

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081519\
 Data File : BG042234.D
 Acq On : 15 Aug 2019 18:17
 Operator : HP/JU
 Sample : K4183-18
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AG8

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 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: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.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	3.141	5	9	24	rVB	443755	750769	35.56%	3.827%
2	3.411	51	55	66	rVB3	7949	14241	0.67%	0.073%
3	4.075	163	168	170	rBV3	2958	4387	0.21%	0.022%
4	4.163	180	183	184	rBV2	2001	2389	0.11%	0.012%
5	5.097	336	342	346	rBV3	2311	3949	0.19%	0.020%
6	5.996	491	495	497	rBV2	1991	2234	0.11%	0.011%
7	6.102	510	513	518	rBV3	1361	2407	0.11%	0.012%
8	6.402	555	564	567	rVB5	1282	2649	0.13%	0.014%
9	6.701	609	615	631	rVB	200272	347349	16.45%	1.771%
10	7.007	662	667	673	rBV3	9308	15815	0.75%	0.081%
11	7.183	690	697	713	rBV2	134535	290173	13.74%	1.479%
12	7.342	716	724	737	rBV	105330	193153	9.15%	0.985%
13	7.465	740	745	753	rVV5	7320	15049	0.71%	0.077%
14	7.553	753	760	772	rVV	154012	289339	13.70%	1.475%
15	8.023	833	840	849	rBV	199224	351585	16.65%	1.792%
16	8.740	954	962	970	rBV	179203	346628	16.42%	1.767%
17	9.193	1031	1039	1051	rBV	138683	258639	12.25%	1.318%
18	9.627	1108	1113	1120	rVB2	4255	6178	0.29%	0.031%
19	9.921	1156	1163	1172	rBV	119831	224813	10.65%	1.146%
20	10.473	1249	1257	1272	rBV	193454	399786	18.93%	2.038%
21	10.849	1314	1321	1334	rBV	283232	513153	24.30%	2.616%
22	10.985	1336	1344	1354	rVV	150358	315045	14.92%	1.606%
23	11.049	1354	1355	1360	rVB3	2064	2543	0.12%	0.013%
24	11.325	1398	1402	1409	rVB2	3960	5930	0.28%	0.030%
25	12.448	1587	1593	1601	rBV2	3045	6629	0.31%	0.034%
26	13.059	1693	1697	1703	rVB3	7311	11024	0.52%	0.056%
27	13.435	1756	1761	1769	rVB5	4842	8554	0.41%	0.044%
28	13.570	1778	1784	1790	rVB	66364	97906	4.64%	0.499%
29	14.016	1853	1860	1866	rVV	792962	1158408	54.86%	5.905%
30	14.087	1866	1872	1877	rVV	386381	656266	31.08%	3.345%
31	14.128	1877	1879	1890	rVB	100587	137578	6.52%	0.701%
32	14.381	1913	1922	1932	rBV2	477762	793526	37.58%	4.045%
33	14.451	1932	1934	1940	rVV3	3549	5363	0.25%	0.027%
34	14.528	1942	1947	1952	rVV3	16104	26170	1.24%	0.133%

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Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
 Title : SVOA CALIBRATION

35	14.592	1952	1958	1960	rVV6	7668	16994	0.80%	0.087%
36	14.651	1960	1968	1971	rVV	438517	722744	34.23%	3.684%
37	14.686	1971	1974	1982	rVV	461079	686254	32.50%	3.498%
38	14.763	1982	1987	1999	rVB7	34228	87781	4.16%	0.447%
39	14.892	2002	2009	2014	rBV2	58518	125672	5.95%	0.641%
40	14.945	2014	2018	2026	rVV2	66659	161139	7.63%	0.821%
41	15.015	2026	2030	2038	rVV2	25271	45328	2.15%	0.231%
42	15.109	2040	2046	2053	rVB6	6547	10003	0.47%	0.051%
43	15.426	2095	2100	2105	rVB	7955	11773	0.56%	0.060%
44	15.532	2114	2118	2121	rBV3	2345	3430	0.16%	0.017%
45	15.567	2121	2124	2128	rBV2	4275	4718	0.22%	0.024%
46	15.626	2128	2134	2137	rBV	40937	59450	2.82%	0.303%
47	15.679	2137	2143	2153	rVB	646247	973338	46.10%	4.962%
48	15.802	2159	2164	2173	rBV2	52819	92094	4.36%	0.469%
49	15.873	2174	2176	2180	rVB2	3136	3224	0.15%	0.016%
50	15.926	2180	2185	2190	rBV3	8246	11573	0.55%	0.059%
51	16.020	2195	2201	2205	rVV4	32769	55322	2.62%	0.282%
52	16.067	2205	2209	2214	rVV2	38214	57028	2.70%	0.291%
53	16.161	2217	2225	2236	rVB	95529	154054	7.30%	0.785%
54	16.255	2236	2241	2247	rBV5	5659	10547	0.50%	0.054%
55	16.372	2257	2261	2271	rBV2	13212	27050	1.28%	0.138%
56	16.466	2274	2277	2281	rVB4	3248	4144	0.20%	0.021%
57	16.537	2285	2289	2298	rBV3	13073	26310	1.25%	0.134%
58	16.672	2307	2312	2319	rVB4	5015	8237	0.39%	0.042%
59	16.831	2335	2339	2342	rVB2	1664	2338	0.11%	0.012%
60	17.183	2398	2399	2403	rVB4	2182	2388	0.11%	0.012%
61	17.442	2436	2443	2452	rBV2	536036	858256	40.65%	4.375%
62	17.536	2452	2459	2471	rVV2	681448	1036713	49.10%	5.285%
63	17.630	2471	2475	2478	rVB3	2353	3008	0.14%	0.015%
64	17.706	2483	2488	2493	rBV5	5579	11606	0.55%	0.059%
65	18.329	2590	2594	2600	rBV5	4485	8801	0.42%	0.045%
66	19.075	2719	2721	2723	rBV3	2490	2460	0.12%	0.013%
67	19.381	2770	2773	2775	rBV4	2618	2758	0.13%	0.014%
68	19.457	2783	2786	2792	rVB2	10036	16631	0.79%	0.085%
69	19.716	2826	2830	2834	rBV2	9425	11363	0.54%	0.058%
70	19.827	2843	2849	2856	rBV	871397	1218908	57.73%	6.214%
71	20.191	2908	2911	2912	rBV2	12680	13881	0.66%	0.071%

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Smoothing : OFF Filtering: 5
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72	20.356	2934	2939	2942	rBV	66154	95145	4.51%	0.485%
73	21.008	3046	3050	3057	rVB	58436	91721	4.34%	0.468%
74	21.179	3073	3079	3089	rVB	1627476	2111400	100.00%	10.763%
75	21.372	3107	3112	3127	rBV5	72712	202507	9.59%	1.032%
76	21.725	3166	3172	3181	rVB2	565738	960491	45.49%	4.896%
77	24.768	3681	3690	3702	rVB	441079	1244328	58.93%	6.343%
78	24.998	3720	3729	3744	rVB2	349744	1103915	52.28%	5.627%

