

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100919\
 Data File : BG043147.D
 Acq On : 10 Oct 2019 21:10
 Operator : JU
 Sample : K5177-15
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BEPK1

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-BG100919MA.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	3.160	6	12	20	rBV	206860	390657	31.93%	2.682%
2	3.237	20	25	38	rVB	421730	639789	52.30%	4.393%
3	3.495	66	69	74	rVB4	6169	9352	0.76%	0.064%
4	3.883	133	135	137	rBV3	2505	1844	0.15%	0.013%
5	4.018	154	158	162	rBV3	6591	10784	0.88%	0.074%
6	4.147	177	180	185	rVV4	3880	5364	0.44%	0.037%
7	5.152	347	351	361	rBV3	6250	11981	0.98%	0.082%
8	5.316	377	379	383	rBV4	1207	1659	0.14%	0.011%
9	5.405	390	394	396	rVB2	1551	1765	0.14%	0.012%
10	5.487	406	408	412	rVB3	1601	1855	0.15%	0.013%
11	6.756	622	624	628	rVB2	1902	1761	0.14%	0.012%
12	7.026	669	670	673	rBV2	1449	1441	0.12%	0.010%
13	7.085	677	680	683	rBV	1373	2267	0.19%	0.016%
14	7.226	697	704	717	rBV2	69453	190543	15.58%	1.308%
15	7.379	724	730	745	rVB	61939	120628	9.86%	0.828%
16	7.590	760	766	778	rBV	93592	183002	14.96%	1.257%
17	7.667	778	779	784	rVB3	2647	3765	0.31%	0.026%
18	7.778	795	798	799	rBV	2668	1992	0.16%	0.014%
19	7.802	799	802	804	rVV	2165	1741	0.14%	0.012%
20	8.054	837	845	856	rBV	146928	272620	22.28%	1.872%
21	8.589	932	936	938	rBV3	966	1436	0.12%	0.010%
22	8.765	960	966	977	rBV3	60077	147488	12.06%	1.013%
23	8.883	984	986	996	rVB4	5356	10466	0.86%	0.072%
24	9.218	1035	1043	1050	rBV	81545	174305	14.25%	1.197%
25	9.641	1110	1115	1121	rVB3	4465	5476	0.45%	0.038%
26	9.940	1157	1166	1172	rBV4	73994	155967	12.75%	1.071%
27	10.070	1186	1188	1192	rVB2	1670	1798	0.15%	0.012%
28	10.111	1192	1195	1197	rBV	1182	1583	0.13%	0.011%
29	10.211	1209	1212	1215	rBV	1486	1693	0.14%	0.012%
