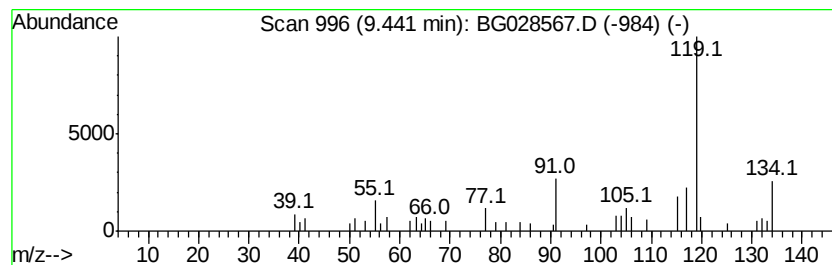
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 p-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.44	2.43 ng/ul	35132	1,4-Dichlorobenzene-d4	8.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	p-Cymene	134	C10H14	000099-87-6	72
2		1,3,8-p-Menthatriene	134	C10H14	018368-95-1	59
3		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	59
4		o-Cymene	134	C10H14	000527-84-4	59
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	58



Data Path : Z:\HPCHEM1\BNA G\DATA\BG083017\
Data File : BG028567.D
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Operator : SJ/JU
Sample : I4917-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B0B51

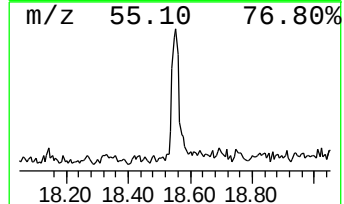
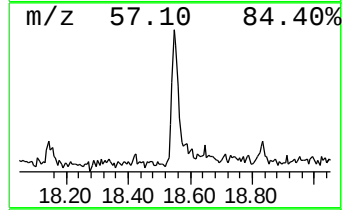
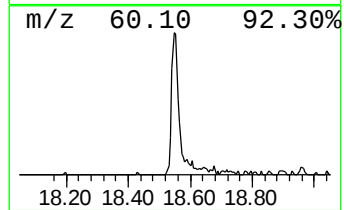
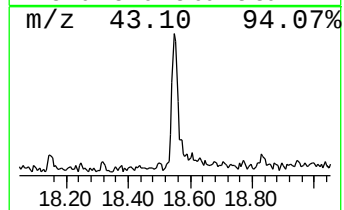
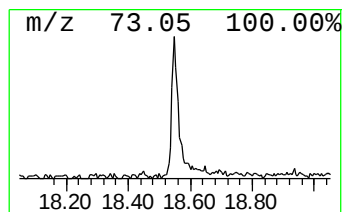
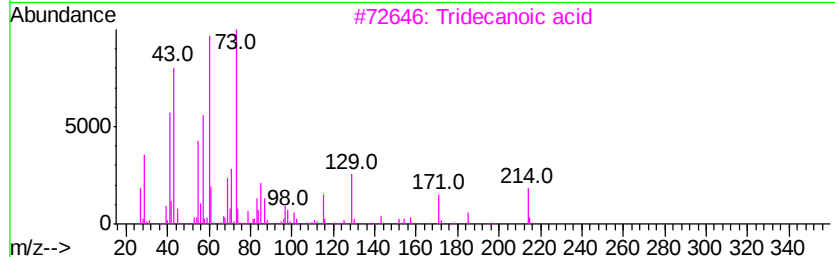
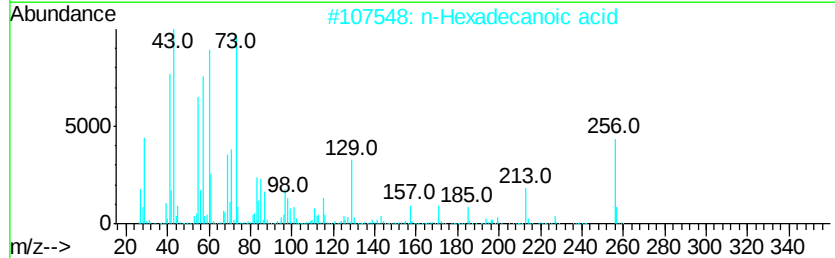
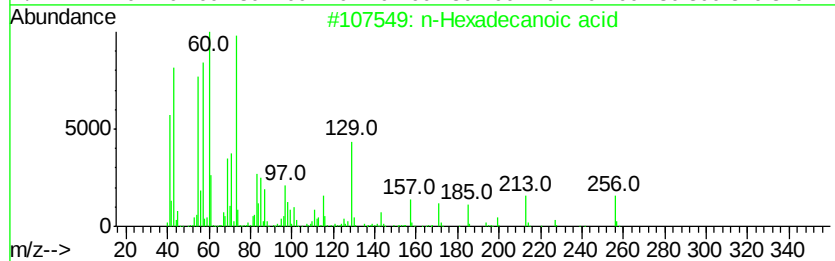
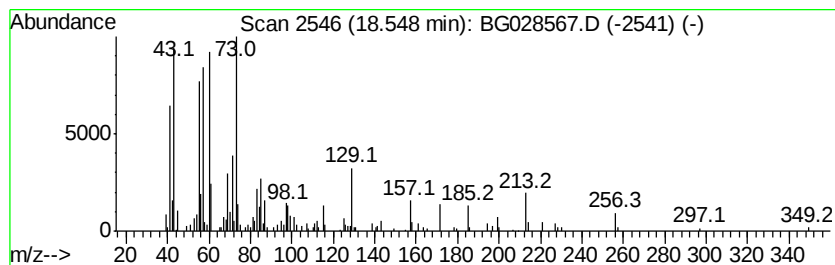
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.55	4.76 ng/ul	176961	Phenanthrene-d10	17.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
3		Tridecanoic acid	214	C13H26O2	000638-53-9	91
4		Tridecanoic acid	214	C13H26O2	000638-53-9	89
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	74