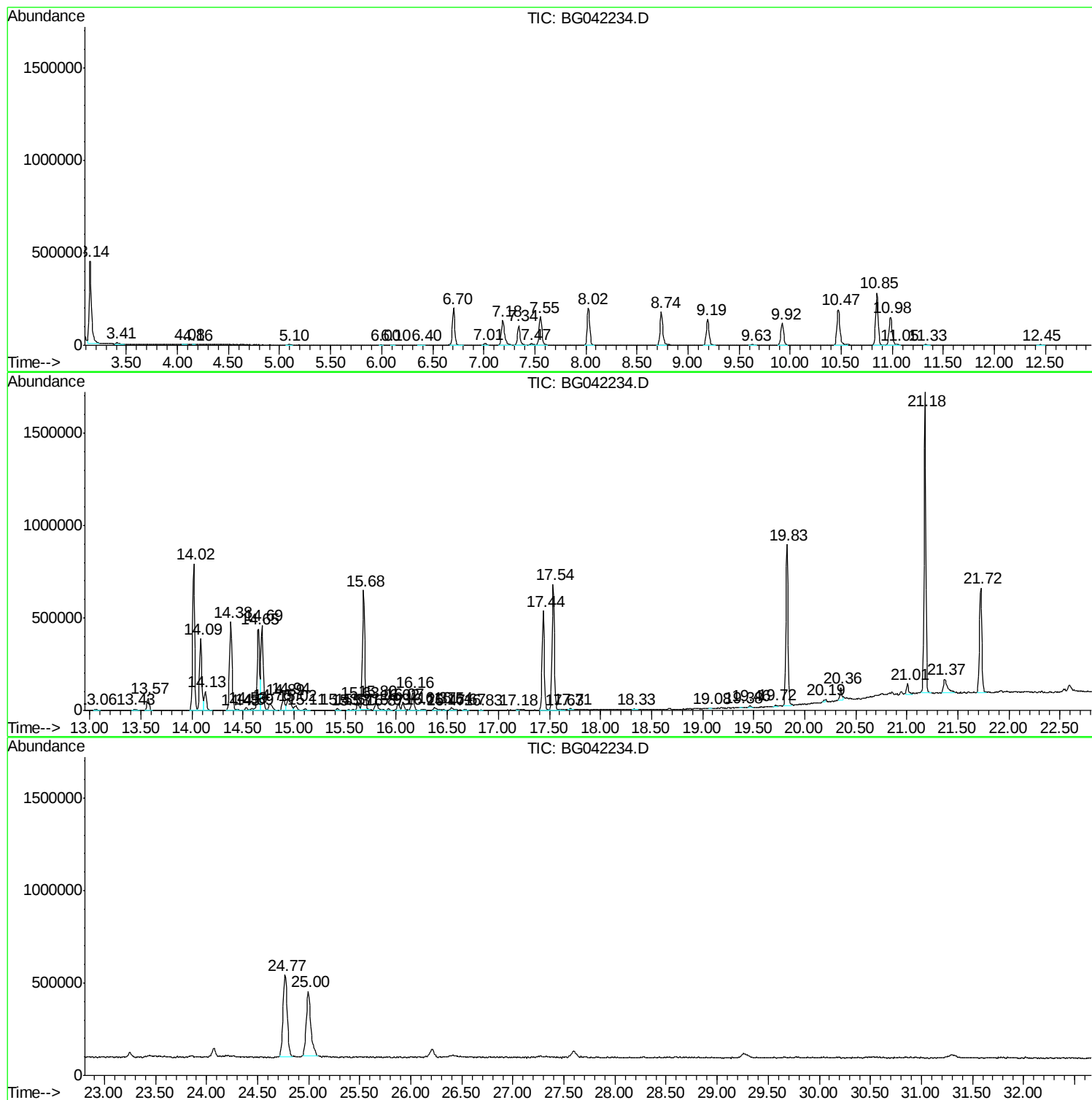
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Instrument :
BNA_G
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
Quant Title : SVOA CALIBRATION