30	10.487	1251	1259	1269	rBV2	91762	224918	18.39%	1.544%
31	10.857	1313	1322	1331	rBV	212928	399534	32.66%	2.743%
32	11.004	1341	1347	1360	rBV4	28646	92152	7.53%	0.633%
33	11.157	1371	1373	1376	rVB2	2527	2850	0.23%	0.020%
34	11.227	1383	1385	1388	rVB	1499	1860	0.15%	0.013%

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 Title : SVOA CALIBRATION

35	11.280	1391	1394	1397	rVB2	1174	1462	0.12%	0.010%
36	11.791	1479	1481	1485	rBV2	1072	1466	0.12%	0.010%
37	12.085	1527	1531	1534	rBV	1063	1472	0.12%	0.010%
38	12.449	1589	1593	1599	rVB3	1927	3586	0.29%	0.025%
39	13.037	1687	1693	1699	rBV	1223	2829	0.23%	0.019%
40	13.184	1713	1718	1720	rBV2	2694	3156	0.26%	0.022%
41	13.389	1748	1753	1756	rBV	1122	1908	0.16%	0.013%
42	13.595	1786	1788	1789	rBV	2923	1765	0.14%	0.012%
43	13.671	1797	1801	1807	rVB4	11504	17934	1.47%	0.123%
44	13.983	1849	1854	1857	rBV2	4662	6520	0.53%	0.045%
45	14.077	1863	1870	1874	rBV	287495	441337	36.08%	3.030%
46	14.124	1874	1878	1890	rVB	102791	175325	14.33%	1.204%
47	14.371	1912	1920	1936	rBV	311773	537855	43.97%	3.693%
48	14.594	1952	1958	1964	rVB3	16199	26176	2.14%	0.180%
49	14.676	1964	1972	1982	rBV	405244	657248	53.73%	4.513%
50	14.764	1983	1987	1992	rVB4	12768	18640	1.52%	0.128%
51	14.911	2003	2012	2028	rBV5	27200	127199	10.40%	0.873%
52	15.176	2054	2057	2064	rBV4	2052	4279	0.35%	0.029%
53	15.422	2095	2099	2100	rBV4	1618	1549	0.13%	0.011%
54	15.469	2104	2107	2112	rVV3	1948	2975	0.24%	0.020%
55	15.528	2113	2117	2119	rVB3	2022	2461	0.20%	0.017%
56	15.587	2124	2127	2130	rVB	1446	1933	0.16%	0.013%
57	15.669	2134	2141	2153	rBV	390090	691591	56.53%	4.749%
58	15.781	2154	2160	2171	rVB	93038	175194	14.32%	1.203%
59	15.910	2179	2182	2186	rVB4	4128	3977	0.33%	0.027%
60	16.268	2240	2243	2247	rBV2	1290	1809	0.15%	0.012%
61	16.345	2254	2256	2259	rVV2	1312	1854	0.15%	0.013%