Data Path : Z:\HPCHEM1\BNA G\DATA\BG083017\
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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B0B51

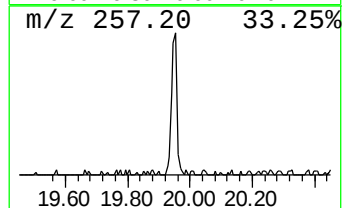
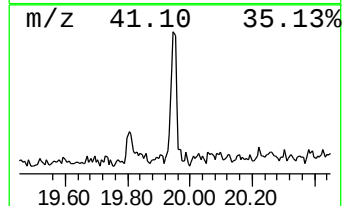
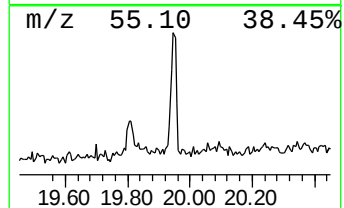
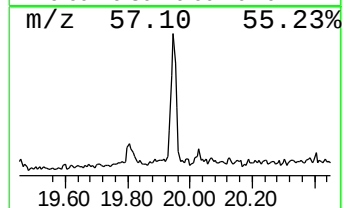
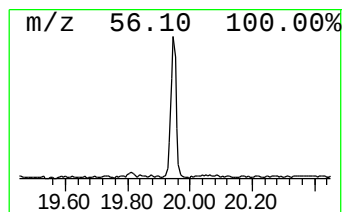
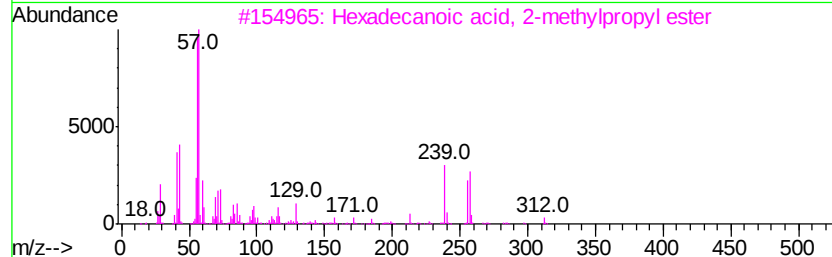
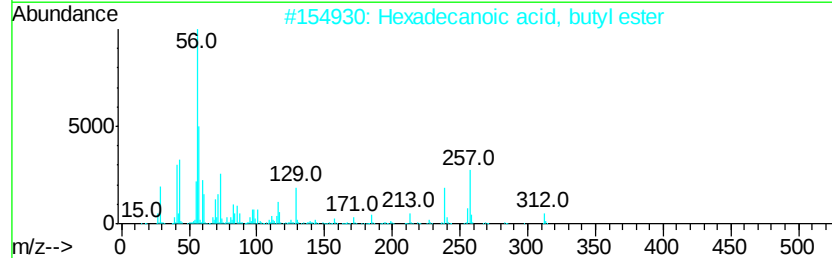
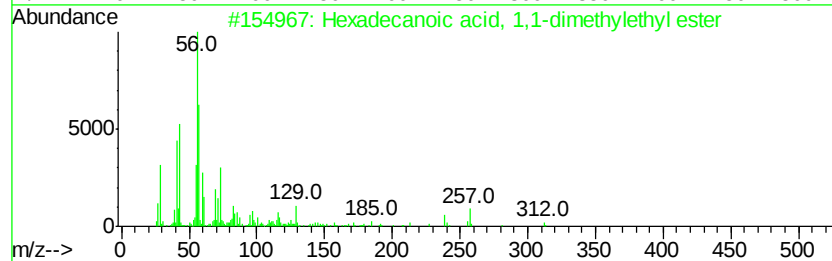
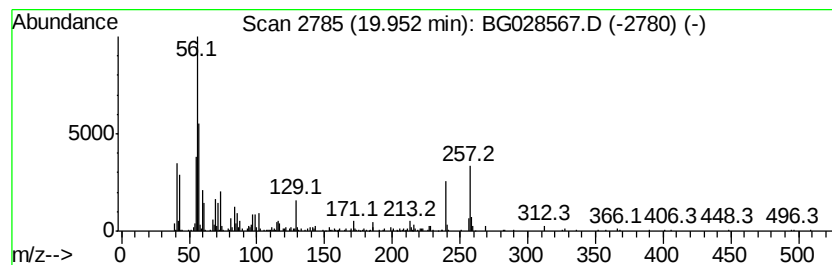
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 Hexadecanoic acid, 1,1-dime... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.95	4.75 ng/ul	214839	Chrysene-d12	22.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	86
2		Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	86
3		Hexadecanoic acid, 2-methylpropyl...	312	C20H40O2	000110-34-9	64
4		Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	55
5		1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	37



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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
B0B51

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Butyn-1-ol	4.76	2.6	ng/ul	37448	1	8.34	289607	20.0
Benzene, 1-ethyl-...	7.42	2.7	ng/ul	39428	1	8.34	289607	20.0
Benzene, 1,3-diet...	8.97	2.4	ng/ul	34109	1	8.34	289607	20.0
p-Cymene	9.44	2.4	ng/ul	35132	1	8.34	289607	20.0
n-Hexadecanoic acid	18.55	4.8	ng/ul	176961	4	17.70	743746	20.0
Hexadecanoic acid...	19.95	4.8	ng/ul	214839	5	22.04	904645	20.0