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



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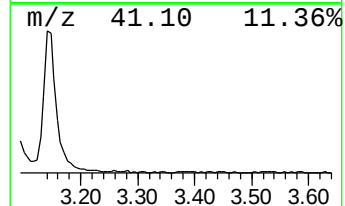
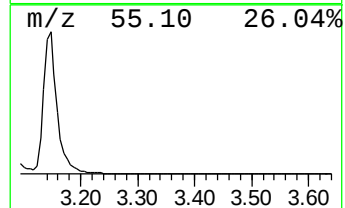
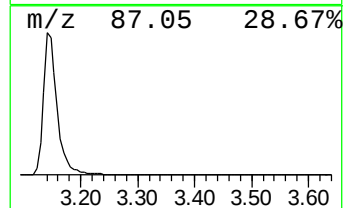
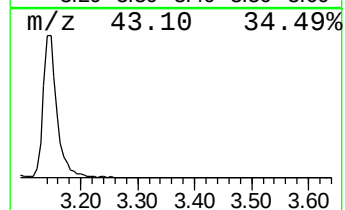
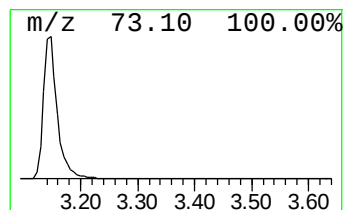
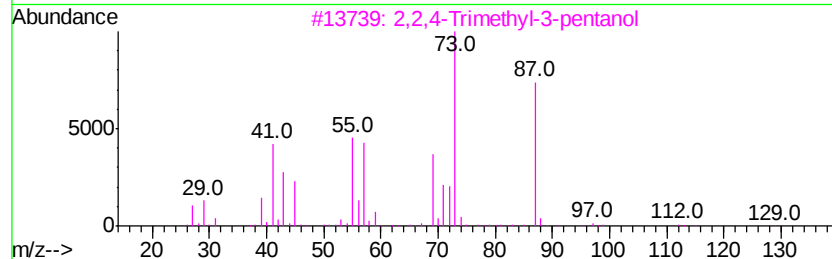
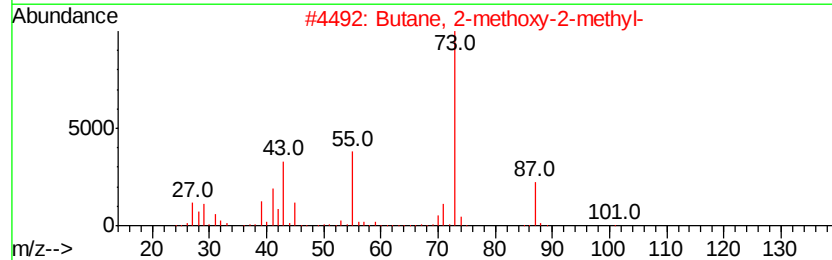
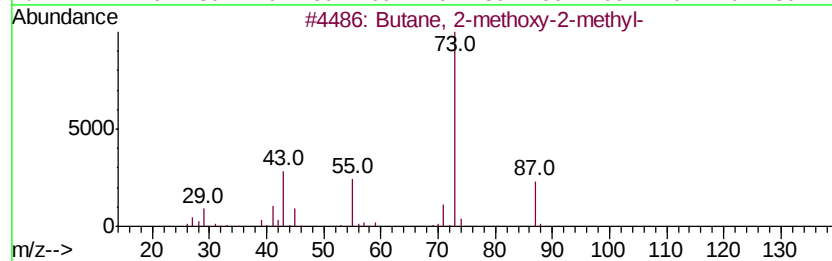
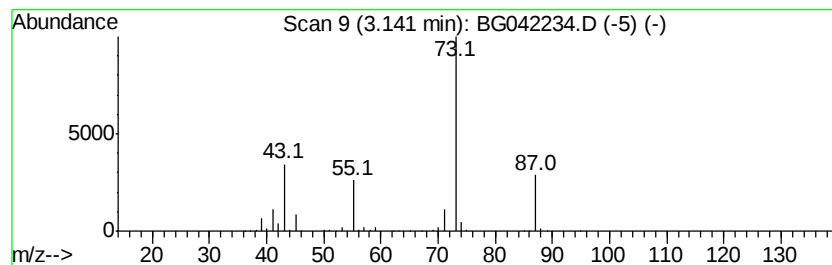
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.14	42.71 ng/ul	750769	1,4-Dichlorobenzene-d4	8.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	25
5		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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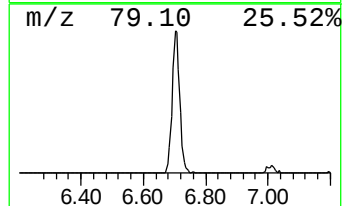
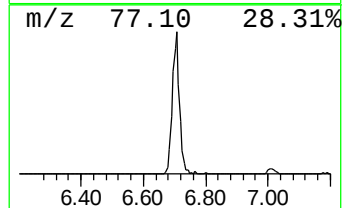
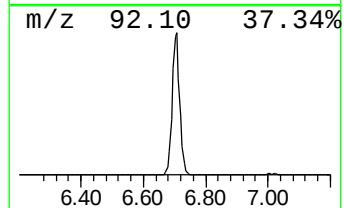
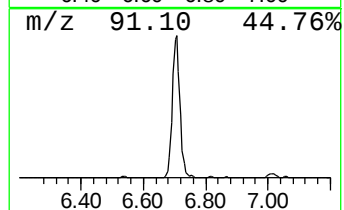
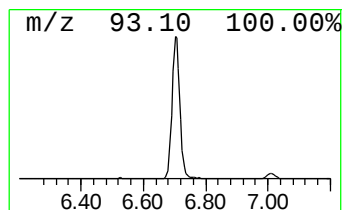
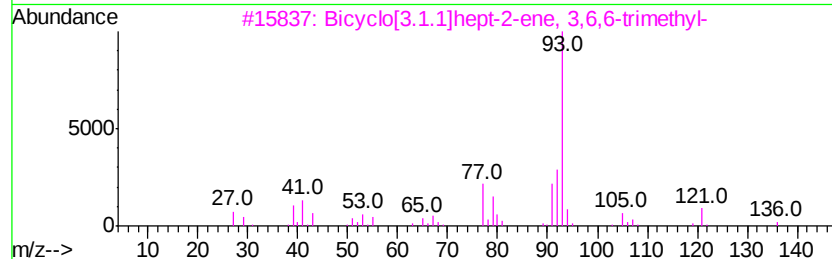
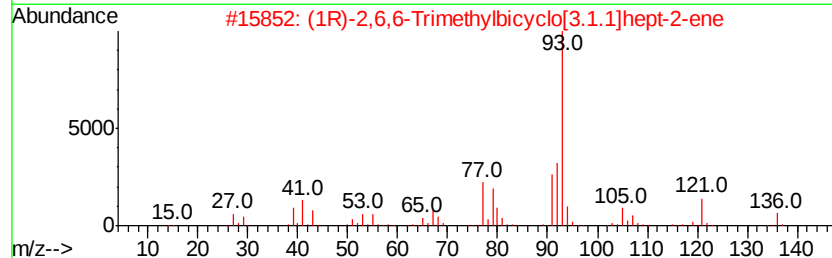
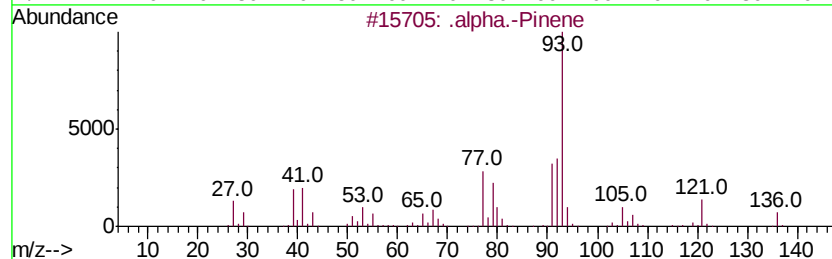
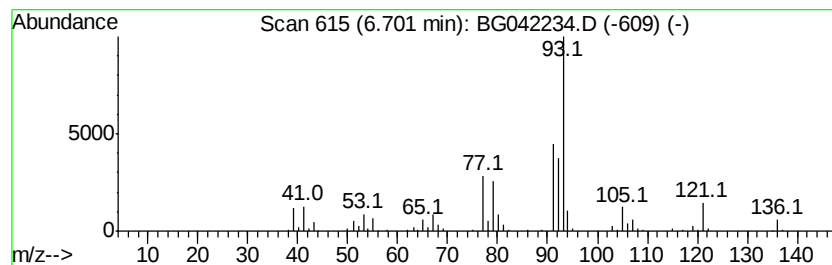
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 .alpha.-Pinene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.70	19.76 ng/ul	347349	1,4-Dichlorobenzene-d4	8.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.alpha.-Pinene	136	C10H16	000080-56-8	96
2			(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	95
3			Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	95
4			(1S)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-26-4	94
5			trans-.beta.-Ocimene	136	C10H16	003779-61-1	91