62	16.374	2259	2261	2265	rVV2	3601	4248	0.35%	0.029%
63	16.968	2358	2362	2363	rBV3	2691	2438	0.20%	0.017%
64	16.997	2363	2367	2368	rBV3	1340	1811	0.15%	0.012%
65	17.173	2393	2397	2400	rBV3	2975	4939	0.40%	0.034%
66	17.214	2400	2404	2407	rVV2	1475	2247	0.18%	0.015%
67	17.285	2412	2416	2417	rBV	1744	1484	0.12%	0.010%
68	17.332	2422	2424	2426	rVV	2350	1537	0.13%	0.011%
69	17.420	2432	2439	2449	rBV	558163	878796	71.84%	6.034%
70	17.520	2450	2456	2466	rVB	427821	704310	57.57%	4.836%
71	17.761	2495	2497	2500	rBV4	1591	1712	0.14%	0.012%

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72	17.802	2500	2504	2506	rVV2	1358	1640	0.13%	0.011%
73	17.908	2518	2522	2526	rBV7	3802	4623	0.38%	0.032%
74	17.949	2526	2529	2530	rBV3	1762	1850	0.15%	0.013%
75	18.078	2548	2551	2553	rBV3	1963	1935	0.16%	0.013%
76	18.131	2559	2560	2562	rBV	2388	1678	0.14%	0.012%
77	18.195	2566	2571	2574	rVV4	4059	5456	0.45%	0.037%
78	18.307	2586	2590	2602	rBV3	40418	85998	7.03%	0.591%
79	18.601	2635	2640	2641	rBV4	4190	4318	0.35%	0.030%
80	18.801	2669	2674	2678	rBV4	3929	7662	0.63%	0.053%
81	19.153	2732	2734	2735	rBV2	2298	1612	0.13%	0.011%
82	19.230	2743	2747	2751	rBV5	11171	14624	1.20%	0.100%
83	19.441	2779	2783	2784	rBV4	9475	11103	0.91%	0.076%
84	19.688	2822	2825	2827	rBV4	4744	7462	0.61%	0.051%
85	19.800	2839	2844	2855	rBV	653771	1030728	84.26%	7.077%
86	20.199	2910	2912	2915	rVB4	5289	3689	0.30%	0.025%
87	20.328	2931	2934	2939	rVB3	30330	35169	2.87%	0.241%
88	20.828	3015	3019	3023	rVB3	24154	28411	2.32%	0.195%
89	21.333	3100	3105	3116	rBV2	184835	395044	32.29%	2.713%
90	21.691	3159	3166	3177	rBV	585762	1072040	87.63%	7.361%
91	21.879	3195	3198	3202	rBV2	31512	43936	3.59%	0.302%
92	22.514	3297	3306	3322	rBV3	176462	634110	51.83%	4.354%
93	23.184	3415	3420	3426	rVB2	55895	102402	8.37%	0.703%