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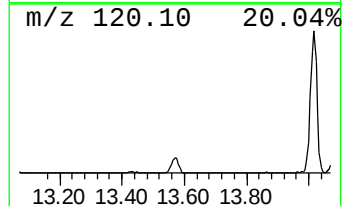
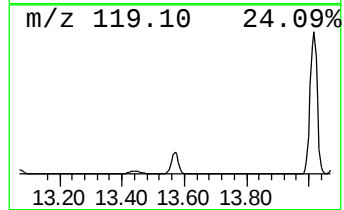
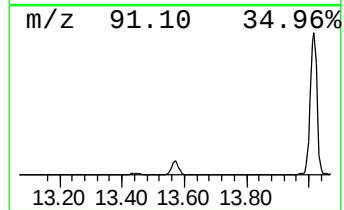
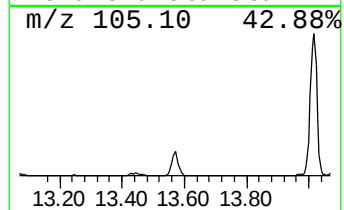
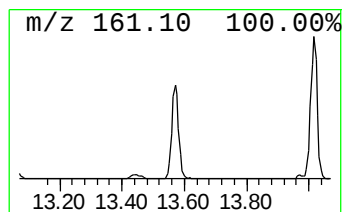
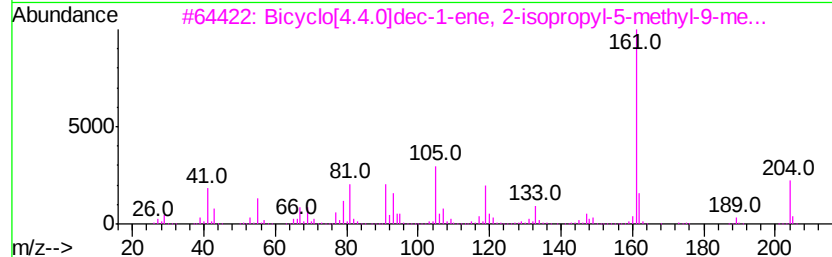
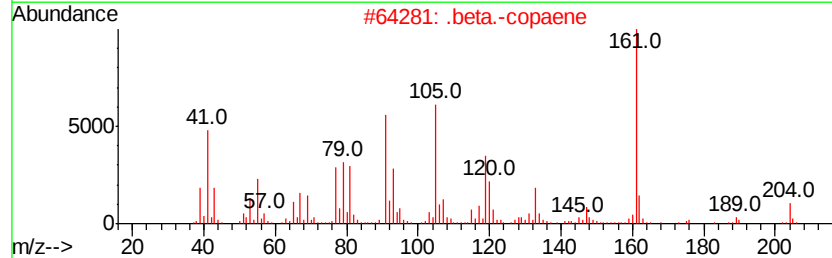
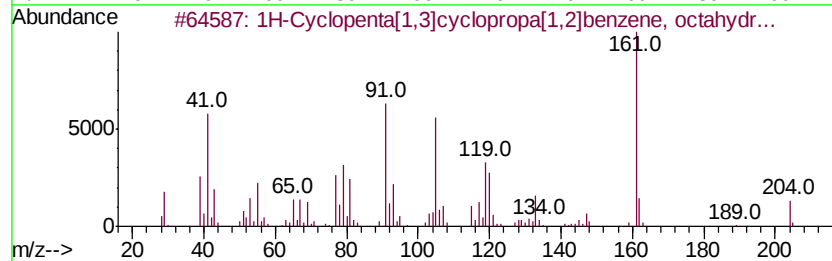
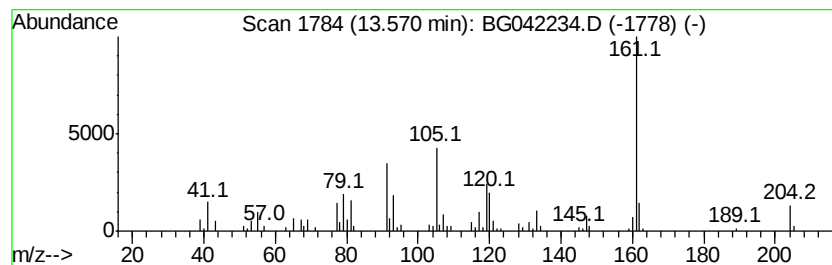
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1H-Cyclopenta[1,3]cycloprop... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.57	2.85 ng/ul	97906	Acenaphthene-d10	14.69
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Cyclopenta[1,3]cyclopropa[1,2...	204 C15H24	013744-15-5 94
2		.beta.-copaene	204 C15H24	1000374-18-9 94
3		Bicyclo[4.4.0]dec-1-ene, 2-isopr...	204 C15H24	150320-52-8 93
4		(+)-epi-Bicyclosesquiphellandrene	204 C15H24	054274-73-6 90
5		1H-Cyclopenta[1,3]cyclopropa[1,2...	204 C15H24	013744-15-5 89



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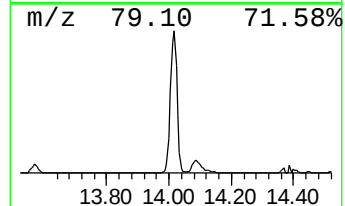
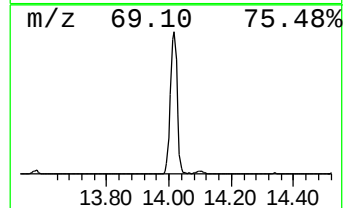
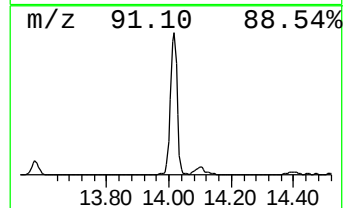
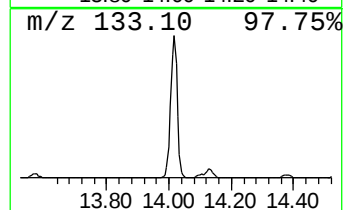
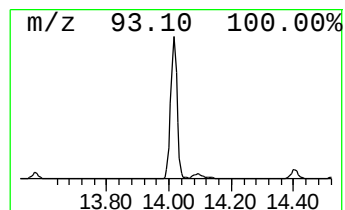
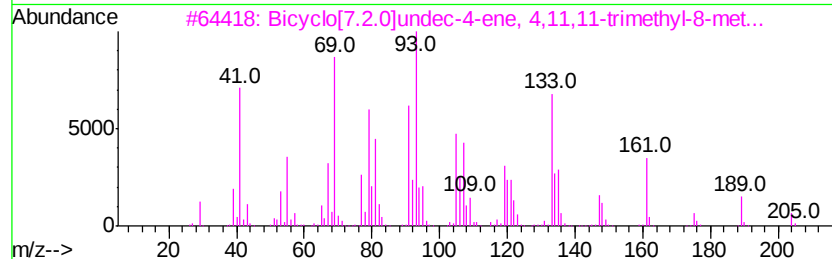
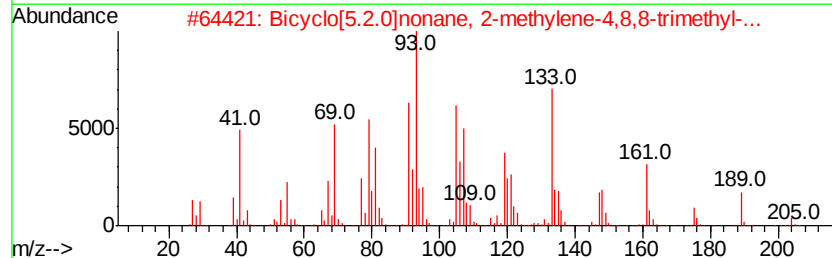
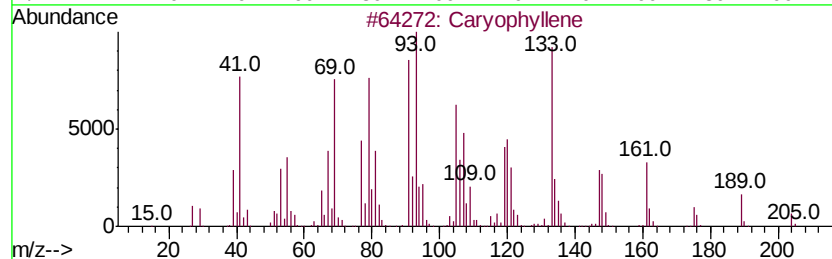
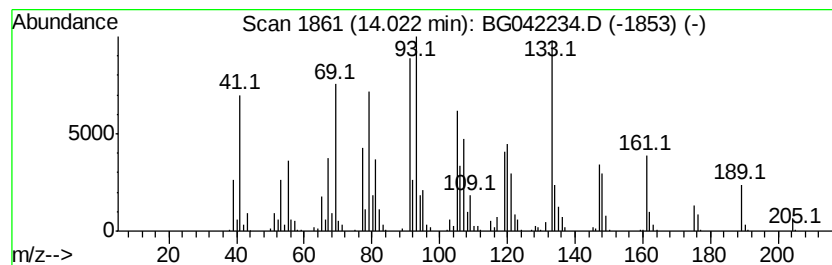
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Caryophyllene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.02	33.76 ng/ul	1158410	Acenaphthene-d10	14.69
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Caryophyllene	204 C15H24	000087-44-5 99
2		Bicyclo[5.2.0]nonane, 2-methylen...	204 C15H24	242794-76-9 94
3		Bicyclo[7.2.0]undec-4-ene, 4,11,...	204 C15H24	013877-93-5 93
4		Bicyclo[7.2.0]undec-4-ene, 4,11,...	204 C15H24	000118-65-0 91
5		Bicyclo[7.2.0]undec-4-ene, 4,11,...	204 C15H24	013877-93-5 89



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081519\
Data File : BG042234.D
Acq On : 15 Aug 2019 18:17
Operator : HP/JU
Sample : K4183-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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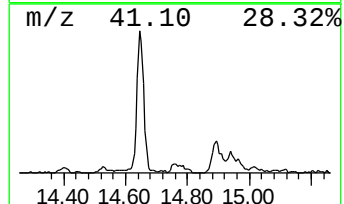
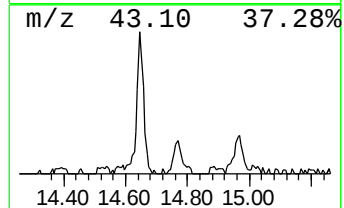
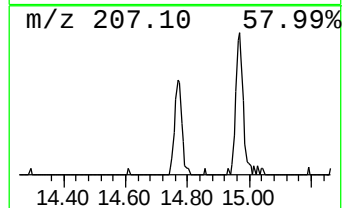
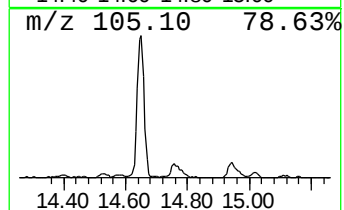
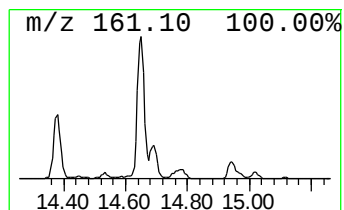
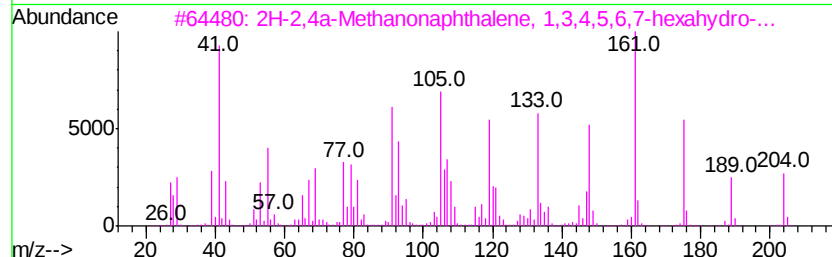
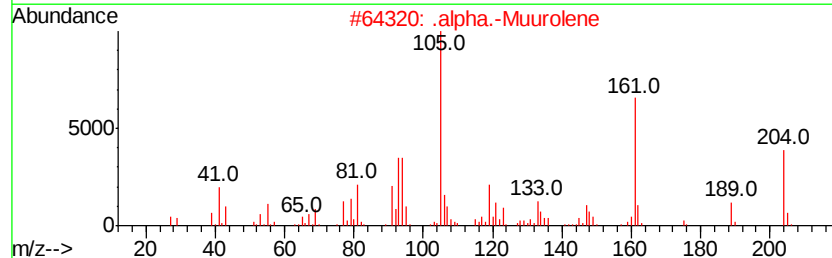
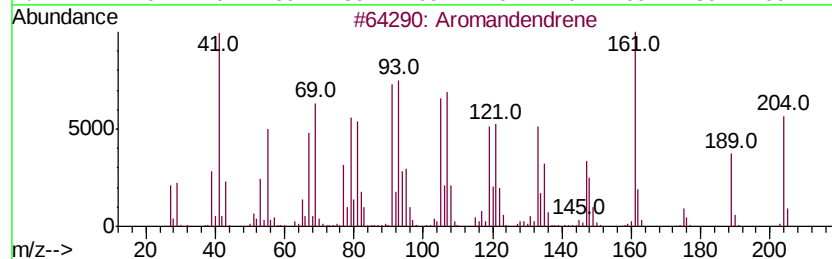
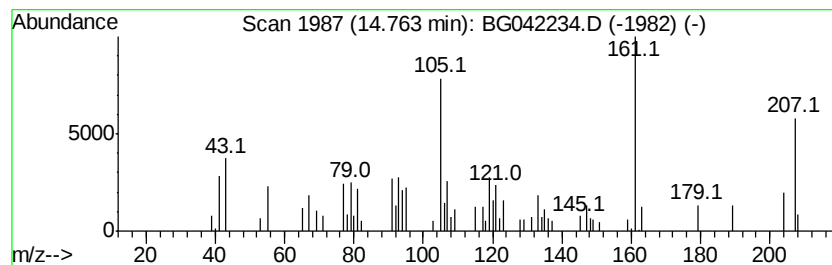
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Aromandendrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	2.56 ng/ul	87781	Acenaphthene-d10	14.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Aromandendrene	204	C15H24	000489-39-4	74
2		.alpha.-Muurolene	204	C15H24	031983-22-9	64
3		2H-2,4a-Methanonaphthalene, 1,3,...	204	C15H24	001135-66-6	56
4		1H-Cyclopropafalnaphthalene, 1a,...	204	C15H24	017334-55-3	55
5		Naphthalene, 1,2,3,4,4a,5,6,8a-o...	204	C15H24	039029-41-9	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081519\
Data File : BG042234.D
Acq On : 15 Aug 2019 18:17
Operator : HP/JU
Sample : K4183-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