94	23.530	3474	3479	3491	rVB9	46947	98722	8.07%	0.678%
95	23.989	3550	3557	3564	rBV2	209246	434620	35.53%	2.984%
96	24.071	3564	3571	3582	rBV7	54267	208973	17.08%	1.435%
97	24.688	3666	3676	3685	rBV	333256	941700	76.98%	6.466%
98	24.905	3704	3713	3727	rVB2	389871	1223343	100.00%	8.400%
99	25.928	3879	3887	3897	rVB4	69064	195143	15.95%	1.340%
100	26.063	3903	3910	3921	rVB3	131385	372159	30.42%	2.555%

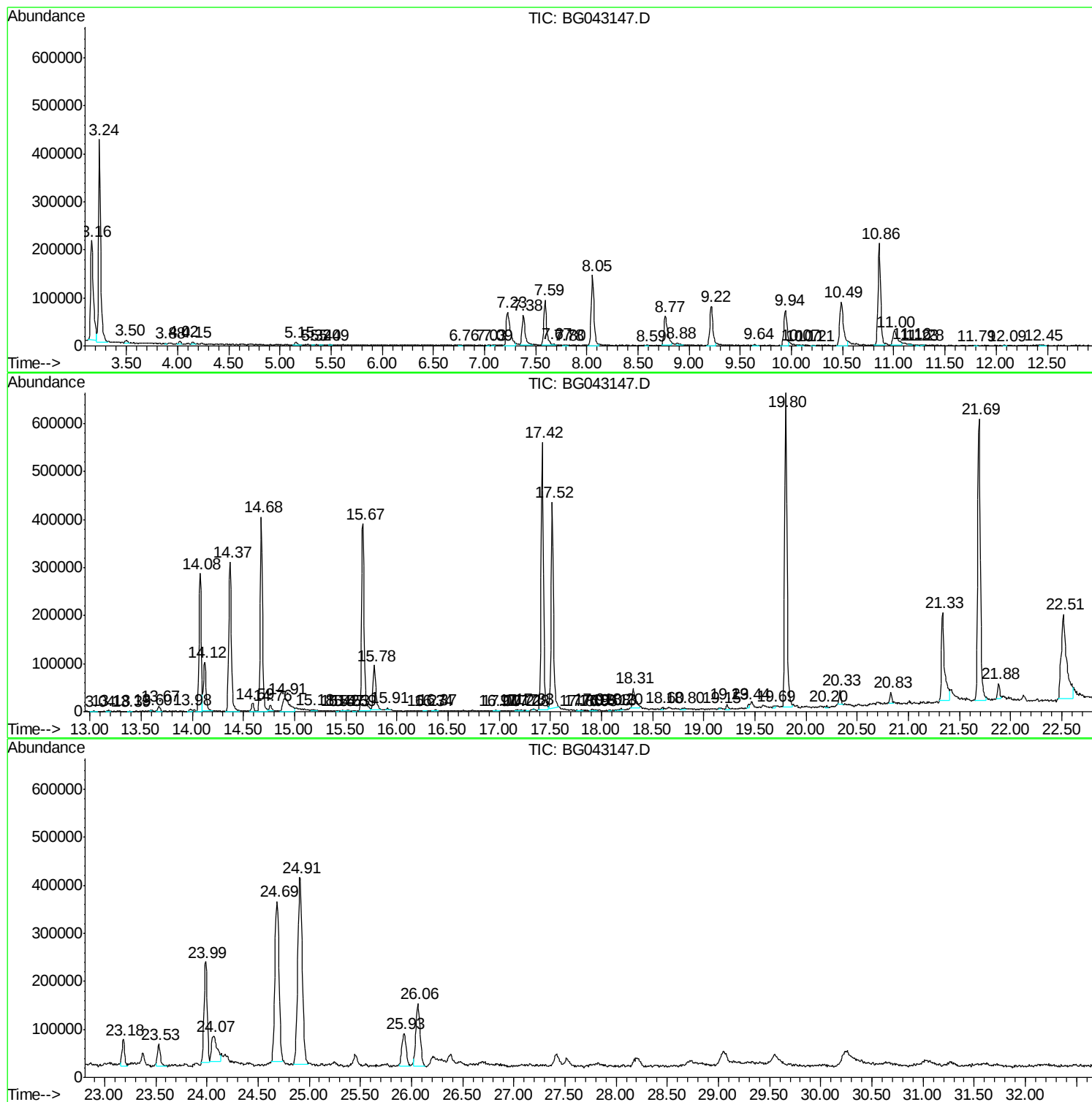
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG100919MA.M
Quant Title : SVOA CALIBRATION

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



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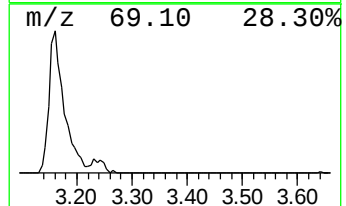
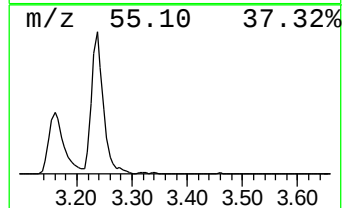
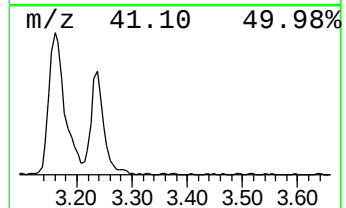
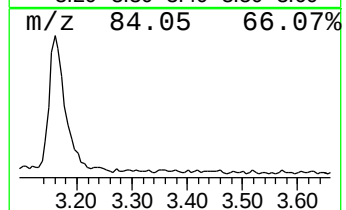
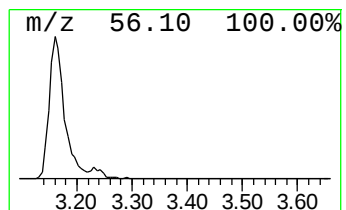
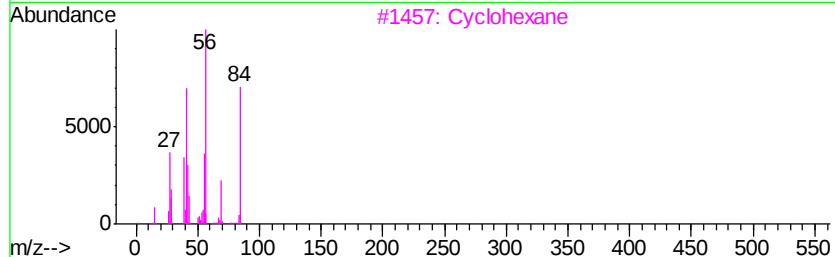
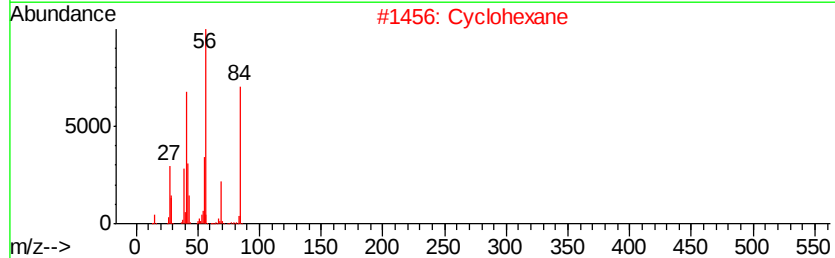
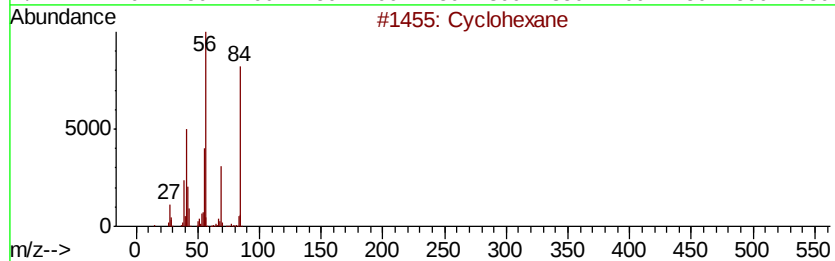
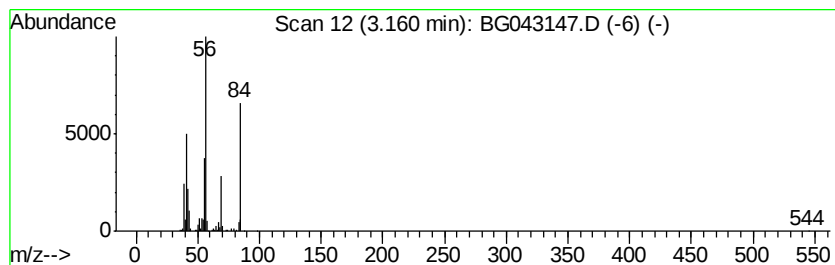
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG100919MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic3.16 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.16	28.66 ng/ul	390657	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane	84	C6H12	000110-82-7	95
2		Cyclohexane	84	C6H12	000110-82-7	91
3		Cyclohexane	84	C6H12	000110-82-7	91
4		Cyclohexane	84	C6H12	000110-82-7	90
5		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	86



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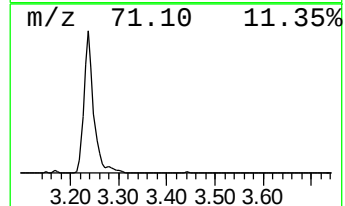
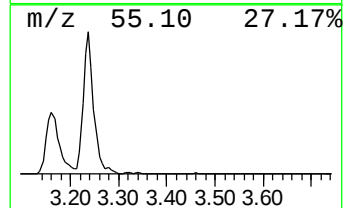
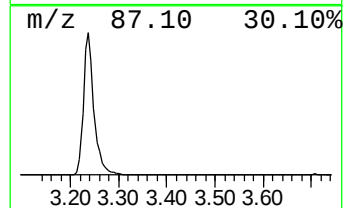
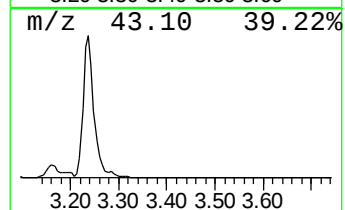
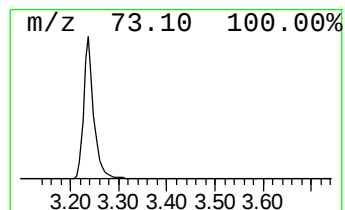
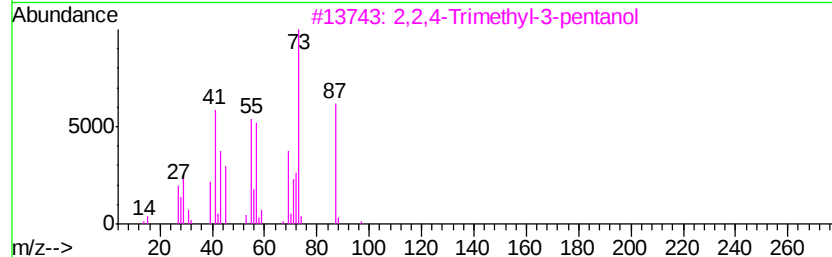
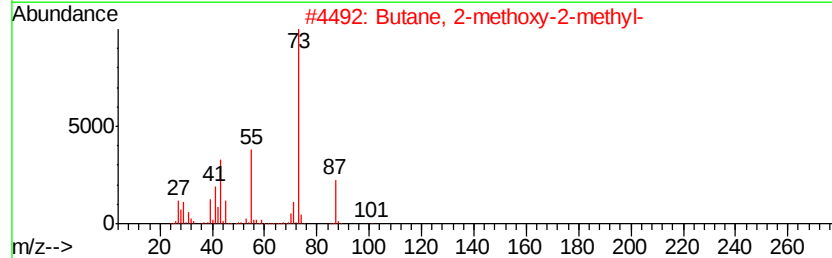
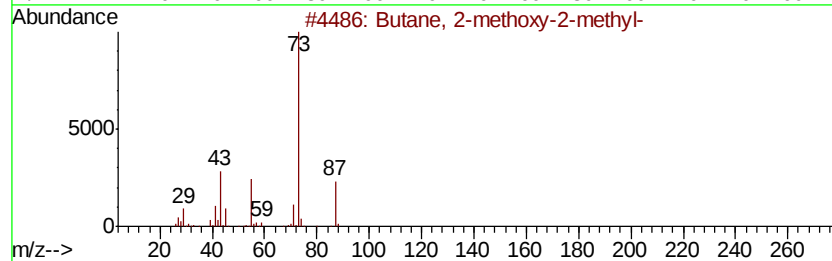
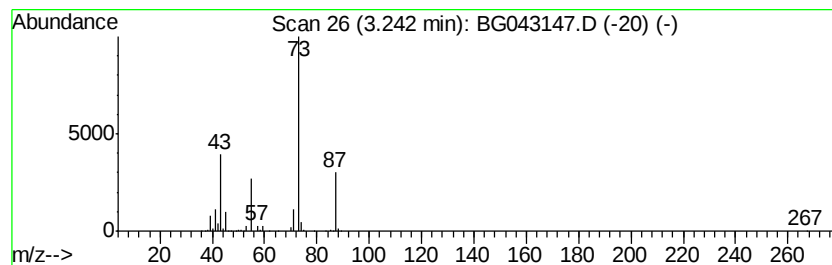
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.24	46.94 ng/ul	639789	1,4-Dichlorobenzene-d4	8.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	56
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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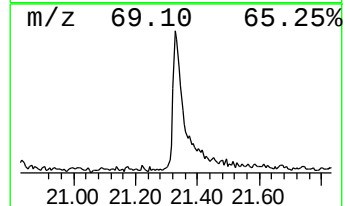
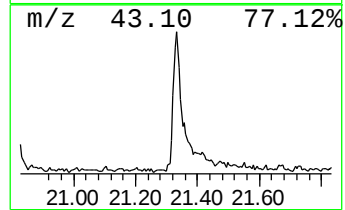
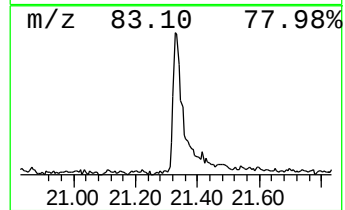
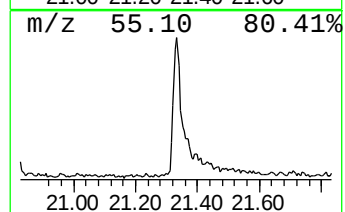
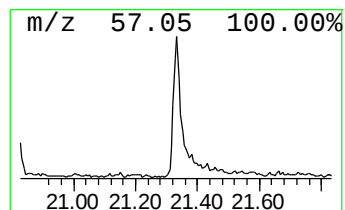
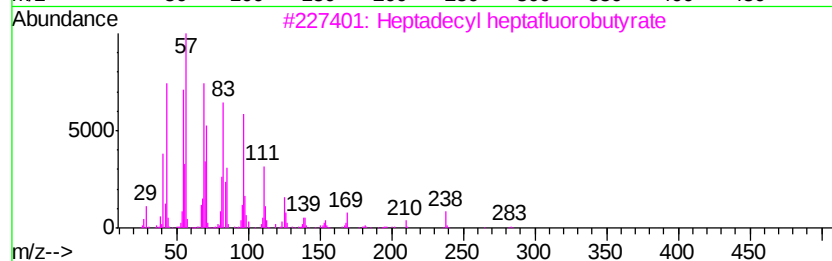
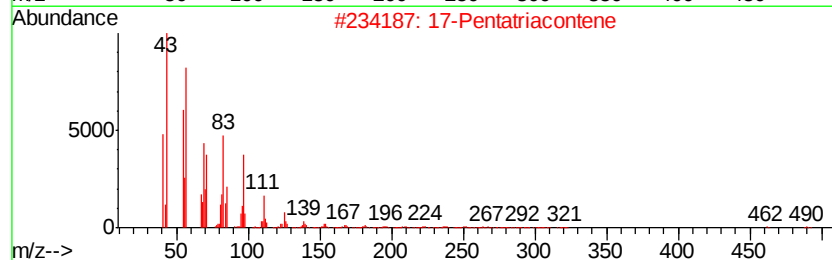
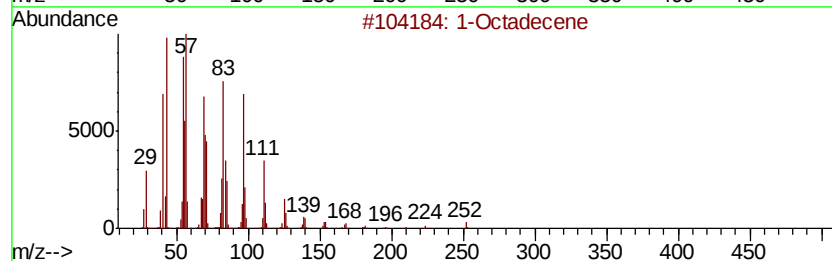
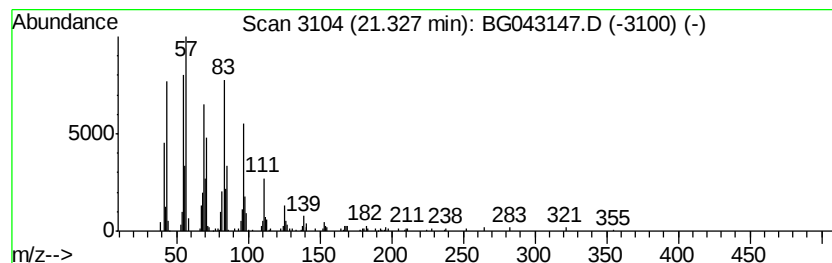
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TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Cyclic21.33 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.33	7.37 ng/ul	395044	Chrysene-d12	21.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Octadecene	252	C18H36	000112-88-9	93
2		17-Pentatriacontene	491	C35H70	006971-40-0	91
3		Heptadecyl heptafluorobutvrate	452	C21H35F7O2	959085-66-6	91
4		Hexacosyl trifluoroacetate	478	C28H53F3O2	1000351-75-0	91
5		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100919\
Data File : BG043147.D
Acq On : 10 Oct 2019 21:10
Operator : JU
Sample : K5177-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BEPK1

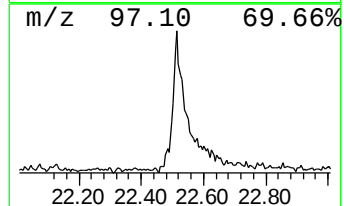
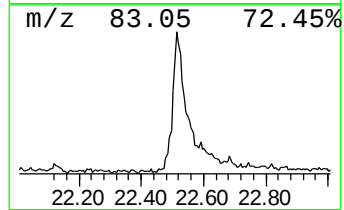
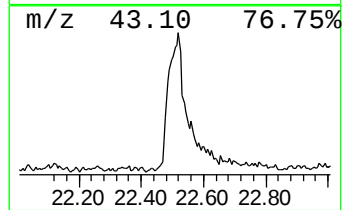
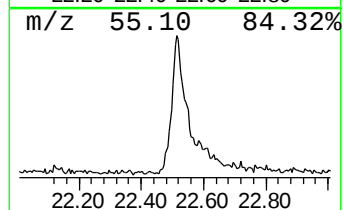
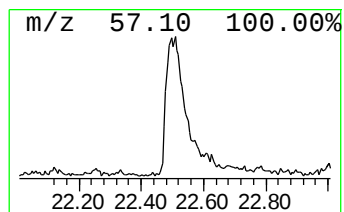
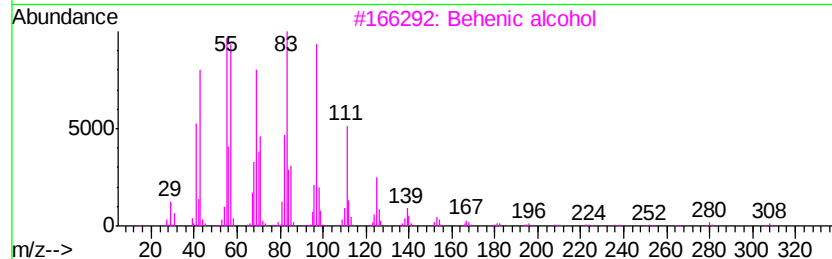
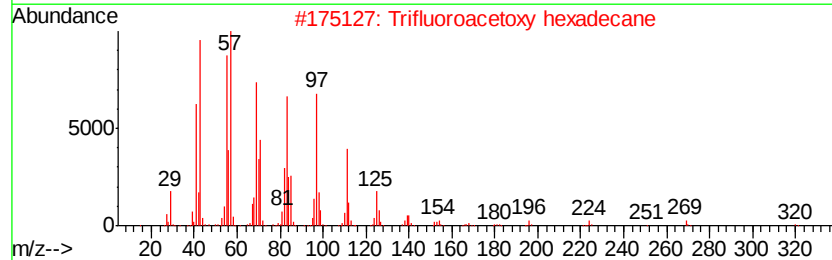
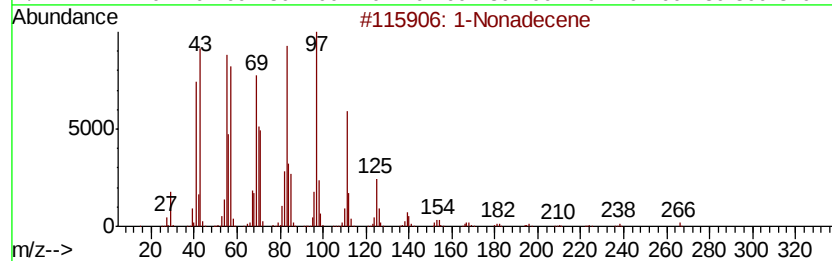