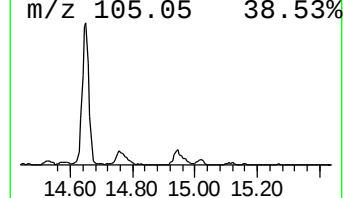
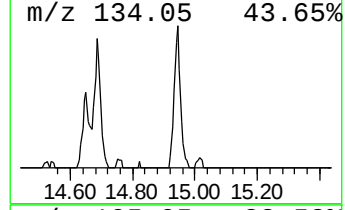
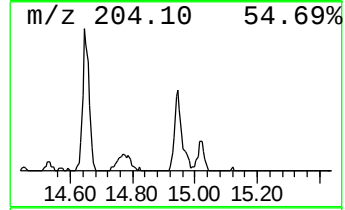
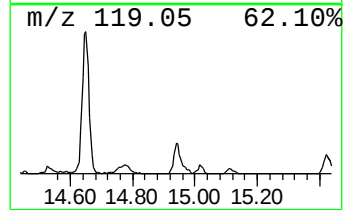
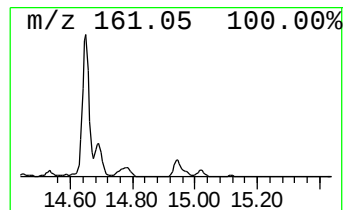
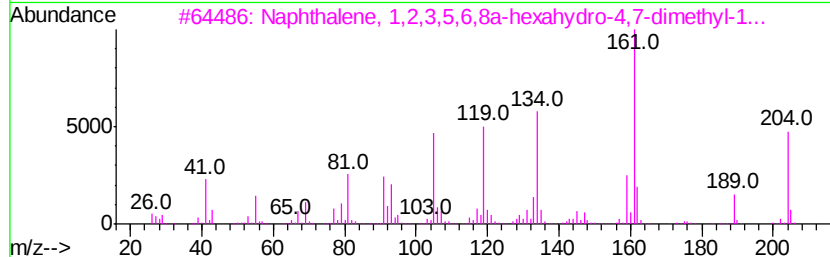
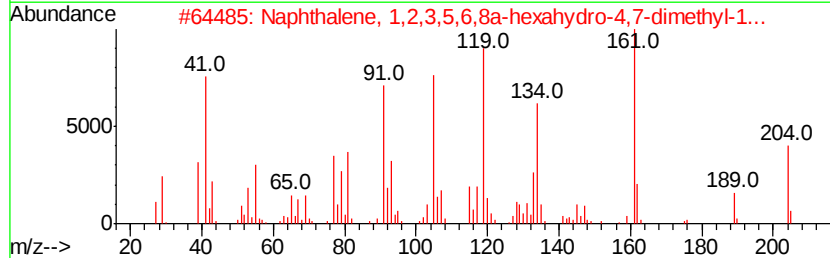
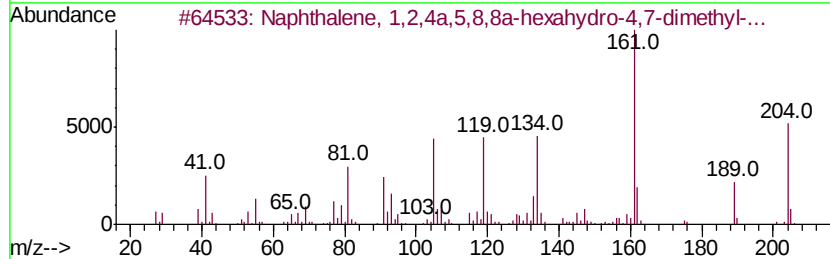
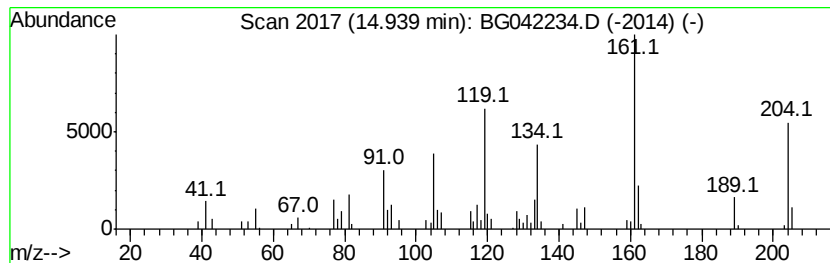
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Naphthalene, 1,2,4a,5,8,8a-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.94	4.70 ng/ul	161139	Acenaphthene-d10	14.69
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Naphthalene, 1,2,4a,5,8,8a-hexah...	204 C15H24	000523-47-7	96
2	Naphthalene, 1,2,3,5,6,8a-hexahy...	204 C15H24	000483-76-1	95
3	Naphthalene, 1,2,3,5,6,8a-hexahv...	204 C15H24	000483-76-1	94
4	Naphthalene, 1,2,3,5,6,8a-hexahy...	204 C15H24	000483-76-1	94
5	Naphthalene, 1,2,3,5,6,8a-hexahy...	204 C15H24	000483-76-1	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081519\
Data File : BG042234.D
Acq On : 15 Aug 2019 18:17
Operator : HP/JU
Sample : K4183-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

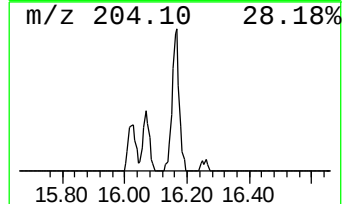
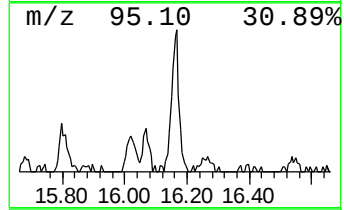
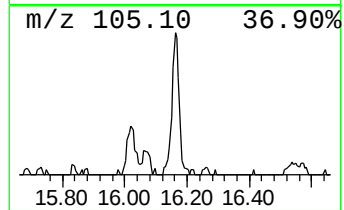
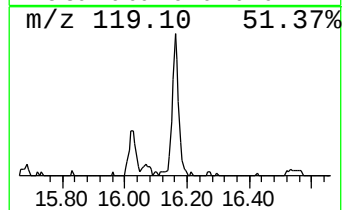
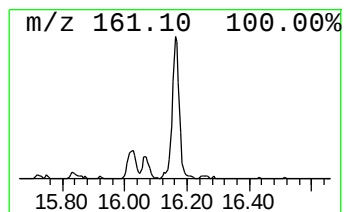
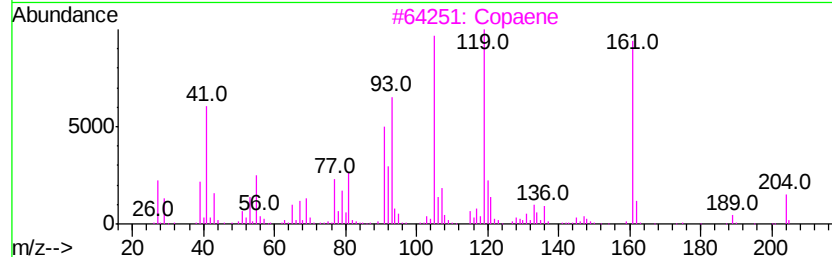
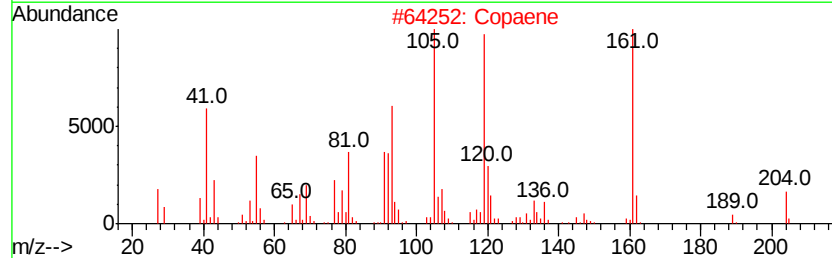
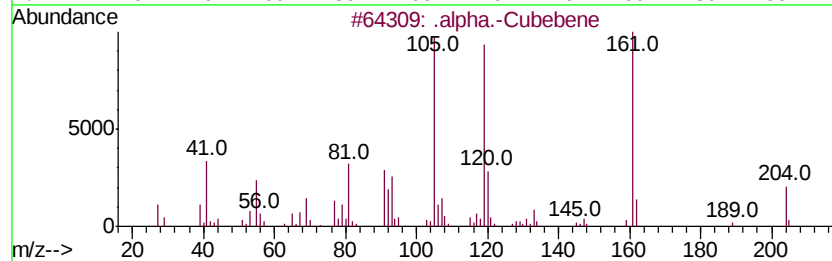
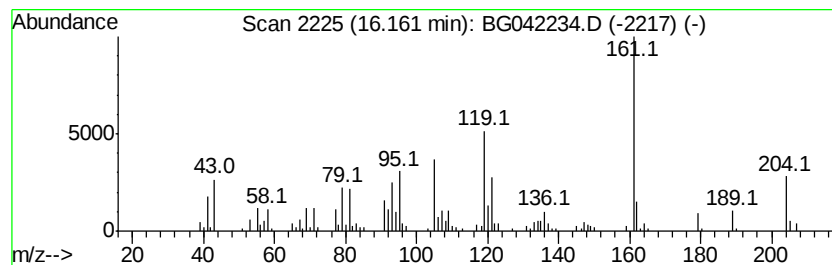
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 .alpha.-Cubebene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
16.16	3.59 ng/ul	154054	Phenanthrene-d10		17.44
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Cubebene	204	C15H24	017699-14-8	95
2	Copaene	204	C15H24	003856-25-5	93
3	Copaene	204	C15H24	003856-25-5	90
4	.gamma.-Muurolene	204	C15H24	030021-74-0	90
5	.alfa.-Copaene	204	C15H24	1000360-33-0	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081519\
Data File : BG042234.D
Acq On : 15 Aug 2019 18:17
Operator : HP/JU
Sample : K4183-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AG8