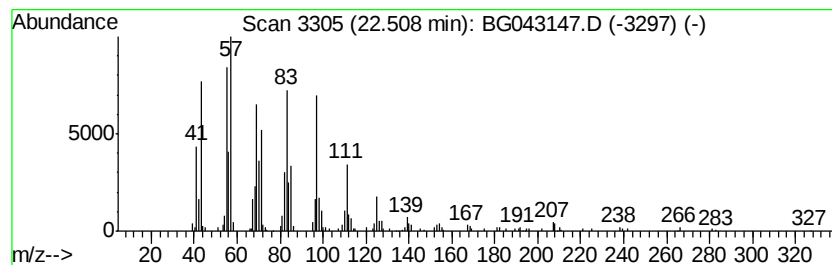
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Cyclic22.51 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.51	11.83 ng/ul	634110	Chrysene-d12	21.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Nonadecene	266	C19H38	018435-45-5	97
2		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
3		Behenic alcohol	326	C22H46O	000661-19-8	93
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	93
5		1-Heneicosanol	312	C21H44O	015594-90-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100919\
Data File : BG043147.D
Acq On : 10 Oct 2019 21:10
Operator : JU
Sample : K5177-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BEPK1

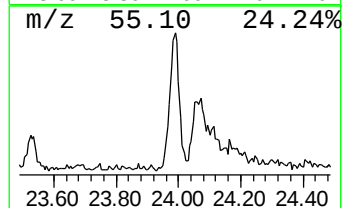
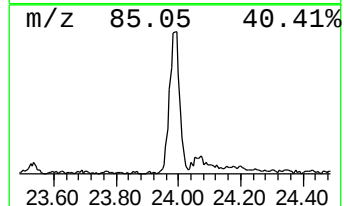
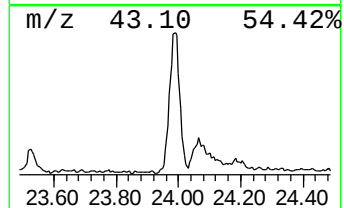
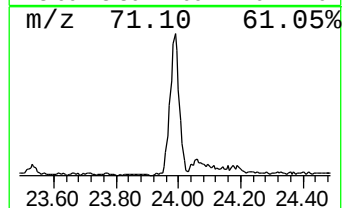
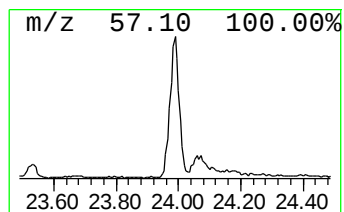
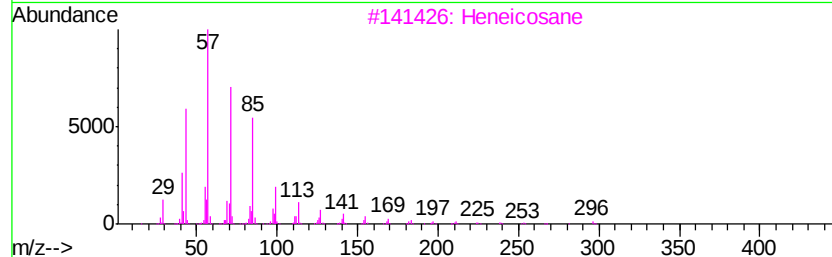
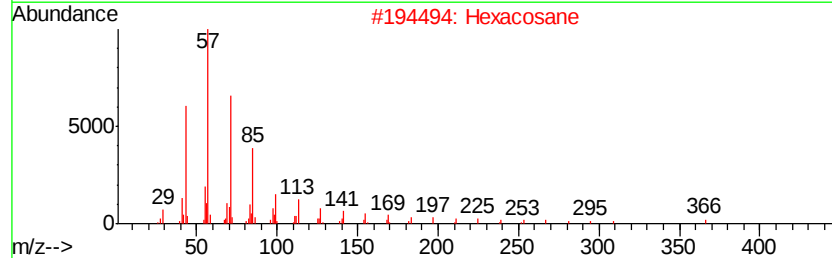
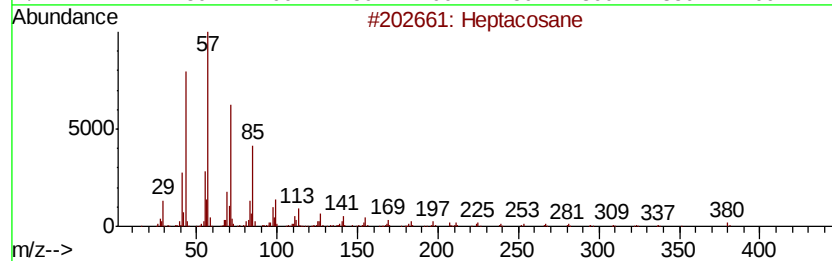
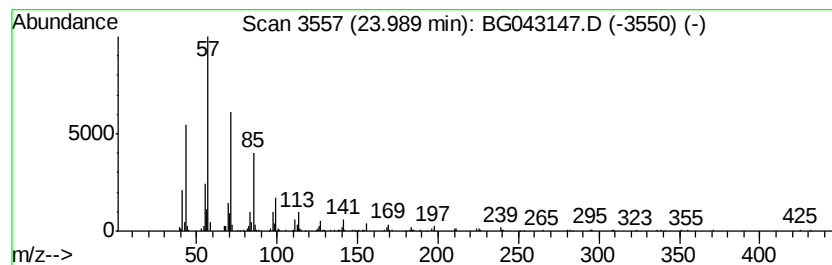
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.99	7.11 ng/ul	434620	Perylene-d12	24.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		Hexacosane	366	C26H54	000630-01-3	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Hentriacontane	437	C31H64	000630-04-6	90
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100919\
Data File : BG043147.D
Acq On : 10 Oct 2019 21:10
Operator : JU
Sample : K5177-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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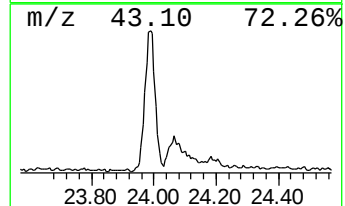
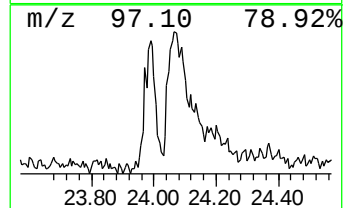
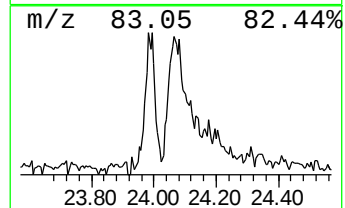
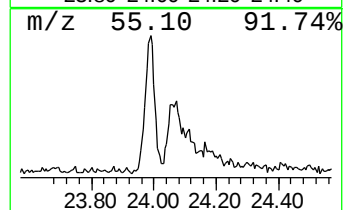
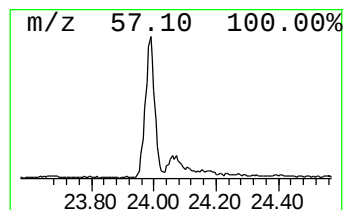
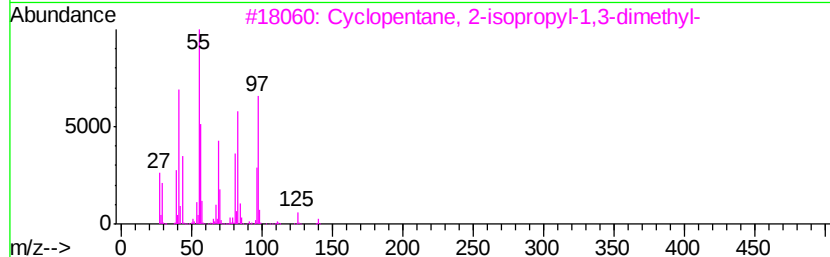
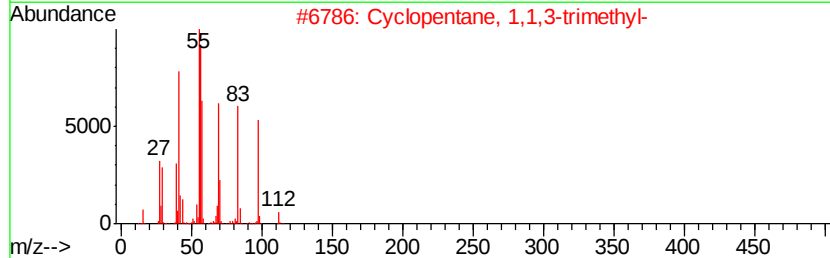
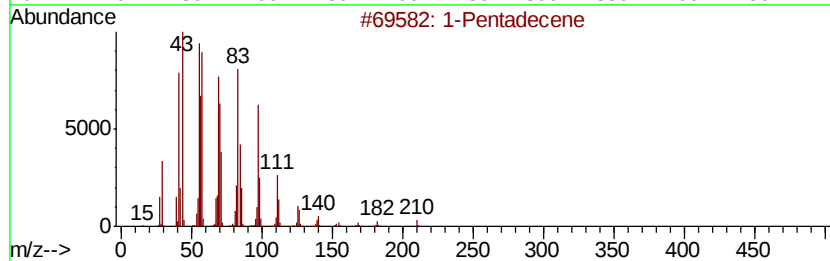
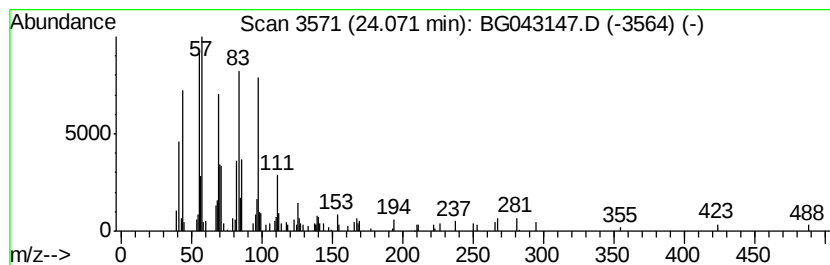
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic24.07 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.07	3.42 ng/ul	208973	Perylene-d12	24.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentadecene	210	C15H30	013360-61-7	83
2		Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	58
3		Cyclopentane, 2-isopropyl-1,3-di...	140	C10H20	032281-85-9	58
4		Octacosyl acetate	452	C30H60O2	018206-97-8	50
5		1-Heptacosanol	396	C27H56O	002004-39-9	45



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100919\
Data File : BG043147.D
Acq On : 10 Oct 2019 21:10
Operator : JU
Sample : K5177-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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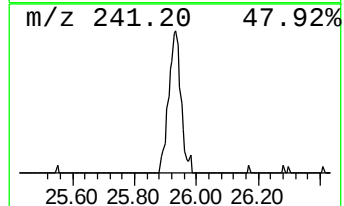
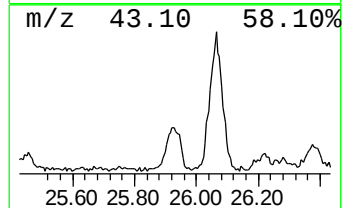
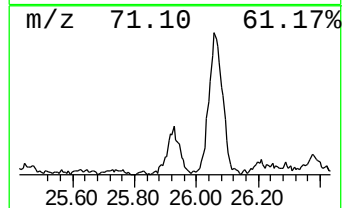
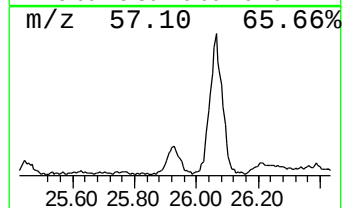
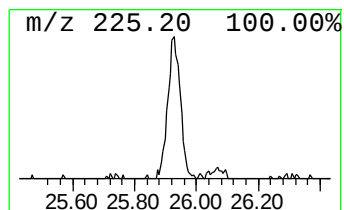
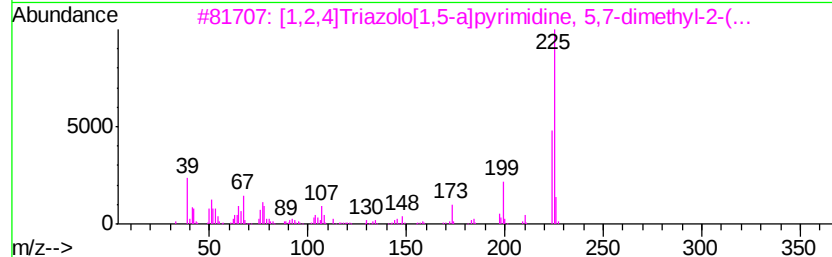
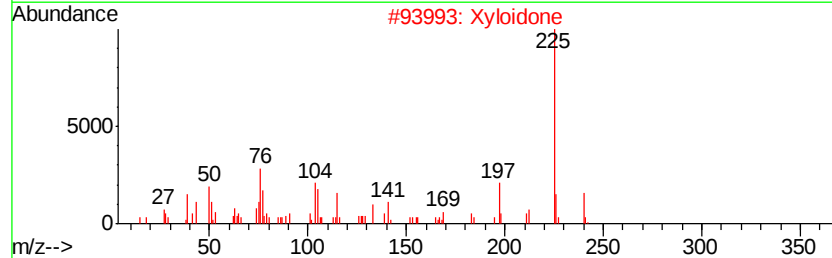
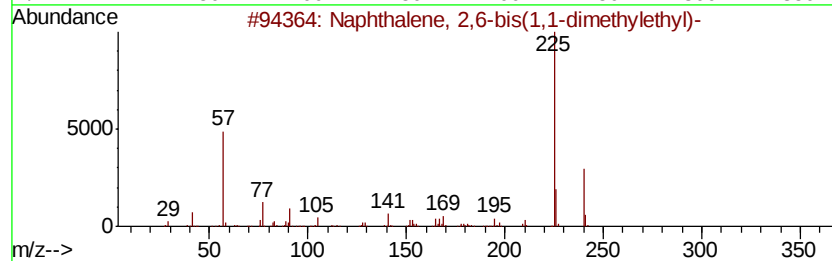
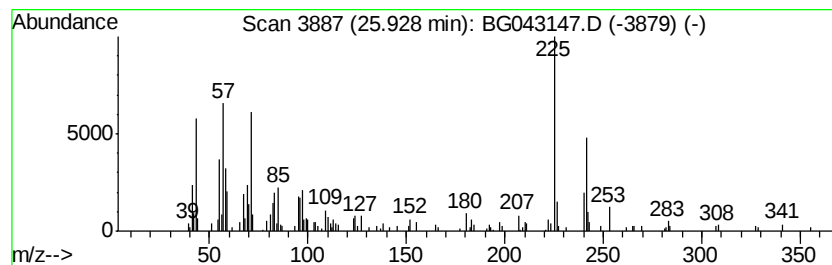
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.93	3.19 ng/ul	195143	Perylene-d12	24.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,6-bis(1,1-dimethyl-...	240	C18H24	003905-64-4	27
2		Xyloidone	240	C15H12O3	015297-92-4	22
3		[1,2,4]Triazolo[1,5-a]pyrimidine...	225	C12H11N5	1000320-01-2	20
4		Propionitrile, 3-(5-methyl-3-phe...	225	C13H11N3O	1000263-77-0	18
5		4-Amino-5,7-dinitrobenzofurazan	225	C6H3N5O5	021123-74-0	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100919\
Data File : BG043147.D
Acq On : 10 Oct 2019 21:10