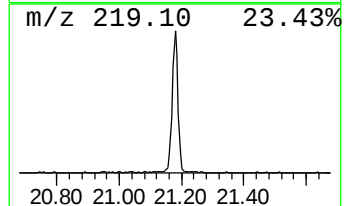
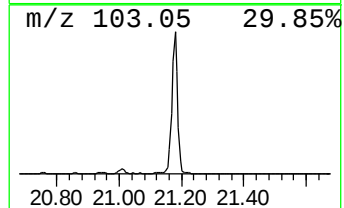
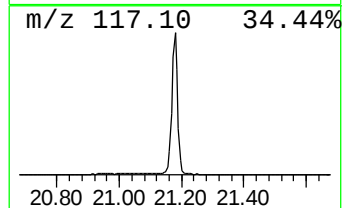
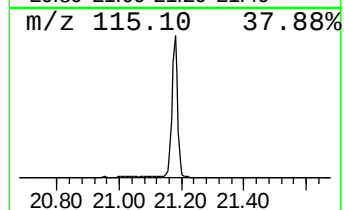
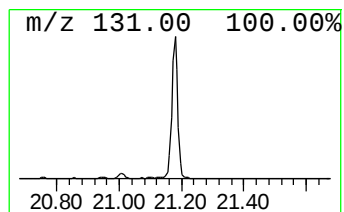
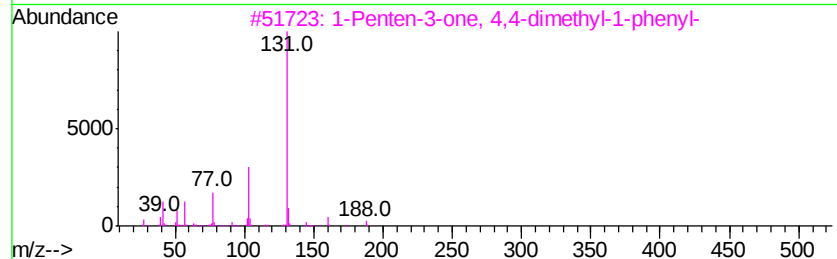
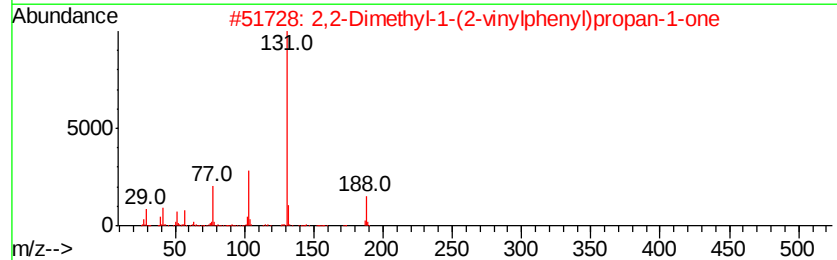
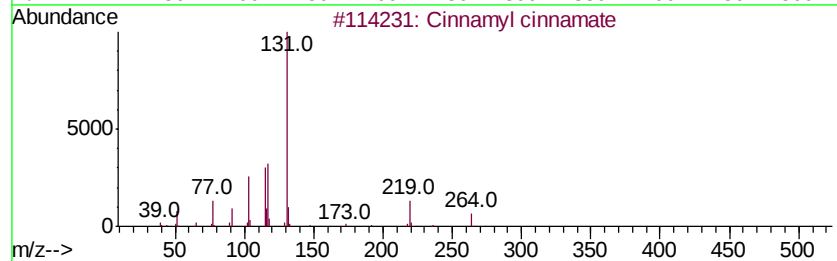
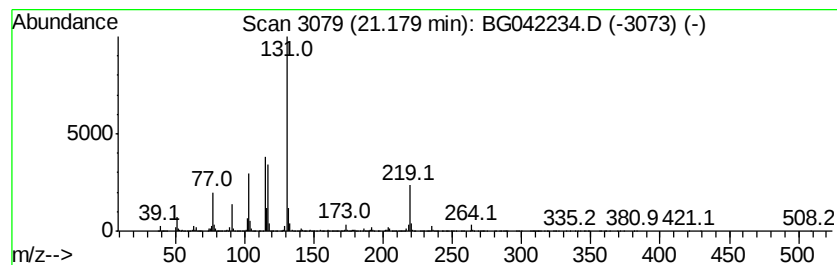
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Cinnamyl cinnamate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.18	43.97 ng/ul	2111400	Chrysene-d12	21.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	72
2		2,2-Dimethyl-1-(2-vinylphenyl)pr...	188	C13H16O	1000210-99-9	49
3		1-Penten-3-one, 4,4-dimethyl-1-p...	188	C13H16O	000538-44-3	47
4		N-(p-Vinylbenzoyl)-l-alanine	219	C12H13NO3	139184-60-4	47
5		2-Iodo-benzoic acid N'-(3-phenyl...	392	C16H13IN2O2	315672-62-9	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081519\
Data File : BG042234.D
Acq On : 15 Aug 2019 18:17
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Sample : K4183-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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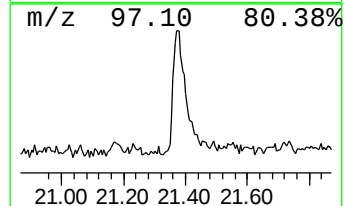
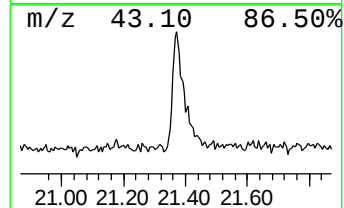
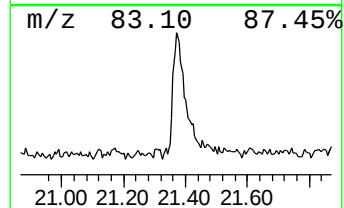
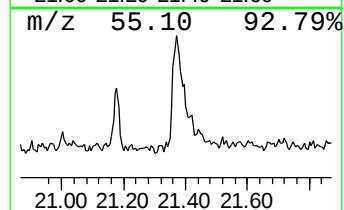
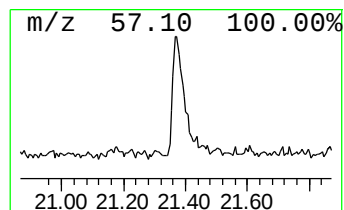
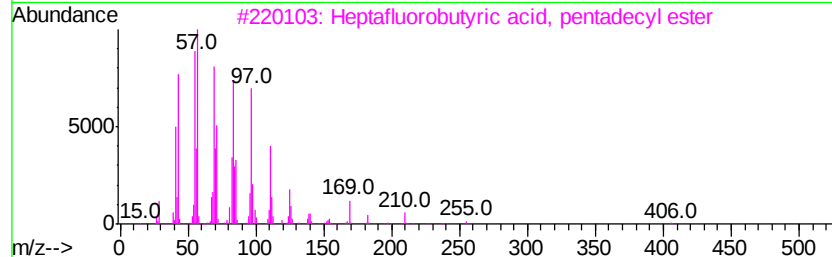
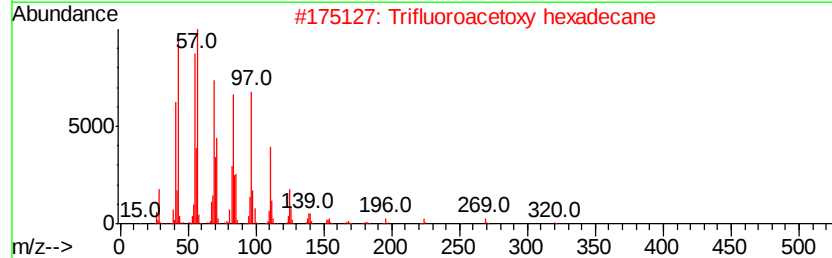
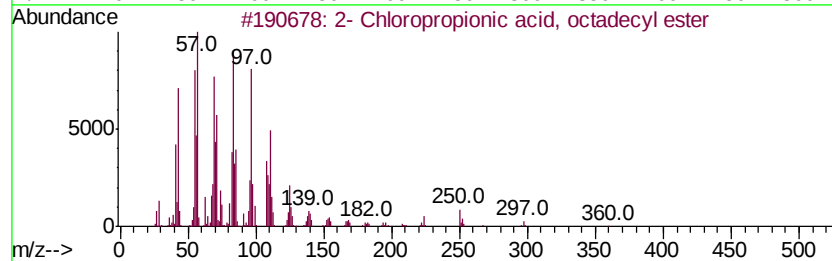
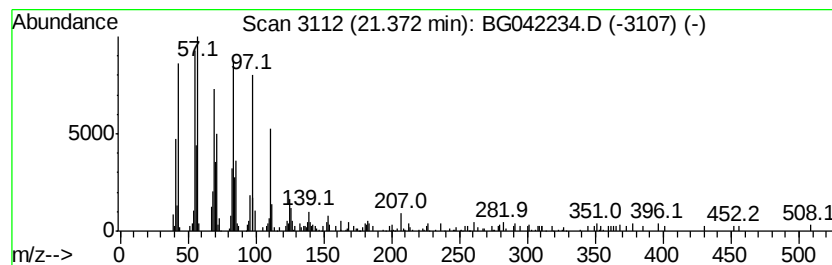
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2- Chloropropionic acid, oc... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.37	4.22 ng/ul	202507	Chrysene-d12	21.73

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-	Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
2	Trifluoroacetoxy	hexadecane	338	C18H33F3O2	006222-03-3	94
3	Heptafluorobutyric	acid, pentade...	424	C19H31F7O2	959261-23-5	93
4	10-Heneicosene	(c,t)	294	C21H42	095008-11-0	93
5	Pentafluoropropionic	acid, penta...	374	C18H31F5O2	959092-08-1	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG081519\
Data File : BG042234.D
Acq On : 15 Aug 2019 18:17
Operator : HP/JU
Sample : K4183-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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
Butane, 2-methoxy...	3.14	42.7	ng/ul	750769	1	8.02	351585	20.0
.alpha.-Pinene	6.70	19.8	ng/ul	347349	1	8.02	351585	20.0
1H-Cyclopenta[1,3...	13.57	2.9	ng/ul	97906	3	14.69	686254	20.0
Caryophyllene	14.02	33.8	ng/ul	1158410	3	14.69	686254	20.0
Aromandendrene	14.76	2.6	ng/ul	87781	3	14.69	686254	20.0
Naphthalene, 1,2,...	14.94	4.7	ng/ul	161139	3	14.69	686254	20.0
.alpha.-Cubebene	16.16	3.6	ng/ul	154054	4	17.44	858256	20.0
Cinnamyl cinnamate	21.18	44.0	ng/ul	2111400	5	21.73	960491	20.0
2- Chloropropioni...	21.37	4.2	ng/ul	202507	5	21.73	960491	20.0