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Sample : K5177-15
Misc :
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Instrument :
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ClientSampleId :
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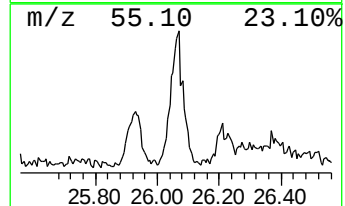
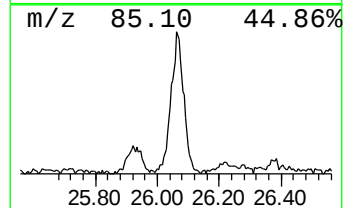
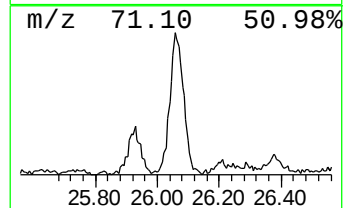
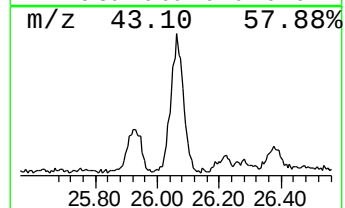
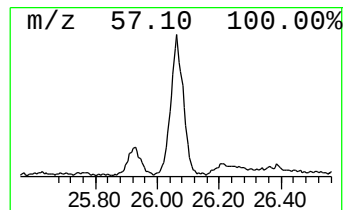
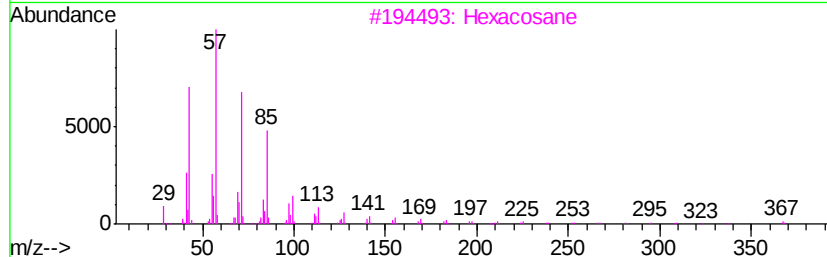
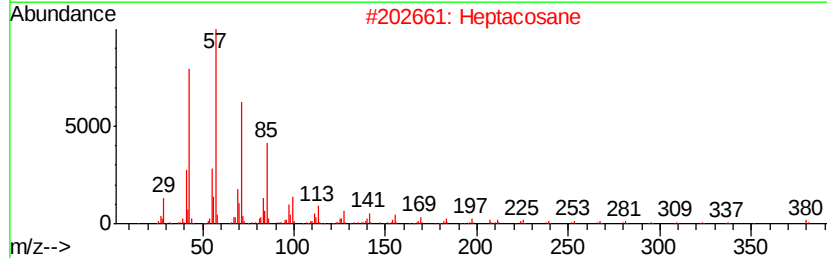
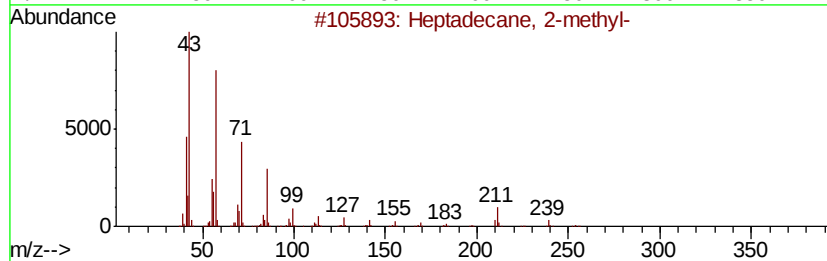
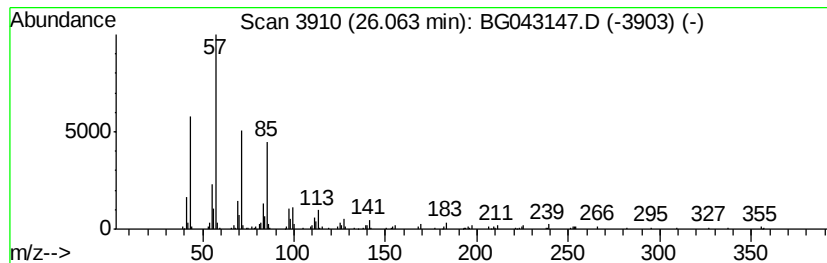
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.06	6.08 ng/ul	372159	Perylene-d12	24.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2-methyl-	254	C18H38	001560-89-0	86
2		Heptacosane	380	C27H56	000593-49-7	86
3		Hexacosane	366	C26H54	000630-01-3	86
4		Nonadecane	268	C19H40	000629-92-5	80
5		Hexadecane	226	C16H34	000544-76-3	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG100919\
Data File : BG043147.D
Acq On : 10 Oct 2019 21:10
Operator : JU
Sample : K5177-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BEPK1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG100919MA.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
(DEL) Alkane: Cyc...	3.16	28.7	ng/ul	390657	1	8.05	272620	20.0
Butane, 2-methoxy...	3.24	46.9	ng/ul	639789	1	8.05	272620	20.0
(DEL) Alkane: Cyc...	21.33	7.4	ng/ul	395044	5	21.69	1072040	20.0
(DEL) Alkane: Cyc...	22.51	11.8	ng/ul	634110	5	21.69	1072040	20.0
(DEL) Alkane: Str...	23.99	7.1	ng/ul	434620	6	24.91	1223340	20.0
(DEL) Alkane: Cyc...	24.07	3.4	ng/ul	208973	6	24.91	1223340	20.0
unknown-01	25.93	3.2	ng/ul	195143	6	24.91	1223340	20.0
(DEL) Alkane: Str...	26.06	6.1	ng/ul	372159	6	24.91	1223340	20.0